(805) 526-7270 fax

Columbia Analytical Services INC.

LABORATORY REPORT

May 28, 2008

Robert Kennedy ENSR 2 Technology Park Drive Westford, MA 01886

RE: Phase B Soil Gas / 04020-023-4311

Dear Robert:

Enclosed are the results of the samples submitted to our laboratory on May 8, 2008. For your reference, these analyses have been assigned our service request number P0801342.

All analyses were performed in accordance with our laboratory's quality assurance program. Results are intended to be considered in their entirety and apply only to the samples analyzed and reported herein. Your report contains **279** pages.

Columbia Analytical Services, Inc. is certified by the California Department of Health Services, NELAP Laboratory Certificate No. 02115CA; Arizona Department of Health Services, Certificate No. AZ0694; Florida Department of Health, NELAP Certification E871020; New Jersey Department of Environmental Protection, NELAP Laboratory Certification ID #CA009; New York State Department of Health, NELAP NY Lab ID No: 11221; Oregon Environmental Laboratory Accreditation Program, NELAP ID: CA20007; The American Industrial Hygiene Association, Laboratory #101661; Department of the Navy (NFESC); Pennsylvania Registration No. 68-03307. Each of the certifications listed above have an explicit Scope of Accreditation that applies to specific matrices/methods/analytes; therefore, please contact me for information corresponding to a particular certification.

If you have any questions, please call me at (805) 526-7161.

Respectfully submitted,

Columbia Analytical Services, Inc.

Keely M Horince

Kelly Horiuchi **Project Manager**

Page 1 of <u>279</u>



Client: ENSR Project: Phase B Soil Gas / 04020-023-4311 CAS Project No: I

P0801342

CASE NARRATIVE

The samples were received intact under chain of custody on May 8, 2008 and were stored in accordance with the analytical method requirements. Please refer to the sample acceptance check form for additional information. The results reported herein are applicable only to the condition of the samples at the time of sample receipt.

Helium Analysis

Two of the samples were analyzed for helium according to modified EPA Method 3C using a gas chromatograph equipped with a thermal conductivity detector (TCD).

Volatile Organic Compound Analysis

All of the samples were analyzed for selected volatile organic compounds. In addition, one of the samples was analyzed for tentatively identified compounds in accordance with EPA Method TO-15 from the Compendium of Methods for the Determination of Toxic Organic Compounds in Ambient Air, Second Edition (EPA/625/R-96/010b), January, 1999. The analytical system was comprised of a gas chromatograph / mass spectrometer (GC/MS) interfaced to a whole-air preconcentrator.

The results of analyses are given in the attached laboratory report. All results are intended to be considered in their entirety, and Columbia Analytical Services, Inc. (CAS) is not responsible for utilization of less than the complete report.

Client: ENSR Project: Phase B Soil Gas 04020-023-4311

Folder: P0801342

Detailed Sample Information

| Order # | 8616 | |
|---|-----------------------------|---|
| FC ID | OA00080 | |
| Order # | 8616 | 8616 F |
| Cont ID | SC00979 | SC00564 SC00791 |
| Pf2 | | |
| <u>Pi2</u> (psig) | | |
| <u>Pi2</u> (Hg) | | 1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1 |
| <u>Pf1</u> | 3.5 | 3.5 3.5 |
| <u>Pi1</u> (psig) | -3.6 | - 3.8 -4.2 |
| <u>Pi1</u> (Hg) | -7.4 | -7.8 -8.5 |
| Container Type | 6.0 L-Summa Canister Source | 6.0 L-Summa Canister Source 6.0 L-Summa Canister Source |
| CAS Sample ID Client Sample ID Container Type | 20801342-001.01 SG83B-05-1 | 20801342-002.01 SG83B-05-3 20801342-003.01 SG83B-05-7 |
| CAS S | P080134 | P080134 P080134 |

Miscellaneous Items - received

C3/8/2008 2:52:06PM

| Columbia Analytical Services ^{we} | An Employee - Owned Company |
|--|-----------------------------|
| N. | An |

אוומווי הו המשיטעו וובהטוח א הוומוזווים הבו גורם וובלתבאו 2655 Park Center Drive, Suite A Simi Valley California 93065

Page _____ of

| Services ^{we} | SILLI VALLEY, CALIFOLITIA 93003 | IIII 200 | 00 | | | | | | | | | |
|---|--|--|----------------------------------|--|---|---|--------------------------------|---|---|-----------------------------|----------------|--------------------------------------|
| An Employee - Owned Company | Phone (805) 526-7161 Fax (805) 526-7270 | 7161 70 | | He - He | quested Jur ay (100%) | naround Time 2 Day (75%) 3 | In Business Day (50%) | Hequested Turnaround Time in Business Days (Surcharges) please circle 1 Day (100%) 2 Day (75%) 3 Day (50%) 4 Day (35%) 5 Day (25%) 10 Day - Standard | jes) please circ Jay (25%) 10 D | :le ay - Standard | CAS Proje | CAS Project No. |
| | | | | | | | | CAS Contact | | | | 1010 |
| Company Name & Address (F | Reporting Information | | Project Name | ne | | | | | | | | - |
| ENSK Provinski da | NA ACO | | Phase | Physe B South | ill Gas | 5 | L | A | Analysis Method and/or Analytes | nd/or Analytes | | - <u>-</u> |
| Camavillo, CA 93012 | 93012 | | Project Number | ect Number $U \neq 0 2 v = 0.2$ | 0.23- | 3-4311 | | | | | | |
| Project Manager MILC FIACK | | <u> </u> | P.O. # / Bill | P.O. # / Billing Information $\vec{E} \wedge$ | uo | | | 5 | 200104-00-00 |)I | | Comments e.g. Actual Preservative |
| Phone 805-388-5775 845.388-3577 | ax 105 - 38 8 - 39 | 116 | | | | | | 1-01 | | <u> </u> | | or specific instructions |
| Email Address for Result Reporting | orting | | Sampler (Print & S I.C. Strue | Sampler (Print & Sign) Ian Stone | Ju | B | | . 4 | шnj | 51- | | |
| Client Sample ID | Laboratory ID Number Co | Date Collected | Time Collected | Sample Type (Air/Tube/ Solid) | Canister ID (Bar Code # - AC, SC, etc.) | Flow Controller (Bar Code - FC #) | Sample Volume | <u>24-1</u> 52 | 3H | Q | | |
| 59838-06-1 | 1/5/ | 517/00/1433 | 1433 | MR | 6160075 | 0800040 | 19 | × | × | | | |
| SG 833-05-3 | | XOLE S | 1221 | Alk | Puzosiy | SLEDST4 UA00080 | 9L | X | × | | | |
| SG833-05-7 | (C) 5/ | 30/6/5 | N 39 | AIR | SLOOPH !! | 08000 YO | ĥι | × | | × | | |
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| Report Tier Levels - please select Tier 1 - (Results/Default if not specified) _ Tier II - (Results + QC) | | Tier III - (Data Validation Tier V - (client specified) | Validation Pack | Tier III - (Data Validation Package) 10% Surcharge- Tier V - (client specified) | large | | EDD required Yes / No Type: | 0N / s | EDD Units: | | Project Re | Project Requirements (MRLs, QAPP) |
| Relinquished by: (Signature) | d | | Date: 1 fo & | Time: | Received by: (Signature) | (Signature) | 1470 | YYYY | 2 por | Lar Tople | 2 | v |
| Relinquished by: (Signature) | | | Date: | | Received by: (Signature) | (Signature) | | | Date: | Time: | Cooler / Blank | ank |
| Relinquished by: (Signature) | | | Date: | Time: | Received by: (Signature) | (Signature) | | | . Date: | Time: | Temnerature | °C °° |

| | | | | bia Analytica de Acceptance | | | | | | |
|-------------------|-------------------------------|---|---------------------|--|---------------------------------------|--------------------------|-----------------------|------------|----------------|--------------|
| Clien! | t: ENSR | | Samb | le Acceptiones | Click I VII | Work order: | P0801342 | | | |
| | | Gas / 04020-023-4311 | 1 | | - | | | | | |
| U | e(s) received on | | | | Date opened: | : 05/08/08 | by: | MZAN | MORA | |
| <u>Note:</u> This | s form is used for <u>all</u> | ll samples received by CAS. | The use of this for | m for custody seals | s is strictly meant | to indicate presence/ | /absence and not as a | n indicati | on of | |
| compliance | e or nonconformity. | Thermal preservation and pl | H will only be eva | duated either at the | request of the cli | ient and/or as require | d by the method/SOI | | Na | NEA |
| 1 | Wore comple | | trad with (| tiont comple II | D 0 | | | <u>Yes</u> | <u>No</u> □ | <u>N/A</u> |
| 1 | - | e containers properly r supplied by CAS? | Markeu wiui e | ment sample it |)(| | | | | |
| 2 | | containers arrive in go | and condition (| n | | | | | | |
| 3 4 | • | of-custody papers used | | | | | | | | |
| 5 | | container labels and/o | | | march | | | | | |
| 6 | - | volume received adequ | | - ~ . | pers: | | | | | |
| 7 | - | within specified holdir | - | 515 : | | | | X | | |
| 8 | - - | emperature (thermal | - | of cooler at rec | ceint adhered | to? | | | | \mathbf{X} |
| C | | Cooler Temperature | preser | | Temperature | | °C | - | Kana | |
| 9 | | lank received? | | •••••••••••••••••••••••••••••••••••••• | * ****P +2 | | | | \mathbf{X} | |
| | - | supplied by CAS: Seri | ial # | | -TB | | | | | - |
| 10 | - | y seals on outside of co | | | • | | | | \mathbf{X} | |
| | Location of | · | | | | | Sealing Lid? | | | \mathbf{X} |
| | Were signat | ture and date included | 1? | •••••••••••••••••••••••••••••••••••••• | | | | | | \times |
| | Were seals i | intact? | | | | | | | | \mathbf{X} |
| | Were custody | v seals on outside of sam | mple containe | r? | | | | | X | |
| | Location of | seal(s)? | | · . | | | Sealing Lid? | | | \mathbf{X} |
| | Were signat | ture and date included | 1? | | | | | | | X |
| | Were seals i | intact? | | | | | | | | X |
| 11 | Do containers | s have appropriate pre | servation, acc | cording to met | nod/SOP or C | Client specified i | information? | | | \mathbf{X} |
| | Is there a clie | ent indication that the | submitted san | nples are pH p | reserved? | | | | | \mathbf{X} |
| | Were <u>VOA v</u> | vials checked for prese | nce/absence o | of air bubbles? | | | | | | \mathbf{X} |
| | Does the clier | nt/method/SOP requir | e that the ana | lyst check the : | sample pH ar | nd <u>if necessary</u> a | lter it? | | | X |
| 12 | Tubes: | Are the tubes cap | ped and intac | t? | | | | | | \mathbf{X} |
| | | Do they contain a | moisture? | | | | | | | \mathbf{X} |
| 13 | Badges: | Are the badges p | roperly cappe | d and intact? | | | | | | \mathbf{X} |
| | | Are dual bed bad | lges separated | and individua | lly capped an | id intact? | | | | X |
| Lab | Sample ID | Container | Required | Received | Adjusted | VOA Headspac | e Receip | ot / Prese | ervation | , I |
| | | Description | pH * | pH | pH | (Presence/Absence | | Commen | | |
| P0801342 | 2-001.01 | 6.0 L Source Can | | | 1 | | Ť | | | |
| P0801342 | 2-002.01 | 6.0 L Source Can | | | · · · · · · · · · · · · · · · · · · · | | | | | |

Required pH: Phenols/COD/NH3/TOC/TOX/NO3+NO2/TKN/T.PHOS, H2SO4 (pH<2); Metals, HNO3 (pH<2); CN (NaOH or NaOH/Asc Acid) (pH>12);

Explain any discrepancies: (include lab sample ID numbers):

6.0 L Source Can

P0801342-003.01

5

RESULTS OF ANALYSIS

Page 1 of 1

| Client: Client Project ID: | ENSR Phase B Soil Gas / 04020-023 | 3-4311 | | CA | S Project ID: P0801342 | |
|--|--|------------------------------|--------------------------------|----------------|---|--|
| | | I | Ielium | | | |
| Fest Code: Instrument ID: Analyst: Sampling Media: Fest Notes: | EPA 3C Modified HP5890 II/GC8/TCD Zheng Wang/Wade Henton/Ch 6.0 L Summa Canister(s) | nris Cornett | | D | (s) Collected: 5/7/08 ate Received: 5/8/08 ate Analyzed: 5/8/08 | |
| Client Sample ID | CAS Sample ID | Injection Volume ml(s) | Canister Dilution Factor | Result ppmV | MRL ppmV | |
| SG83B-05-1 SG83B-05-3 | P0801342-001 P0801342-002 | 1.00 | 1.64 | 2,700 160 | 41 42 | |

1.00 1.00

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

P080508-MB

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

1.00

ND

25

Method Blank

Verified By:

Date:

6

Data Qualifier

RESULTS OF ANALYSIS

Page 1 of 3

ENSR Client: Client Sample ID: SG83B-05-1 Client Project ID: Phase B Soil Gas / 04020-023-4311

Test Code: EPA TO-15 Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 Instrument ID: Rusty Bravo Sampling Media: 6.0 L Summa Canister Test Notes:

SC00979 Container ID:

Analyst:

Initial Pressure (psig):

-3.6

Final Pressure (psig): 3.5

Canister Dilution Factor: 1.64

0.010 Liter(s) 0.0010 Liter(s)

CAS Project ID: P0801342 CAS Sample ID: P0801342-001

Date Collected: 5/7/08

Date Received: 5/8/08

Date Analyzed: 5/8/08

Volume(s) Analyzed:

| CAS # | Compound | Result | MRL | MDL | Result | MRL | MDL | Data |
|-----------|--|--------|-------------------|-------------------|--------|------|------|---|
| | | μg/m³ | μg/m ³ | μg/m ³ | ppbV | ppbV | ppbV | Qualifier |
| 75-71-8 | Dichlorodifluoromethane (CFC 12) | ND | 82 | 8.2 | ND | 17 | 1.7 | |
| 74-87-3 | Chloromethane | ND | 16 | 8.2 | ND | 7.9 | 4.0 | |
| 76-14-2 | 1,2-Dichloro-1,1,2,2- tetrafluoroethane (CFC 114) | ND | 82 | 8.2 | ND | 12 | 1.2 | |
| 75-01-4 | Vinyl Chloride | ND | 16 | 8.2 | ND | 6.4 | 3.2 | |
| 74-83-9 | Bromomethane | ND | 16 | 8.2 | ND | 4.2 | 2.1 | |
| 75-00-3 | Chloroethane | ND | 16 | 8.2 | ND | 6.2 | 3.1 | |
| 64-17-5 | Ethanol | 12 | 820 | 8.2 | 6.6 | 440 | 4.4 | \mathbf{J} |
| 67-64-1 | Acetone | 150 | 820 | 12 | 64 | 350 | 5.0 | J, B |
| 75-69-4 | Trichlorofluoromethane | 1,500 | 16 | 8.2 | 280 | 2.9 | 1.5 | |
| 107-13-1 | Acrylonitrile | ND | 82 | 11 | ND | 38 | 5.3 | |
| 75-35-4 | 1,1-Dichloroethene | ND | 16 | 8.2 | ND | 4.1 | 2.1 | |
| 75-65-0 | 2-Methyl-2-Propanol (tert-Butyl Alcohol) | ND | 82 | 12 | ND | 27 | 4.0 | |
| 75-09-2 | Methylene Chloride | ND | 82 | 8.2 | ND | 24 | 2.4 | |
| 107-05-1 | 3-Chloro-1-propene (Allyl Chloride) | ND | 16 | 8.2 | ND | 5.2 | 2.6 | |
| 76-13-1 | Trichlorotrifluoroethane | ND | 16 | 9.2 | ND | 2.1 | 1.2 | |
| 75-15-0 | Carbon Disulfide | ND | 82 | 20 | ND | 26 | 6.3 | energy and a second s |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 16 | 8.2 | ND | 4.1 | 2.1 | |
| 75-34-3 | 1,1-Dichloroethane | ND | 16 | 8.2 | ND | 4.1 | 2.0 | |
| 1634-04-4 | Methyl tert-Butyl Ether | ND | 16 | .8.2 | ND | 4.6 | 2.3 | |
| 108-05-4 | Vinyl Acetate | ND | 820 | 26 | ND | 230 | 7.5 | |
| 78-93-3 | 2-Butanone (MEK) | 26 | 82 | 8.2 | 8.7 | 28 | 2.8 | J |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 16 | 8.2 | ND | 4.1 | 2.1 | |
| 108-20-3 | Diisopropyl Ether | ND | 82 | 9.7 | ND | 20 | 2.3 | |
| 67-66-3 | Chloroform | 52,000 | 16 | 9.7 | 11,000 | 3.4 | 2.0 | |

JD = Compound was analyzed for, but not detected above the laboratory detection limit.

ARL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

= The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

3 = Analyte was found in the method blank.

Verified By: Ru- Date: 5 9 08

7

RESULTS OF ANALYSIS

Page 2 of 3

Client:ENSRClient Sample ID:SG83B-05-1Client Project ID:Phase B Soil Gas / 04020-023-4311

Fest Code:EPA TO-15Instrument ID:Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13Analyst:Rusty BravoSampling Media:6.0 L Summa CanisterFest Notes:SC00979

CAS Project ID: P0801342 CAS Sample ID: P0801342-001

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

3.5

Initial Pressure (psig):

-3.6 Final Pressure (psig):

Canister Dilution Factor: 1.64

| CAS # | Compound | Result µg/m³ | MRL µg/m³ | MDL µg/m³ | Result ppbV | MRL ppbV | MDL ppbV | Data Qualifier |
|------------|---------------------------|-----------------|--------------|--------------|----------------|-------------|-------------|-------------------|
| 637-92-3 | Ethyl tert-Butyl Ether | ND | 82 | 8.4 | ND | 20 | 2.0 | |
| 107-06-2 | 1,2-Dichloroethane | ND | 16 | 8.2 | ND | 4.1 | 2.0 | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 16 | 8.2 | ND | 3.0 | 1.5 | |
| 71-43-2 | Benzene | 100 | 16 | 8.2 | 31 | 5.1 | 2.6 | |
| 56-23-5 | Carbon Tetrachloride | 12,000 | 16 | 8.2 | 1,900 | 2.6 | 1.3 | |
| 994-05-8 | tert-Amyl Methyl Ether | ND | 82 | 8.2 | ND | 20 | 2.0 | |
| 78-87-5 | 1,2-Dichloropropane | ND | 16 | 8.2 | ND | 3.5 | 1.8 | |
| 75-27-4 | Bromodichloromethane | ND | 16 | 8.2 | ND | 2.4 | 1.2 | |
| 79-01-6 | Trichloroethene | 16 | 16 | 8.2 | 3.0 | 3.1 | 1.5 | J |
| 123-91-1 | 1,4-Dioxane | ND | 82 | 10 | ND | 23 | 2.8 | |
| 80-62-6 | Methyl Methacrylate | ND | 82 | 12 | ND | 20 | 3.0 | |
| 142-82-5 | n-Heptane | ND | 82 | 10 | ND | 20 | 2.6 | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 82 | 8.5 | ND | 18 | 1.9 | |
| 108-10-1 | 4-Methyl-2-pentanone | ND | 82 | 9.2 | ND | 20 | 2.2 | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 82 | 10 | ND | 18 | 2.3 | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 16 | 8.2 | ND | 3.0 | 1.5 | |
| 108-88-3 | Toluene | 13 | 82 | 8.2 | 3.6 | 22 | 2.2 | J |
| 591-78-6 | 2-Hexanone | ND | 82 | 12 | ND | 20 | 3.0 | |
| 124-48-1 | Dibromochloromethane | ND | 16 | 11 | ND | 1.9 | 1.3 | |
| 106-93-4 | 1,2-Dibromoethane | ND | 16 | 8.9 | ND | 2.1 | 1.2 | |
| 111-65-9 | n-Octane | ND | 82 | 8.2 | ND | 18 | 1.8 | |
| 127-18-4 | Tetrachloroethene | 110 | 16 | 8.2 | 16 | 2.4 | 1.2 | |
| 108-90-7 | Chlorobenzene | 120 | 16 | 8.4 | 26 | 3.6 | 1.8 | |

VD = Compound was analyzed for, but not detected above the laboratory detection limit.

ARL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

= The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

RESULTS OF ANALYSIS

Page 3 of 3

ENSR Client: Client Sample ID: SG83B-05-1 Client Project ID: Phase B Soil Gas / 04020-023-4311

EPA TO-15 Test Code: Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 Instrument ID: Analyst: Rusty Bravo 6.0 L Summa Canister Sampling Media: Test Notes:

SC00979 Container ID:

> Initial Pressure (psig): Final Pressure (psig): 3.5 -3.6

> > Canister Dilution Factor: 1.64

0.010 Liter(s)

0.0010 Liter(s)

| | | Result | MRL | MDL | Result | MRL | MDL | Data |
|-------------|-------------------------------|-------------|-------|-------|--------|------|------|-------------------------|
| CAS # | Compound | $\mu g/m^3$ | μg/m³ | μg/m³ | ppbV | ppbV | ppbV | Qualifier |
| 100-41-4 | Ethylbenzene | ND | 82 | 10 | ND | 19 | 2.3 | |
| 179601-23-1 | m,p-Xylenes | ND | 82 | 21 | ND | 19 | 4.9 | |
| 75-25-2 | Bromoform | ND | 82 | 12 | ND | 7.9 | 1.2 | |
| 100-42-5 | Styrene | ND | 82 | 12 | ND | 19 | 2.9 | |
| 95-47-6 | o-Xylene | ND | 82 | 10 | ND | 19 | 2.4 | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 16 | 10 | ND | 2.4 | 1.5 | |
| 98-82-8 | Cumene | 9.7 | 82 | 9.2 | 2.0 | 17 | 1.9 | J, B |
| 103-65-1 | n-Propylbenzene | ND | 82 | 8.5 | ND | 17 | 1.7 | |
| 622-96-8 | 4-Ethyltoluene | 12 | 82 | 9.3 | 2.4 | 17 | 1.9 | J |
| 108-67-8 | 1,3,5-Trimethylbenzene | ND | 82 | 9.8 | ND | 17 | 2.0 | |
| 98-83-9 | alpha-Methylstyrene | ND | 82 | 12 | ND | 17 | 2.5 | |
| 95-63-6 | 1,2,4-Trimethylbenzene | ND | 82 | 11 | ND | 17 | 2.3 | |
| 100-44-7 | Benzyl Chloride | ND | 16 | 14 | ND | 3.2 | 2.7 | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | -16 | 10 | ND | 2.7 | 1.7 | |
| 106-46-7 | 1,4-Dichlorobenzene | 17 | .16 | 9.2 | 2.9 | 2.7 | 1.5 | |
| 135-98-8 | sec-Butylbenzene | ND | 82 | 9.5 | ND | 15 | 1.7 | |
| 99-87-6 | 4-Isopropyltoluene (p-Cymene) | ND | 82 | 11 | ND | 15 | 1.9 | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 16 | 11 | ND | 2.7 | 1.8 | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 82 | 12 | ND | 8.5 | 1.3 | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 16 | 12 | ND | 2.2 | 1.7 | |
| 91-20-3 | Naphthalene | 28 | 33 | 12 | 5.3 | 6.3 | 2.3 | J,B |
| 87-68-3 | Hexachlorobutadiene | ND | 16 | 15 | ND | 1.5 | 1.4 | |
| 98-06-6 | tert-Butylbenzene | ND | 33 | 8.2 | ND | 6.0 | 1.5 | |
| 104-51-8 | n-Butylbenzene | ND | 33 | 8.2 | ND | 6.0 | 1.5 | |

VD = Compound was analyzed for, but not detected above the laboratory detection limit.

VIRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method. = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

Verified By: RC-

3 = Analyte was found in the method blank.

P0801342 TO15 0805091251 SS.xls - Sample

CAS Project ID: P0801342

Date Collected: 5/7/08

Date Received: 5/8/08

Date Analyzed: 5/8/08

Volume(s) Analyzed:

CAS Sample ID: P0801342-001

Date: 5 9 00 000 - PageNo.:

RESULTS OF ANALYSIS

Page 1 of 3

Client: ENSR Client Sample ID: SG83B-05-3 Client Project ID: Phase B Soil Gas / 04020-023-4311

EPA TO-15 Test Code: Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 Rusty Bravo 6.0 L Summa Canister SC00564

CAS Project ID: P0801342 CAS Sample ID: P0801342-002

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

Instrument ID: Analyst: Sampling Media: Test Notes: Container ID:

Initial Pressure (psig):

-3.8

Final Pressure (psig): 3.5

Canister Dilution Factor: 1.67

| CAS # | Compound | Result | MRL | MDL | Result | MRL | MDL | Data |
|-----------|--|-------------------|-------------------------|-------------|--------|------|------|-----------|
| | | μg/m ³ | <u>μg/m³</u> | $\mu g/m^3$ | ppbV | ppbV | ppbV | Qualifier |
| 75-71-8 | Dichlorodifluoromethane (CFC 12) | ND | 84 | 8.4 | ND | 17 | 1.7 | |
| 74-87-3 | Chloromethane | ND | 17 | 8.4 | ND | 8.1 | 4.0 | |
| 76-14-2 | 1,2-Dichloro-1,1,2,2- tetrafluoroethane (CFC 114) | ND | 84 | 8.4 | ND | 12 | 1.2 | |
| 75-01-4 | Vinyl Chloride | ND | 17 | 8.4 | ND | 6.5 | 3.3 | |
| 74-83-9 | Bromomethane | ND | 17 | 8.4 | ND | 4.3 | 2.2 | |
| 75-00-3 | Chloroethane | ND | 17 | 8.4 | ND | 6.3 | 3.2 | |
| 64-17-5 | Ethanol | ND | 840 | 8.4 | ND | 440 | 4.4 | |
| 67-64-1 | Acetone | 130 | 840 | 12 | 55 | 350 | 5.1 | J, B |
| 75-69-4 | Trichlorofluoromethane | 1,500 | 17 | 8.4 | 270 | 3.0 | 1.5 | |
| 107-13-1 | Acrylonitrile | ND | 84 | 12 | ND | 38 | 5.4 | |
| 75-35-4 | 1,1-Dichloroethene | ND | 17 | 8.4 | ND | 4.2 | 2.1 | |
| 75-65-0 | 2-Methyl-2-Propanol (tert-Butyl Alcohol) | ND | 84 | 12 | ND | 28 | 4.1 | |
| 75-09-2 | Methylene Chloride | ND | 84 | 8.4 | ND | 24 | 2.4 | |
| 107-05-1 | 3-Chloro-1-propene (Allyl Chloride) | ND | 17 | 8.4 | ND | 5.3 | 2.7 | |
| 76-13-1 | Trichlorotrifluoroethane | ND | 17 | 9.4 | ND | 2.2 | 1.2 | |
| 75-15-0 | Carbon Disulfide | ND | 84 | 20 | ND | 27 | 6.4 | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 17 | 8.4 | ND | 4.2 | 2.1 | |
| 75-34-3 | 1,1-Dichloroethane | ND | 17 | 8.4 | ND | 4.1 | 2.1 | |
| 1634-04-4 | Methyl tert-Butyl Ether | ND | 17 | 8.4 | ND | 4.6 | 2.3 | |
| 108-05-4 | Vinyl Acetate | ND | 840 | 27 | ND | 240 | 7.6 | |
| 78-93-3 | 2-Butanone (MEK) | 24 | 84 | 8.4 | 8.1 | 28 | 2.8 | J |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 17 | 8.4 | ND | 4.2 | 2.1 | |
| 108-20-3 | Diisopropyl Ether | ND | 84 | 9.9 | ND | 20 | 2.4 | |
| 67-66-3 | Chloroform | 49,000 | 17 | 9.9 | 10,000 | 3.4 | 2.0 | |

ID = Compound was analyzed for, but not detected above the laboratory detection limit.

IRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

= The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

= Analyte was found in the method blank.

Verified By: Re-

RESULTS OF ANALYSIS

Page 2 of 3

Client:ENSRClient Sample ID:SG83B-05-3Client Project ID:Phase B Soil Gas / 04020-023-4311

CAS Project ID: P0801342 CAS Sample ID: P0801342-002

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

Test Code: Instrument ID: Analyst: Sampling Media: Test Notes: Container ID:

SC00564

EPA TO-15

Rusty Bravo

6.0 L Summa Canister

Initial Pressure (psig):

-3.8

Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13

Final Pressure (psig): 3.5

Canister Dilution Factor: 1.67

| CAS # | Compound | Result μg/m³ | MRL µg/m³ | MDL µg/m³ | Result ppbV | MRL ppbV | MDL ppbV | Data Qualifier |
|------------|---------------------------|-----------------|--------------|--------------|----------------|-------------|-------------|-------------------|
| 637-92-3 | Ethyl tert-Butyl Ether | ND | 84 | 8.5 | ND | 20 | 2.0 | |
| 107-06-2 | 1,2-Dichloroethane | ND | 17 | 8.4 | ND | 4.1 | 2.1 | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 17 | 8.4 | ND | 3.1 | 1.5 | |
| 71-43-2 | Benzene | 100 | 17 | 8.4 | 32 | 5.2 | 2.6 | |
| 56-23-5 | Carbon Tetrachloride | 12,000 | 17 | 8.4 | 1,900 | 2.7 | 1.3 | |
| 994-05-8 | tert-Amyl Methyl Ether | ND | 84 | 8.4 | ND | 20 | 2.0 | |
| 78-87-5 | 1,2-Dichloropropane | ND | 17 | 8.4 | ND | 3.6 | 1.8 | |
| 75-27-4 | Bromodichloromethane | ND | 17 | 8.4 | ND | 2.5 | 1.2 | |
| 79-01-6 | Trichloroethene | 16 | 17 | 8.4 | 3.0 | 3.1 | 1.6 | J |
| 123-91-1 | 1,4-Dioxane | ND | 84 | 10 | ND | 23 | 2.8 | |
| 80-62-6 | Methyl Methacrylate | ND | 84 | 13 | ND | 20 | 3.1 | |
| 142-82-5 | n-Heptane | ND | 84 | 11 | ND | 20 | 2.6 | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 84 | 8.7 | ND | 18 | 1.9 | |
| 108-10-1 | 4-Methyl-2-pentanone | ND | 84 | 9.4 | ND | 20 | 2.3 | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 84 | 11 | ND | 18 | 2.3 | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 17 | 8.4 | ND | 3.1 | 1.5 | |
| 108-88-3 | Toluene | 8.5 | 84 | 8.4 | 2.3 | 22 | 2.2 | \mathbf{J} |
| 591-78-6 | 2-Hexanone | ND | 84 | 13 | ND | 20 | 3.1 | |
| 124-48-1 | Dibromochloromethane | ND | 17 | 11 | ND | 2.0 | 1.3 | |
| 106-93-4 | 1,2-Dibromoethane | ND | 17 | 9.0 | ND | 2.2 | 1.2 | |
| 111-65-9 | n-Octane | ND | 84 | 8.4 | ND | 18 | 1.8 | |
| 127-18-4 | Tetrachloroethene | 130 | 17 | 8.4 | 19 | 2.5 | 1.2 | |
| 108-90-7 | Chlorobenzene | 160 | 17 | 8.5 | 36 | 3.6 | 1.9 | |

JD = Compound was analyzed for, but not detected above the laboratory detection limit.

ARL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

= The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

Verified By:

Date: 519(08 TO15SCAN.XLT - Tronox - Henderson - PageNo.:

RESULTS OF ANALYSIS

Page 3 of 3

-3.8

Client:ENSRClient Sample ID:SG83B-05-3Client Project ID:Phase B Soil Gas / 04020-023-4311

CAS Project ID: P0801342 CAS Sample ID: P0801342-002

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

3.5

Initial Pressure (psig):

Final Pressure (psig):

Canister Dilution Factor: 1.67

| | | Result | MRL | MDL | Result | MRL | MDL | Data |
|-------------|-------------------------------|-------------------|-------|-------------|--------|------|------|-----------|
| CAS # | Compound | μg/m ³ | μg/m³ | $\mu g/m^3$ | ppbV | ppbV | ppbV | Qualifier |
| 100-41-4 | Ethylbenzene | ND | 84 | 10 | ND | 19 | 2.4 | |
| 179601-23-1 | m,p-Xylenes | ND | 84 | 22 | ND | 19 | 5.0 | |
| 75-25-2 | Bromoform | ND | 84 | 13 | ND | 8.1 | 1.2 | |
| 100-42-5 | Styrene | ND | 84 | 13 | ND | 20 | 3.0 | |
| 95-47-6 | o-Xylene | ND | 84 | 11 | ND | 19 | 2.4 | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 17 | 11 | ND | 2.4 | 1.6 | |
| 98-82-8 | Cumene | ND | 84 | 9.4 | ND | 17 | 1.9 | |
| 103-65-1 | n-Propylbenzene | ND | 84 | 8.7 | ND | 17 | 1.8 | |
| 622-96-8 | 4-Ethyltoluene | ND | 84 | 9.5 | ND | 17 | 1.9 | |
| 108-67-8 | 1,3,5-Trimethylbenzene | ND | 84 | 10 | ND | 17 | 2.0 | |
| 98-83-9 | alpha-Methylstyrene | ND | 84 | 12 | ND | 17 | 2.5 | |
| 95-63-6 | 1,2,4-Trimethylbenzene | ND | 84 | 12 | ND | 17 | 2.3 | |
| 100-44-7 | Benzyl Chloride | ND | 17 | 14 | ND | 3.2 | 2.8 | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 17 | 10 | ND | 2.8 | 1.7 | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 17 | 9.4 | ND | 2.8 | 1.6 | |
| 135-98-8 | sec-Butylbenzene | ND | 84 | 9.7 | ND | 15 | 1.8 | |
| 99-87-6 | 4-Isopropyltoluene (p-Cymene) | ND | 84 | 11 | ND | 15 | 2.0 | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 17 | 11 | ND | 2.8 | 1.8 | |
| 96-12-8 | 1.2-Dibromo-3-chloropropane | ND | 84 | 13 | ND | 8.6 | 1.3 | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 17 | 13 | ND | 2.3 | 1.7 | |
| 91-20-3 | Naphthalene | ND | 33 | 12 | ND | 6.4 | 2.4 | |
| 87-68-3 | Hexachlorobutadiene | ND | 17 | 15 | ND | 1.6 | 1.4 | |
| 98-06-6 | tert-Butylbenzene | ND | 33 | 8.4 | ND | 6.1 | 1.5 | |
| 104-51-8 | n-Butylbenzene | ND | 33 | 8.4 | ND | 6.1 | 1.5 | |

VD = Compound was analyzed for, but not detected above the laboratory detection limit.

ARL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

12

RESULTS OF ANALYSIS

Page 1 of 4

Client:ENSRClient Sample ID:SG83B-05-7Client Project ID:Phase B Soil Gas / 04020-023-4311

Test Code:EPA TO-15Instrument ID:Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13Analyst:Rusty BravoSampling Media:6.0 L Summa CanisterTest Notes:Sc00791

CAS Project ID: P0801342 CAS Sample ID: P0801342-003

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

Initial Pressure (psig):

(psig): -4.2 Fina

Final Pressure (psig): 3.5

Canister Dilution Factor: 1.73

| CAS # | Compound | Result μg/m³ | MRL μg/m³ | MDL µg/m³ | Result ppbV | MRL ppbV | MDL ppbV | Data Qualifier |
|-----------|--|-----------------|--------------|--------------|----------------|-------------|-------------|-------------------|
| 75-71-8 | Dichlorodifluoromethane (CFC 12) | ND | 87 | 8.7 | ND | 18 | 1.8 | |
| 74-87-3 | Chloromethane | ND | 17 | 8.7 | ND | 8.4 | 4.2 | |
| 76-14-2 | 1,2-Dichloro-1,1,2,2- tetrafluoroethane (CFC 114) | ND | 87 | 8.7 | ND | 12 | 1.2 | |
| 75-01-4 | Vinyl Chloride | ND | 17 | 8.7 | ND | 6.8 | 3.4 | |
| 74-83-9 | Bromomethane | ND | 17 | 8.7 | ND | 4.5 | 2.2 | |
| 75-00-3 | Chloroethane | ND | 17 | 8.7 | ND | 6.6 | 3.3 | |
| 64-17-5 | Ethanol | ND | 870 | 8.7 | ND | 460 | 4.6 | |
| 67-64-1 | Acetone | 110 | 870 | 13 | 48 | 360 | 5.3 | J, B |
| 75-69-4 | Trichlorofluoromethane | 1,500 | 17 | 8.7 | 270 | 3.1 | 1.5 | |
| 107-13-1 | Acrylonitrile | ND | 87 | 12 | ND | 40 | 5.6 | |
| 75-35-4 | 1,1-Dichloroethene | ND | 17 | 8.7 | ND | 4.4 | 2.2 | |
| 75-65-0 | 2-Methyl-2-Propanol (tert-Butyl Alcohol) | ND | 87 | 13 | ND | 29 | 4.2 | |
| 75-09-2 | Methylene Chloride | 9.3 | 87 | 8.7 | 2.7 | 25 | 2.5 | J |
| 107-05-1 | 3-Chloro-1-propene (Allyl Chloride) | ND | 17 | 8.7 | ND | 5.5 | 2.8 | |
| 76-13-1 | Trichlorotrifluoroethane | ND | 17 | 9.7 | ND | 2.3 | 1.3 | |
| 75-15-0 | Carbon Disulfide | ND | 87 | 21 | ND | 28 | 6.7 | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 17 | 8.7 | ND | 4.4 | 2.2 | |
| 75-34-3 | 1,1-Dichloroethane | ND | 17 | 8.7 | ND | 4.3 | 2.1 | |
| 1634-04-4 | Methyl tert-Butyl Ether | ND | 17 | 8.7 | ND | 4.8 | 2.4 | |
| 108-05-4 | Vinyl Acetate | ND | 870 | 28 | ND | 250 | 7.9 | |
| 78-93-3 | 2-Butanone (MEK) | 23 | 87 | 8.7 | 7.8 | 29 | 2.9 | J |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 17 | 8.7 | ND | 4.4 | 2.2 | |
| 108-20-3 | Diisopropyl Ether | ND | 87 | 10 | ND | 21 | 2.4 | |
| 67-66-3 | Chloroform | 54,000 | 17 | 10 | 11,000 | 3.5 | 2.1 | |

VD = Compound was analyzed for, but not detected above the laboratory detection limit.

ARL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

= The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

3 = Analyte was found in the method blank.

Verified By: <u>Rc-</u>

RESULTS OF ANALYSIS

Page 2 of 4

Client:ENSRClient Sample ID:SG83B-05-7Client Project ID:Phase B Soil Gas / 04020-023-4311

6.0 L Summa Canister

CAS Project ID: P0801342 CAS Sample ID: P0801342-003

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

Fest Code: Instrument ID: Analyst: Sampling Media: Fest Notes: Container ID:

SC00791

EPA TO-15

Rusty Bravo

Initial Pressure (psig): -4.2

Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13

Final Pressure (psig): 3.5

Canister Dilution Factor: 1.73

| CAS # | Compound | Result μg/m³ | MRL µg/m³ | MDL µg/m³ | Result ppbV | MRL ppbV | MDL ppbV | Data Qualifier |
|------------|---------------------------|-----------------|--------------|--------------|----------------|-------------|-------------|-------------------|
| 637-92-3 | Ethyl tert-Butyl Ether | ND | 87 | 8.8 | ND | 21 | 2.1 | |
| 107-06-2 | 1,2-Dichloroethane | ND | 17 | 8.7 | ND | 4.3 | 2.1 | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 17 | 8.7 | ND | 3.2 | 1.6 | |
| 71-43-2 | Benzene | 100 | 17 | 8.7 | 32 | 5.4 | 2.7 | |
| 56-23-5 | Carbon Tetrachloride | 12,000 | 17 | 8.7 | 1,900 | 2.8 | 1.4 | |
| 994-05-8 | tert-Amyl Methyl Ether | ND | 87 | 8.7 | ND | 21 | 2.1 | |
| 78-87-5 | 1,2-Dichloropropane | ND | 17 | 8.7 | ND | 3.7 | 1.9 | |
| 75-27-4 | Bromodichloromethane | ND | 17 | 8.7 | ND | 2.6 | 1.3 | |
| 79-01-6 | Trichloroethene | 11 | 17 | 8.7 | 2.1 | 3.2 | 1.6 | J |
| 123-91-1 | 1,4-Dioxane | ND | 87 | 11 | ND | 24 | 2.9 | |
| 80-62-6 | Methyl Methacrylate | ND | 87 | 13 | ND | 21 | 3.2 | |
| 142-82-5 | n-Heptane | ND | 87 | 11 | ND | 21 | 2.7 | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 87 | 9.0 | ND | 19 | 2.0 | |
| 108-10-1 | 4-Methyl-2-pentanone | ND | 87 | 9.7 | ND | 21 | 2.4 | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 87 | 11 | ND | 19 | 2.4 | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 17 | 8.7 | ND | 3.2 | 1.6 | |
| 108-88-3 | Toluene | ND | 87 | 8.7 | ND | 23 | 2.3 | |
| 591-78-6 | 2-Hexanone | ND | 87 | 13 | ND | 21 | 3.2 | |
| 124-48-1 | Dibromochloromethane | ND | 17 | 12 | ND | 2.0 | 1.4 | |
| 106-93-4 | 1,2-Dibromoethane | ND | 17 | 9.3 | ND | 2.3 | 1.2 | |
| 111-65-9 | n-Octane | ND | 87 | 8.7 | ND | 19 | 1.9 | |
| 127-18-4 | Tetrachloroethene | 130 | 17 | 8.7 | 19 | 2.6 | 1.3 | |
| 108-90-7 | Chlorobenzene | 180 | 17 | 8.8 | 39 | 3.8 | 1.9 | |

JD = Compound was analyzed for, but not detected above the laboratory detection limit.

*A*RL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

= The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

Verified By: R_C_

14

Date: 5/4/58 TO15SCAN.XLT - Tronox - Henderson - PageNo.:

RESULTS OF ANALYSIS

Page 3 of 4

Client:ENSRClient Sample ID:SG83B-05-7Client Project ID:Phase B Soil Gas / 04020-023-4311Fest Code:EPA TO-15

6.0 L Summa Canister

Rusty Bravo

SC00791

nstrument ID:

Sampling Media:

Analyst:

[est Notes:

Container ID:

CAS Project ID: P0801342 CAS Sample ID: P0801342-003

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

Initial Pressure (psig):

Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13

-4.2 Final Pressure (psig): 3.5

Canister Dilution Factor: 1.73

| | | Result | MRL | MDL | Result | MRL | MDL Data |
|-------------|-------------------------------|--------|-------|-------|--------|------|----------------|
| CAS # | Compound | μg/m³ | μg/m³ | μg/m³ | ppbV | ppbV | ppbV Qualifier |
| 100-41-4 | Ethylbenzene | ND | 87 | 11 | ND | 20 | 2.5 |
| 179601-23-1 | m,p-Xylenes | ND | 87 | 22 | ND | 20 | 5.2 |
| 75-25-2 | Bromoform | ND | 87 | 13 | ND | 8.4 | 1.3 |
| 100-42-5 | Styrene | ND | 87 | 13 | ND | 20 | 3.1 |
| 95-47-6 | o-Xylene | ND | 87 | 11 | ND | 20 | 2.5 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 17 | 11 | ND | 2.5 | 1.6 |
| 98-82-8 | Cumene | ND | 87 | 9.7 | ND | 18 | 2.0 |
| 103-65-1 | n-Propylbenzene | ND | 87 | 9.0 | ND | 18 | 1.8 |
| 622-96-8 | 4-Ethyltoluene | ND | 87 | 9.9 | ND | 18 | 2.0 |
| 108-67-8 | 1,3,5-Trimethylbenzene | ND | 87 | 10 | ND | 18 | 2.1 |
| 98-83-9 | alpha-Methylstyrene | ND | 87 | 13 | ND | 18 | 2.6 |
| 95-63-6 | 1,2,4-Trimethylbenzene | ND | 87 | 12 | ND | 18 | 2.4 |
| 100-44-7 | Benzyl Chloride | ND | 17 | 15 | ND | 3.3 | 2.9 |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 17 | 11 | ND | 2.9 | 1.8 |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 17 | 9.7 | ND | 2.9 | 1.6 |
| 135-98-8 | sec-Butylbenzene | ND | 87 | - 10 | ND | 16 | 1.8 |
| 99-87-6 | 4-Isopropyltoluene (p-Cymene) | ND | 87 | 11 | ND | 16 | 2.0 |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 17 | 11 | ND | 2.9 | 1.9 |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 87 | 13 | ND | 9.0 | 1.4 |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 17 | 13 | ND | 2.3 | 1.8 |
| 91-20-3 | Naphthalene | ND | 35 | 13 | ND | 6.6 | 2.4 |
| 87-68-3 | Hexachlorobutadiene | ND | 17 | 16 | ND | 1.6 | 1.5 |
| 98-06-6 | tert-Butylbenzene | ND | 35 | 8.7 | ND | 6.3 | 1.6 |
| 104-51-8 | n-Butylbenzene | ND | 35 | 8.7 | ND | 6.3 | 1.6 |

JD = Compound was analyzed for, but not detected above the laboratory detection limit.

ARL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

Verified By: Rer

RESULTS OF ANALYSIS Page 4 of 4

Client: ENSR CAS Project ID: P0801342 Client Sample ID: SG83B-05-7 CAS Sample ID: P0801342-003 Client Project ID: Phase B Soil Gas / 04020-023-4311 **Tentatively Identified Compounds** Date Collected: 5/7/08 EPA TO-15 Test Code: Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 Date Received: 5/8/08 instrument ID: Date Analyzed: 5/8/08 Rusty Bravo Analyst: 6.0 L Summa Canister 0.010 Liter(s) Sampling Media: Volume(s) Analyzed: 0.0010 Liter(s) **Fest Notes:** Container ID: SC00791 Initial Pressure (psig): -4.2 Final Pressure (psig): 3.5 Canister Dilution Factor: 1.73

| GC/MS | Compound Identification | Concentration | Data |
|----------------|-------------------------|---------------|-----------|
| Retention Time | | $\mu g/m^3$ | Qualifier |
| | No Compounds Detected | | |

P0801342_TO15_0805091251_SS.xls - TIC (3)

RESULTS OF ANALYSIS

Page 1 of 4

Client:ENSRClient Sample ID:Method BlankClient Project ID:Phase B Soil Gas / 04020-023-4311

EPA TO-15

Rusty Bravo

6.0 L Summa Canister

Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13

Test Code:

Analyst:

[nstrument ID:

Sampling Media: Test Notes:

CAS Project ID: P0801342 CAS Sample ID: P080508-MB

Date Collected: NA Date Received: NA Date Analyzed: 5/8/08 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

| CAS # | Compound | Result µg/m³ | MRL µg/m³ | MDL µg/m³ | Result ppbV | MRL ppbV | MDL ppbV | Data Qualifier |
|-----------|--|-----------------|--------------|--------------|----------------|-------------|-------------|-------------------------------------|
| 75-71-8 | Dichlorodifluoromethane (CFC 12) | ND | 0.50 | 0.050 | ND | 0.10 | 0.010 | |
| 74-87-3 | Chloromethane | ND | 0.10 | 0.050 | ND | 0.048 | 0.024 | |
| 76-14-2 | 1,2-Dichloro-1,1,2,2- tetrafluoroethane (CFC 114) | ND | 0.50 | 0.050 | ND | 0.072 | 0.0072 | |
| 75-01-4 | Vinyl Chloride | ND | 0.10 | 0.050 | ND | 0.039 | 0.020 | |
| 74-83-9 | Bromomethane | ND | 0.10 | 0.050 | ND | 0.026 | 0.013 | |
| 75-00-3 | Chloroethane | ND | 0.10 | 0.050 | ND | 0.038 | 0.019 | |
| 64-17-5 | Ethanol | ND | 5.0 | 0.050 | ND | 2.7 | 0.027 | |
| 67-64-1 | Acetone | 0.42 | 5.0 | 0.073 | 0.18 | 2.1 | 0.031 | J |
| 75-69-4 | Trichlorofluoromethane | ND | 0.10 | 0.050 | ND | 0.018 | 0.0089 | |
| 107-13-1 | Acrylonitrile | ND | 0.50 | 0.070 | ND | 0.23 | 0.032 | |
| 75-35-4 | 1,1-Dichloroethene | ND | 0.10 | 0.050 | ND | 0.025 | 0.013 | |
| 75-65-0 | 2-Methyl-2-Propanol (tert-Butyl Alcohol) | ND | 0.50 | 0.074 | ND | 0.17 | 0.024 | |
| 75-09-2 | Methylene Chloride | ND | 0.50 | 0.050 | ND | 0.14 | 0.014 | |
| 107-05-1 | 3-Chloro-1-propene (Allyl Chloride) | ND | 0.10 | 0.050 | ND | 0.032 | 0.016 | |
| 76-13-1 | Trichlorotrifluoroethane | ND | 0.10 | 0.056 | ND | 0.013 | 0.0073 | |
| 75-15-0 | Carbon Disulfide | ND | 0.50 | 0.12 | ND | 0.16 | 0.039 | n ala se anna a se a bR alanta e Re |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 0.10 | 0.050 | ND | 0.025 | 0.013 | |
| 75-34-3 | 1,1-Dichloroethane | ND | 0.10 | 0.050 | ND | 0.025 | 0.012 | |
| 1634-04-4 | Methyl tert-Butyl Ether | ND | 0.10 | 0.050 | ND | 0.028 | 0.014 | |
| 108-05-4 | Vinyl Acetate | ND | 5.0 | 0.16 | ND | 1.4 | 0.045 | |
| 78-93-3 | 2-Butanone (MEK) | ND | 0.50 | 0.050 | ND | 0.17 | 0.017 | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 0.10 | 0.050 | ND | 0.025 | 0.013 | |
| 108-20-3 | Diisopropyl Ether | ND | 0.50 | 0.059 | ND | 0.12 | 0.014 | |
| 67-66-3 | Chloroform | ND | 0.10 | 0.059 | ND | 0.020 | 0.012 | |

JD = Compound was analyzed for, but not detected above the laboratory detection limit.

ARL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.
 The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

Verified By: Re-

17

RESULTS OF ANALYSIS

Page 2 of 4

| Client: | ENSR |
|---------------------------|-----------------------------------|
| Client Sample ID: | Method Blank |
| Client Project ID: | Phase B Soil Gas / 04020-023-4311 |

EPA TO-15

Rusty Bravo

6.0 L Summa Canister

Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13

Fest Code:

Analyst:

[est Notes:

nstrument ID:

Sampling Media:

CAS Project ID: P0801342 CAS Sample ID: P080508-MB

Date Collected: NA Date Received: NA Date Analyzed: 5/8/08 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

| Compound | Result | MRL | MDL | Result | MRL | MDL | Data |
|---------------------------|---|---|---|--|---|---|--|
| - | μg/m³ | µg∕m³ | $\mu g/m^3$ | ppbV | ppbV | ppbV | Qualifier |
| Ethyl tert-Butyl Ether | ND | 0.50 | 0.051 | ND | 0.12 | 0.012 | |
| 1,2-Dichloroethane | ND | 0.10 | 0.050 | ND | 0.025 | 0.012 | |
| 1,1,1-Trichloroethane | ND | 0.10 | 0.050 | ND | 0.018 | 0.0092 | |
| Benzene | ND | 0.10 | 0.050 | ND | 0.031 | 0.016 | |
| Carbon Tetrachloride | ND | 0.10 | 0.050 | ND | 0.016 | 0.0080 | |
| tert-Amyl Methyl Ether | ND | 0.50 | 0.050 | ND | 0.12 | 0.012 | |
| 1,2-Dichloropropane | ND | 0.10 | 0.050 | ND | 0.022 | 0.011 | |
| Bromodichloromethane | ND | 0.10 | 0.050 | ND | 0.015 | 0.0075 | |
| Trichloroethene | ND | 0.10 | 0.050 | ND | 0.019 | 0.0093 | |
| 1,4-Dioxane | ND | 0.50 | 0.061 | ND | 0.14 | 0.017 | |
| Methyl Methacrylate | ND | 0.50 | 0.075 | ND | 0.12 | 0.018 | |
| n-Heptane | ND | 0.50 | 0.064 | ND | 0.12 | 0.016 | |
| cis-1,3-Dichloropropene | ND | 0.50 | 0.052 | ND | 0.11 | 0.011 | |
| 4-Methyl-2-pentanone | ND | 0.50 | 0.056 | ND | 0.12 | 0.014 | |
| trans-1,3-Dichloropropene | ND | 0.50 | 0.063 | ND | 0.11 | 0.014 | |
| 1,1,2-Trichloroethane | ND | 0.10 | 0.050 | ND | 0.018 | 0.0092 | |
| Toluene | ND | 0.50 | 0.050 | ND | 0.13 | 0.013 | |
| 2-Hexanone | ND | 0.50 | 0.076 | ND | 0.12 | 0.019 | |
| Dibromochloromethane | ND | 0.10 | 0.068 | ND | 0.012 | 0.0080 | |
| 1,2-Dibromoethane | ND | 0.10 | 0.054 | ND | 0.013 | 0.0070 | |
| n-Octane | ND | 0.50 | 0.050 | ND | 0.11 | 0.011 | |
| Tetrachloroethene | ND | 0.10 | 0.050 | ND | 0.015 | 0.0074 | |
| Chlorobenzene | ND | 0.10 | 0.051 | ND | 0.022 | 0.011 | |
| | Ethyl tert-Butyl Ether 1,2-Dichloroethane 1,1,1-Trichloroethane Benzene Carbon Tetrachloride tert-Amyl Methyl Ether 1,2-Dichloropropane Bromodichloromethane Trichloroethene 1,4-Dioxane Methyl Methacrylate n-Heptane cis-1,3-Dichloropropene 4-Methyl-2-pentanone trans-1,3-Dichloropropene 1,1,2-Trichloroethane Toluene 2-Hexanone Dibromochloromethane 1,2-Dibromoethane n-Octane Tetrachloroethene | μg/m³Ethyl tert-Butyl EtherND1,2-DichloroethaneND1,1,1-TrichloroethaneNDBenzeneNDCarbon TetrachlorideNDtert-Amyl Methyl EtherND1,2-DichloropropaneNDBromodichloromethaneNDTrichloroetheneND1,4-DioxaneNDMethyl MethacrylateNDn-HeptaneNDcis-1,3-DichloropropeneND4-Methyl-2-pentanoneNDtrans-1,3-DichloropropeneND1,1,2-TrichloroethaneND2-HexanoneNDDibromochloromethaneND1,2-DibromoethaneND1,2-DibromoethaneND1,2-DibromoethaneND1,2-DibromoethaneND1,2-DibromoethaneND1,2-DibromoethaneNDTetrachloroetheneNDTetrachloroetheneND | $\mu g/m^3$ $\mu g/m^3$ Ethyl tert-Butyl EtherND0.501,2-DichloroethaneND0.101,1,1-TrichloroethaneND0.10BenzeneND0.10Carbon TetrachlorideND0.10tert-Amyl Methyl EtherND0.10BromodichloropropaneND0.10BromodichloromethaneND0.10TrichloroetheneND0.101,4-DioxaneND0.50Methyl MethacrylateND0.50n-HeptaneND0.50cis-1,3-DichloropropeneND0.501,1,2-TrichloroethaneND0.501,1,2-TrichloroethaneND0.501,1,2-TrichloroethaneND0.502-HexanoneND0.50DibromochloromethaneND0.101,2-DibromoethaneND0.101,2-DibromoethaneND0.101,2-DibromoethaneND0.101,2-DibromoethaneND0.101,2-DibromoethaneND0.101,2-DibromoethaneND0.101,2-DibromoethaneND0.101,2-DibromoethaneND0.101,2-DibromoethaneND0.101,2-DibromoethaneND0.101,2-DibromoethaneND0.101,2-DibromoethaneND0.101,2-DibromoethaneND0.101,2-DibromoethaneND0.101,2-DibromoethaneND0.10 | $\mu g/m^3$ $\mu g/m^3$ $\mu g/m^3$ $\mu g/m^3$ Ethyl tert-Butyl EtherND 0.50 0.051 1,2-DichloroethaneND 0.10 0.050 1,1,1-TrichloroethaneND 0.10 0.050 BenzeneND 0.10 0.050 Carbon TetrachlorideND 0.10 0.050 tert-Amyl Methyl EtherND 0.10 0.050 1,2-DichloropropaneND 0.10 0.050 BromodichloromethaneND 0.10 0.050 TrichloroetheneND 0.10 0.050 1,4-DioxaneND 0.50 0.061 Methyl MethacrylateND 0.50 0.075 n-HeptaneND 0.50 0.052 4-Methyl-2-pentanoneND 0.50 0.050 1,1,2-TrichloroethaneND 0.50 0.050 1,1,2-TrichloroethaneND 0.50 0.050 1,1,2-TrichloroethaneND 0.50 0.050 2-HexanoneND 0.50 0.076 DibromochloromethaneND 0.10 0.054 1,2-DibromoethaneND 0.10 0.054 1,2-DibromoethaneND 0.10 0.054 1,2-DibromoethaneND 0.10 0.054 1,2-DibromoethaneND 0.10 0.050 1,1,2-TrichloroethaneND 0.10 0.054 1,2-DibromoethaneND 0.10 0.054 1,2-DibromoethaneND 0.10 0.050 1,3-Dichl | $\mu g/m^3$ $\mu g/m^3$ $\mu g/m^3$ $\mu g/m^3$ $p p b V$ Ethyl tert-Butyl EtherND 0.50 0.051 ND $1,2$ -DichloroethaneND 0.10 0.050 ND $1,1$ -TrichloroethaneND 0.10 0.050 NDBenzeneND 0.10 0.050 NDCarbon TetrachlorideND 0.10 0.050 NDtert-Amyl Methyl EtherND 0.10 0.050 ND1,2-DichloropropaneND 0.10 0.050 NDBromodichloromethaneND 0.10 0.050 NDTrichloroetheneND 0.10 0.050 ND1,4-DioxaneND 0.50 0.061 NDMethyl MethacrylateND 0.50 0.052 NDn-HeptaneND 0.50 0.056 NDtrans-1,3-DichloropropeneND 0.50 0.056 ND1,1,2-TrichloroethaneND 0.50 0.050 ND1,1,2-TrichloroethaneND 0.50 0.050 ND2-HexanoneND 0.50 0.050 NDDibromochloromethaneND 0.10 0.054 ND1,2-DibromoethaneND 0.10 0.054 NDn-OctaneND 0.50 0.050 ND | $\mu g/m^3$ $\mu g/m^3$ $\mu g/m^3$ $p p b V$ $p p b V$ Ethyl tert-Butyl EtherND0.500.051ND0.121,2-DichloroethaneND0.100.050ND0.0251,1,1-TrichloroethaneND0.100.050ND0.018BenzeneND0.100.050ND0.031Carbon TetrachlorideND0.100.050ND0.016tert-Amyl Methyl EtherND0.100.050ND0.121,2-DichloropropaneND0.100.050ND0.022BromodichloromethaneND0.100.050ND0.015TrichloroetheneND0.100.050ND0.0191,4-DioxaneND0.500.061ND0.12n-HeptaneND0.500.064ND0.12cis-1,3-DichloropropeneND0.500.052ND0.114-Methyl-2-pentanoneND0.500.063ND0.111,1,2-TrichloroethaneND0.500.063ND0.132-HexanoneND0.500.050ND0.018TolueneND0.500.076ND0.121,2-DibromochloromethaneND0.500.056ND0.132-HexanoneND0.500.050ND0.018TolueneND0.500.056ND0.12DibromochloromethaneND0.100.054ND0.012 <t< td=""><td>$\mu g/m^3$$\mu g/m^3$$\mu g/m^3$$ppbV$$ppbV$$ppbV$Ethyl tert-Butyl EtherND0.500.051ND0.120.0121,2-DichloroethaneND0.100.050ND0.0250.0121,1,1-TrichloroethaneND0.100.050ND0.0180.0092BenzeneND0.100.050ND0.0160.0080tert-Amyl Methyl EtherND0.100.050ND0.0120.0121,2-DichloropropaneND0.100.050ND0.0220.011BromodichloromethaneND0.100.050ND0.0220.011BromodichloromethaneND0.100.050ND0.0150.0075TrichloroetheneND0.100.050ND0.0120.018n-HeptaneND0.500.061ND0.120.018n-HeptaneND0.500.056ND0.120.014trist-1,3-DichloropropeneND0.500.056ND0.120.014trast-1,3-DichloropropeneND0.500.056ND0.120.014trast-1,3-DichloropropeneND0.500.050ND0.120.014trast-1,3-DichloropropeneND0.500.056ND0.120.014trast-1,3-DichloropropeneND0.500.056ND0.120.014trast-1,3-DichloropropeneND0.500.050ND0.130.0</td></t<> | $\mu g/m^3$ $\mu g/m^3$ $\mu g/m^3$ $ppbV$ $ppbV$ $ppbV$ Ethyl tert-Butyl EtherND0.500.051ND0.120.0121,2-DichloroethaneND0.100.050ND0.0250.0121,1,1-TrichloroethaneND0.100.050ND0.0180.0092BenzeneND0.100.050ND0.0160.0080tert-Amyl Methyl EtherND0.100.050ND0.0120.0121,2-DichloropropaneND0.100.050ND0.0220.011BromodichloromethaneND0.100.050ND0.0220.011BromodichloromethaneND0.100.050ND0.0150.0075TrichloroetheneND0.100.050ND0.0120.018n-HeptaneND0.500.061ND0.120.018n-HeptaneND0.500.056ND0.120.014trist-1,3-DichloropropeneND0.500.056ND0.120.014trast-1,3-DichloropropeneND0.500.056ND0.120.014trast-1,3-DichloropropeneND0.500.050ND0.120.014trast-1,3-DichloropropeneND0.500.056ND0.120.014trast-1,3-DichloropropeneND0.500.056ND0.120.014trast-1,3-DichloropropeneND0.500.050ND0.130.0 |

ID = Compound was analyzed for, but not detected above the laboratory detection limit.

IRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS Page 3 of 4

Client:

Test Code:

Analyst:

Test Notes:

Instrument ID:

ENSR Client Sample ID: Method Blank Client Project ID: Phase B Soil Gas / 04020-023-4311

CAS Project ID: P0801342 CAS Sample ID: P080508-MB

EPA TO-15 Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 Rusty Bravo Sampling Media: 6.0 L Summa Canister

Date Collected: NA Date Received: NA Date Analyzed: 5/8/08 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

| | | Result | MRL | MDL | Result | MRL | MDL | Data |
|-------------|-------------------------------|-------------|-------------|-------------|--------|--------|--------|-----------|
| CAS # | Compound | $\mu g/m^3$ | $\mu g/m^3$ | $\mu g/m^3$ | ppbV | ppbV | ppbV | Qualifier |
| 100-41-4 | Ethylbenzene | ND | 0.50 | 0.062 | ND | 0.12 | 0.014 | |
| 179601-23-1 | m,p-Xylenes | ND | 0.50 | 0.13 | ND | 0.12 | 0.030 | |
| 75-25-2 | Bromoform | ND | 0.50 | 0.076 | ND | 0.048 | 0.0074 | |
| 100-42-5 | Styrene | ND | 0.50 | 0.076 | ND | 0.12 | 0.018 | |
| 95-47-6 | o-Xylene | ND | 0.50 | 0.063 | ND | 0.12 | 0.015 | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 0.10 | 0.064 | ND | 0.015 | 0.0093 | |
| 98-82-8 | Cumene | 0.060 | 0.50 | 0.056 | 0.012 | 0.10 | 0.011 | J |
| 103-65-1 | n-Propylbenzene | ND | 0.50 | 0.052 | ND | 0.10 | 0.011 | |
| 622-96-8 | 4-Ethyltoluene | ND | 0.50 | 0.057 | ND | 0.10 | 0.012 | |
| 108-67-8 | 1,3,5-Trimethylbenzene | ND | 0.50 | 0.060 | ND | 0.10 | 0.012 | |
| 98-83-9 | alpha-Methylstyrene | ND | 0.50 | 0.073 | ND | 0.10 | 0.015 | |
| 95-63-6 | 1,2,4-Trimethylbenzene | ND | 0.50 | 0.069 | ND | 0.10 | 0.014 | |
| 100-44-7 | Benzyl Chloride | ND | 0.10 | 0.086 | ND | 0.019 | 0.017 | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 0.10 | 0.062 | ND | 0.017 | 0.010 | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 0.10 | 0.056 | ND | 0.017 | 0.0093 | |
| 135-98-8 | sec-Butylbenzene | ND | 0.50 | 0.058 | ND | 0.091 | 0.011 | |
| 99-87-6 | 4-Isopropyltoluene (p-Cymene) | ND | 0.50 | 0.065 | ND | 0.091 | 0.012 | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 0.10 | 0.066 | ND | 0.017 | 0.011 | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 0.50 | 0.076 | ND | 0.052 | 0.0079 | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 0.10 | 0.076 | ND | 0.013 | 0.010 | |
| 91-20-3 | Naphthalene | 0.10 | 0.20 | 0.074 | 0.020 | 0.038 | 0.014 | J |
| 87-68-3 | Hexachlorobutadiene | ND | 0.10 | 0.090 | ND | 0.0094 | 0.0084 | |
| 98-06-6 | tert-Butylbenzene | ND | 0.20 | 0.050 | ND | 0.036 | 0.0091 | |
| 104-51-8 | n-Butylbenzene | ND | 0.20 | 0.050 | ND | 0.036 | 0.0091 | |

JD = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method. = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

Verified By: R_ Date: 5/9/38 TOI5SCAN.XLT - Tronox - Henderson - PageNo.:

RESULTS OF ANALYSIS Page 4 of 4

| Client: | ENSR | |
|--------------------------|--|-----------------------------------|
| Client Sample ID: | Method Blank | CAS Project ID: P0801342 |
| Client Project ID: | Phase B Soil Gas / 04020-023-4311 | CAS Sample ID: P080508-MB |
| | Tentatively Identified Co | ompounds |
| Fest Code: | EPA TO-15 | Date Collected: NA |
| nstrument ID: | Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 | Date Received: NA |
| Analyst: | Rusty Bravo | Date Analyzed: 5/8/08 |
| Sampling Media: | 6.0 L Summa Canister | Volume(s) Analyzed: 1.00 Liter(s) |
| Fest Notes: | | |

Canister Dilution Factor: 1.00

| GC/MS | Compound Identification | Concentration | Data |
|----------------|-------------------------|---------------|-----------|
| Retention Time | | μg/m³ | Qualifier |
| | No Compounds Detected | | |

Verified By: Re.

SURROGATE SPIKE RECOVERY RESULTS

Page 1 of 1

ENSR Client: Phase B Soil Gas / 04020-023-4311 Client Project ID:

EPA TO-15

Rusty Bravo

CAS Project ID: P0801342

[est Code: nstrument ID: \nalyst: Sampling Media: [est Notes:

Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 6.0 L Summa Canister(s)

Date(s) Collected: 5/7/08 Date(s) Received: 5/8/08 Date(s) Analyzed: 5/8/08

| | | 1,2-Dichlor | oethane-d4 | Tolue | ene-d8 | Bromofluo | robenzene | |
|--------------------|-----------------|-------------|------------|-----------|------------|-----------|------------|-----------|
| Client Sample ID | CAS Sample ID | % | Acceptance | % | Acceptance | % | Acceptance | Data |
| * | | Recovered | Limits | Recovered | Limits | Recovered | Limits | Qualifier |
| Method Blank | P080508-MB | 89 | 70-130 | 103 | 70-130 | 100 | 70-130 | |
| Lab Control Sample | P080508-LCS | 91 | 70-130 | 103 | 70-130 | 101 | 70-130 | |
| SG83B-05-1 | P0801342-001 | 92 | 70-130 | 102 | 70-130 | 101 | 70-130 | |
| SG83B-05-3 | P0801342-002 | 91 | 70-130 | 106 | 70-130 | 102 | 70-130 | ~ |
| SG83B-05-7 | P0801342-003 | 91 | 70-130 | 102 | 70-130 | 98 | 70-130 | |
| SG83B-05-7 | P0801342-003DUP | 91 | 70-130 | 103 | 70-130 | 99 | 70-130 | - |

21

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 3

Client:ENSRClient Sample ID:Lab Control SampleClient Project ID:Phase B Soil Gas / 04020-023-4311

CAS Project ID: P0801342 CAS Sample ID: P080508-LCS

Fest Code:EPA TO-15Date Collected: NAInstrument ID:Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13Date Received: NAAnalyst:Rusty BravoDate Analyzed: 5/08/08Sampling Media:6.0 L Summa CanisterVolume(s) Analyzed:Fest Notes:Fest Notes:NA Liter(s)

| CAS # | Compound | Spike Amount | Result ng | % Recovery | CAS Acceptance Limits | Data Qualifier |
|-----------|--|--------------|--------------|------------|-----------------------------|-------------------|
| 75-71-8 | Dichlorodifluoromethane (CFC 12) | 25.5 | 21.9 | 86 | 69-117 | |
| 74-87-3 | Chloromethane | 24.5 | 20.8 | 85 | 53-131 | |
| 76-14-2 | 1,2-Dichloro-1,1,2,2- tetrafluoroethane (CFC 114) | 26.0 | 22.6 | 87 | 58-133 | |
| 75-01-4 | Vinyl Chloride | 24.8 | 20.7 | 83 | 61-127 | |
| 74-83-9 | Bromomethane | 25.0 | 23.9 | 96 | 67-124 | |
| 75-00-3 | Chloroethane | 25.0 | 23.2 | 93 | 69-123 | |
| 64-17-5 | Ethanol | 23.8 | 20.6 | 87 | 56-137 | |
| 67-64-1 | Acetone | 26.8 | 24.9 | 93 | 63-116 | |
| 75-69-4 | Trichlorofluoromethane | 26.3 | 24.4 | 93 | 71-120 | |
| 107-13-1 | Acrylonitrile | 25.5 | 25.5 | 100 | 74-129 | |
| 75-35-4 | 1,1-Dichloroethene | 27.8 | 25.8 | 93 | 77-116 | |
| 75-65-0 | 2-Methyl-2-Propanol (tert-Butyl Alcohol) | 25.8 | 24.4 | 95 | 35-141 | |
| 75-09-2 | Methylene Chloride | 27.8 | 24.0 | 86 | 71-113 | |
| 107-05-1 | 3-Chloro-1-propene (Allyl Chloride) | 26.8 | 29.6 | 110 | 75-127 | |
| 76-13-1 | Trichlorotrifluoroethane | 27.8 | 25.4 | 91 | 63-129 | |
| 75-15-0 | Carbon Disulfide | 25.0 | 23.1 | 92 | 72-122 | |
| 156-60-5 | trans-1,2-Dichloroethene | 26.5 | 24.5 | 92 | 74-118 | |
| 75-34-3 | 1,1-Dichloroethane | 26.8 | 24.6 | 92 | 74-118 | |
| 1634-04-4 | Methyl tert-Butyl Ether | 26.8 | 24.2 | 90 | 72-119 | |
| 108-05-4 | Vinyl Acetate | 25.3 | 28.0 | 111 | 32-163 | |
| 78-93-3 | 2-Butanone (MEK) | 27.0 | 27.0 | 100 | 71-122 | - · · · · |
| 156-59-2 | cis-1,2-Dichloroethene | 27.0 | 24.4 | 90 | 74-117 | |
| 108-20-3 | Diisopropyl Ether | 26.3 | 22.8 | 87 | 70-131 | |
| 67-66-3 | Chloroform | 29.8 | 28.3 | 95 | 72-113 | |

Date: 5663 TOI5SCAN.XLT - Tronox - Henderson - PageNo.:

22

LABORATORY CONTROL SAMPLE SUMMARY Page 2 of 3

ENSR

Client: Client Sample ID: Lab Control Sample Client Project ID: Phase B Soil Gas / 04020-023-4311

CAS Project ID: P0801342 CAS Sample ID: P080508-LCS

EPA TO-15 Test Code: Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 Instrument ID: Analyst: Rusty Bravo 6.0 L Summa Canister Sampling Media: Test Notes:

Date Collected: NA Date Received: NA Date Analyzed: 5/08/08 Volume(s) Analyzed: NA Liter(s)

| | | | | | CAS | |
|------------|---------------------------|--------------|--------|------------|------------|--|
| CAS # | Compound | Spike Amount | Result | % Recovery | Acceptance | Data |
| | | ng | ng | | Limits | Qualifier |
| 637-92-3 | Ethyl tert-Butyl Ether | 26.0 | 24.0 | 92 | 74-123 | |
| 107-06-2 | 1,2-Dichloroethane | 26.3 | 23.6 | 90 | 72-117 | |
| 71-55-6 | 1,1,1-Trichloroethane | 26.8 | 25.5 | 95 | 78-114 | |
| 71-43-2 | Benzene | 27.0 | 24.9 | 92 | 73-111 | |
| 56-23-5 | Carbon Tetrachloride | 26.0 | 27.2 | 105 | 78-126 | |
| 994-05-8 | tert-Amyl Methyl Ether | 26.0 | 24.9 | 96 | 81-118 | |
| 78-87-5 | 1,2-Dichloropropane | 26.5 | 24.1 | 91 | 78-117 | |
| 75-27-4 | Bromodichloromethane | 27.8 | 26.8 | 96 | 77-120 | |
| 79-01-6 | Trichloroethene | 27.3 | 26.4 | 97 | 80-116 | |
| 123-91-1 | 1,4-Dioxane | 27.5 | 27.8 | 101 | 79-122 | |
| 80-62-6 | Methyl Methacrylate | 25.8 | 26.3 | 102 | 79-128 | |
| 142-82-5 | n-Heptane | 26.8 | 24.3 | 91 | 77-117 | |
| 10061-01-5 | cis-1,3-Dichloropropene | 25.0 | 25.9 | 104 | 78-112 | |
| 108-10-1 | 4-Methyl-2-pentanone | 27.5 | 24.6 | 89 | 78-128 | |
| 10061-02-6 | trans-1,3-Dichloropropene | 28.0 | 29.8 | 106 | 81-121 | |
| 79-00-5 | 1,1,2-Trichloroethane | 26.3 | 24.7 | 94 | 80-117 | Antonio Martinio Mantenanti en ante la los a de las endes districtivas en entre ent |
| 108-88-3 | Toluene | 26.5 | 27.0 | 102 | 76-116 | |
| 591-78-6 | 2-Hexanone | 26.3 | 25.9 | 98 | 69-131 | |
| 124-48-1 | Dibromochloromethane | 27.0 | 30.2 | 112 | 80-128 | |
| 106-93-4 | 1,2-Dibromoethane | 26.3 | 29.6 | 113 | 79-122 | |
| 111-65-9 | n-Octane | 26.0 | 26.2 | 101 | 78-122 | |
| 127-18-4 | Tetrachloroethene | 26.0 | 27.1 | 104 | 77-118 | |
| 108-90-7 | Chlorobenzene | 26.5 | 27.0 | 102 | 78-117 | |

RG

23

LABORATORY CONTROL SAMPLE SUMMARY

Page 1 of 3

| Client: | ENSR |
|--------------------|-----------------------------------|
| Client Sample ID: | Lab Control Sample |
| Client Project ID: | Phase B Soil Gas / 04020-023-4311 |

Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13

EPA TO-15

Rusty Bravo

6.0 L Summa Canister

Test Code:

Analyst:

Test Notes:

Instrument ID:

Sampling Media:

CAS Project ID: P0801342 CAS Sample ID: P080508-LCS

Date Collected: NA Date Received: NA Date Analyzed: 5/08/08 Volume(s) Analyzed: NA Liter(s)

| | | | | | CAS | |
|-----------|--|--------------|--------|------------|------------|-----------|
| CAS # | Compound | Spike Amount | Result | % Recovery | Acceptance | Data |
| | | ng | ng | | Limits | Qualifier |
| 75-71-8 | Dichlorodifluoromethane (CFC 12) | 25.5 | 21.9 | 86 | 69-117 | |
| 74-87-3 | Chloromethane | 24.5 | 20.8 | 85 | 53-131 | |
| 76-14-2 | 1,2-Dichloro-1,1,2,2- tetrafluoroethane (CFC 114) | 26.0 | 22.6 | 87 | 58-133 | |
| 75-01-4 | Vinyl Chloride | 24.8 | 20.7 | 83 | 61-127 | |
| 74-83-9 | Bromomethane | 25.0 | 23.9 | 96 | 67-124 | |
| 75-00-3 | Chloroethane | 25.0 | 23.2 | 93 | 69-123 | |
| 64-17-5 | Ethanol | 23.8 | 20.6 | 87 | 56-137 | |
| 67-64-1 | Acetone | 26.8 | 24.9 | 93 | 63-116 | |
| 75-69-4 | Trichlorofluoromethane | 26.3 | 24.4 | 93 | 71-120 | |
| 107-13-1 | Acrylonitrile | 25.5 | 25.5 | 100 | 74-129 | |
| 75-35-4 | 1,1-Dichloroethene | 27.8 | 25.8 | 93 | 77-116 | |
| 75-65-0 | 2-Methyl-2-Propanol (tert-Butyl Alcohol) | 25.8 | 24.4 | 95 | 35-141 | |
| 75-09-2 | Methylene Chloride | 27.8 | 24.0 | 86 | 71-113 | |
| 107-05-1 | 3-Chloro-1-propene (Allyl Chloride) | 26.8 | 29.6 | 110 | 75-127 | |
| 76-13-1 | Trichlorotrifluoroethane | 27.8 | 25.4 | 91 | 63-129 | |
| 75-15-0 | Carbon Disulfide | 25.0 | 23.1 | 92 | 72-122 | |
| 156-60-5 | trans-1,2-Dichloroethene | 26.5 | 24.5 | 92 | 74-118 | |
| 75-34-3 | 1,1-Dichloroethane | 26.8 | 24.6 | 92 | 74-118 | |
| 1634-04-4 | Methyl tert-Butyl Ether | 26.8 | 24.2 | 90 | 72-119 | |
| 108-05-4 | Vinyl Acetate | 25.3 | 28.0 | 111 | 32-163 | |
| 78-93-3 | 2-Butanone (MEK) | 27.0 | 27.0 | 100 | 71-122 | |
| 156-59-2 | cis-1,2-Dichloroethene | 27.0 | 24.4 | 90 | 74-117 | |
| 108-20-3 | Diisopropyl Ether | 26.3 | 26.8 | 102 | 70-131 | |
| 67-66-3 | Chloroform | 26.8 | 28.7 | 107 | 72-113 | |

Date: 5900 24

LABORATORY DUPLICATE SUMMARY RESULTS

Page 1 of 3

Client:ENSRClient Sample ID:SG83B-05-7Client Project ID:Phase B Soil Gas / 04020-023-4311

SC00791

Container ID:

Test Code:EPA TO-15Instrument ID:Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13Analyst:Rusty BravoSampling Media:6.0 L Summa CanisterFest Notes:Fest Notes:

CAS Project ID: P0801342 CAS Sample ID: P0801342-003DUP

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

Initial Pressure (psig): -4.2

Final Pressure (psig): 3.5

| | | | | Canis | ter Dilutio | n Factor: | 1.73 |
|--------|---|---|---|---|--|--|--|
| | | | | | | | |
| Sample | Result | Sample | Result | Average | % RPD | RPD | Data |
| μg/m³ | ppbV | µg/m³ | ppbV | µg/m³ | | Limit | Qualifier |
| ND | ND | ND | ND | - | - | 25 | |
| ND | ND | ND | ND | | - | 25 | |
| ND | ND | ND | ND | - | - | 25 | |
| ND | ND | ND | ND | - | - | 25 | |
| ND | ND | ND | ND | - | - | . 25 | |
| ND | ND | ND | ND | - | - | 25 | |
| ND | ND | ND | ND | - | · · - | | |
| 114 | 48.1 | 107 | 45.2 | 110.5 | 6 | | J, B |
| 1,520 | 271 | 1,530 | 273 | 1525 | 0.7 | 25 | |
| ND | ND | ND | ND | | N | 25 | |
| ND | ND | ND | ND | - | - | 25 | |
| ND | ND | ND | ND | - | - | 25 | |
| 9.34 | 2.69 | ND | ND | - | - | 25 | J |
| ND | ND | ND | ND | - | - | 25 | |
| ND | ND | ND | ND | | - | 25 | |
| ND | ND | ND | ND | - | - | 25 | |
| ND | ND | ND | ND | | - | 25 | |
| ND | ND | ND | ND | - | - | 25 | |
| ND | ND | ND | ND | - | - | 25 | |
| ND | ND | ND | ND | | - | 25 | |
| 23.0 | 7.80 | 18.7 | 6.34 | 20.85 | 21 | 25 | J |
| ND | ND | ND | ND | - | - | 25 | |
| ND | ND | ND | ND | - | - | 25 | |
| 54,300 | 11,100 | 45,700 | 9,370 | 50000 | 17 | 25 | |
| | μg/m ³ ND ND ND ND ND ND 114 1,520 ND ND ND ND ND ND ND ND ND ND ND ND ND | ND ND 114 48.1 1,520 271 ND ND ND | Sample Result Sample $\mu g/m^3$ $\mu g/m^3$ $ppbV$ $\mu g/m^3$ ND ND ND ND ND </td <td>μg/m³ ppbV μg/m³ ppbV ND ND ND ND ND ND</td> <td>DuplicationSample FesultSample FesultAverageµg/m³ppbVµg/m³ppbVµg/m³ND1,5202711,5302731525ND<</td> <td>DuplicitsSample FundSample FundAverage% RPDµg/m³pbVµg/m³pbVµg/m³% RPDµg/m³pbVµg/m³µg/m³µg/m³NDANDNDNDNDNDANDNDNDNDNDANDNDNDNDNDANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDND</td> <td>Sample ResultSample ResultAverage% RPDRPD$\mu g/m^3$ppbV$\mu g/m^3$LimitNDNDNDND-25NDN</td> | μg/m³ ppbV μg/m³ ppbV ND ND ND ND ND ND | DuplicationSample FesultSample FesultAverageµg/m³ppbVµg/m³ppbVµg/m³ND1,5202711,5302731525ND< | DuplicitsSample FundSample FundAverage% RPDµg/m³pbVµg/m³pbVµg/m³% RPDµg/m³pbVµg/m³µg/m³µg/m³NDANDNDNDNDNDANDNDNDNDNDANDNDNDNDNDANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDNDAANDNDNDND | Sample ResultSample ResultAverage% RPDRPD $\mu g/m^3$ ppbV $\mu g/m^3$ LimitNDNDNDND-25NDN |

JD = Compound was analyzed for, but not detected above the laboratory detection limit.

= The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

Verified By: <u>RG</u>

3 = Analyte was found in the method blank.

P0801342_TO15_0805091251_SS.xls - Dup (3)

LABORATORY DUPLICATE SUMMARY RESULTS

Page 2 of 3

| Client: | ENSR |
|--------------------|--|
| Client Sample ID: | SG83B-05-7 |
| Client Project ID: | Phase B Soil Gas / 04020-023-4311 |
| | |
| Fest Code: | EPA TO-15 |
| instrument ID: | Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 |
| Analyst: | Rusty Bravo |
| Sampling Media: | 6.0 L Summa Canister |

SC00791

Fest Notes: Container ID:

CAS Project ID: P0801342 CAS Sample ID: P0801342-003DUP

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

Initial Pressure (psig): -4.2 Final Pressure (psig): 3.5

| | | | | | Canis | ter Dilutio | n Factor: | 1.73 |
|---------------------------|-------------|--------|--------|--------|---------|-------------|-----------|--|
| | | | Dupli | cate | | | | |
| Compound | Sample | Result | Sample | Result | Average | % RPD | RPD | Data |
| * | $\mu g/m^3$ | ppbV | µg/m³ | ppbV | μg/m³ | | Limit | Qualifier |
| Ethyl tert-Butyl Ether | ND | ND | ND | ND | - | - | 25 | |
| 1,2-Dichloroethane | ND | ND | ND | ND | - | - | 25 | |
| 1,1,1-Trichloroethane | ND | ND | ND | ND | - | - | 25 | |
| Benzene | 101 | 31.7 | 102 | 32.1 | 101.5 | 1 | 25 | |
| Carbon Tetrachloride | 12,000 | 1,910 | 12,300 | 1,950 | 12150 | 2 | 25 | |
| tert-Amyl Methyl Ether | ND | ND | ND | ND | - | - | 25 | |
| 1,2-Dichloropropane | ND | ND | ND | ND | - | - | 25 | |
| Bromodichloromethane | ND | ND | ND | ND | - | - | 25 | |
| Trichloroethene | 11.4 | 2.13 | 11.9 | 2.22 | 11.65 | 4 | 25 | \mathbf{J} |
| 1,4-Dioxane | ND | ND | ND | ND | * | | 25 | anna gu an ann an airtean airte an chuileachta dan bha a |
| Methyl Methacrylate | ND | ND | ND | ND | - | - | 25 | |
| n-Heptane | ND | ND | ND - | ND | - | - | 25 | |
| cis-1,3-Dichloropropene | ND | ND | ND | ND | - | - | 25 | |
| 4-Methyl-2-pentanone | ND | ND | ND | ND | - | - | 25 | |
| trans-1,3-Dichloropropene | ND | ND | ND | ND | | | 25 | - |
| 1,1,2-Trichloroethane | ND | ND | ND | , ND | - | - | 25 | |
| Toluene | ND | ND | ND | ND | - | - | 25 | |
| 2-Hexanone | ND | ND | ND | ND | - | - | 25 | |
| Dibromochloromethane | ND | ND | ND | ND | - | - | 25 | |
| 1,2-Dibromoethane | ND | ND | ND | ND | | - | 25 | |
| n-Octane | ND | ND | ND | ND | - | - | 25 | |
| Tetrachloroethene | 127 | 18.7 | 124 | 18.3 | 125.5 | 2 | 25 | |
| Chlorobenzene | 180 | 39.1 | 197 | 42.7 | 188.5 | 9 | 25 | |

JD = Compound was analyzed for, but not detected above the laboratory detection limit.

= The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

Verified By: <u>Rc-</u>

Date: 56608 TO15SCAN.XLT - Tronox - Henderson - PageNo.:

26

LABORATORY DUPLICATE SUMMARY RESULTS

Page 3 of 3

| Client: Client Sample ID: Client Project ID: | ENSR SG83B-05-7 Phase B Soil Gas / 04020-023-4311 |
|--|---|
| Fest Code: | EPA TO-15 |
| nstrument ID: | Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 |
| Analyst: | Rusty Bravo |
| Sampling Media: | 6.0 L Summa Canister |
| Fest Notes: | |
| Container ID: | SC00791 |

CAS Project ID: P0801342 CAS Sample ID: P0801342-003DUP

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

Initial Pressure (psig): -4.2

Canister Dilution Factor: 1.73

27

Date: 5/0/05 TO15SCAN.XLT - Tronox - Henderson - PageNo.:

Final Pressure (psig): 3.5

| | | | Dupli | icate | Carris | ster Difutio. | | 1.75 |
|-------------------------------|--------|--------|--------|-------|---------|---------------|-------|-----------|
| Compound | Sample | Result | Sample | | Average | % RPD | RPD | Data |
| | μg/m³ | ppbV | µg/m³ | ppbV | μg/m³ | | Limit | Qualifier |
| Ethylbenzene | ND | ND | ND | NI |) - | - | 25 | |
| m,p-Xylenes | ND | ND | ND | NI |) - | - | 25 | |
| Bromoform | ND | ND | ND | NI |) - | - | 25 | |
| Styrene | ND | ND | ND | NI |) - | - | 25 | |
| o-Xylene | ND | ND | ND | NI |) - | - . | 25 | |
| 1,1,2,2-Tetrachloroethane | ND | ND | ND | NI |) - | - | 25 | |
| Cumene | ND | ND | ND | NI |) - | - | 25 | |
| n-Propylbenzene | ND | ND | ND | NI |) - | - | 25 | |
| 4-Ethyltoluene | ND | ND | ND | NI |) | - | 25 | |
| 1,3,5-Trimethylbenzene | ND | ND | ND | NI |) - | - | 25 | |
| alpha-Methylstyrene | ND | ND | ND | NE |) - | - | 25 | |
| 1,2,4-Trimethylbenzene | ND | ND | ND | NE |) - | - | 25 | |
| Benzyl Chloride | ND | ND | ND | NI |) - | - | 25 | |
| 1,3-Dichlorobenzene | ND | ND | ND | NI |) - | - | 25 | |
| 1,4-Dichlorobenzene | ND | ND | ND | NI |) _ ' | - | 25 | |
| sec-Butylbenzene | ND | ND | ND | NE |) - | - | 25 | |
| 4-Isopropyltoluene (p-Cymene) | ND | ND | ND | NI |) - | - | 25 | |
| 1,2-Dichlorobenzene | ND | ND | ND | NE |) - | - | 25 | |
| 1,2-Dibromo-3-chloropropane | ND | ND | ND | NE |) _ ' | - | 25 | |
| 1,2,4-Trichlorobenzene | ND | ND | ND | NE |) - | - | 25 | - |
| Naphthalene | ND | ND | ND | NI |) - | - | 25 | |
| Hexachlorobutadiene | ND | ND | ND | NE |) - | - | 25 | |
| tert-Butylbenzene | ND | ND | ND | NE |) _ | - | 25 | |
| n-Butylbenzene | ND | ND | ND | NI |) _ | | 25 | |

Verified By:___

JD = Compound was analyzed for, but not detected above the laboratory detection limit.

RESULTS OF ANALYSIS

Page 1 of 1

Client: ENSR Client Project ID: Phase B Soil Gas / 04020-023-4311

CAS Project ID: P0801342

Internal Standard Area and RT Summary

| Test Code: | EPA TO-15 | |
|-----------------|--|-------------------------|
| Instrument ID: | Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 | Lab File ID: 05080801.D |
| Analyst: | Rusty Bravo | Date Analyzed: 5/8/08 |
| Sampling Media: | 6.0 L Summa Canister(s) | Time Analyzed: 08:41 |
| Test Notes: | | |

| | IS1 (BCM) | IS2 (DFB) | | | IS3 (CBZ) | |
|------------------|-----------|-----------|---------|-------|-----------|-------|
| | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 24 Hour Standard | 242477 | 12.59 | 1048411 | 15.52 | 500730 | 21.35 |
| Upper Limit | 339468 | 12.92 | 1467775 | 15.85 | 701022 | 21.68 |
| Lower Limit | 145486 | 12.26 | 629047 | 15.19 | 300438 | 21.02 |

| | Client Sample ID | | | | | | |
|-----|---------------------------------------|--------|-------|---------|-------|--------|-------|
| 01 | Method Blank | 241346 | 12.58 | 1037647 | 15.51 | 489334 | 21.35 |
| 02 | Lab Control Sample | 244114 | 12.59 | 1061640 | 15.52 | 502772 | 21.35 |
| 03 | SG83B-05-1 | 242853 | 12.58 | 1040889 | 15.51 | 494646 | 21.35 |
| 04 | SG83B-05-3 | 231941 | 12.58 | 1004035 | 15.51 | 467452 | 21.35 |
| 05 | SG83B-05-7 | 226199 | 12.58 | 971717 | 15.51 | 470342 | 21.35 |
| 06 | SG83B-05-1 (Dilution) | 219916 | 12.58 | 959711 | 15.51 | 446346 | 21.35 |
| 07 | SG83B-05-3 (Dilution) | 213718 | 12.58 | 929659 | 15.51 | 442512 | 21.35 |
| 08 | SG83B-05-7 (Dilution) | 205516 | 12.58 | 899772 | 15.51 | 428319 | 21.35 |
| 09 | SG83B-05-7 (Lab Duplicate - Dilution) | 202052 | 12.58 | 880501 | 15.51 | 420894 | 21.35 |
| 10 | SG83B-05-7 (Lab Duplicate) | 199552 | 12.58 | 853249 | 15.51 | 416006 | 21.35 |
| 1 1 | · · · · · · | | | | | | |

- 11 12 13 14 15 16
- 10 17 18
- 19
- 20

IS1 (BCM) = Bromochloromethane

IS2 (DFB) = 1,4-Difluorobenzene

IS3 (CBZ) = Chlorobenzene-d5

AREA UPPER LIMIT = 140% of internal standard area AREA LOWER LIMIT = 60% of internal standard area RT UPPER LIMIT = 0.33 minutes of internal standard RT RT LOWER LIMIT = 0.33 minutes of internal standard RT

Column used to flag values outside QC limits with an asterisk.

* Values outside of QC limits.

Verified By: R.

Date: 59.668 TO15SCAN.XLT - Tronox - Henderson - PageNo.:

RESULTS OF HELIUM ANALYSIS

RESULTS OF ANALYSIS

Page 1 of 1

| Client: Client Project ID: | ENSR : Phase B Soil Gas / 04020-023-4311 | | | CA | S Project ID: P08013 | 42 |
|--|--|------------------------------|--------------------------------|----------------|---|-------------------|
| | | I | Helium | | | |
| Test Code: Instrument ID: Analyst: Sampling Media: Test Notes: | EPA 3C Modified HP5890 II/GC8/TCD Zheng Wang/Wade Henton/Cl 6.0 L Summa Canister(s) | hris Cornett | | Da | (s) Collected: 5/7/08 ate Received: 5/8/08 ate Analyzed: 5/8/08 | |
| | | | | , | | |
| Client Sample ID | CAS Sample ID | Injection Volume ml(s) | Canister Dilution Factor | Result ppmV | MRL ppmV | Data Qualifier |
| SG83B-05-1 | P0801342-001 | 1.00 | 1.64 | 2,700 | 41 | |
| SG83B-05-3 | P0801342-002 | 1.00 | 1.67 | 160 | 42 | |
| Method Blank | P080508-MB | 1.00 | 1.00 | ND | 25 | |

ND = Compound was analyzed for, but not detected above the laboratory reporting limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

P0801342_3CHE_0805081513_SS - Helium

Verified By: <u>RG</u>



RESULTS OF VOLATILE ORGANIC ANALYSIS

RESULTS OF ANALYSIS

Page 1 of 3

| Client: | ENSR |
|---------------------------|-----------------------------------|
| Client Sample ID: | SG83B-05-1 |
| Client Project ID: | Phase B Soil Gas / 04020-023-4311 |
| | |

Test Code:EPA TO-15Instrument ID:Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13Analyst:Rusty BravoSampling Media:6.0 L Summa CanisterTest Notes:Container ID:SC00979

CAS Project ID: P0801342 CAS Sample ID: P0801342-001

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

3.5

Initial Pressure (psig):

ig): -3.6 Final Pressure (psig):

Canister Dilution Factor: 1.64

| CAS # | Compound | Result | MRL | MDL | Result | MRL | MDL | Data |
|-----------|-------------------------------------|--------|-------|-------------|--------|------|------|-----------|
| | - | μg/m³ | μg/m³ | $\mu g/m^3$ | ppbV | ppbV | ppbV | Qualifier |
| 75-71-8 | Dichlorodifluoromethane (CFC 12) | ND | 82 | 8.2 | ND | 17 | 1.7 | |
| 74-87-3 | Chloromethane | ND | 16 | 8.2 | ND | 7.9 | 4.0 | |
| 76-14-2 | 1,2-Dichloro-1,1,2,2- | ND 82 | 80 | 82 8.2 | ND | 12 | 1.2 | |
| /0-14-2 | tetrafluoroethane (CFC 114) | ND | 02 | | | | | |
| 75-01-4 | Vinyl Chloride | ND | 16 | 8.2 | ND | 6.4 | 3.2 | |
| 74-83-9 | Bromomethane | ND | 16 | 8.2 | ND | 4.2 | 2.1 | |
| 75-00-3 | Chloroethane | ND | 16 | 8.2 | ND | 6.2 | 3.1 | |
| 64-17-5 | Ethanol | 12 | 820 | 8.2 | 6.6 | 440 | 4.4 | J |
| 67-64-1 | Acetone | 150 | 820 | 12 | 64 | 350 | 5.0 | J, B |
| 75-69-4 | Trichlorofluoromethane | 1,500 | 16 | 8.2 | 280 | 2.9 | 1.5 | |
| 107-13-1 | Acrylonitrile | ND | 82 | 11 | ND | 38 | 5.3 | |
| 75-35-4 | 1,1-Dichloroethene | ND | 16 | 8.2 | ND | 4.1 | 2.1 | |
| 75-65-0 | 2-Methyl-2-Propanol | ND | 82 | 12 | ND | 27 | 4.0 | |
| | (tert-Butyl Alcohol) | ND | | | ND | | | |
| 75-09-2 | Methylene Chloride | ND | 82 | 8.2 | ND | 24 | 2.4 | |
| 107-05-1 | 3-Chloro-1-propene (Allyl Chloride) | ND | 16 | 8.2 | ND | 5.2 | 2.6 | |
| 76-13-1 | Trichlorotrifluoroethane | ND | 16 | 9.2 | ND | 2.1 | 1.2 | |
| 75-15-0 | Carbon Disulfide | ND | 82 | 20 | ND | 26 | 6.3 | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 16 | 8.2 | ND | 4.1 | 2.1 | |
| 75-34-3 | 1,1-Dichloroethane | ND | 16 | 8.2 | ND | 4.1 | 2.0 | |
| 1634-04-4 | Methyl tert-Butyl Ether | ND | 16 | 8.2 | ND | 4.6 | 2.3 | |
| 108-05-4 | Vinyl Acetate | ND | 820 | 26 | ND | 230 | 7.5 | |
| 78-93-3 | 2-Butanone (MEK) | 26 | 82 | 8.2 | 8.7 | 28 | 2.8 | J |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 16 | 8.2 | ND | 4.1 | 2.1 | |
| 108-20-3 | Diisopropyl Ether | ND | 82 | 9.7 | ND | 20 | 2.3 | |
| 67-66-3 | Chloroform | 52,000 | 16 | 9.7 | 11,000 | 3.4 | 2.0 | |

ND = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method. J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

P0801342_TO15_0805091251_SS - Sample

Verified By: <u>Rc-</u>

32

RESULTS OF ANALYSIS

Page 2 of 3

ENSR Client: Client Sample ID: SG83B-05-1 Client Project ID: Phase B Soil Gas / 04020-023-4311

6.0 L Summa Canister

EPA TO-15

Rusty Bravo

SC00979

Test Code:

Analyst:

Instrument ID:

Container ID:

Sampling Media: Test Notes:

CAS Project ID: P0801342 CAS Sample ID: P0801342-001

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

Initial Pressure (psig):

-3.6

Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13

Final Pressure (psig): 3.5

Canister Dilution Factor: 1.64

| CAS # | Compound | Result μg/m³ | MRL µg/m³ | MDL µg/m³ | Result ppbV | MRL ppbV | MDL ppbV | Data Qualifier |
|------------|---------------------------|-----------------|--------------|--------------|----------------|-------------|-------------|---|
| 637-92-3 | Ethyl tert-Butyl Ether | ND | 82 | 8.4 | ND | 20 | 2.0 | |
| 107-06-2 | 1,2-Dichloroethane | ND | 16 | 8.2 | ND | 4.1 | 2.0 | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 16 | 8.2 | ND | 3.0 | 1.5 | |
| 71-43-2 | Benzene | 100 | 16 | 8.2 | 31 | 5.1 | 2.6 | |
| 56-23-5 | Carbon Tetrachloride | 12,000 | 16 | 8.2 | 1,900 | 2.6 | 1.3 | |
| 994-05-8 | tert-Amyl Methyl Ether | ND | 82 | 8.2 | ND | 20 | 2.0 | |
| 78-87-5 | 1,2-Dichloropropane | ND | 16 | 8.2 | ND | 3.5 | 1.8 | |
| 75-27-4 | Bromodichloromethane | ND | 16 | 8.2 | ND | 2.4 | 1.2 | |
| 79-01-6 | Trichloroethene | 16 | 16 | 8.2 | 3.0 | 3.1 | 1.5 | J |
| 123-91-1 | 1,4-Dioxane | ND | 82 | 10 | ND | 23 | 2.8 | na adama kana yang ting a di atau kanang tana mang gana na mana ngan na |
| 80-62-6 | Methyl Methacrylate | ND | 82 | 12 | ND | 20 | 3.0 | |
| 142-82-5 | n-Heptane | ND | 82 | 10 | ND | 20 | 2.6 | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 82 | 8.5 | ND | 18 | 1.9 | |
| 108-10-1 | 4-Methyl-2-pentanone | ND | 82 | 9.2 | ND | 20 | 2.2 | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 82 | 10 | ND | 18 | 2.3 | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 16 | 8.2 | ND | 3.0 | 1.5 | |
| 108-88-3 | Toluene | 13 | 82 | 8.2 | 3.6 | 22 | 2.2 | J |
| 591-78-6 | 2-Hexanone | ND | 82 | 12 | ND | 20 | 3.0 | |
| 124-48-1 | Dibromochloromethane | ND | 16 | 11 | ND | 1.9 | 1.3 | |
| 106-93-4 | 1,2-Dibromoethane | ND | 16 | 8.9 | ND | 2.1 | 1.2 | |
| 111-65-9 | n-Octane | ND | 82 | 8.2 | ND | 18 | 1.8 | |
| 127-18-4 | Tetrachloroethene | 110 | 16 | 8.2 | 16 | 2.4 | 1.2 | |
| 108-90-7 | Chlorobenzene | 120 | 16 | 8.4 | 26 | 3.6 | 1.8 | |

ND = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

P0801342_TO15_0805091251_SS - Sample

Verified By: <u>RG</u>

RESULTS OF ANALYSIS

Page 3 of 3

Client:ENSRClient Sample ID:SG83B-05-1Client Project ID:Phase B Soil Gas / 04020-023-4311

Test Code:EPA TO-15Instrument ID:Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13Analyst:Rusty BravoSampling Media:6.0 L Summa CanisterTest Notes:Container ID:SC00979

CAS Project ID: P0801342 CAS Sample ID: P0801342-001

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

3.5

Initial Pressure (psig):

-3.6 Final Press

Final Pressure (psig):

Canister Dilution Factor: 1.64

| | | Result | MRL | MDL | Result | MRL | MDL | Data |
|-------------|-------------------------------|--------|-------------|-------------|--------|------|------|-----------|
| CAS # | Compound | μg/m³ | $\mu g/m^3$ | $\mu g/m^3$ | ppbV | ppbV | ppbV | Qualifier |
| 100-41-4 | Ethylbenzene | ND | 82 | 10 | ND | 19 | 2.3 | |
| 179601-23-1 | m,p-Xylenes | ND | 82 | 21 | ND | 19 | 4.9 | |
| 75-25-2 | Bromoform | ND | 82 | 12 | ND | 7.9 | 1.2 | |
| 100-42-5 | Styrene | ND | 82 | 12 | ND | 19 | 2.9 | |
| 95-47-6 | o-Xylene | ND | 82 | 10 | ND | 19 | 2.4 | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 16 | 10 | ND | 2.4 | 1.5 | |
| 98-82-8 | Cumene | 9.7 | 82 | 9.2 | 2.0 | 17 | 1.9 | J, B |
| 103-65-1 | n-Propylbenzene | ND | 82 | 8.5 | ND | 17 | 1.7 | |
| 622-96-8 | 4-Ethyltoluene | 12 | 82 | 9.3 | 2.4 | 17 | 1.9 | J |
| 108-67-8 | 1,3,5-Trimethylbenzene | ND | 82 | 9.8 | ND | 17 | 2.0 | |
| 98-83-9 | alpha-Methylstyrene | ND | 82 | 12 | ND | 17 | 2.5 | |
| 95-63-6 | 1,2,4-Trimethylbenzene | ND | 82 | 11 | ND | 17 | 2.3 | |
| 100-44-7 | Benzyl Chloride | ND | 16 | 14 | ND | 3.2 | 2.7 | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 16 | 10 | ND | 2.7 | 1.7 | |
| 106-46-7 | 1,4-Dichlorobenzene | 17 | 16 | 9.2 | 2.9 | 2.7 | 1.5 | |
| 135-98-8 | sec-Butylbenzene | ND | 82 | 9.5 | ND | 15 | 1.7 | |
| 99-87-6 | 4-Isopropyltoluene (p-Cymene) | ND | 82 | 11 | ND | 15 | 1.9 | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 16 | 11 | ND | 2.7 | 1.8 | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 82 | 12 | ND | 8.5 | 1.3 | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 16 | 12 | ND | 2.2 | 1.7 | |
| 91-20-3 | Naphthalene | 28 | 33 | 12 | 5.3 | 6.3 | 2.3 | J, B |
| 87-68-3 | Hexachlorobutadiene | ND | 16 | 15 | ND | 1.5 | 1.4 | |
| 98-06-6 | tert-Butylbenzene | ND | 33 | 8.2 | ND | 6.0 | 1.5 | |
| 104-51-8 | n-Butylbenzene | ND | 33 | 8.2 | ND | 6.0 | 1.5 | |

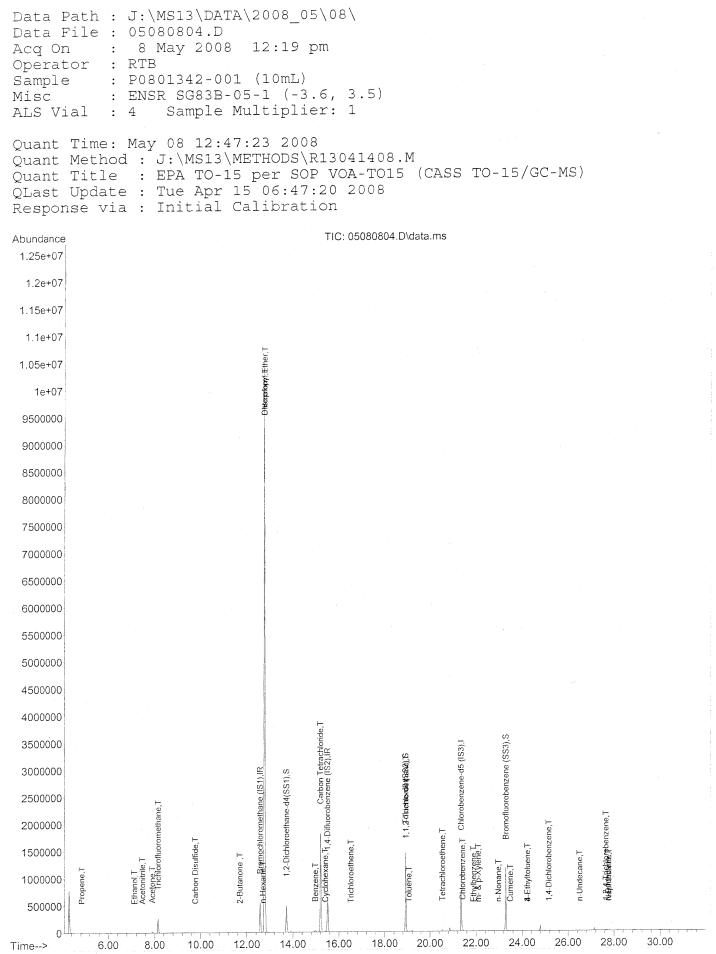
ND = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

34



R13041408.M Fri May 09 14:48:30 2008

| | - | | | | |
|--|----------------------|---------------|-------------|------------------------|--------------------|
| Data Path : J:\MS13\DATA\2008_0 | 5\08\ | | | | |
| Data File : 05080804.D | | | | | |
| Acq On : 8 May 2008 12:19 j | pm | | | | |
| Operator : RTB Sample · P0801342-001 (10mL) | | | | | |
| Misc : ENSR SG83B-05-1 (-3 | .6, 3.5) | | | | |
| Sample : P0801342-001 (10mL) Misc : ENSR SG83B-05-1 (-3 ALS Vial : 4 Sample Multiplie | er: 1 | | | | |
| | | | | | |
| Quant Time: May 08 12:47:23 200 | 8 . | o M | | | |
| Quant Method : J:\MS13\METHODS\ Quant Title : EPA TO-15 per SO | RI304140 P VOD-TO | 8.M 15 (C) | ASS TO-15/0 | C-MS) | |
| QLast Update : Tue Apr 15 06:47 | $\cdot 20 2008$ | 10 (01 | 400 10 10/ | | |
| Response via : Initial Calibrat. | ion | | | | |
| - | | | | ~ | |
| Internal Standards | R.T. | QION | Response | Conc Units | ; Dev(Min) |
| 1) Bromochloromethane (IS1) | 12.58 | 130 | 242853 | 25.000 ng | -0.03 |
| 1) Bromochloromethane (IS1) 37) 1,4-Difluorobenzene (IS2) 56) Chlorobenzene-d5 (IS3) | 15.51 | 114 | 1040889 | 25.000 ng | -0.02 |
| 56) Chlorobenzene-d5 (IS3) | 21.35 | 82 | 494646 | 25.000 ng | -0.01 |
| | | | | | |
| System Monitoring Compounds 33) 1,2-Dichloroethane-d4(| 13 72 | 65 | 448165 | 23.011 ng | -0.03 |
| spiked Amount 25 000 | | | Recove | erv = 92 | 2.048 / |
| 57) Toluene-d8 (SS2) | 18.92 | 98 | 1129319 | 25.472 ng | -0.01 |
| Spiked Amount 25.000 | | | Recove | ery = 101 | |
| 73) Bromofluorobenzene (SS3) | 23.29 | 174 | 384401 | 25.194 ng ery = 100 | 0.00 |
| Spiked Amount 25.000 | | | Recove | ery = 100 | . / 0 % |
| Target Compounds | | | | | Qvalue |
| 2) Propene | 4.84 | 42 | 3109 | 0.155 ng | # 63 |
| 3) Dichlorodifluoromethane | 4.99 | 85 | 488 | N.D. 🗸 | |
| <pre>Target Compounds 2) Propene 3) Dichlorodifluoromethane 4) Chloromethane 5) Freon 114 6) Vinyl Chloride 7) 1,3-Butadiene 8) Bromomethane 0) Chloroothane</pre> | 0.00 | 50 | 0 | N.D. | |
| 5) Freon 114 | 0.00 | 135 62 | 0 | | |
| 6) VINYI CHIOFIde 7) 1 2-Dutadiene | 6.00 | 54 | 661 | N.D. | |
| 8) Bromomethane | 6.51 | 94 | 57 | N.D. | |
| 9) Chloroethane | 0.00 | 64 | 0 | | |
| 10) Ethanol | 7.13 | 45 | 1035 < | 0.076 ng> | # 52 |
| 11) Acetonitrile | 7.47 | 41 | 4208 | 0.117 ng | 81 |
| 12) Acrolein | 7.66 | 56 > 58 | 70 12415 | N.D. (0.924 ng | # 66 |
| 13) Acetone 14) Trichlorofluoromethane | 8.14 |)101 | 271272 | 9.422 ng | π <u>99</u> |
| 14) Trichlorofluoromethane 15) Isopropanol | 8.36 | 45 | 568 | N.D. | |
| 16) Acrylonitrile | 8.69 | 53 | 78 | N.D. | |
| 17) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | |
| 18) tert-Butanol | 9.32 | 59 | 1244 | N.D. | |
| 19) Methylene Chloride | 9.35 | 84 | 524 56 | N.D. / N.D. / | |
| 20) Allyl Chloride | 9.54 0.00 | 41 151 | | N.D. | |
| 21) Trichlorotrifluoroethane 22) Carbon Disulfide | 9.77 | 76 | 2481 | -0.043 ng | 82 |
| 23) trans-1,2-Dichloroethene | 0.00 | 61 | 0 | N.D. | |
| 24) 1,1-Dichloroethane | 11.10 | 63 | 1027 | N.D. | |
| 25) Methyl tert-Butyl Ether | 0.00 | 73 | 0 | N.D. | |
| 26) Vinyl Acetate | 0.00 | 86 | 0 | N.Đ. | # 50 |
| 27) 2-Butanone (| 11.70 |) 72 61 | 1470 0 | (0.156 ng) N.D./ | # 50 |
| 28) cis-1,2-Dichloroethene 29) Diisopropyl Ether | 0.00 | 61 87 | 991565 | <u>80.106 ng</u> N | K # 1 |
| 30) Ethyl Acetate | 12.74 | 61 | 77 | N.D. | l |
| 31) n-Hexane | 12.70 | 57 | 1641 | 0.054 ng | # ⁷¹ 36 |
| | | | Į. | Inglio | |
| 3041408.M Fri May 09 14:48:30 2 | 008 | | 70 | 8-11-0 | Page: 1 |
| | | | | | |
| | | | , | | |

R13041408.M Fri May 09 14:48:30 2008

Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080804.D Acq On : 8 May 2008 12:19 pm Operator : RTB Sample : P0801342-001 (10mL) Misc : ENSR SG83B-05-1 (-3.6, 3.5) ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 12:47:23 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration Internal StandardsR.T. QIon Response Conc Units Dev (Min)32) Chloroform12.78839317625410.934 ng9934) Tetrahydrofuran0.00720N.D.35) Ethyl tert-Butyl Ether0.00870N.D.36) 1,2-Dichloroethane13.7262177N.D.38) 1,1,1-Trichloroethane0.00970N.D.39) Isopropyl Acetate14.9761213N.D.40) 1-Butanol14.9878336940.610 ng9841) Benzene15.21117133008272.830 ng9843) Cyclohexane15.418468840.337 ng#44) tert-Amyl Methyl Ether0.00730N.D.45) 1,2-Dichloropropane0.00630N.D.46) Bromodichloromethane16.488369N.D.47) Trichloroethene16.6057464N.D.48) 1,4-Dioxane0.001000N.D.50) Methyl Methacrylate0.001000N.D.51) n-Heptane16.9871586N.D.52) cis-1,3-Dichloropropene0.00750N.D.53) ND100750N.D. R.T. QIon Response Conc Units Dev(Min) Internal Standards 49)Isooctane16.6057464N.D.50)Methyl Methacrylate0.001000N.D.,51)n-Heptane16.9871586N.D.,52)cis-1,3-Dichloropropene0.00750N.D.,53)4-Methyl-2-pentanone17.795865N.D.,54)trans-1,3-Dichloropropene18.457588N.D.,55)1,1,2-Trichloroethane19.069145820.082 ng58)Toluene19.4043792N.D.,60)Dibromochloromethane0.001070N.D.,61)1,2-Dibromoethane0.001070N.D.,62)Butyl Acetate20.194351N.D.63)n-Octane20.36573930.657 ng64)Tetrachloroethene21.41112248720.721 ng65)Chlorobenzene21.899125030.040 ng67)m-& p-Xylene22.099147470.1144 ng 7 91 94 99 66)Ethylbenzene21.899125030.040 ng67)m- & p-Xylene22.099147470.114 ng#68)Bromoform0.001730N.D.69)Styrene22.58104973N.D.70)o-Xylene22.72911739N.D.71)n-Nonane22.984318450.051 ng72)1,1,2,2-Tetrachloroethane22.708356N.D.74)Cumene23.4810533220.059 ng#75)alpha-Pinene23.959370N.D.76)n-Propylbenzene24.11911200N.D.77)3-Ethyltoluene24.2310541850.068 ng#78)4-Ethyltoluene24.381051648N.D.#79)1,3,5-Trimethylbenzene24.381051648N.D. 66 # 68 88 49 0.068 ng # 50 52 N.D. ĵ Page: 2 Fostoglog

R13041408.M Fri May 09 14:48:30 2008

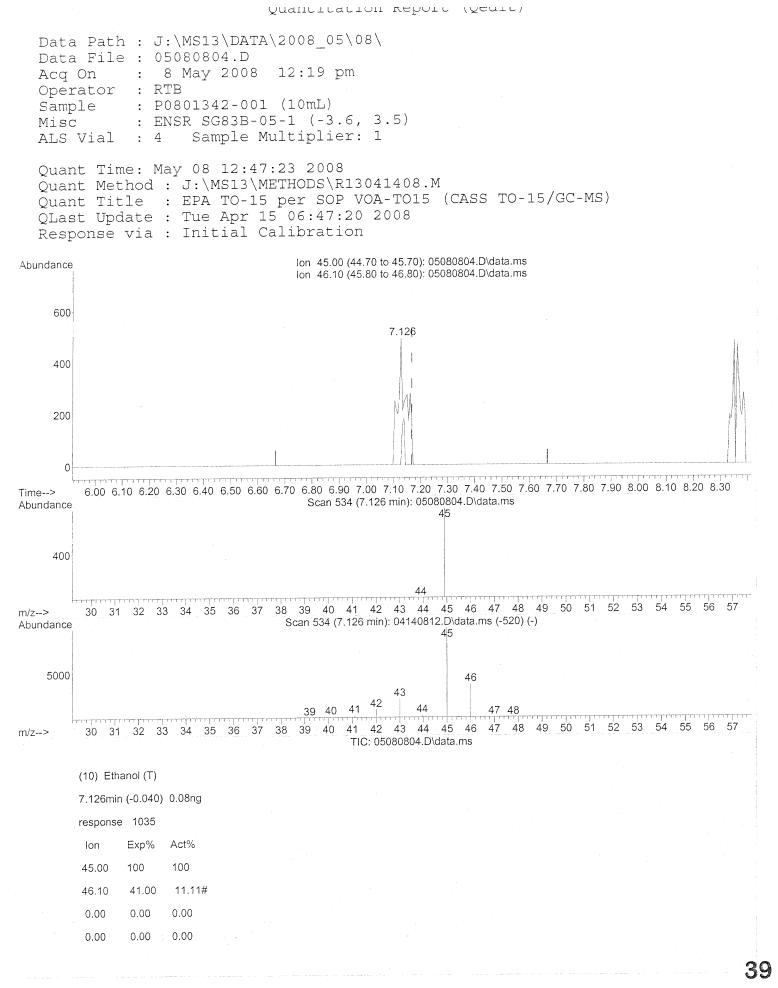
37

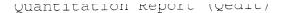
| Data Path : J:\MS13\DATA\2008_0 Data File : 05080804.D Acq On : 8 May 2008 12:19 Operator : RTB Sample : P0801342-001 (10mL) Misc : ENSR SG83B-05-1 (-3 ALS Vial : 4 Sample Multipli | pm .6, 3.5) | | | | | | | |
|--|---|--|---|--|----------------------|--|--|--|
| Quant Time: May 08 12:47:23 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration | | | | | | | | |
| Internal Standards | R.T. | QIon | Response | Conc Units | Dev(Min) | | | |
| <pre>80) alpha-Methylstyrene 81) 2-Ethyltoluene 82) 1,2,4-Trimethylbenzene 83) n-Decane 84) Benzyl Chloride 85) 1,3-Dichlorobenzene 86) 1,4-Dichlorobenzene 87) sec-Butylbenzene 88) p-Isopropyltoluene 89) 1,2,3-Trimethylbenzene 90) 1,2-Dichlorobenzene 91) d-Limonene 92) 1,2-Dibromo-3-Chloropr</pre> | 25.06 25.08 25.16 25.21 25.41 25.41 25.57 25.57 26.11 | 105 57 91 146 105 105 146 105 146 68 157 | 241 766 1849 1071 1051 829 3129 359 1171 1178 524 59 54 | N.D./ N.D. N.D./ N.D. N.D./ N.D. N.D./ N.D./ N.D./ N.D./ N.D./ | 96 | | | |
| 93) n-Undecane 94) 1,2,4-Trichlorobenzene 95) Naphthalene 96) n-Dodecane 97) Hexachloro-1,3-butadiene | 26.50 27.64 27.77 27.74 28.18 | 180 >128 57 | 1790 1427 10630 1927 81 | 0.054 ng 0.073 ng 0.168 ng 0.057 ng N.D. | 92 87 88 92 | | | |
| | | ~~~~~~~ | | | | | | |

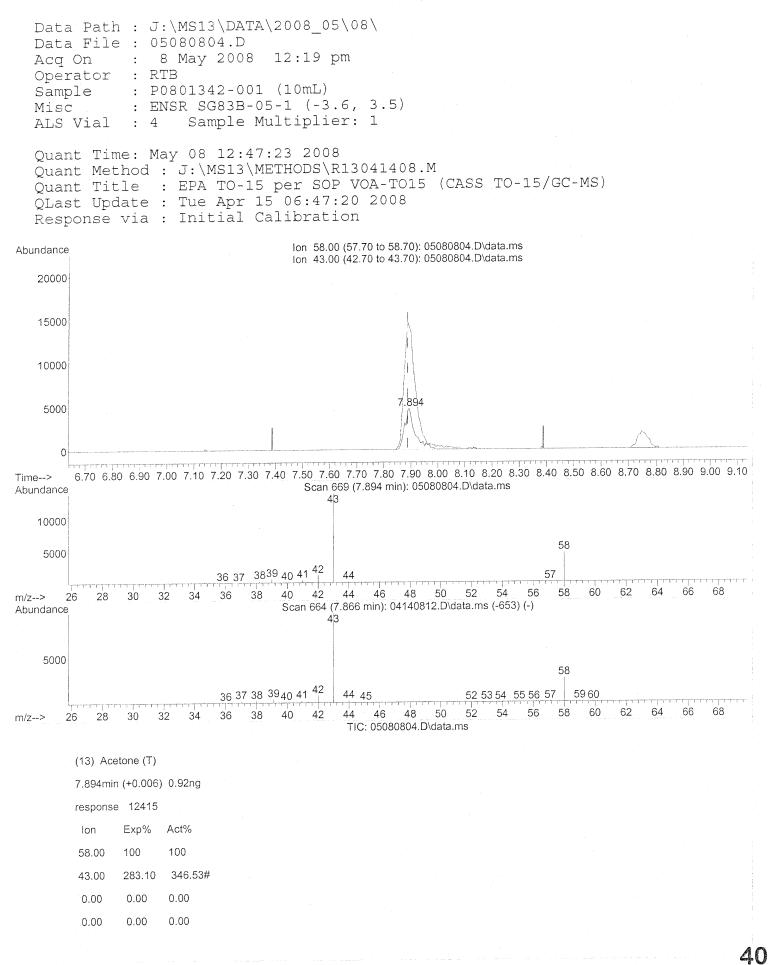
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Fostoglog

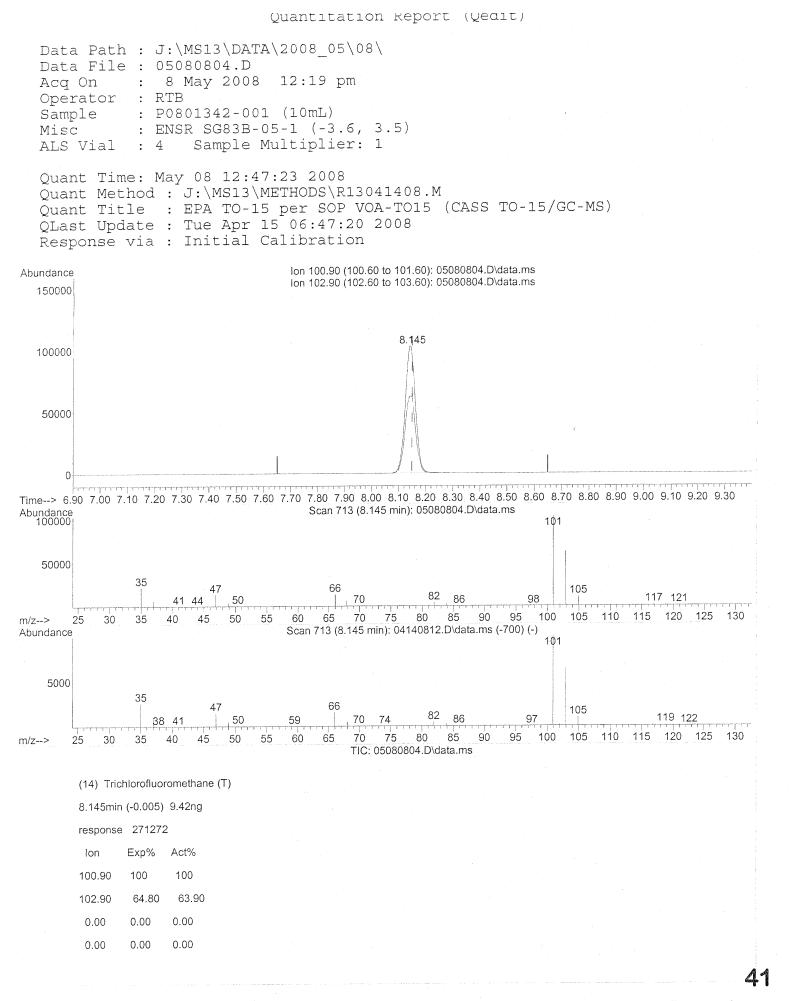
38

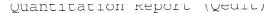


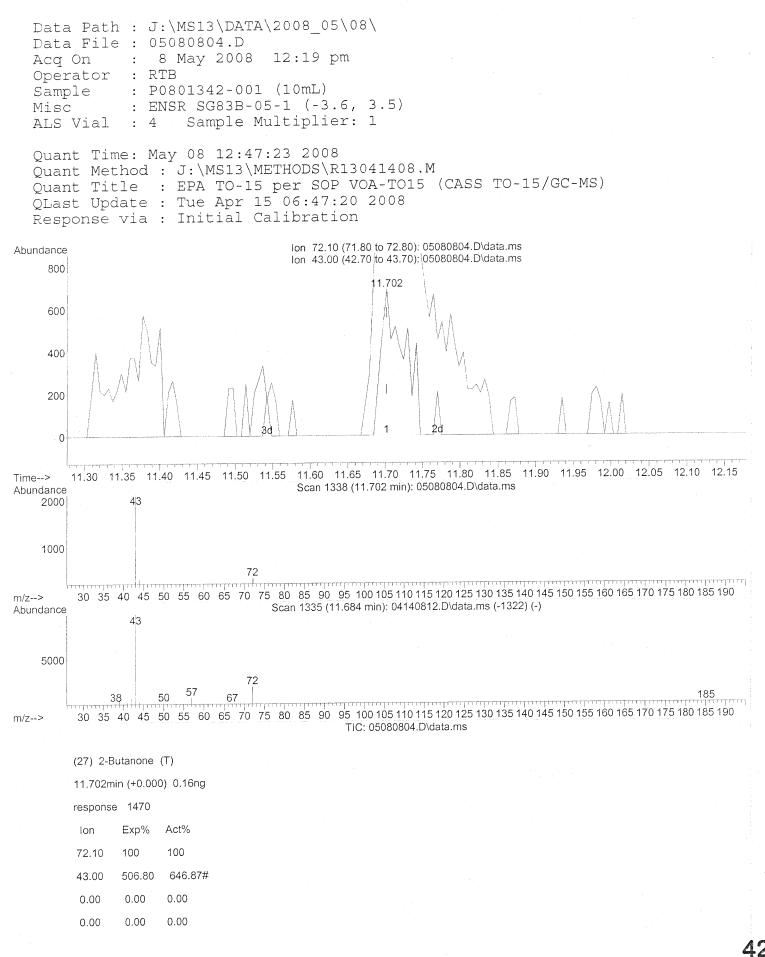




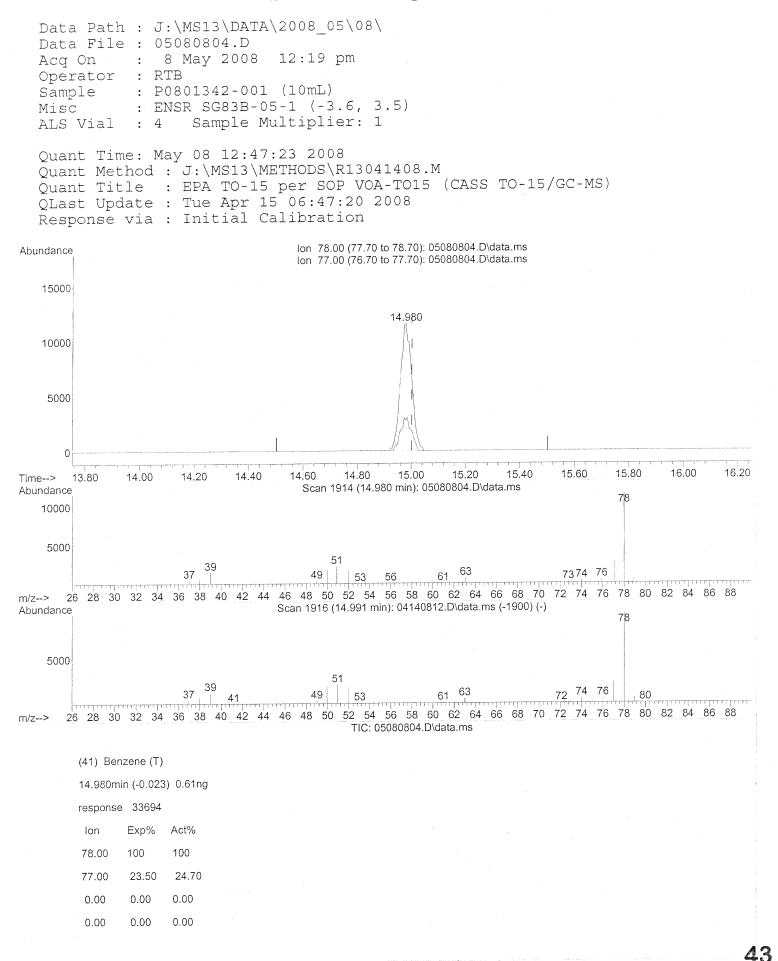
R13041408.M Thu May 08 16:23:52 2008



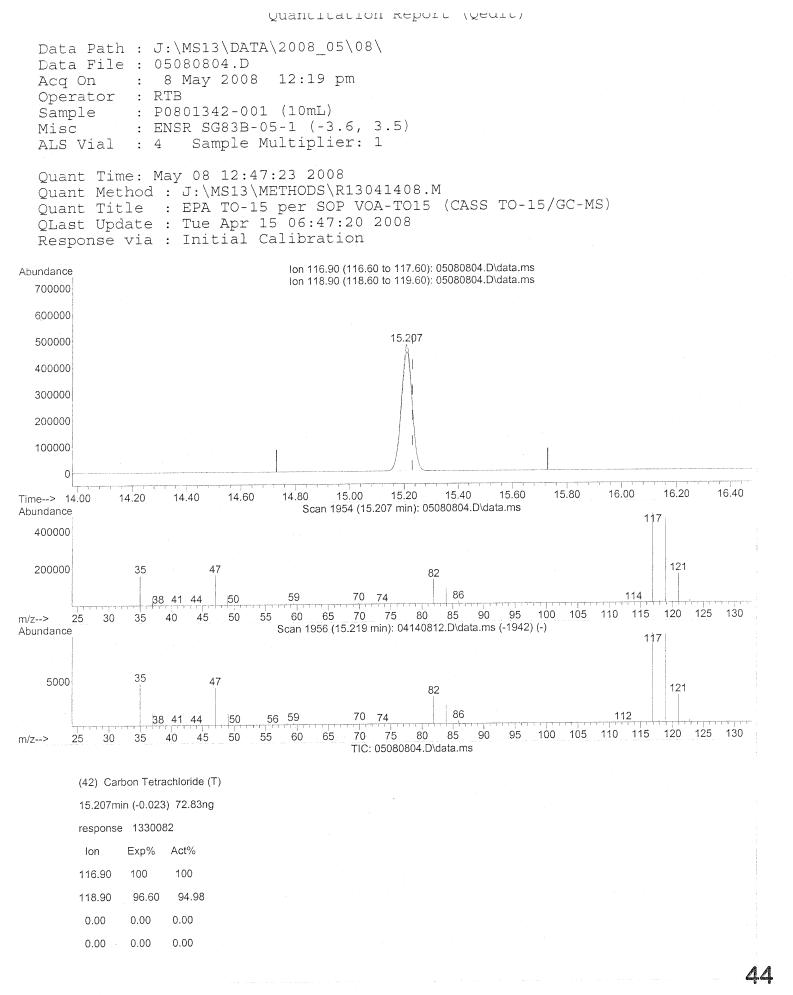


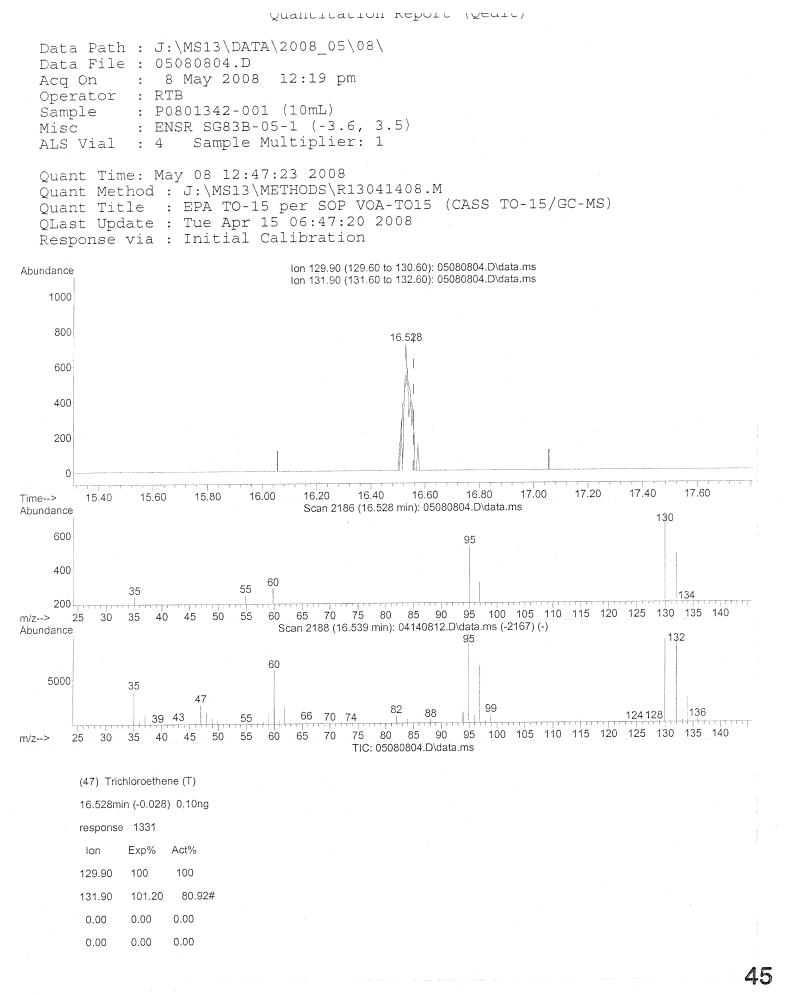


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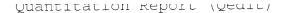


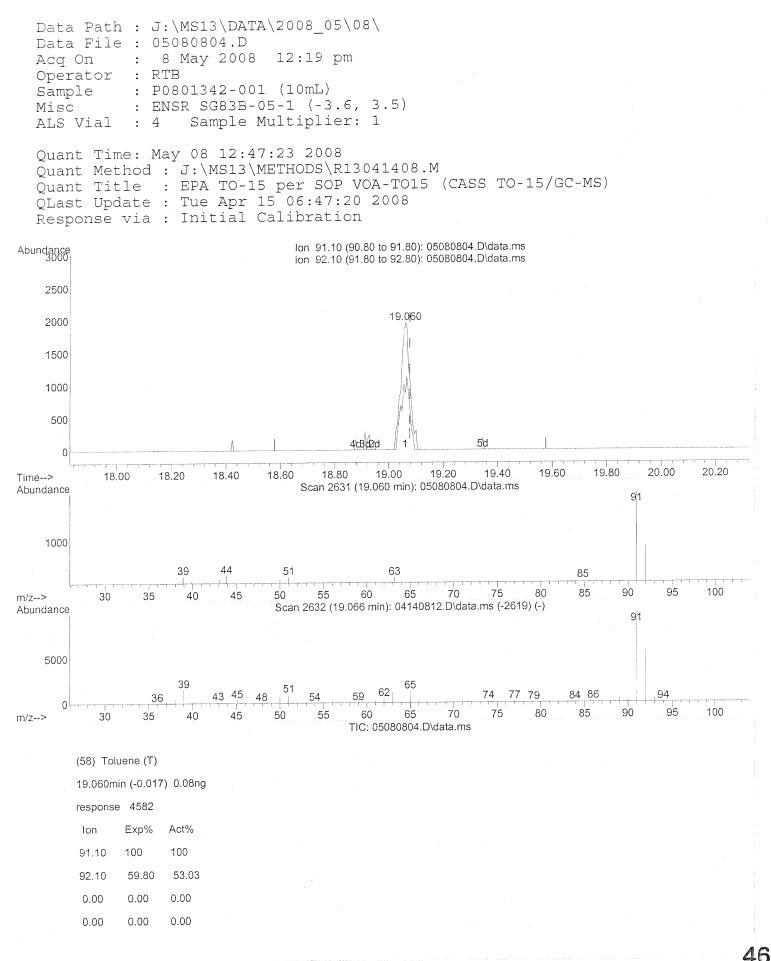
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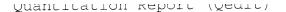


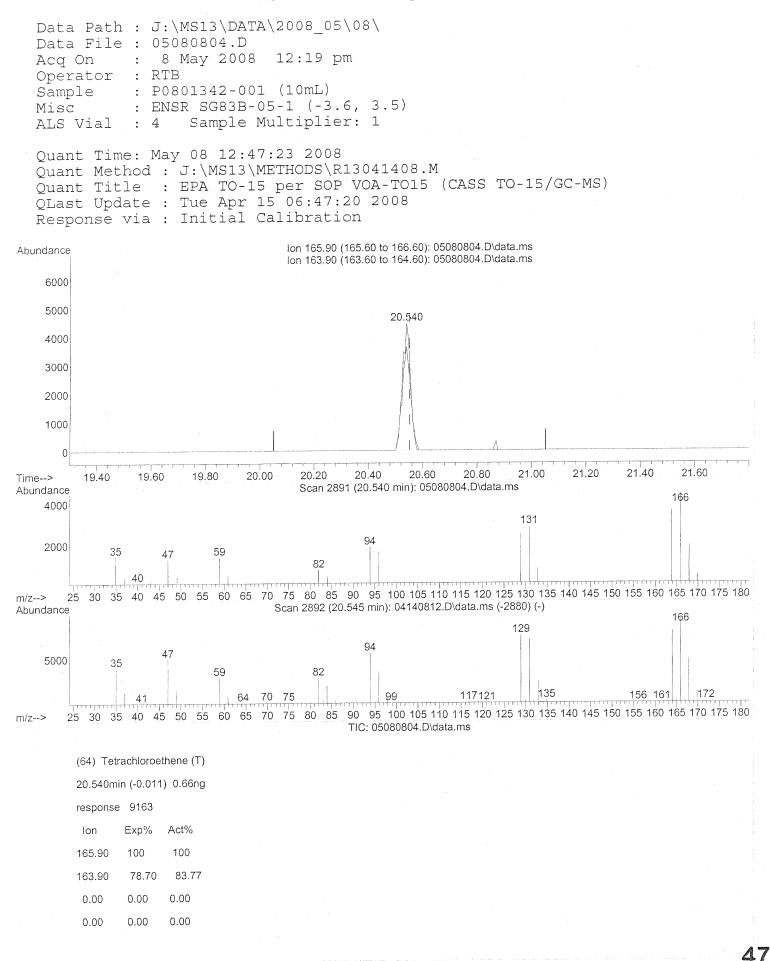
R13041408.M Fri May 09 14:52:13 2008



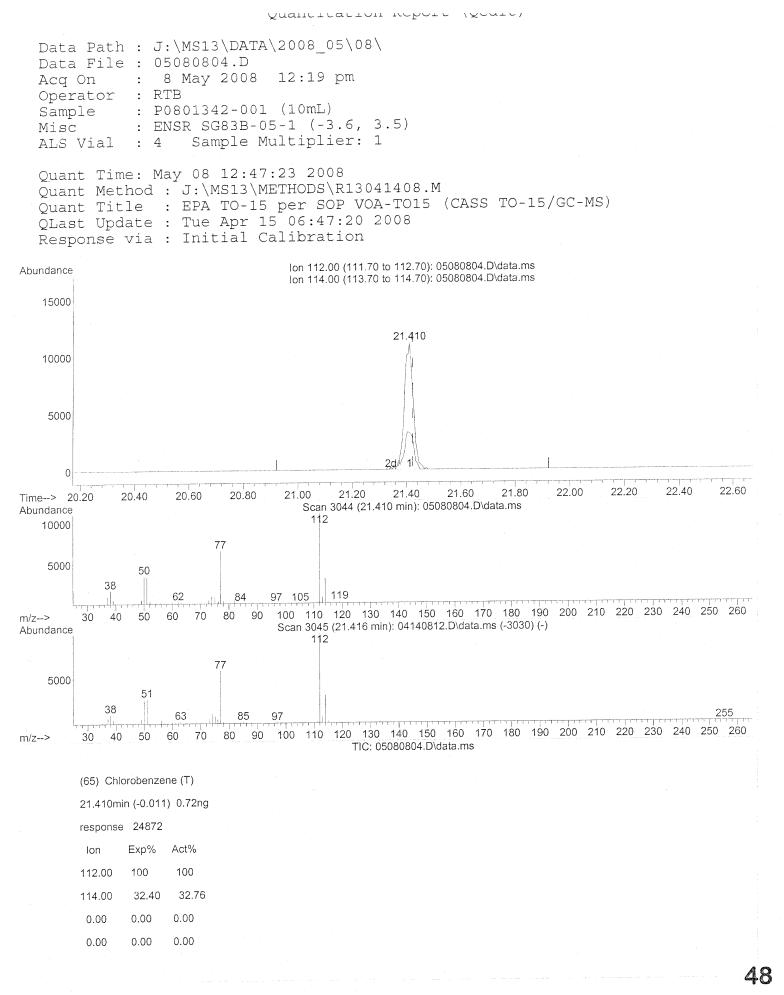


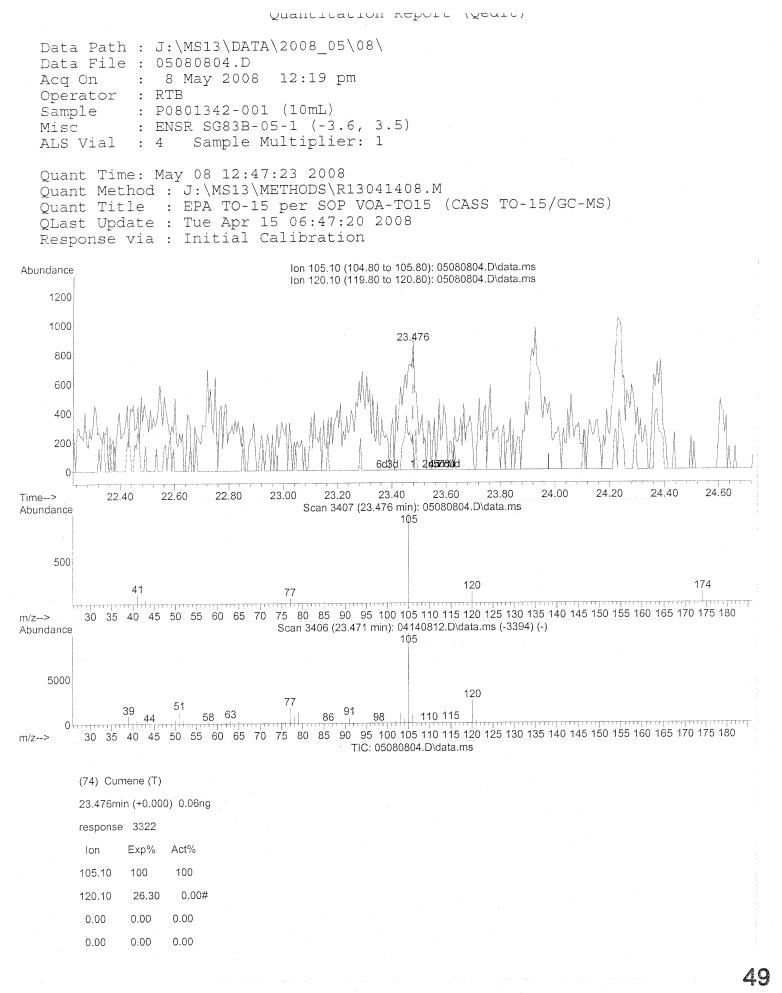
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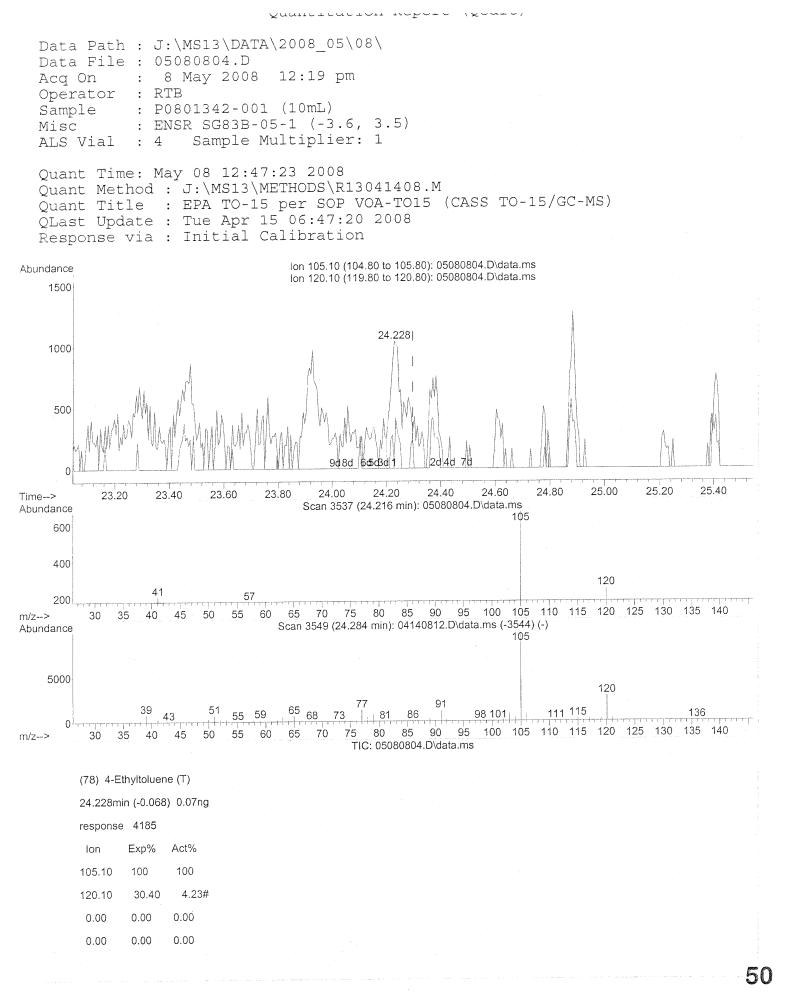


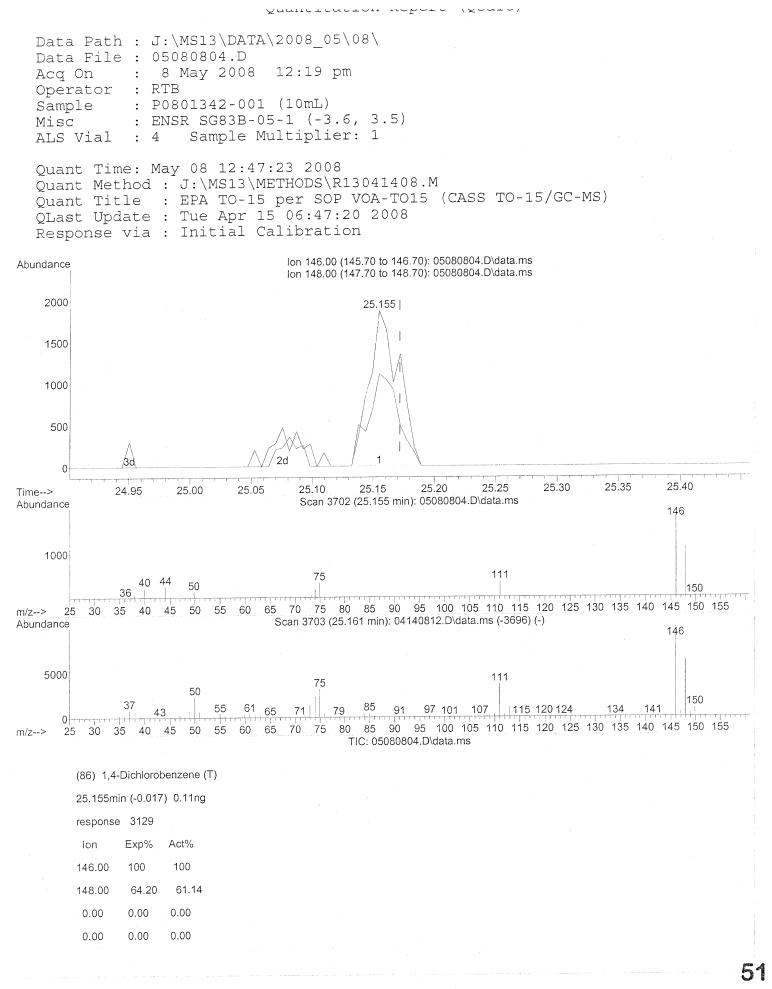


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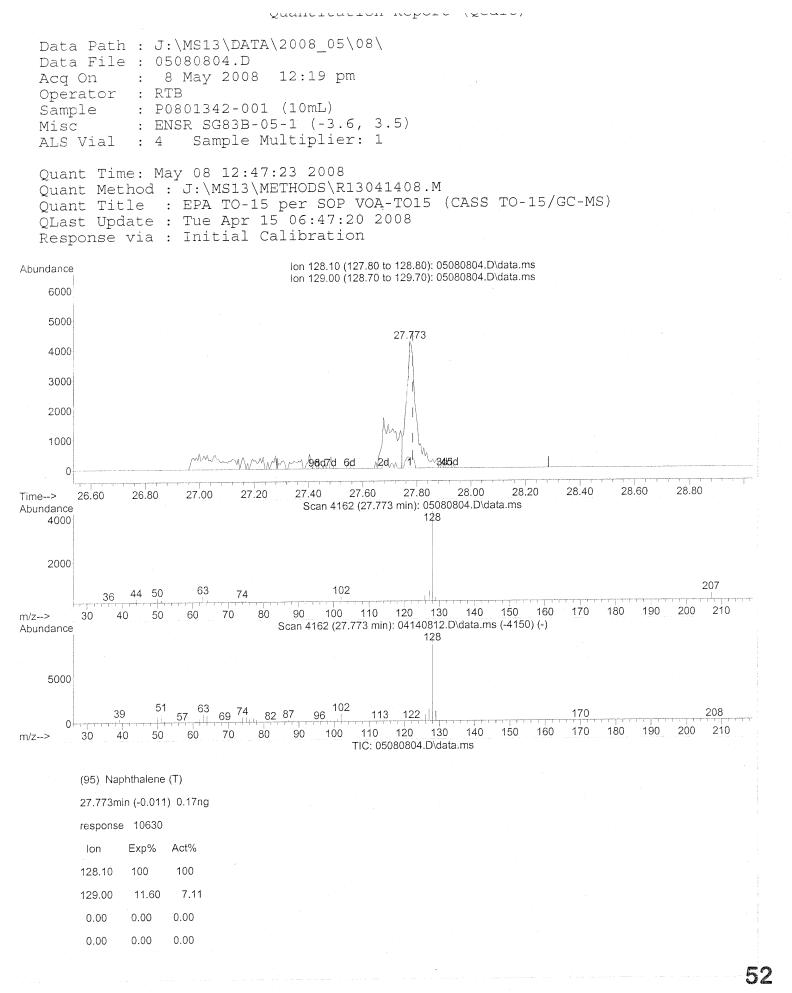


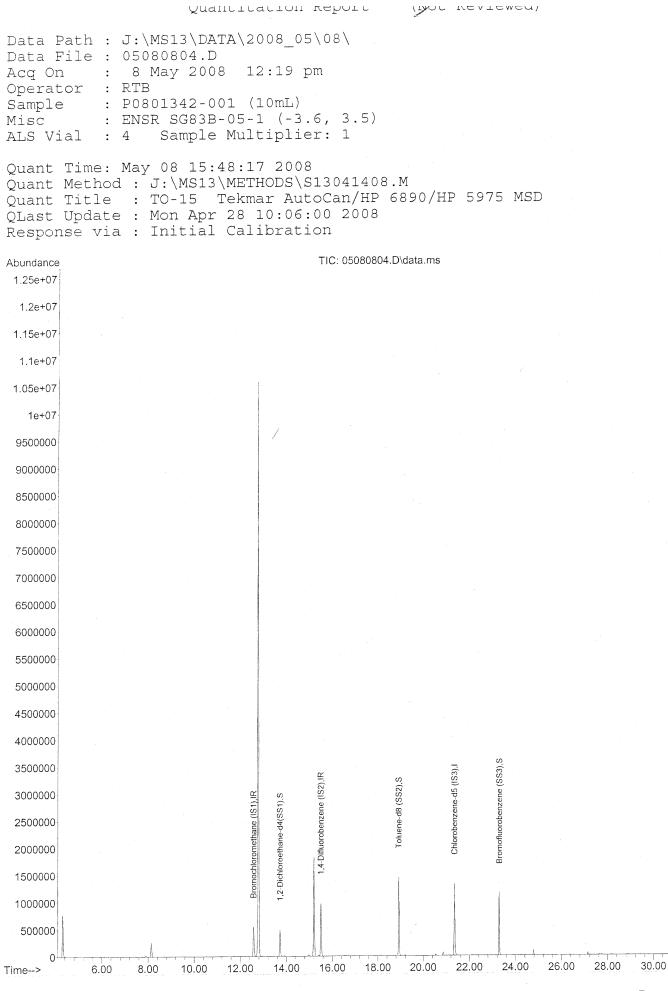






R13041408.M Thu May 08 16:32:02 2008





Page: 2

53

S13041408.M Thu May 08 15:54:35 2008

Quantitation Report (INCL INCATONICA) Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080804.D Acq On : 8 May 2008 12:19 pm Operator : RTB Sample : P0801342-001 (10mL) Misc : ENSR SG83B-05-1 (-3.6, 3.5) ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 15:48:17 2008 Quant Method : J:\MS13\METHODS\S13041408.M Ouant Title : TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD QLast Update : Mon Apr 28 10:06:00 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards _____ 1) Bromochloromethane (IS1) 12.58 130 242853 25.000 ng -0.03 3) 1,4-Difluorobenzene (IS2)15.51114104088925.000 ng-0.024) Chlorobenzene-d5 (IS3)21.358249464625.000 ng-0.01 System Monitoring Compounds 2) 1,2-Dichloroethane-d4(... 13.72 65 448165 23.011 ng -0.03 Recovery = 92.04% 🗸

 Spiked Amount
 25.000
 Recovery
 = 92.04%/

 5) Toluene-d8 (SS2)
 18.92
 98
 1129319
 25.472 ng
 -0.01

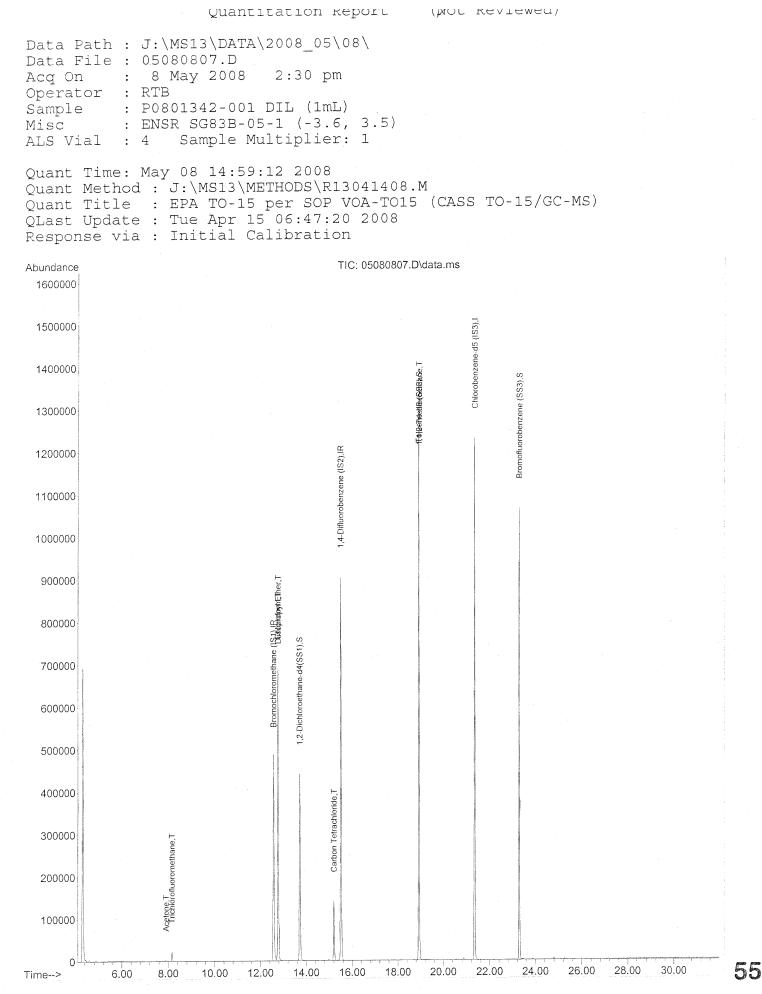
 Spiked Amount
 25.000
 Recovery
 = 101.88%,

 6) Bromofluorobenzene (SS3)
 23.29
 174
 384401
 25.194 ng
 0.00

 Recovery = 100.76% 🗸 Spiked Amount 25.000 Qvalue Target Compounds 7) tert-Butylbenzene24.881193108) n-Butylbenzene25.9191875 N.D. 🗸 N.D. / _______ _____

(#) = qualifier out of range (m) = manual integration (+) = signals summed

605/08/08



R13041408.M Thu May 08 15:05:30 2008

| Quantitation | keport | | DL KEVLEWED | 1) | | |
|---|------------------------------|-----------------------|---|--|-----------------------------|---|
| Data Path : J:\MS13\DATA\2008_05 Data File : 05080807.D Acq On : 8 May 2008 2:30 p Operator : RTB Sample : P0801342-001 DIL (1m Misc : ENSR SG83B-05-1 (-3. ALS Vial : 4 Sample Multiplie | | | | | | |
| Quant Time: May 08 14:59:12 2008 Quant Method : J:\MS13\METHODS\R Quant Title : EPA TO-15 per SOP QLast Update : Tue Apr 15 06:47: Response via : Initial Calibrati | 1304140 VOA-TO 20 2008 | 15 (CA | ASS TO-15/G | GC-MS) | | |
| Internal Standards | R.T. | QIon | Response | Conc Units | s Dev(Min |) |
| 1) Bromochloromethane (IS1) 37) 1,4-Difluorobenzene (IS2) 56) Chlorobenzene-d5 (IS3) | 15.51 | 130 114 | 219916 959711 | 25.000 ng 25.000 ng | -0.02 | |
| System Monitoring Compounds 33) 1,2-Dichloroethane-d4(Spiked Amount 25.000 57) Toluene-d8 (SS2) | 13.72 18.92 | 65 98 | Recove 1054023 | 23.432 ng ery = 93 26.346 ng | 3.72%∕ -0.01 | |
| Spiked Amount 25.000 73) Bromofluorobenzene (SS3) Spiked Amount 25.000 | | | Recove 345507 | ery = 105 | 0.00 | |
| | 0.00 0.00 0.00 | 85 50 135 62 | $\begin{array}{c} 0\\ 0\\ 0\\ 0\\ 0\\ 196\\ 0\\ 2047\\ 21764\\ 0\\ 0\\ 0\\ 51\\ 121\\ 75\\ 0\\ 295\\ 0\\ 56\\ 0\\ 56\\ 0\\ 67\\ 0\end{array}$ | N.D. N.D. N.D. N.D. N.D. N.D. N.D. N.D. | Qvalue # 74 98 # 1 | |
| 13041408.M Thu May 08 15:05:30 20 | 08 | | Ŧ | 5-108/00 | Page: 1 | |

| Quantitation | Report | | t Keviewed | .) | | |
|---|---|---|--|---|--------|------|
| Data Path : J:\MS13\DATA\2008_05 Data File : 05080807.D Acq On : 8 May 2008 2:30 p Operator : RTB Sample : P0801342-001 DIL (1r Misc : ENSR SG83B-05-1 (-3 ALS Vial : 4 Sample Multiplie | om nL) .6, 3.5) | / | | | | |
| Quant Time: May 08 14:59:12 2008 Quant Method : J:\MS13\METHODS\H Quant Title : EPA TO-15 per SOH QLast Update : Tue Apr 15 06:47 Response via : Initial Calibrat | R1304140 P VOA-TC :20 2008 | 15 (CA | SS TO-15/G | C-MS) | | |
| Internal Standards | R.T. | QIon | Response | Conc Units | Dev (| Min) |
| <pre>32) Chloroform 34) Tetrahydrofuran 35) Ethyl tert-Butyl Ether 36) 1,2-Dichloroethane 38) 1,1,1-Trichloroethane 39) Isopropyl Acetate 40) 1-Butanol 41) Benzene 42) Carbon Tetrachloride 43) Cyclohexane 44) tert-Amyl Methyl Ether 45) 1,2-Dichloropropane 46) Bromodichloromethane 47) Trichloroethene 48) 1,4-Dioxane 49) Isooctane 50) Methyl Methacrylate 51) n-Heptane 52) cis-1,3-Dichloropropene 53) 4-Methyl-2-pentanone 54) trans-1,3-Dichloropropene 55) 1,1,2-Trichloroethane 50) Dibromochloromethane 61) 1,2-Dibromoethane 62) Butyl Acetate 63) n-Octane 64) Tetrachloroethene 65) Chlorobenzene 66) Ethylbenzene 67) m- & p-Xylene 68) Bromoform 69) Styrene 70) o-Xylene 71) n-Nonane 72) 1,1,2,2-Tetrachloroethane 73) alpha-Pinene 76) alpha-Pinene 77) 3-Ethyltoluene</pre> | 0.00 0.00 0.00 0.00 0.00 18.93 19.06 0.00 20.34 0.00 20.34 0.00 20.55 21.40 22.13 0.00 22.75 23.16 0.00 23.48 0.00 24.24 | $\begin{array}{c} 87\\ 62\\ 97\\ 61\\ 56\\ 78\\ 117\\ 83\\ 63\\ 83\\ 130\\ 83\\ 130\\ 71\\ 75\\ 58\\ 75\\ 97\\ 129\\ 107\\ 43\\ 107\\ 43\\ 107\\ 43\\ 107\\ 166\\ 112\\ 91\\ 173\\ 104\\ 93\\ 105\\ 93\\ 105\\ 105\\ \end{array}$ | $\begin{array}{c} 0 \\ 0 \\ 74 \\ 0 \\ 0 \\ 0 \\ 3472 \\ 101995 \\ 438 \\ 0 \\ 0 \\ 0 \\ 0 \\ 0 \\ 0 \\ 0 \\ 0 \\ 0 \\ $ | 6.057 ng N.D. N.D. N.D. N.D. N.D. N.D. N.D. N.D | * * | 99 |
| 78) 4-Ethyltoluene 79) 1,3,5-Trimethylbenzene | 24.24 24.38 | 105 105 | 195 61 | N.D. N.D. | | 5 |
| 13041408.M Thu May 08 15:05:30 20 | 008 | | 7=0 | 5/68/68 | Page: | |

57

(NOL REVIEWEU) Quantitation keporu Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080807.D Acq On : 8 May 2008 2:30 pm Operator : RTB Sample : P0801342-001 DIL (1mL) Misc : ENSR SG83B-05-1 (-3.6, 3.5) ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 14:59:12 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards

 80) alpha-Methylstyrene
 0.00
 118
 0
 N.D.

 81) 2-Ethyltoluene
 24.38
 105
 61
 N.D.

 82) 1,2,4-Trimethylbenzene
 0.00
 105
 0
 N.D.

 83) n-Decane
 25.25
 57
 2095
 N.D.

 84) Benzyl Chloride
 0.00
 91
 0
 N.D.

 85) 1,3-Dichlorobenzene
 25.15
 146
 55
 N.D.

 86) 1,4-Dichlorobenzene
 25.41
 105
 187
 N.D.

 87) sec-Butylbenzene
 25.41
 105
 187
 N.D.

 89) 1,2,3-Trimethylbenzene
 25.41
 105
 187
 N.D.

 90) 1,2-Dichlorobenzene
 25.18
 146
 53
 N.D.

 90) 1,2-Dichlorobenzene
 25.41
 105
 187
 N.D.

 90) 1,2-Dichlorobenzene
 25.18
 146
 53
 N.D.

 91) d-Limonene
 0.00
 68
 0
 N.D.

 92) 1,2-Dibromo-3-Chloropr...
 0.00
 157
 0
 N.D.

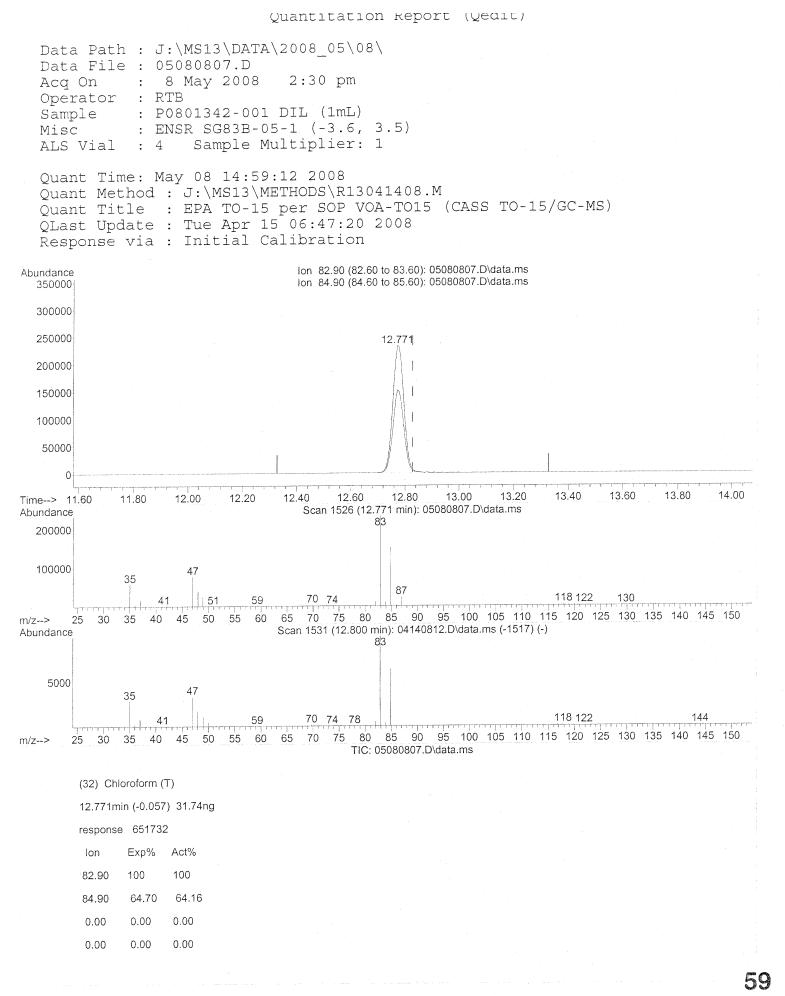
 93) n-Undecane
 26.50
 57
 196
 N.D.

 94) 1,2,4-T _____

(#) = qualifier out of range (m) = manual integration (+) = signals summed

R13041408.M Thu May 08 15:05:30 2008

705/08/08



R13041408.M Thu May 08 16:36:16 2008

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Page 1 of 3

-3.8

| Client: Client Sample ID: Client Project ID: | ENSR SG83B-05-3 Phase B Soil Gas / 04020-023-4311 |
|--|---|
| Test Code: | EPA TO-15 |
| Instrument ID: | Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 |
| Analyst: | Rusty Bravo |
| Sampling Media: | 6.0 L Summa Canister |

CAS Project ID: P0801342 CAS Sample ID: P0801342-002

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

Test Notes: Container ID:

SC00564

Initial Pressure (psig):

Final Pressure (psig): 3.5

Canister Dilution Factor: 1.67

| CAS # | Compound | Result | MRL | MDL | Result | MRL | MDL | Data |
|-----------|--|--------------|-------------|-------------|--------|------|------|-----------|
| | | <u>μg/m³</u> | $\mu g/m^3$ | $\mu g/m^3$ | ppbV | ppbV | ppbV | Qualifier |
| 75-71-8 | Dichlorodifluoromethane (CFC 12) | ND | 84 | 8.4 | ND | 17 | 1.7 | |
| 74-87-3 | Chloromethane | ND | 17 | 8.4 | ND | 8.1 | 4.0 | |
| 76-14-2 | 1,2-Dichloro-1,1,2,2- tetrafluoroethane (CFC 114) | ND | 84 | 8.4 | ND | 12 | 1.2 | |
| 75-01-4 | Vinyl Chloride | ND | 17 | 8.4 | ND | 6.5 | 3.3 | |
| 74-83-9 | Bromomethane | ND | 17 | 8.4 | ND | 4.3 | 2.2 | |
| 75-00-3 | Chloroethane | ND | 17 | 8.4 | ND | 6.3 | 3.2 | |
| 64-17-5 | Ethanol | ND | 840 | 8.4 | ND | 440 | 4.4 | |
| 67-64-1 | Acetone | 130 | 840 | 12 | 55 | 350 | 5.1 | J, B |
| 75-69-4 | Trichlorofluoromethane | 1,500 | 17 | 8.4 | 270 | 3.0 | 1.5 | |
| 107-13-1 | Acrylonitrile | ND | 84 | 12 | ND | 38 | 5.4 | |
| 75-35-4 | 1,1-Dichloroethene | ND | 17 | 8.4 | ND | 4.2 | 2.1 | |
| 75-65-0 | 2-Methyl-2-Propanol (tert-Butyl Alcohol) | ND | 84 | 12 | ND | 28 | 4.1 | |
| 75-09-2 | Methylene Chloride | ND | 84 | 8.4 | ND | 24 | 2.4 | |
| 107-05-1 | 3-Chloro-1-propene (Allyl Chloride) | ND | 17 | 8.4 | ND | 5.3 | 2.7 | |
| 76-13-1 | Trichlorotrifluoroethane | ND | 17 | 9.4 | ND | 2.2 | 1.2 | |
| 75-15-0 | Carbon Disulfide | ND | 84 | 20 | ND | 27 | 6.4 | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 17 | 8.4 | ND | 4.2 | 2.1 | |
| 75-34-3 | 1,1-Dichloroethane | ND | 17 | 8.4 | ND | 4.1 | 2.1 | |
| 1634-04-4 | Methyl tert-Butyl Ether | ND | 17 | 8.4 | ND | 4.6 | 2.3 | |
| 108-05-4 | Vinyl Acetate | ND | 840 | 27 | ND | 240 | 7.6 | |
| 78-93-3 | 2-Butanone (MEK) | 24 | 84 | 8.4 | 8.1 | 28 | 2.8 | J |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 17 | 8.4 | ND | 4.2 | 2.1 | |
| 108-20-3 | Diisopropyl Ether | ND | 84 | 9.9 | ND | 20 | 2.4 | |
| 67-66-3 | Chloroform | 49,000 | 17 | 9.9 | 10,000 | 3.4 | 2.0 | |

ND = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.



COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Page 2 of 3

| Client: | ENSR |
|---------------------------|-----------------------------------|
| Client Sample ID: | SG83B-05-3 |
| Client Project ID: | Phase B Soil Gas / 04020-023-4311 |
| | |
| Test Code: | EPA TO-15 |

6.0 L Summa Canister

CAS Project ID: P0801342 CAS Sample ID: P0801342-002

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

Instrument ID: Analyst: Sampling Media: Test Notes: Container ID:

SC00564

Rusty Bravo

Initial Pressure (psig): -3.8

Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13

Final Pressure (psig):

3.5

Canister Dilution Factor: 1.67

| CAS # | Compound | Result µg/m³ | MRL µg/m³ | MDL µg/m³ | Result ppbV | MRL ppbV | MDL ppbV | Data Qualifier |
|------------|---------------------------|-----------------|--------------|--------------|----------------|-------------|-------------|-------------------|
| 637-92-3 | Ethyl tert-Butyl Ether | ND | 84 | 8.5 | ND | 20 | 2.0 | |
| 107-06-2 | 1,2-Dichloroethane | ND | 17 | 8.4 | ND | 4.1 | 2.1 | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 17 | 8.4 | ND | 3.1 | 1.5 | |
| 71-43-2 | Benzene | 100 | 17 | 8.4 | 32 | 5.2 | 2.6 | |
| 56-23-5 | Carbon Tetrachloride | 12,000 | 17 | 8.4 | 1,900 | 2.7 | 1.3 | |
| 994-05-8 | tert-Amyl Methyl Ether | ND | 84 | 8.4 | ND | 20 | 2.0 | |
| 78-87-5 | 1,2-Dichloropropane | ND | 17 | 8.4 | ND | 3.6 | 1.8 | |
| 75-27-4 | Bromodichloromethane | ND | 17 | 8.4 | ND | 2.5 | 1.2 | |
| 79-01-6 | Trichloroethene | 16 | 17 | 8.4 | 3.0 | 3.1 | 1.6 | J |
| 123-91-1 | 1,4-Dioxane | ND | 84 | 10 | ND | 23 | 2.8 | |
| 80-62-6 | Methyl Methacrylate | ND | 84 | 13 | ND | 20 | 3.1 | |
| 142-82-5 | n-Heptane | ND | 84 | 11 | ND | 20 | 2.6 | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 84 | 8.7 | ND | 18 | 1.9 | |
| 108-10-1 | 4-Methyl-2-pentanone | ND | 84 | 9.4 | ND | 20 | 2.3 | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 84 | 11 | ND | 18 | 2.3 | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 17 | 8.4 | ND | 3.1 | 1.5 | |
| 108-88-3 | Toluene | 8.5 | 84 | 8.4 | 2.3 | 22 | 2.2 | J |
| 591-78-6 | 2-Hexanone | ND | 84 | 13 | ND | 20 | 3.1 | |
| 124-48-1 | Dibromochloromethane | ND | 17 | 11 | ND | 2.0 | 1.3 | |
| 106-93-4 | 1,2-Dibromoethane | ND | 17 | 9.0 | ND | 2.2 | 1.2 | |
| 111-65-9 | n-Octane | ND | 84 | 8.4 | ND | 18 | 1.8 | |
| 127-18-4 | Tetrachloroethene | 130 | 17 | 8.4 | 19 | 2.5 | 1.2 | |
| 108-90-7 | Chlorobenzene | 160 | 17 | 8.5 | 36 | 3.6 | 1.9 | ***** |

ND = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

Verified By: <u>Re-</u>Date: <u>519(08</u> TO15SCAN.XLT - Tronox - Henderson - PageNo.:

COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Page 3 of 3

Client: ENSR Client Sample ID: SG83B-05-3 Client Project ID: Phase B Soil Gas / 04020-023-4311

CAS Project ID: P0801342 CAS Sample ID: P0801342-002

3.5

Date Collected: 5/7/08 EPA TO-15 Test Code: Date Received: 5/8/08 Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 Instrument ID: Date Analyzed: 5/8/08 Analyst: Rusty Bravo Volume(s) Analyzed: 6.0 L Summa Canister 0.010 Liter(s) Sampling Media: 0.0010 Liter(s) Test Notes: SC00564 Container ID:

-3.8

Initial Pressure (psig):

Final Pressure (psig):

Canister Dilution Factor: 1.67

| | | Result | MRL | MDL | Result | MRL | MDL | Data |
|-------------|-------------------------------|-------------|-------|-------|--------|------|------|-----------|
| CAS # | Compound | $\mu g/m^3$ | μg/m³ | μg/m³ | ppbV | ppbV | ppbV | Qualifier |
| 100-41-4 | Ethylbenzene | ND | 84 | 10 | ND | 19 | 2.4 | |
| 179601-23-1 | m,p-Xylenes | ND | 84 | 22 | ND | 19 | 5.0 | |
| 75-25-2 | Bromoform | ND | 84 | 13 | ND | 8.1 | 1.2 | |
| 100-42-5 | Styrene | ND | 84 | 13 | ND | 20 | 3.0 | |
| 95-47-6 | o-Xylene | ND | 84 | 11 | ND | 19 | 2.4 | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 17 | 11 | ND | 2.4 | 1.6 | |
| 98-82-8 | Cumene | ND | 84 | 9.4 | ND | 17 | 1.9 | |
| 103-65-1 | n-Propylbenzene | ND | 84 | 8.7 | ND | 17 | 1.8 | |
| 622-96-8 | 4-Ethyltoluene | ND | 84 | 9.5 | ND | 17 | 1.9 | |
| 108-67-8 | 1,3,5-Trimethylbenzene | ND | 84 | 10 | ND | 17 | 2.0 | |
| 98-83-9 | alpha-Methylstyrene | ND | 84 | 12 | ND | 17 | 2.5 | |
| 95-63-6 | 1,2,4-Trimethylbenzene | ND | 84 | 12 | ND | 17 | 2.3 | |
| 100-44-7 | Benzyl Chloride | ND | 17 | 14 | ND | 3.2 | 2.8 | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 17 | 10 | ND | 2.8 | 1.7 | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 17 | 9.4 | ND | 2.8 | 1.6 | |
| 135-98-8 | sec-Butylbenzene | ND | 84 | 9.7 | ND | 15 | 1.8 | |
| 99-87-6 | 4-Isopropyltoluene (p-Cymene) | ND | 84 | 11 | ND | 15 | 2.0 | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 17 | 11 | ND | 2.8 | 1.8 | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 84 | 13 | ND | 8.6 | 1.3 | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 17 | 13 | ND | 2.3 | 1.7 | |
| 91-20-3 | Naphthalene | ND | 33 | 12 | ND | 6.4 | 2.4 | |
| 87-68-3 | Hexachlorobutadiene | ND | 17 | 15 | ND | 1.6 | 1.4 | |
| 98-06-6 | tert-Butylbenzene | ND | 33 | 8.4 | ND | 6.1 | 1.5 | |
| 104-51-8 | n-Butylbenzene | ND | 33 | 8.4 | ND | 6.1 | 1.5 | |

ND = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

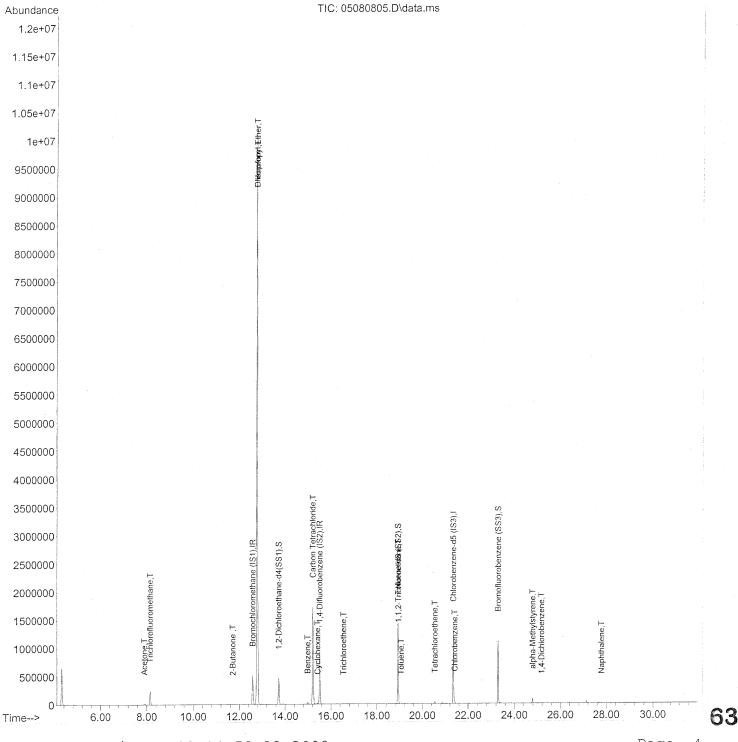
Verified By: <u>Rc-</u>

62 59108 TO15SCAN.XLT - Tronox - Henderson - PageNo.

Date:

| Data Path | : | J:\MS13\DATA\2008_05\08\ | |
|-----------|---|---------------------------|----|
| Data File | : | 05080805.D | |
| Acq On | : | 8 May 2008 1:04 pm | |
| Operator | | | |
| Sample | | P0801342-002 (10mL) | |
| Misc | | ENSR SG83B-05-3 (-3.8, 3. | 5) |
| ALS Vial | : | 4 Sample Multiplier: 1 | |
| | | | |

Quant Time: May 08 13:34:04 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration



R13041408.M Fri May 09 14:58:29 2008



Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080805.D Acq On : 8 May 2008 1:04 pm Operator : RTB Sample : P0801342-002 (10mL) Misc : ENSR SG83B-05-3 (-3.8, 3.5) ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 13:34:04 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards 1) Bromochloromethane (IS1)12.5813023194125.000 ng-0.0237) 1,4-Difluorobenzene (IS2)15.51114100403525.000 ng-0.02 21.35 82 467452 25.000 ng 0.00 56) Chlorobenzene-d5 (IS3) System Monitoring Compounds 13.72 65 424227 22.807 ng -0.03 33) 1,2-Dichloroethane-d4(... Recovery = 91.24% / Spiked Amount 25.000 18.93 98 1110032 26.493 ng 0.00 57) Toluene-d8 (SS2) Recovery = 105.96% / Spiked Amount 25.000 23.29 174 366213 25.399 ng 0.00 73) Bromofluorobenzene (SS3) Recovery = 101.60% Spiked Amount 25.000 Qvalue Target Compounds N.D. 332 4.83 42 2) Propene 353 N.D./ 4.98 85 3) Dichlorodifluoromethane 0 0.00 50 N.D./ 4) Chloromethane 0 N.D.V 0.00 135 5) Freon 114 0 N.D./ 0.00 62 6) Vinyl Chloride 0.00 54 0.00 94 0 N.D. 7) 1,3-Butadiene 0 0 N.D./ 8) Bromomethane N.D./ 0.00 64 9) Chloroethane N.D.V 7.15 45 234 10) Ethanol 7.46 41 413 N.D. 11) Acetonitrile 7<u>69</u>56 /0 7<u>89</u>58 9997 N.D. 12) Acrolein (0.779 ng # 65 13) Acetone 8.14 101 (9.171 ng 100 252177 14) Trichlorofluoromethane 8.37 45 462 N.D. 15) Isopropanol N.D./ 0.00 53 0 16) Acrylonitrile 0.00 96 0 N.D. 🗸 17) 1,1-Dichloroethene 9.32 59 9.36 84 53 N.D./ 18) tert-Butanol N.D. 🗸 414 19) Methylene Chloride 0 0.00 41 N.D./ 20) Allyl Chloride 0.00 151 0 N.D./ 21) Trichlorotrifluoroethane 9.77 76 22) Carbon Disulfide 1415 N.D./ N.D.V 0.00 61 0 23) trans-1,2-Dichloroethene 492 11.11 63 N.D./ 24) 1,1-Dichloroethane 0.00 73 0 N.D.V 25) Methyl tert-Butyl Ether 0.00 86 0 N.D.K 26) Vinyl Acetate 11.71)72 0.143 ng 60 1292 27) 2-Butanone N.D. 0 -0.00 61 28) cis-1,2-Dichloroethene 12.78 87 12.78 61 959634 81.174 ng 1 29) Diisopropyl Ether 60 N.D. 30) Ethyl Acetate N.D. 12.70 57 632 31) n-Hexane

R13041408.M Fri May 09 14:58:29 2008

05/09/08

64

| Data Acq C Opera Sampl Misc | | pm .8, 3.5) | | | | | |
|---|--|---|------------|-------------|-----------------------------|-------------------|----------|
| Quant Quant OLast | Time: May 08 13:34:04 200 Method : J:\MS13\METHODS\ Title : EPA TO-15 per SO Update : Tue Apr 15 06:47 onse via : Initial Calibrat | R1304140 P VOA-TC :20 2008 ion |)15 (C | | | | |
| Inte | rnal Standards | R.T. | QIon | Respons | e Conc Unit | s Dev | /(Min) |
| | Chloroform | 12.78 | 83 | 9063536 | 4 <u>18</u> .5 <u>34</u> ng | Ea a | 100 |
| 34) | Tetrahydrofuran | 0.00 | 72 | 0 | N.D. | | |
| | Ethyl tert-Butyl Ether | 0.00 | | 0 306 | N.D. | | |
| | 1,2-Dichloroethane 1,1,1-Trichloroethane | 13.73 0.00 | | 306 0 | N.D. 🗸 N.D. 🗸 | | |
| | Isopropyl Acetate | 14.97 | | 74 | N.D. | | |
| | 1-Butanol | 14.97 | | 84 | N.D. | | |
| 41) | Benzene | 14.98 | | 32830 | | > | 100 |
| | Carbon Tetrachloride | 15.21 | | 1267210 | (71.934 ng | > # | 99 84 |
| | Cyclohexane | 15.42 | 84 73 | 6181 0 | 0.314 ng N.D. | ++ | 04 |
| | tert-Amyl Methyl Ether 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D.V | | |
| | Bromodichloromethane | 16,48 | | 623 | N.D.V | | |
| | Trichloroethene | 16.55 | | 1273 | 0.097 ng | > # | 77 |
| | 1,4-Dioxane | 0.00 | 88 | 0 | N.D.V | | |
| | Isooctane | 0.00 | 57 | 0 | N.D. | | |
| | Methyl Methacrylate | 0.00 | | 0 214 | N.D. | | |
| | n-Heptane | 16.98 0.00 | 71 75 | 214 | N.D. 🗸 N.D. 🗸 | | |
| | cis-1,3-Dichloropropene 4-Methyl-2-pentanone | 0.00 | 58 | 0 | N.D. | | |
| | trans-1,3-Dichloropropene | | | 0 | / | ID | |
| | 1,1,2-Trichloroethane | 18_94 | 97 | 100833 | 7.861=19 | 14-# | 8 |
| | Toluene | (19.06 | > 91 | 2698 | 0.051 ng | \mathcal{Y}_{n} | 82 |
| | 2-Hexanone | 19.54 | 43 | 54 | N.D. | | |
| | Dibromochloromethane | 0.00 | 129 | 0 | N.D./ N.D./ | | |
| | 1,2-Dibromoethane | 0.00 20.35 | 107 43 | 743 | N.D. | | |
| | Butyl Acetate n-Octane | 20.35 | _ 57 | 159 | N.D. | | |
| | Tetrachloroethene | 20.55 | $>_{166}$ | 10046 | 0.762 ng | > | 100 |
| | Chlorobenzene | 21.41 | ∕112 | 32038 | 0.983 ng | \sum | 100 |
| | Ethylbenzene | 21.89 | 91 | 817 | N.D. | | |
| | m- & p-Xylene | 22.17 | 91 | 82 | N.D.V | | |
| | Bromoform | 0.00 22.58 | 173 104 | 0 109 | N.D. 🗸 N.D. 🗸 | | |
| | Styrene o-Xylene | 22.58 | 104 91 | 536 | N.D.V | | |
| | n-Nonane | 22.98 | 43 | 685 | N.D. | | |
| | 1,1,2,2-Tetrachloroethane | | 83 | 78 | N.D./ | | |
| | Cumene | 23.48 | | 885 | N.D. 🗸 | | |
| | alpha-Pinene | 0.00 | 93 | 0 | N.D. | | |
| | n-Propylbenzene | 24.12 | 91 | 159 1668 | N.D. N.D. | | |
| | 3-Ethyltoluene | 24.23 24.30 | | 475 | N.D.V | | |
| | 4-Ethyltoluene 1,3,5-Trimethylbenzene | 24.30 | 105 | 589 | N.D. | | C |
| 121 | $\pm, 5, 5$ $\pm \pm 1$ meetry ± 5 cm2 cm2 | _1.00 | | | RII- | | 6 |
| 130414 | 08.M Fri May 09 14:58:29 2 | 008 | | | 205109108 | Page | 2: 2 |
| | | | | | | | |

⁶⁵

Data Path : J:\MS13\DATA\2008_05\08\ Data File : 05080805.D Acq On : 8 May 2008 1:04 pm Operator : RTB Sample : P0801342-002 (10mL) Misc : ENSR SG83B-05-3 (-3.8, 3.5) ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 13:34:04 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards

 80) alpha-Methylstyrene
 24.79
 118
 1393
 0.055 ng
 #
 27

 81) 2-Ethyltoluene
 24.60
 105
 237
 N.D.

 82) 1,2,4-Trimethylbenzene
 24.89
 105
 1088
 N.D.

 83) n-Decane
 24.98
 57
 244
 N.D.

 84) Benzyl Chloride
 25.05
 91
 221
 N.D.

 85) 1,3-Dichlorobenzene
 25.08
 146
 138
 N.D.

 86) 1,4-Dichlorobenzene
 25.17
 146
 1203
 0.043 ng
 82

 87) sec-Butylbenzene
 25.20
 105
 62
 N.D.
 82

 89) 1,2,3-Trimethylbenzene
 25.41
 119
 600
 N.D.
 89
 1,2-Dichlorobenzene
 25.59
 146
 128
 N.D.

 90) 1,2-Dichlorobenzene
 25.59
 146
 128
 N.D.
 91
 1.2-Dibromo-3-Chloropr...
 0.00
 157
 0
 N.D.

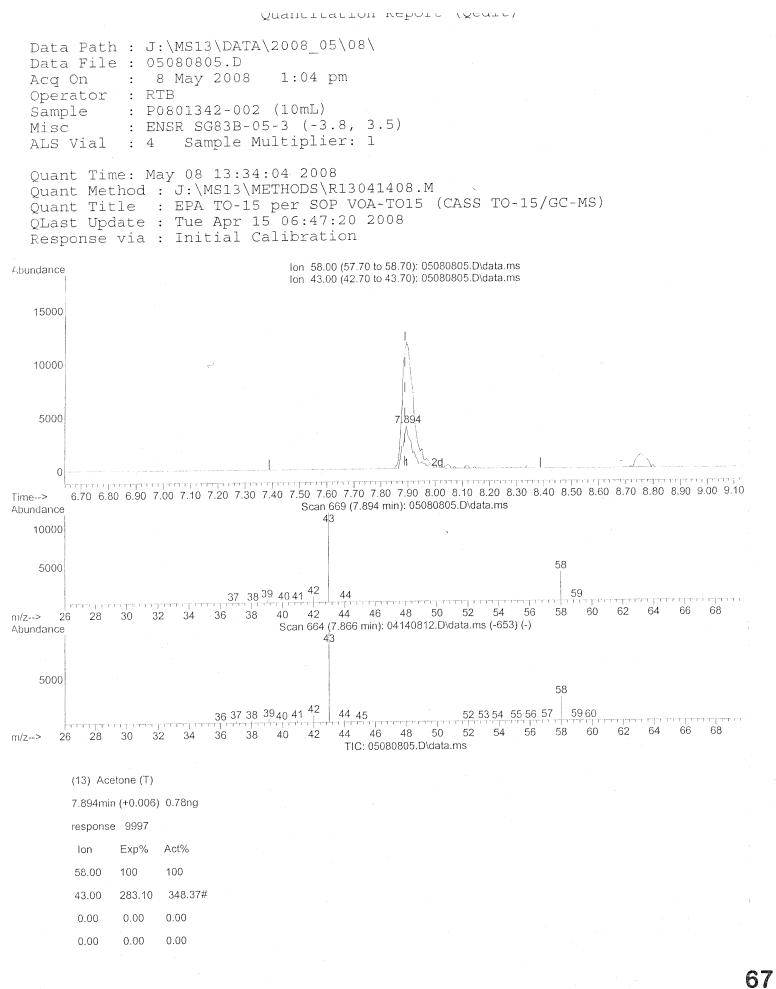
 91) d-Limonene
 26.51
 57
 915
 N.D.
 91
 1,2,4-Trichlorobenzene
 27.64
 180
 229
 N.D.

 94) 1,2,4-Trichlorobenzene
 27.64
 _____ _____

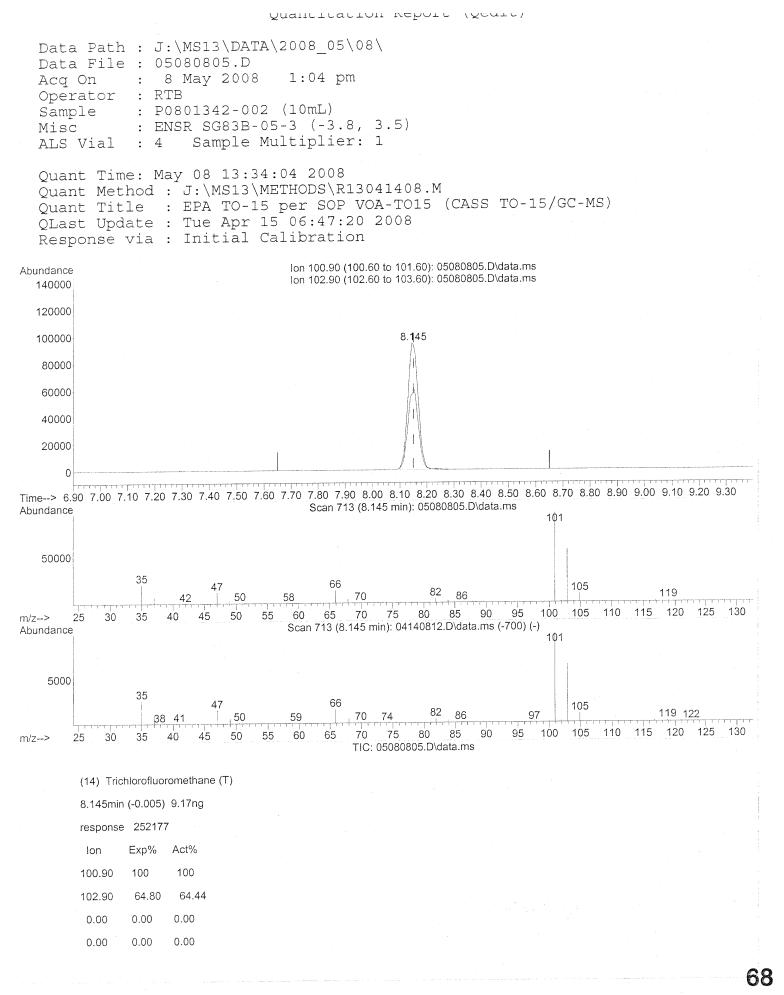
(#) = qualifier out of range (m) = manual integration (+) = signals summed

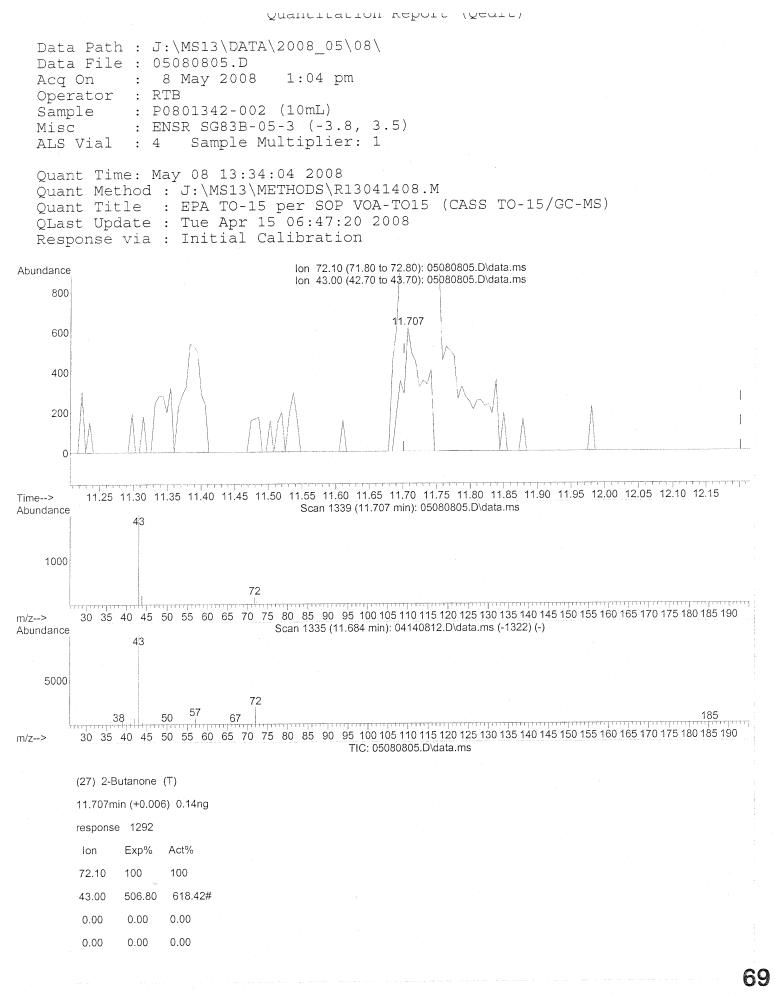
505/09/15

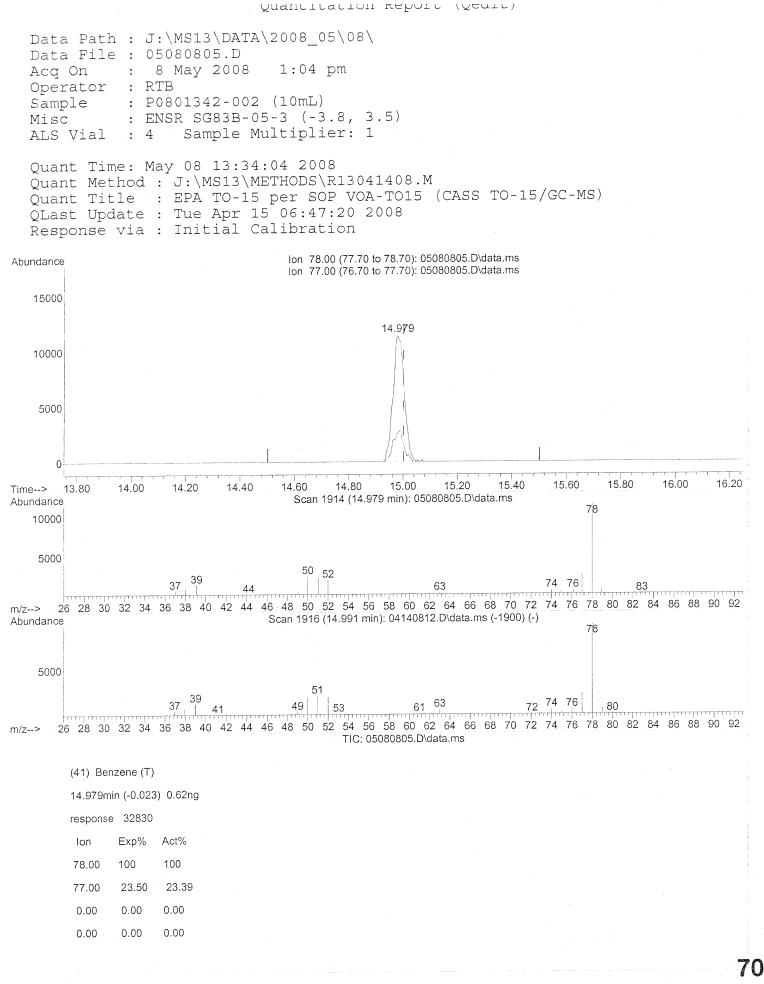
R13041408.M Fri May 09 14:58:29 2008

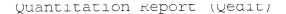


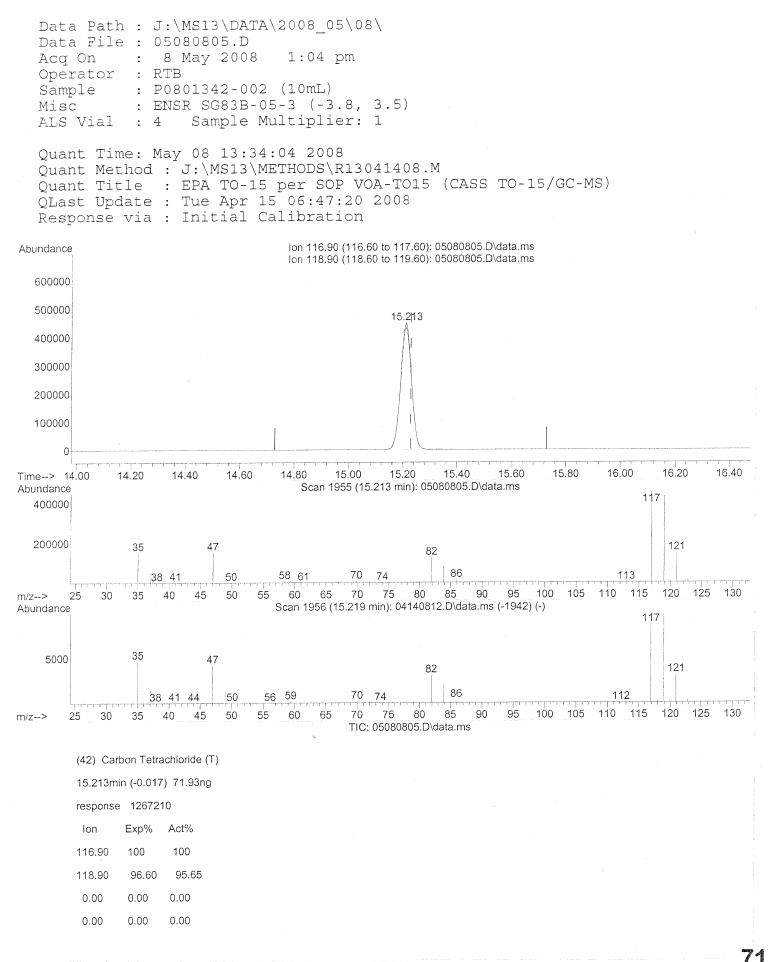
R13041408.M Thu May 08 16:38:23 2008

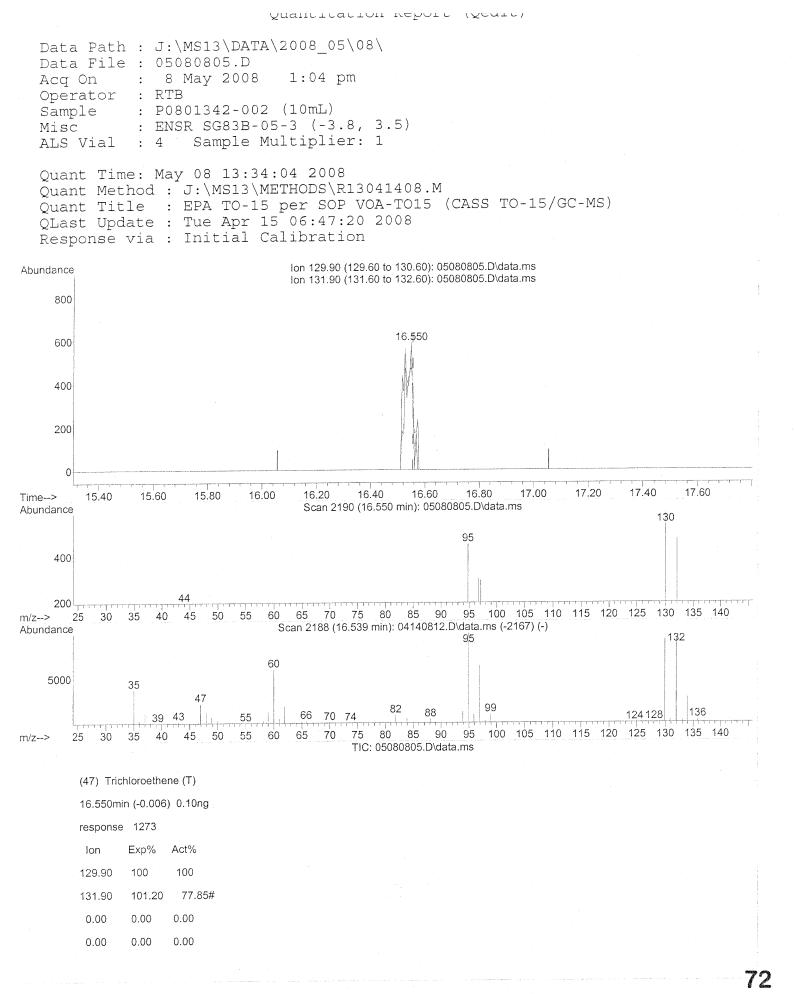




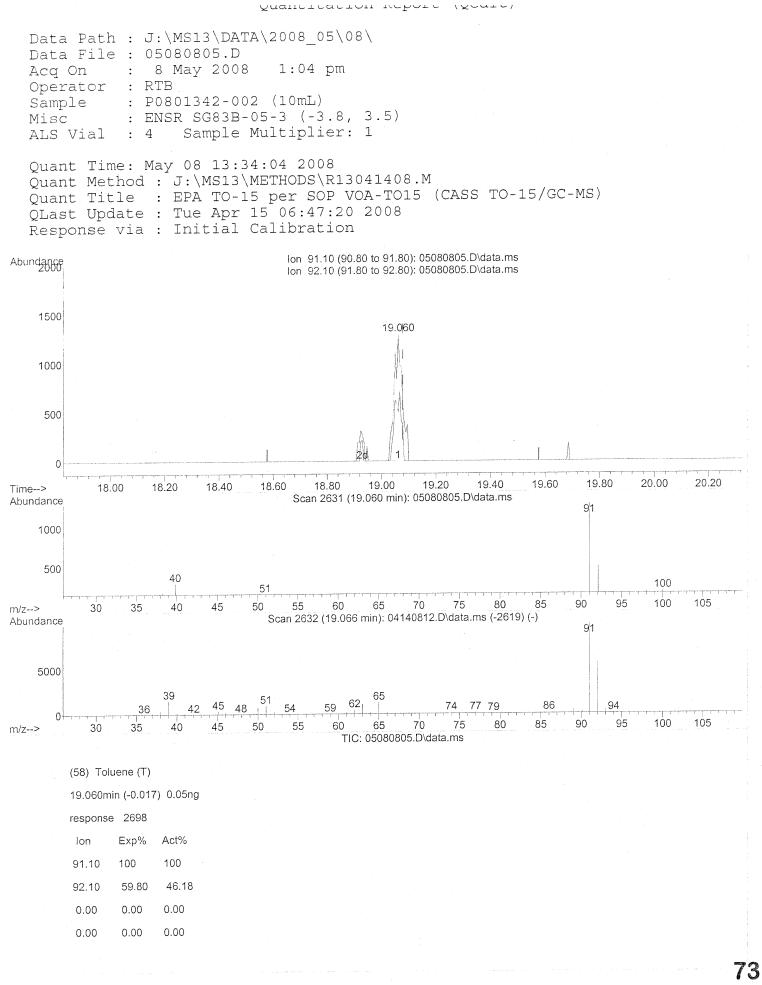


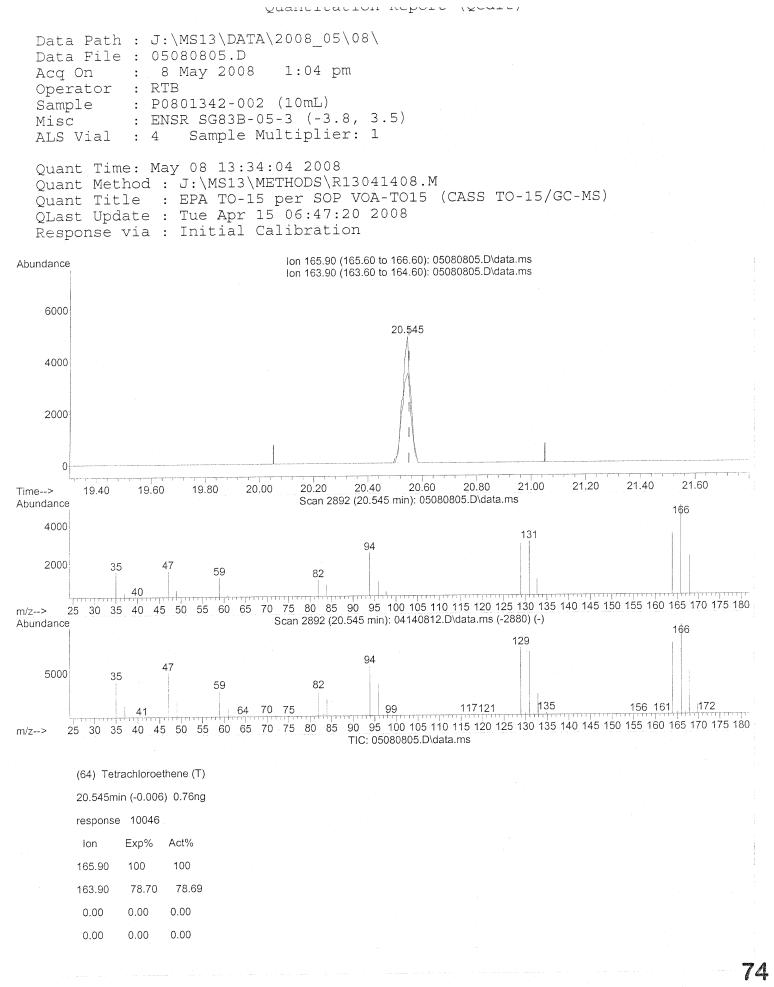


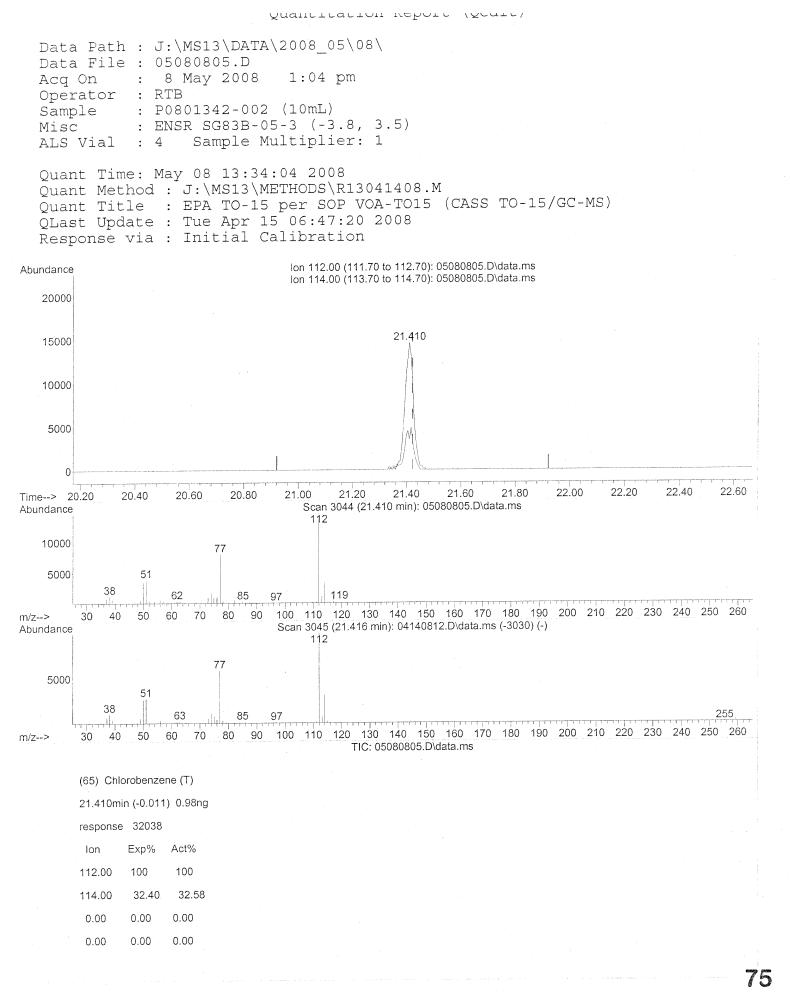


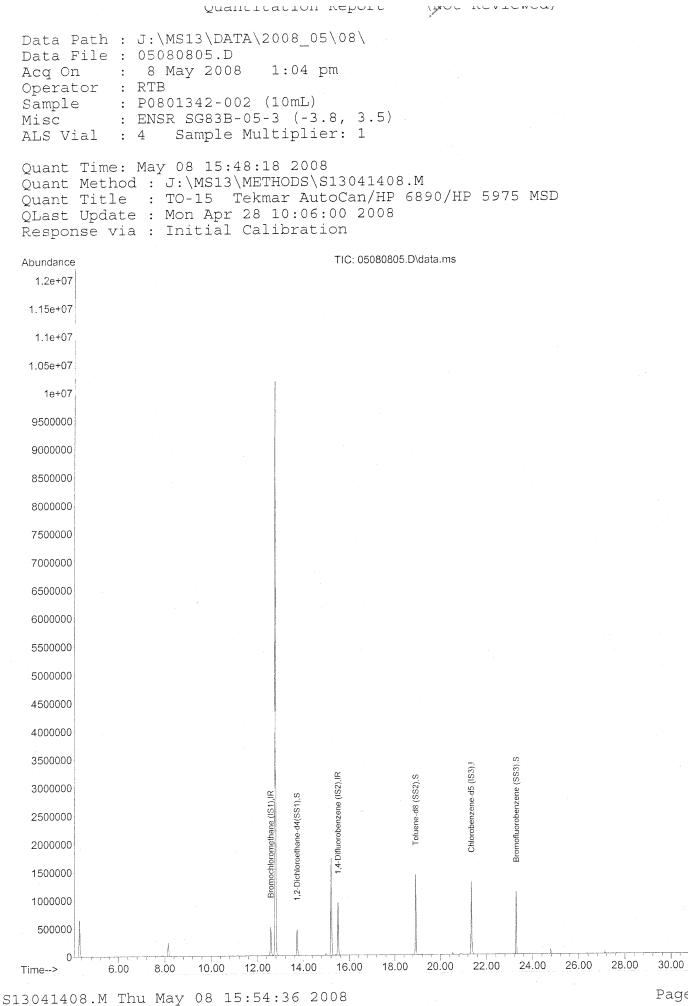


R13041408.M Fri May 09 15:19:21 2008









Page: 2

Quantitation Report

(INOL REVIEWEN)

Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080805.D Acq On : 8 May 2008 1:04 pm Operator : RTB Sample : P0801342-002 (10mL) Misc : ENSR SG83B-05-3 (-3.8, 3.5) ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 15:48:18 2008 Quant Method : J:\MS13\METHODS\S13041408.M Quant Title : TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD OLast Update : Mon Apr 28 10:06:00 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards 1) Bromochloromethane (IS1)12.5813023194125.000 ng-0.023) 1,4-Difluorobenzene (IS2)15.51114100403525.000 ng-0.024) Chlorobenzene-d5 (IS3)21.358246745225.000 ng0.00 System Monitoring Compounds 2) 1,2-Dichloroethane-d4(... 13.72 65 424227 22.807 ng -0.03

 Spiked Amount
 25.000
 Recovery = 91.24%

 5) Toluene-d8 (SS2)
 18.93
 98
 1110032
 26.493 ng
 0.00

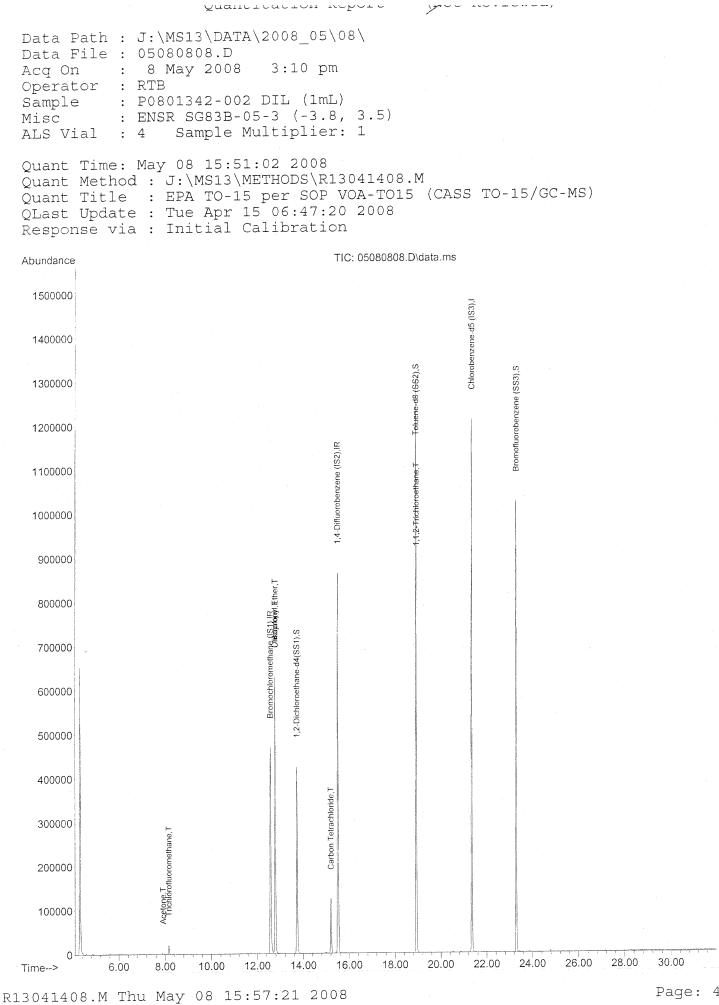
 Spiked Amount
 25.000
 Recovery = 105.96%
 105.96%

 6) Bromofluorobenzene (SS3)
 23.29
 174
 366213
 25.399 ng
 0.00

 Recovery = 101.60% Spiked Amount 25.000 Qvalue Target Compounds 7) tert-Butylbenzene24.8811952N.D.↓8) n-Butylbenzene25.919152N.D.↓

(#) = qualifier out of range (m) = manual integration (+) = signals summed

105/08/08



Page: 4

| Quality cattor | NCPUL C | \-7 | | ur / | |
|---|------------------------------|---|---|------------------------------------|---------------------------|
| Data Path : J:\MS13\DATA\2008_05 Data File : 05080808.D Acq On : 8 May 2008 3:10 p Operator : RTB Sample : P0801342-002 DIL (1m Misc : ENSR SG83B-05-3 (-3. ALS Vial : 4 Sample Multiplie | m L) 8, 3.5) | | | | |
| Quant Time: May 08 15:51:02 2008 Quant Method : J:\MS13\METHODS\R Quant Title : EPA TO-15 per SOP QLast Update : Tue Apr 15 06:47: Response via : Initial Calibratio | 1304140 VOA-TO 20 2008 | 15 (C2 | ASS TO-15/0 | GC-MS) | |
| Internal Standards | R.T. | QIon | Response | Conc Units | Dev(Min) |
| 1) Bromochloromethane (IS1) 37) 1,4-Difluorobenzene (IS2) 56) Chlorobenzene-d5 (IS3) | 15.51 | 114 | 929659 | 25.000 ng | -0.02 |
| System Monitoring Compounds 33) 1,2-Dichloroethane-d4(Spiked Amount 25.000 | | | Recove | 23.206 ng ery = 92 25.854 ng | .84% |
| 57) Toluene-d8 (SS2) Spiked Amount 25.000 73) Bromofluorobenzene (SS3) Spiked Amount 25.000 | | | Recove 334113 | ery = 103 | .40%/ 0,00 |
| <pre>Target Compounds 2) Propene 3) Dichlorodifluoromethane 4) Chloromethane 5) Freon 114 6) Vinyl Chloride 7) 1,3-Butadiene 8) Bromomethane 9) Chloroethane 10) Ethanol 11) Acetonitrile 12) Acrolein 13) Acetone 14) Trichlorofluoromethane 15) Isopropanol 16) Acrylonitrile 17) 1,1-Dichloroethene 18) tert-Butanol 19) Methylene Chloride 20) Allyl Chloride 21) Trichlorotrifluoroethane 22) Carbon Disulfide 23) trans-1,2-Dichloroethene 24) 1,1-Dichloroethane 25) Methyl tert-Butyl Ether 26) Vinyl Acetate 27) 2-Butanone 28) cis-1,2-Dichloroethene 29) Diisopropyl Ether 30) Ethyl Acetate</pre> | 0.00 9.37 | 85 50 135 62 54 94 45 41 58 101 53 69 41 156 63 73 62 61 87 | 200 0 0 0 0 0 0 0 0 2566 20884 0 0 0 0 0 0 0 220 0 0 0 0 0 0 0 0 0 0 0 0 0 | N.D. N.D. N.D. N.D. | Qvalue 97 98 # 1 |
| 30) Ethyl Acetate 31) n-Hexane 13041408.M Thu May 08 15:57:20 20 | 12.69 | | 53 | N.D. | 7 Page: 1 |

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| QUAILILALION | керотс | (1) | IL REVIEW | | |
|---|--|-------------------------------|-------------------------------------|--|---------------------|
| Data Path : J:\MS13\DATA\2008_0 Data File : 05080808.D Acq On : 8 May 2008 3:10 Operator : RTB Sample : P0801342-002 DIL (1 Misc : ENSR SG83B-05-3 (-3 ALS Vial : 4 Sample Multipli | pm mL) .8, 3.5) | | | | |
| Quant Time: May 08 15:51:02 200 Quant Method : J:\MS13\METHODS\ Quant Title : EPA TO-15 per SC QLast Update : Tue Apr 15 06:47 Response via : Initial Calibrat | R1304140 P VOA-TO :20 2008 ion | 15 (CA | | | |
| Internal Standards | R.T. | QIon | Response | e Conc Unit | s Dev(Min) |
| 32) Chloroform 34) Tetrahydrofuran 35) Ethyl tert-Butyl Ether 36) 1,2-Dichloroethane 38) 1,1,1-Trichloroethane 39) Isopropyl Acetate | 12.78 0.00 0.00 13.72 0.00 0.00 | 62 97 | 587649 0 324 0 0 | 29.450 ng N.D. N.D. N.D. N.D. N.D. N.D. |) 100 |
| 40) 1-Butanol 41) Benzene 42) Carbon Tetrachloride 43) Cyclohexane 44) tert-Amyl Methyl Ether 45) 1,2-Dichloropropane | 15.41 | 78 117 84 73 | 0 3172 91943 146 0 0 | 5.637 ng | 99 |
| <pre>46) Bromodichloromethane 47) Trichloroethene 48) 1,4-Dioxane 49) Isooctane 50) Methyl Methacrylate 51) n-Heptane 52) cis-1,3-Dichloropropene 53) 4 Methacl 2 mentanene</pre> | 0.00 0.00 0.00 0.00 0.00 0.00 0.00 | 130 88 57 | | N.D. N.D. N.D. N.D. N.D. N.D. N.D. N.D. | |
| 53) 4-Methyl-2-pentanone 54) trans-1,3-Dichloropropene 55) 1,1,2-Trichloroethane 58) Toluene 59) 2-Hexanone 60) Dibromochloromethane | | 75 97 91 43 129 | 0 93721 226 0 0 | N.D. 7.891 ng N.D. N.D. N.D. N.D. | # 8 |
| 61) 1,2-Dibromoethane 62) Butyl Acetate 63) n-Octane 64) Tetrachloroethene 65) Chlorobenzene | 0.00 20.34 0.00 20.53 21.42 | 107 43 57 166 112 | 0 129 0 722 2785 | N.D. N.D. N.D. N.D. N.D. | |
| 66) Ethylbenzene 67) m- & p-Xylene 68) Bromoform 69) Styrene 70) o-Xylene | 21.92 22.12 0.00 0.00 22.73 | 91 91 173 104 91 | 117 63 0 125 | N.D. N.D. N.D. N.D. N.D. | |
| 71) n-Nonane 72) 1,1,2,2-Tetrachloroethane 74) Cumene 75) alpha-Pinene 76) n-Propylbenzene | 23.31 0.00 0.00 | 43 83 105 93 91 | 60 0 788 0 0 | N.D. N.D. N.D. N.D. N.D. | |
| 77) 3-Ethyltoluene 78) 4-Ethyltoluene 79) 1,3,5-Trimethylbenzene 13041408.M Thu May 08 15:57:20 2 | | | 201 201 201 | N.D. N.D. N.D. Fostoslos | 8 Page: 2 |
| 13041400.m illu may oo i3:37.20 Z | | | | ~ [• | |

QUALICICALIUM REPORT (IVEL ILLVILLVILL) Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080808.D Acq On : 8 May 2008 3:10 pm Operator : RTB Sample : P0801342-002 DIL (1mL) Misc : ENSR SG83B-05-3 (-3.8. : ENSR SG83B-05-3 (-3.8, 3.5) ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 15:51:02 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards _____

 80) alpha-Methylstyrene
 0.00
 118
 0
 N.D.

 81) 2-Ethyltoluene
 24.23
 105
 201
 N.D.

 82) 1,2,4-Trimethylbenzene
 0.00
 105
 0
 N.D.

 N.D. 83) n-Decane 84) Benzyl Chloride N.D.

 84)
 Benzyl Chloride
 0.00
 91

 85)
 1,3-Dichlorobenzene
 25.16
 146

 86)
 1,4-Dichlorobenzene
 25.16
 146

 87)
 sec-Butylbenzene
 0.00
 105

 N.D. N.D. N.D. N.D. 55 55 55

 86)
 1,4-Dichlorobenzene
 0.00
 105

 87)
 sec-Butylbenzene
 0.00
 105

 88)
 p-Isopropyltoluene
 25.41
 119

 89)
 1,2,3-Trimethylbenzene
 0.00
 105

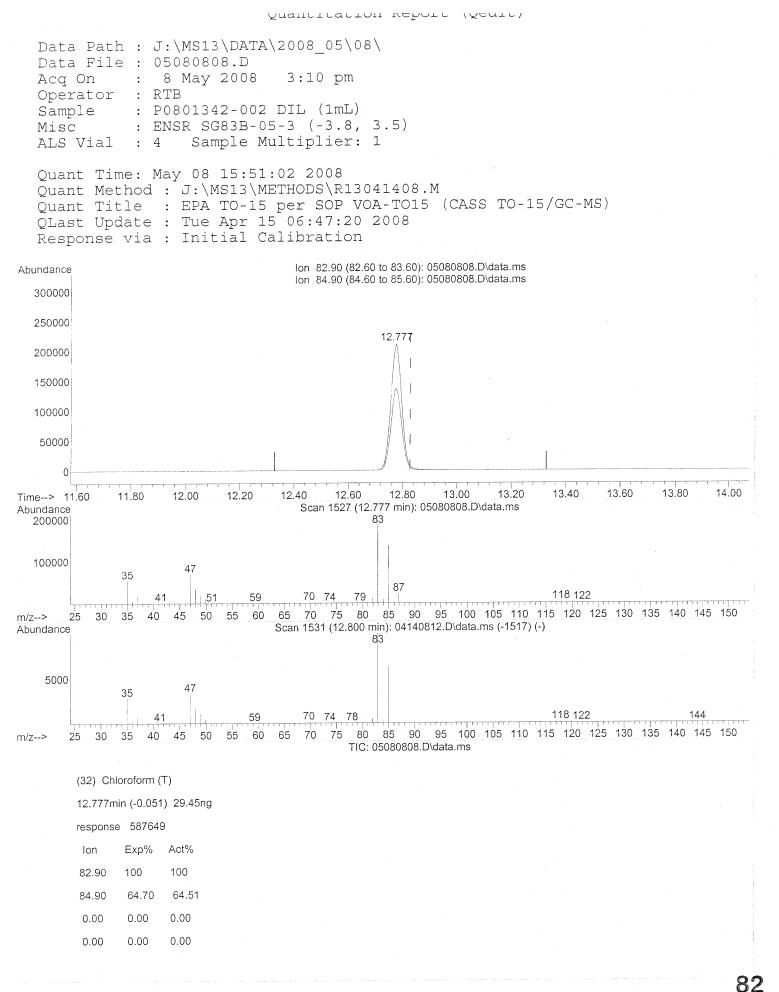
 1.2
 Dichlorobenzene
 25.16
 146

 2.20
 68

 0 171 0 N.D. N.D. 55 0 0 0.00 68 N.D. 91) d-Limonene N.D. N.D. N.D. N.D. 92) 1,2-Dibromo-3-Chloropr... 0.00 157 93) n-Undecane26.375794) 1,2,4-Trichlorobenzene0.0018095) Naphthalene27.82128 160 0 94) 1,2,1 12-5 95) Naphthalene 489 96) n-Dodecane 27.75 57 97) Hexachloro-1,3-butadiene 0.00 225 75 N.D. 0 N.D. _____

(#) = qualifier out of range (m) = manual integration (+) = signals summed

R13041408.M Thu May 08 15:57:20 2008



R13041408.M Thu May 08 16:42:44 2008

RESULTS OF ANALYSIS

Page 1 of 4

| Client: | ENSR |
|--------------------|-----------------------------------|
| Client Sample ID: | SG83B-05-7 |
| Client Project ID: | Phase B Soil Gas / 04020-023-4311 |
| | |
| Test Code | FPA TO-15 |

-

| Test Code. | EFA 10-15 |
|-----------------|--|
| Instrument ID: | Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 |
| Analyst: | Rusty Bravo |
| Sampling Media: | 6.0 L Summa Canister |
| Test Notes: | |
| Container ID: | SC00791 |

CAS Project ID: P0801342 CAS Sample ID: P0801342-003

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

3.5

Initial Pressure (psig):

-4.2 Final Pressure (psig):

Canister Dilution Factor: 1.73

| CAS # | Compound | Result µg/m³ | MRL µg/m³ | MDL µg/m³ | Result ppbV | MRL ppbV | MDL ppbV | Data Qualifier |
|-----------|--|-----------------|--------------|--------------|----------------|-------------|----------------|------------------------------------|
| 75-71-8 | Dichlorodifluoromethane (CFC 12) | ND | . 87 | 8.7 | ND | 18 | <u></u> 1.8 | <u></u> |
| 74-87-3 | Chloromethane | ND | 17 | 8.7 | ND | 8.4 | 4.2 | |
| 76-14-2 | 1,2-Dichloro-1,1,2,2- tetrafluoroethane (CFC 114) | ND | 87 | 8.7 | ND | 12 | 1.2 | |
| 75-01-4 | Vinyl Chloride | ND | 17 | 8.7 | ND | 6.8 | 3.4 | |
| 74-83-9 | Bromomethane | ND | 17 | 8.7 | ND | 4.5 | 2.2 | |
| 75-00-3 | Chloroethane | ND | 17 | 8.7 | ND | 6.6 | 3.3 | |
| 64-17-5 | Ethanol | ND | 870 | 8.7 | ND | 460 | 4.6 | |
| 67-64-1 | Acetone | 110 | 870 | 13 | 48 | 360 | 5.3 | J, B |
| 75-69-4 | Trichlorofluoromethane | 1,500 | 17 | 8.7 | 270 | 3.1 | 1.5 | |
| 107-13-1 | Acrylonitrile | ND | 87 | 12 | ND | 40 | 5.6 | |
| 75-35-4 | 1,1-Dichloroethene | ND | 17 | 8.7 | ND | 4.4 | 2.2 | |
| 75-65-0 | 2-Methyl-2-Propanol (tert-Butyl Alcohol) | ND | 87 | 13 | ND | 29 | 4.2 | |
| 75-09-2 | Methylene Chloride | 9.3 | 87 | 8.7 | 2.7 | 25 | 2.5 | J |
| 107-05-1 | 3-Chloro-1-propene (Allyl Chloride) | ND | 17 | 8.7 | ND | 5.5 | 2.8 | |
| 76-13-1 | Trichlorotrifluoroethane | ND | 17 | 9.7 | ND | 2.3 | 1.3 | |
| 75-15-0 | Carbon Disulfide | ND | 87 | 21 | ND | 28 | 6.7 | and a contract of the marked dates |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 17 | 8.7 | ND | 4.4 | 2.2 | |
| 75-34-3 | 1,1-Dichloroethane | ND | 17 | 8.7 | ND | 4.3 | 2.1 | |
| 1634-04-4 | Methyl tert-Butyl Ether | ND | 17 | 8.7 | ND | 4.8 | 2.4 | |
| 108-05-4 | Vinyl Acetate | ND | 870 | 28 | ND | 250 | 7.9 | |
| 78-93-3 | 2-Butanone (MEK) | 23 | 87 | 8.7 | 7.8 | 29 | 2.9 | J |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 17 | 8.7 | ND | 4.4 | 2.2 | |
| 108-20-3 | Diisopropyl Ether | ND | 87 | 10 | ND | 21 | 2.4 | |
| 67-66-3 | Chloroform | 54,000 | 17 | 10 | 11,000 | 3.5 | 2.1 | |

ND = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method. J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

Verified By: Rc-Date: 59108 TOISSCAN.XLT - Tronox - Henderson - PageNo.:

RESULTS OF ANALYSIS

Page 2 of 4

| Client: | ENSR |
|---------------------------|--|
| Client Sample ID: | SG83B-05-7 |
| Client Project ID: | Phase B Soil Gas / 04020-023-4311 |
| | |
| Test Code: | EPA TO-15 |
| Instrument ID: | Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 |

6.0 L Summa Canister

Rusty Bravo

SC00791

Analyst:

Sampling Media: Test Notes:

Container ID:

CAS Project ID: P0801342 CAS Sample ID: P0801342-003

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

Initial Pressure (psig):

e (psig): -4.2

Final Pressure (psig):

3.5

Canister Dilution Factor: 1.73

| CAS # | Compound | Result µg/m³ | MRL µg/m³ | MDL µg/m³ | Result ppbV | MRL ppbV | MDL ppbV | Data Qualifier |
|------------|---------------------------|-----------------|--------------|--------------|----------------|-------------|-------------|-------------------|
| 637-92-3 | Ethyl tert-Butyl Ether | ND | 87 | 8.8 | ND | 21 | 2.1 | |
| 107-06-2 | 1,2-Dichloroethane | ND | 17 | 8.7 | ND | 4.3 | 2.1 | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 17 | 8.7 | ND | 3.2 | 1.6 | |
| 71-43-2 | Benzene | 100 | 17 | 8.7 | 32 | 5.4 | 2.7 | |
| 56-23-5 | Carbon Tetrachloride | 12,000 | 17 | 8.7 | 1,900 | 2.8 | 1.4 | |
| 994-05-8 | tert-Amyl Methyl Ether | ND | 87 | 8.7 | ND | 21 | 2.1 | |
| 78-87-5 | 1,2-Dichloropropane | ND | 17 | 8.7 | ND | 3.7 | 1.9 | |
| 75-27-4 | Bromodichloromethane | ND | 17 | 8.7 | ND | 2.6 | 1.3 | |
| 79-01-6 | Trichloroethene | 11 | 17 | 8.7 | 2.1 | 3.2 | 1.6 | J |
| 123-91-1 | 1,4-Dioxane | ND | 87 | 11 | ND | 24 | 2.9 | |
| 80-62-6 | Methyl Methacrylate | ND | 87 | 13 | ND | 21 | 3.2 | |
| 142-82-5 | n-Heptane | ND | 87 | 11 | ND | 21 | 2.7 | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 87 | 9.0 | ND | 19 | 2.0 | |
| 108-10-1 | 4-Methyl-2-pentanone | ND | 87 | 9.7 | ND | 21 | 2.4 | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 87 | 11 | ND | 19 | 2.4 | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 17 | 8.7 | ND | 3.2 | 1.6 | |
| 108-88-3 | Toluene | ND | 87 | 8.7 | ND | 23 | 2.3 | |
| 591-78-6 | 2-Hexanone | ND | 87 | 13 | ND | 21 | 3.2 | |
| 124-48-1 | Dibromochloromethane | ND | 17 | 12 | ND | 2.0 | 1.4 | |
| 106-93-4 | 1,2-Dibromoethane | ND | 17 | 9.3 | ND | 2.3 | 1.2 | |
| 111-65-9 | n-Octane | ND | 87 | 8.7 | ND | 19 | 1.9 | |
| 127-18-4 | Tetrachloroethene | 130 | 17 | 8.7 | 19 | 2.6 | 1.3 | |
| 108-90-7 | Chlorobenzene | 180 | 17 | 8.8 | 39 | 3.8 | 1.9 | |

ND = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

Date: 5/4/58 TO15SCAN.XLT - Tronox - Henderson - PageNo.:

RESULTS OF ANALYSIS

Page 3 of 4

| Client: Client Sample ID: Client Project ID: | ENSR SG83B-05-7 Phase B Soil Gas / 04020-023-4311 | |
|--|---|--|
| Test Code: | EPA TO-15 | |
| Instrument ID: | Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 | |
| Analyst: | Rusty Bravo | |

CAS Project ID: P0801342 CAS Sample ID: P0801342-003

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 0.010 Liter(s) Volume(s) Analyzed: 0.0010 Liter(s)

Test Notes: Container ID:

Sampling Media:

SC00791

6.0 L Summa Canister

Initial Pressure (psig): -4.2 Final Pressure (psig):

3.5

Canister Dilution Factor: 1.73

85

Date: 5 (9) 38 TO15SCAN.XLT - Tronox - Henderson - PageNo.:

| | | Result | MRL | MDL | Result | MRL | MDL Data |
|-------------|-------------------------------|--------|-------|-------|--------|------|----------------|
| CAS # | Compound | μg/m³ | μg/m³ | μg/m³ | ppbV | ppbV | ppbV Qualifier |
| 100-41-4 | Ethylbenzene | ND | 87 | 11 | ND | 20 | 2.5 |
| 179601-23-1 | m,p-Xylenes | ND | 87 | 22 | ND | 20 | 5.2 |
| 75-25-2 | Bromoform | ND | 87 | 13 | ND | 8.4 | 1.3 |
| 100-42-5 | Styrene | ND | 87 | 13 | ND | 20 | 3.1 |
| 95-47-6 | o-Xylene | ND | 87 | 11 | ND | 20 | 2.5 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 17 | 11 | ND | 2.5 | 1.6 |
| 98-82-8 | Cumene | ND | 87 | 9.7 | ND | 18 | 2.0 |
| 103-65-1 | n-Propylbenzene | ND | 87 | 9.0 | ND | 18 | 1.8 |
| 622-96-8 | 4-Ethyltoluene | ND | 87 | 9.9 | ND | 18 | 2.0 |
| 108-67-8 | 1,3,5-Trimethylbenzene | ND | 87 | 10 | ND | 18 | 2.1 |
| 98-83-9 | alpha-Methylstyrene | ND | 87 | 13 | ND | 18 | 2.6 |
| 95-63-6 | 1,2,4-Trimethylbenzene | ND | 87 | 12 | ND | 18 | 2.4 |
| 100-44-7 | Benzyl Chloride | ND | 17 | 15 | ND | 3.3 | 2.9 |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 17 | 11 | ND | 2.9 | 1.8 |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 17 | 9.7 | ND | 2.9 | 1.6 |
| 135-98-8 | sec-Butylbenzene | ND | 87 | 10 | ND | 16 | 1.8 |
| 99-87-6 | 4-Isopropyltoluene (p-Cymene) | ND | 87 | 11 | ND | 16 | 2.0 |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 17 | 11 | ND | 2.9 | 1.9 |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 87 | 13 | ND | 9.0 | 1.4 |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 17 | 13 | ND | 2.3 | 1.8 |
| 91-20-3 | Naphthalene | ND | 35 | 13 | ND | 6.6 | 2.4 |
| 87-68-3 | Hexachlorobutadiene | ND | 17 | 16 | ND | 1.6 | 1.5 |
| 98-06-6 | tert-Butylbenzene | ND | 35 | 8.7 | ND | 6.3 | 1.6 |
| 104-51-8 | n-Butylbenzene | ND | 35 | 8.7 | ND | 6.3 | 1.6 |

ND = Compound was analyzed for, but not detected above the laboratory detection limit.

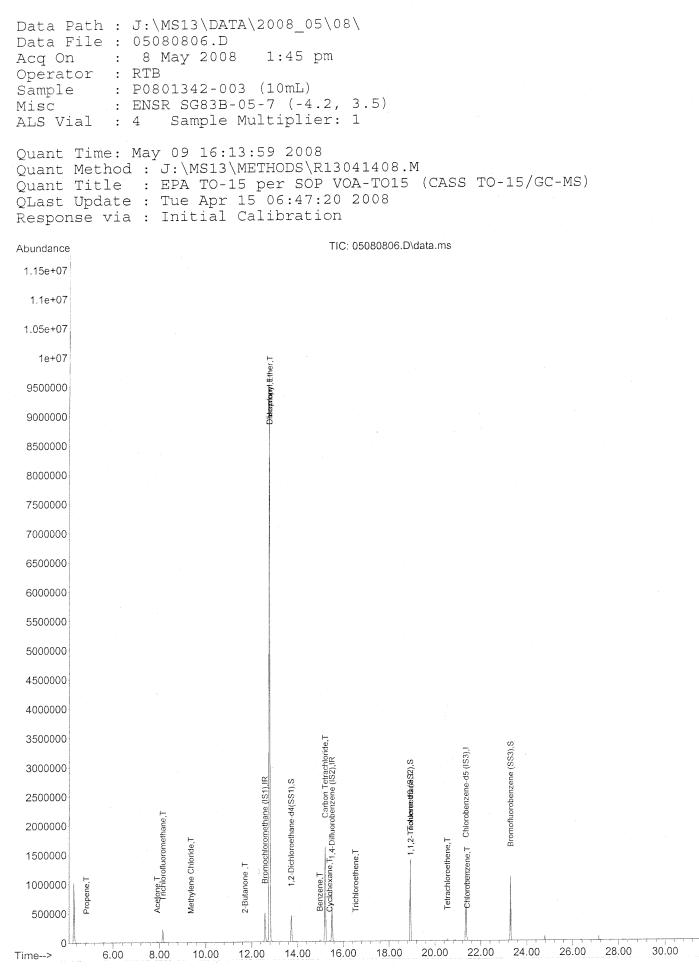
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

RESULTS OF ANALYSIS

Page 4 of 4

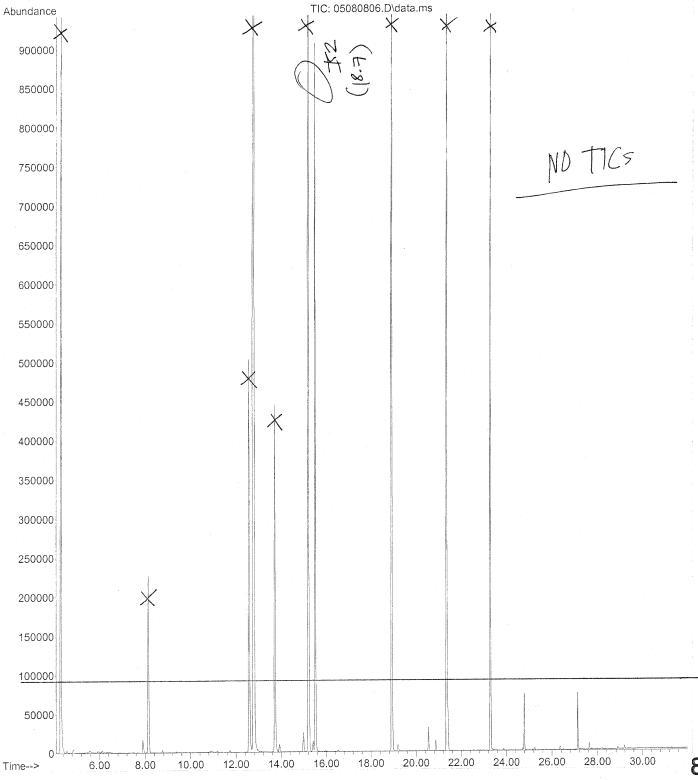
| Client: | ENSR | | |
|---------------------------|--|-------------------------|-------------------|
| Client Sample ID: | SG83B-05-7 | CAS Project ID: P080 | 1342 |
| Client Project ID: | Phase B Soil Gas / 04020-023-4311 | CAS Sample ID: P080 | 1342-003 |
| | Tentatively Identified Compo | unds | |
| Test Code: | EPA TO-15 | Date Collected: 5/7/0 | 8 |
| Instrument ID: | Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 | Date Received: 5/8/0 | 8 |
| Analyst: | Rusty Bravo | Date Analyzed: 5/8/0 | 8 |
| Sampling Media: | 6.0 L Summa Canister | Volume(s) Analyzed: 0.4 | 010 Liter(s) |
| Test Notes: | | 0.0 | 010 Liter(s) |
| Container ID: | SC00791 | | |
| | Initial Pressure (psig): -4.2 Final Pressur | re (psig): 3.5 | |
| | | Canister Dilu | tion Factor: 1.73 |
| GC/MS Retention Time | Compound Identification | Concentration µg/m³ | Data Qualifier |

No Compounds Detected



R13041408.M Fri May 09 16:16:03 2008

```
File :J:\MS13\DATA\2008_05\08\05080806.D
Operator : RTB
Acquired : 8 May 2008 1:45 pm using AcqMethod TO15.M
Instrument : GCMS13
Sample Name: P0801342-003 (10mL)
Misc Info : ENSR SG83B-05-7 (-4.2, 3.5)
Vial Number: 4
```



Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080806.D Acq On : 8 May 2008 1:45 pm Operator : RTB Sample : P0801342-003 (10mL) Misc : ENSR SG83B-05-7 (-4.2, 3.5) ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 09 16:13:59 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards 1) Bromochloromethane (IS1)12.5813022619925.000 ng-0.0237) 1,4-Difluorobenzene (IS2)15.5111497171725.000 ng-0.02 56) Chlorobenzene-d5 (IS3) 21.35 82 470342 25.000 ng 0.00 System Monitoring Compounds 33) 1,2-Dichloroethane-d4(... 13.73 65 414789 22.866 ng -0.02 Recovery = 91.48% Spiked Amount 25.000 18.93 98 1080045 25.619 ng 0.00 57) Toluene-d8 (SS2) Spiked Amount 25.000 Recovery = 102.48% 73) Bromofluorobenzene (SS3) 23.29 174 355347 24.493 ng 0.00 Recovery = 97.96% Spiked Amount 25.000 Qvalue Target Compounds 1167 0.062 ng # 22 4.85 42 2) Propene 3) Dichlorodifluoromethane 4.99 85 477 Ν.Ο. 0 0.00 50 N.D. 4) Chloromethane N.D. 0.00 135 0 5) Freon 114 N.D. 6) Vinyl Chloride N.D. 7) 1,3-Butadiene N.D. N.D. N.D. N.D. N.D. 8) Bromomethane 9) Chloroethane 10) Ethanol
 7.48
 41
 922

 7.69
 56
 232

 7.91
 58
 8266

 8.15
 01
 235861
 922 11) Acetonitrile 12) Acrolein 55 0.660 nq# 13) Acetone (8.795 ng) 99 14) Trichlorofluoromethane \langle 8.15 <u>101</u> 8.37 45 0.00 53 0.00 96 N.D. 221 15) Isopropanol 0 N.D. 16) Acrylonitrile N.D. 0 17) 1,1-Dichloroethene 9.46 59 9.36 84 18) tert-Butanol 156 N.D. 779 (0.054 ng) 97 19) Methylene Chloride 0 0 0.00 41 N.D. 20) Allyl Chloride N.D. 21) Trichlorotrifluoroethane 0.00 151 9.79 76 0.00 61 862 N.D. 22) Carbon Disulfide 0 N.D. 23) trans-1,2-Dichloroethene 24) 1,1-Dichloroethane 11.10 63 997 N.D. 25) Methyl tert-Butyl Ether 0.00 73 0 N.D. 0.00 86 0 11.71 72 1164m 0.00 61 0 12.78 87 904230 N.D. 26) Vinyl Acetate 1164m (0.133 ng 0 N.D. 27) 2-Butanone 28) cis-1,2-Dichloroethene 78.429 ng 1 29) Diisopropyl Ether 12.79 61 96 N.D. 30) Ethyl Acetate 12.69 57 162 N.D. 31) n-Hexane

R13041408.M Fri May 09 16:16:03 2008

Page: 1

| Quantitatio | | | |
|--|---|--|--|
| Data Path : J:\MS13\DATA\2008_ Data File : 05080806.D Acq On : 8 May 2008 1:45 Operator : RTB Sample : P0801342-003 (10mL Misc : ENSR SG83B-05-7 (- ALS Vial : 4 Sample Multipl | pm) 4.2, 3.5) | | |
| Quant Time: May 09 16:13:59 20 Quant Method : J:\MS13\METHODS Quant Title : EPA TO-15 per S QLast Update : Tue Apr 15 06:4 Response via : Initial Calibra | \R13041408.M OP VOA-TO15 (C 7:20 2008 tion | | |
| Internal Standards | R.T. QIor | n Response Cond | |
| 32) Chloroform 34) Tetrahydrofuran 35) Ethyl tert-Butyl Ether 36) 1,2-Dichloroethane 38) 1,1,1-Trichloroethane 39) Isopropyl Acetate | 13.74 62 0.00 97 0.00 61 | 0 N. 0 N. 250 N. 0 N. 0 N. | D. D. D. D. |
| 40) 1-Butanol 41) Benzene 42) Carbon Tetrachloride 43) Cyclohexane 44) tert-Amyl Methyl Ether 45) 1,2-Dichloropropane 46) Bromodichloromethane | $ \begin{array}{r} 14.96 \\ 56 \\ 14.98 \\ 78 \\ 15.21 \\ 15.41 \\ 84 \\ 0.00 \\ 73 \\ 0.00 \\ 63 \\ 16.46 \\ 83 \\ \end{array} $ | 30187 0.58 1183939 69.44 5856 0.30 0 N. 0 N. | - |
| 46) Bromodicinoronale finale 47) Trichloroethene 48) 1,4-Dioxane 49) Isooctane 50) Methyl Methacrylate 51) n-Heptane 52) cis-1,3-Dichloropropene 53) 4-Methyl-2-pentanone 54) trans-1,3-Dichloropropen | 16.54 130 0.00 88 0.00 57 0.00 100 16.98 71 0.00 75 0.00 58 | 837 0.06 0 N. 0 N. 0 N. 58 N. 0 N. 0 N. | 56 ng 94 D. D. D. D. D. D. D. D. |
| 54) trans-1, 3-Dichiolopiopen 55) 1,1,2-Trichloroethane 58) Toluene 59) 2-Hexanone 60) Dibromochloromethane 61) 1,2-Dibromoethane 62) Butyl Acetate 63) n-Octane | $ \begin{array}{cccccccccccccccccccccccccccccccccccc$ | 98611 7.94 1751 N. 88 N. 0 N. 0 N. 327 N. 0 N. | <u>4 ng</u> N # 8 D. D. D. D. D. D. |
| 64) Tetrachloroethene 65) Chlorobenzene 66) Ethylbenzene 67) m- & p-Xylene 68) Bromoform 69) Styrene 70) o-Xylene | 20.54 166 21.41 112 21.90 91 22.10 91 0.00 173 0.00 104 22.71 91 22.98 43 | 34136 402 1008 0 N 469 N | 95 0 ng 95 0 ng 97 0. 0. 0. 0. 0. 0. 0. 0. |
| <pre>71) n-Nonane 72) 1,1,2,2-Tetrachloroethan 74) Cumene 75) alpha-Pinene 76) n-Propylbenzene 77) 3-Ethyltoluene 78) 4-Ethyltoluene 79) 1,3,5-Trimethylbenzene</pre> | | 54 N. 105 N. 0 N. 58 N. 566 N. 67 N. | D. D. D. D. D. D. D. D. 90 |
| 13041408.M Fri May 09 16:16:03 | 2008 | Rosfiglos | Page: 2 |

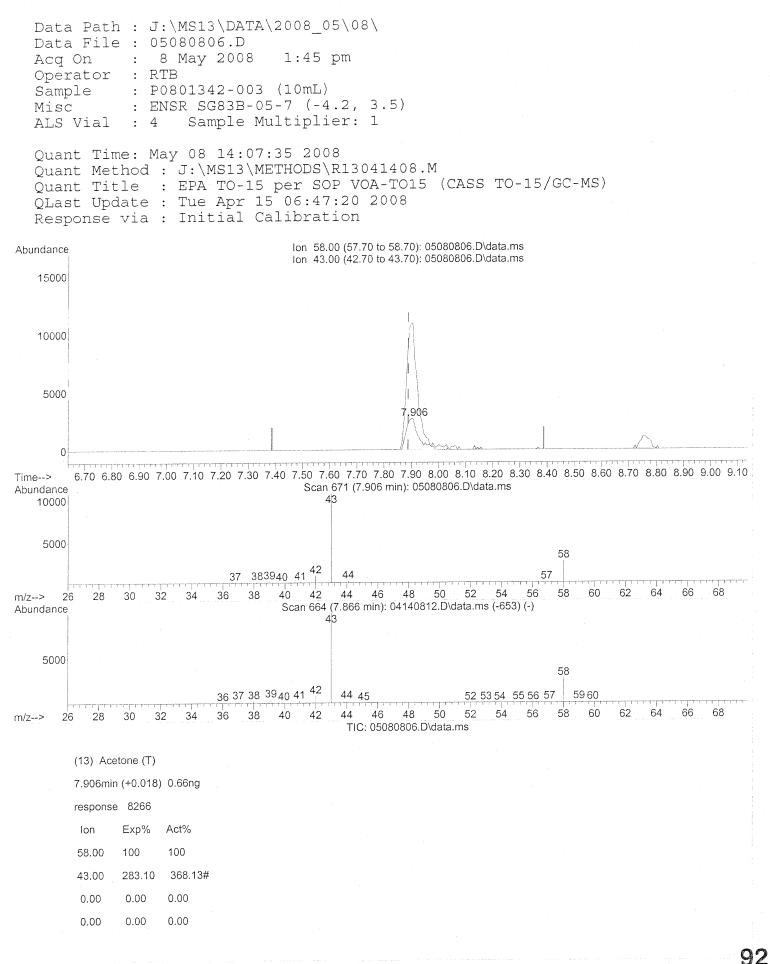
Quantituacion Report (Xr Review,

Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080806.D Acq On : 8 May 2008 1:45 pm Operator : RTB Sample : P0801342-003 (10mL) Misc : ENSR SG83B-05-7 (-4 : ENSR SG83B-05-7 (-4.2, 3.5) ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 09 16:13:59 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards 80) alpha-Methylstyrene24.77118713N.D.81) 2-Ethyltoluene24.6210555N.D.82) 1,2,4-Trimethylbenzene24.88105655N.D.83) n-Decane25.0057289N.D.84) Benzyl Chloride0.00910N.D.85) 1,3-Dichlorobenzene25.17146777N.D.86) 1,4-Dichlorobenzene25.17146777N.D.87) sec-Butylbenzene25.40105129N.D.88) p-Isopropyltoluene25.41119148N.D.90) 1,2-Dichlorobenzene25.17146777N.D.91) d-Limonene25.586856N.D. N.D. N.D. N.D. N.D. N.D. N.D. N.D. 90) 1,2-Dichlorobenzene 56 25.58 68 91) d-Limonene 92) 1,2-Dibromo-3-Chloropr... 0.00 157 0 N.D. 93) n-Undecane26.505794) 1,2,4-Trichlorobenzene0.0018095) Naphthalene27.80128 N.D. N.D. N.D. 661 0909 95) Naphthalene 96) n-Dodecane 27.74 57 696 97) Hexachloro-1,3-butadiene 0.00 225 0 696 N.D. N.D. ------

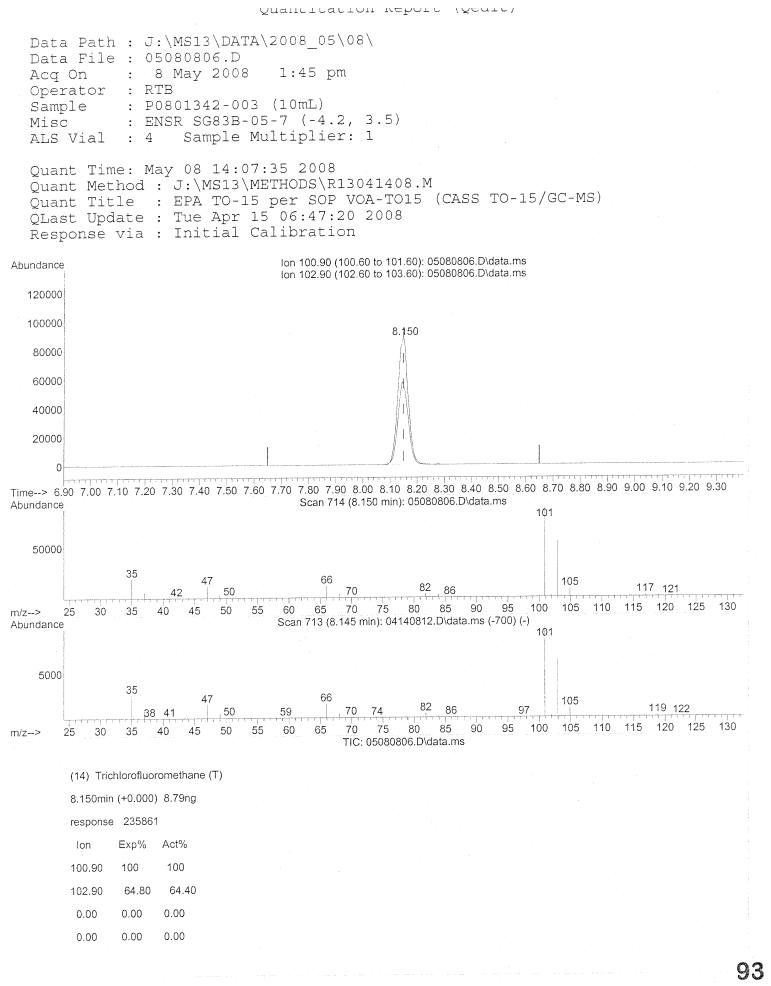
(#) = qualifier out of range (m) = manual integration (+) = signals summed

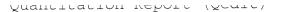
50 stoglor

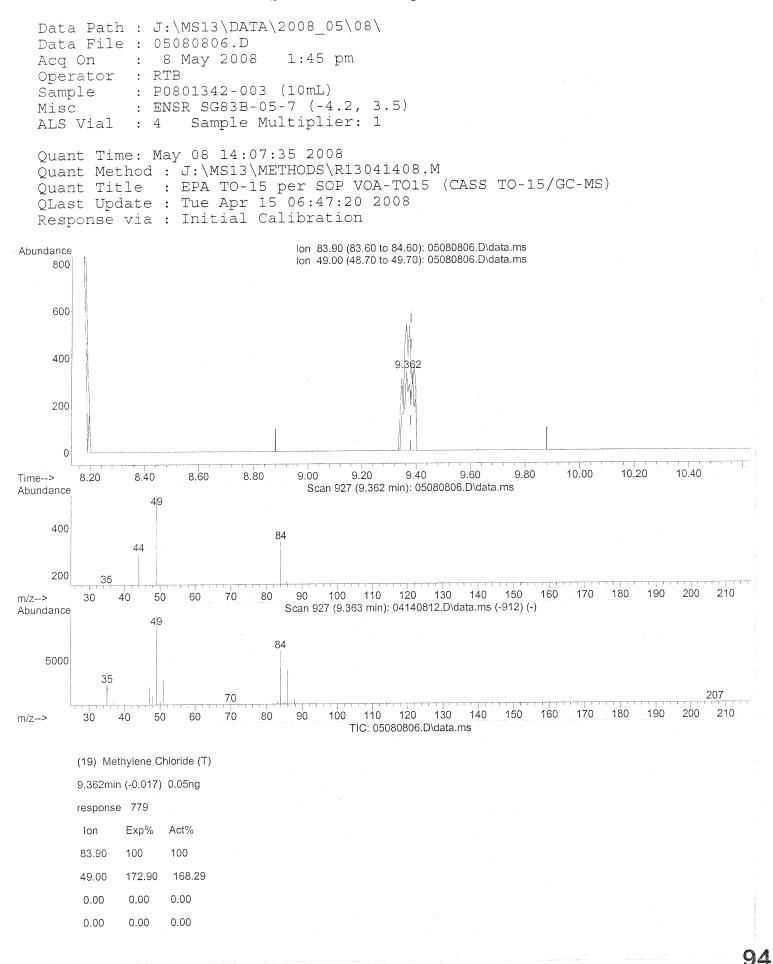




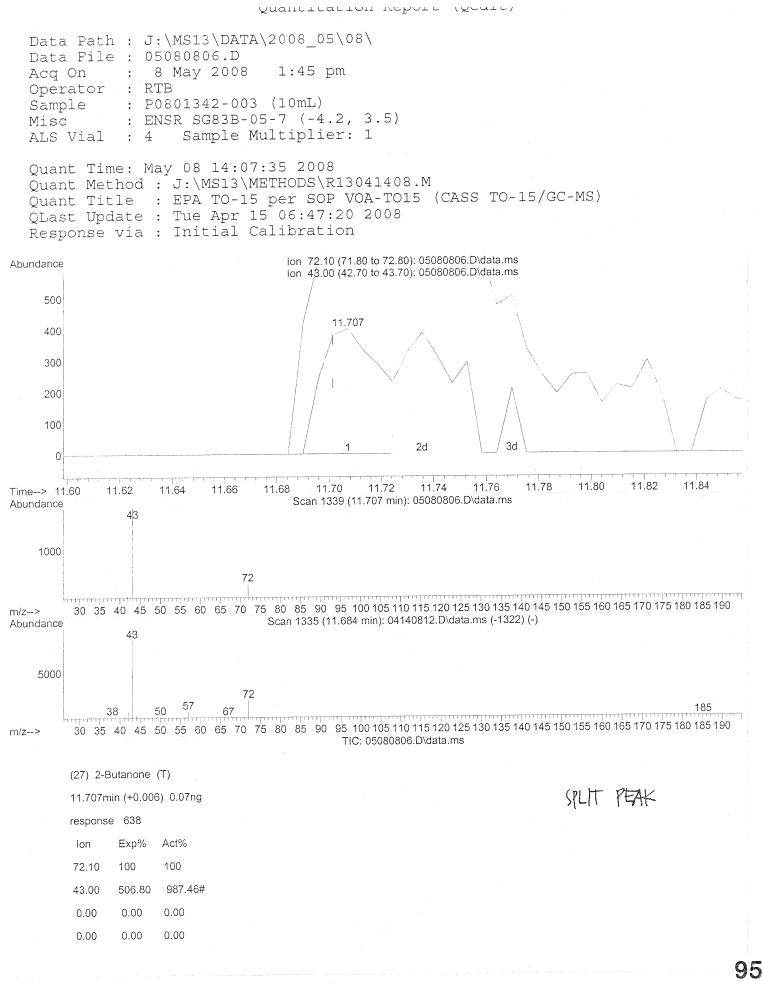
R13041408.M Thu May 08 16:44:44 2008

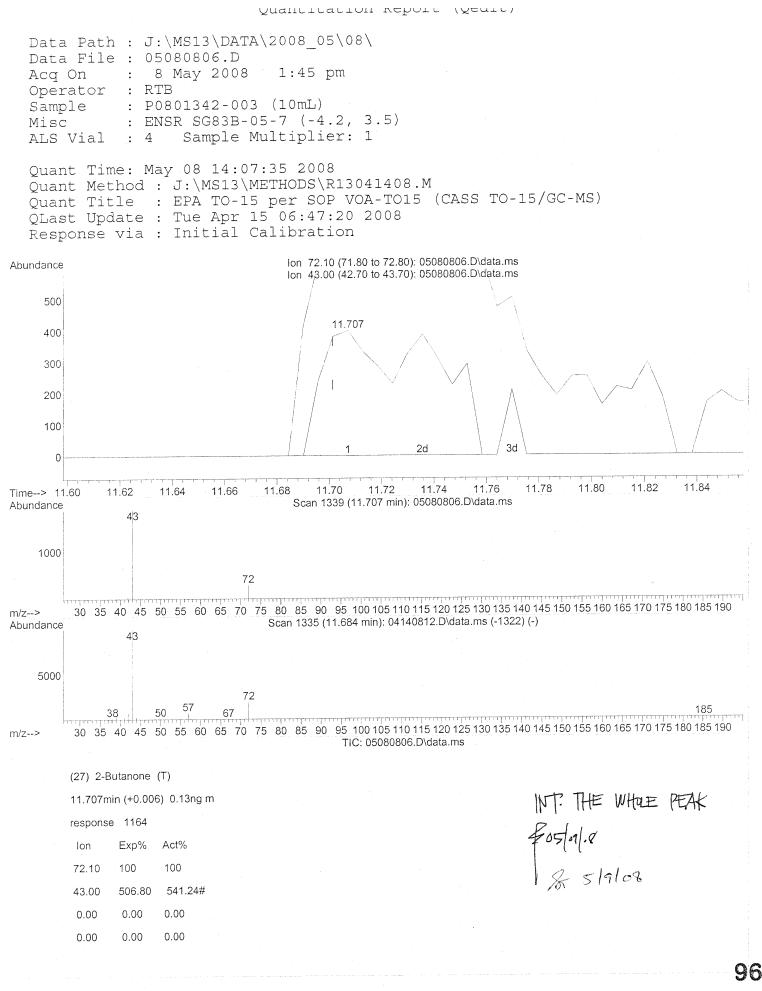


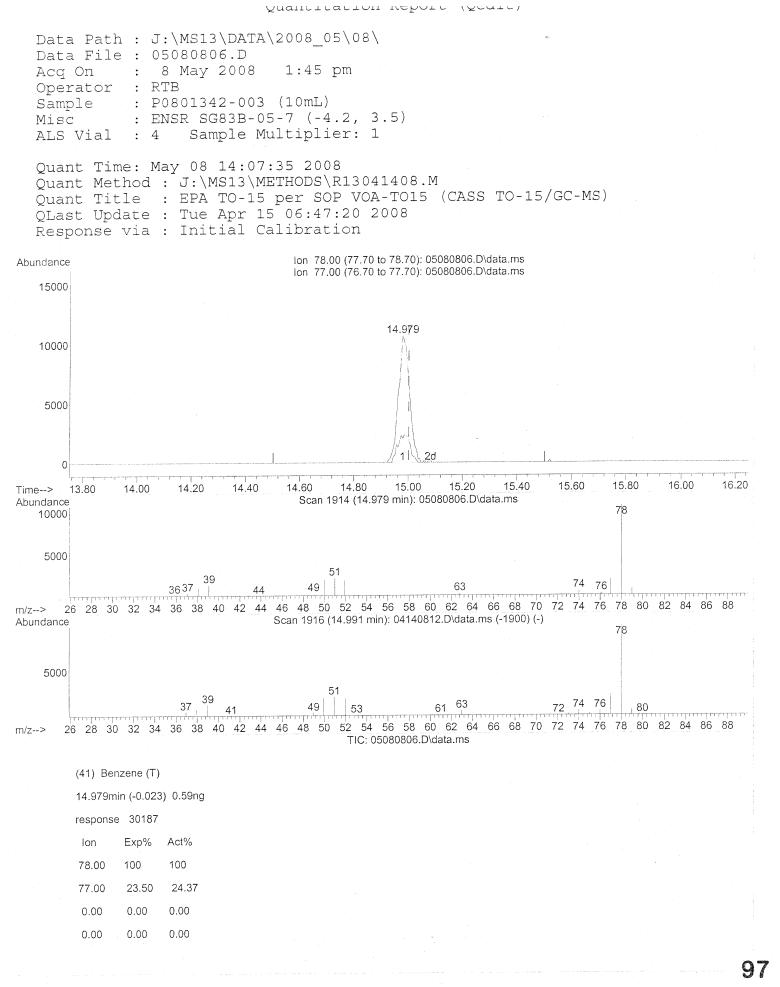




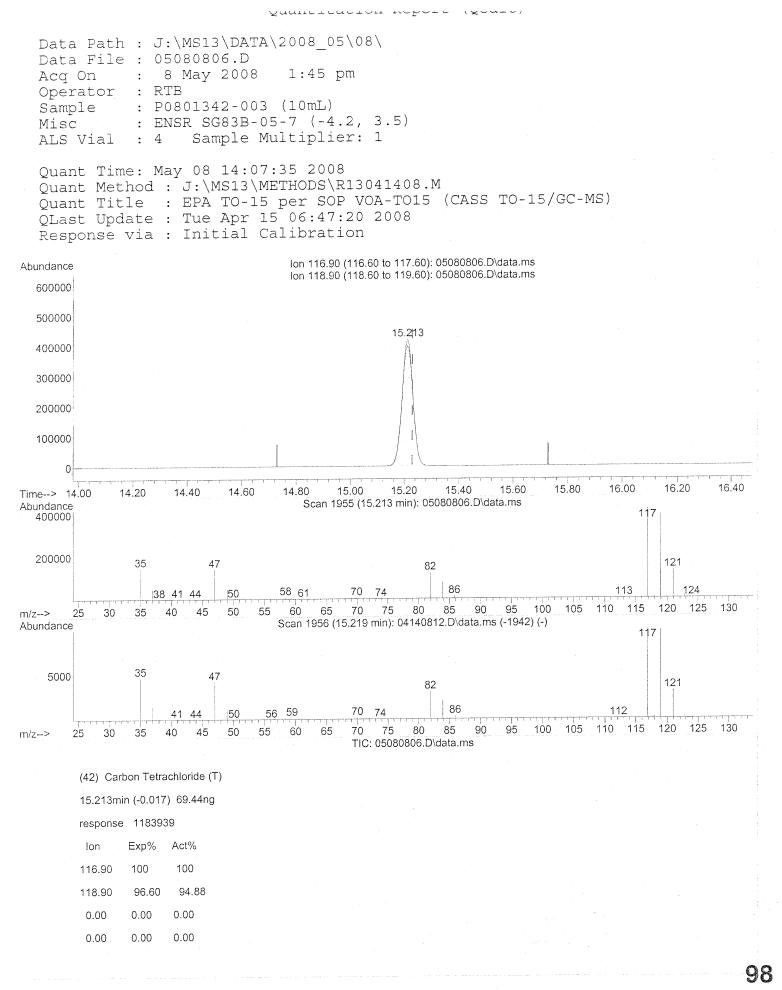
R13041408.M Fri May 09 15:25:06 2008

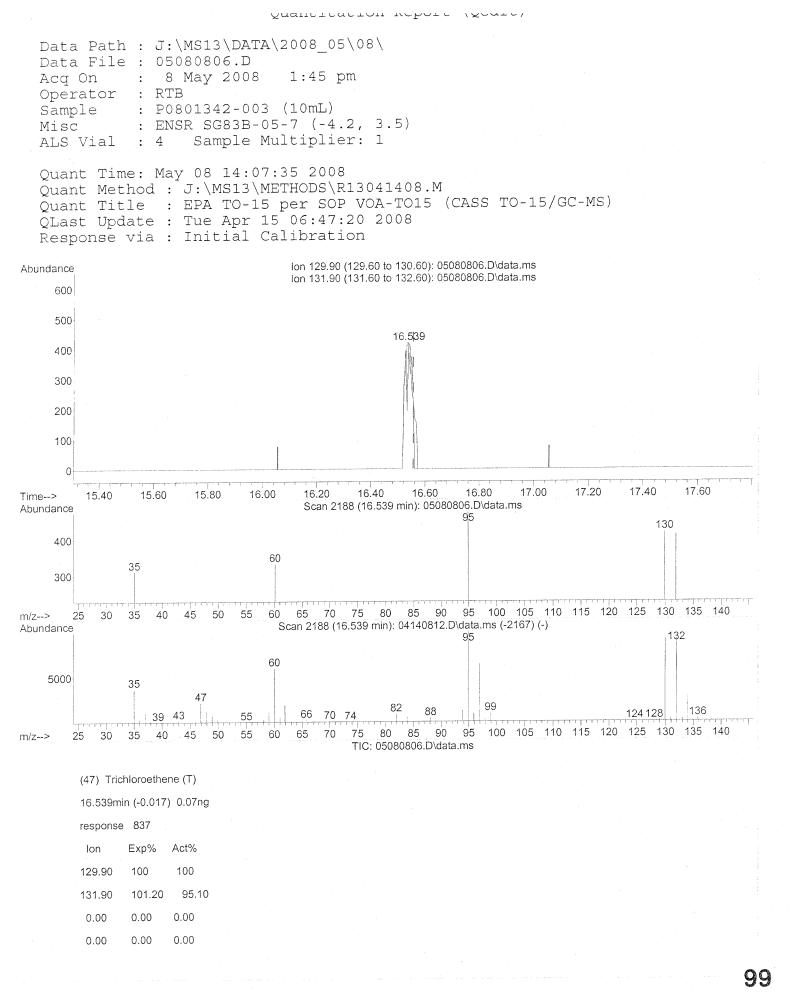




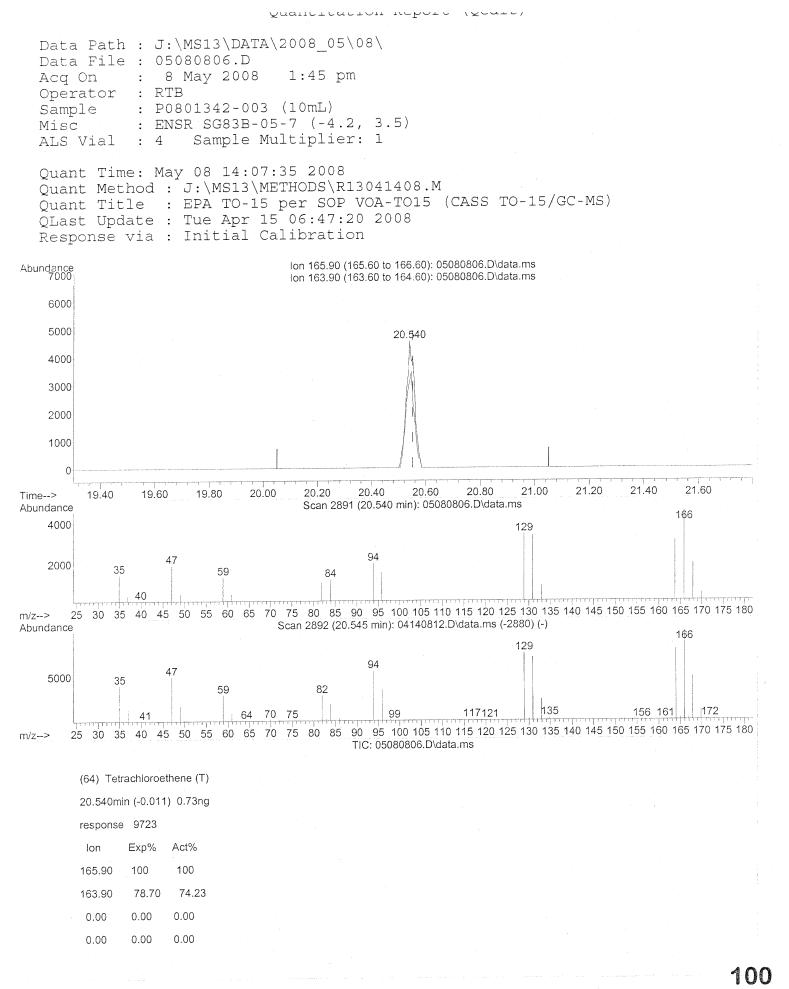


R13041408.M Thu May 08 16:45:54 2008



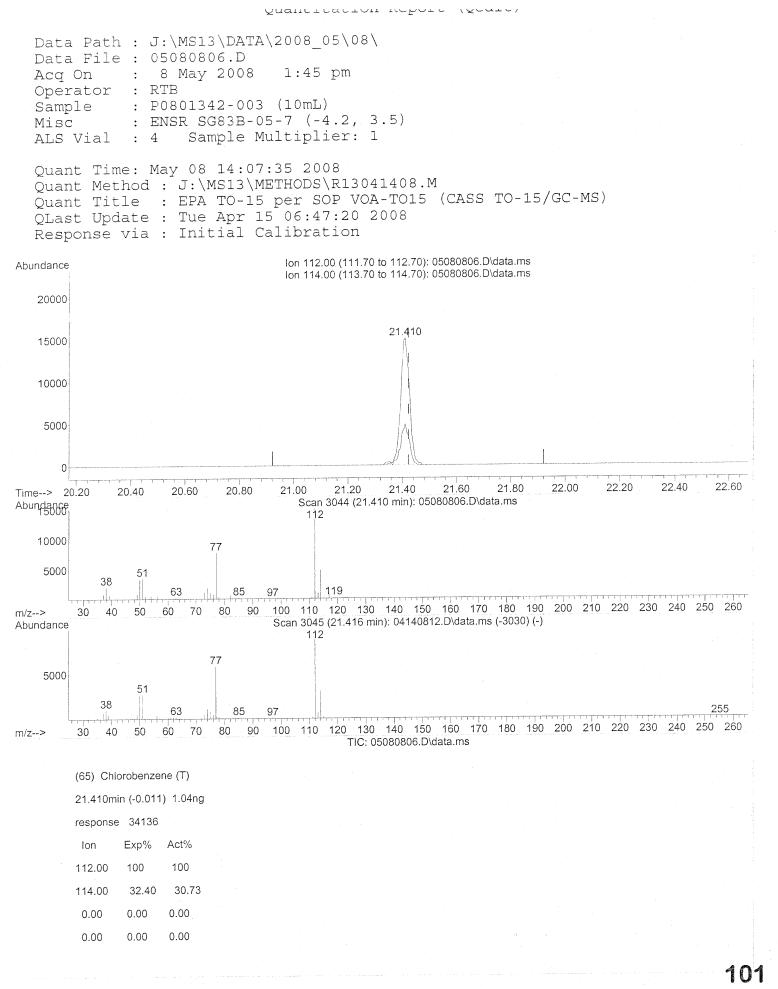


R13041408.M Fri May 09 15:25:51 2008

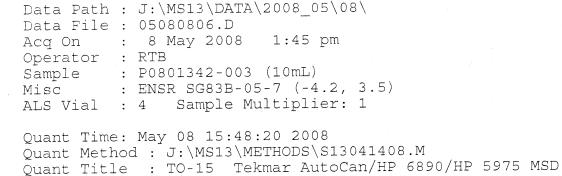


R13041408.M Thu May 08 16:46:40 2008

Page: 1



(INON VENTEMEN)



QLast Update : Mon Apr 28 10:06:00 2008

Response via : Initial Calibration

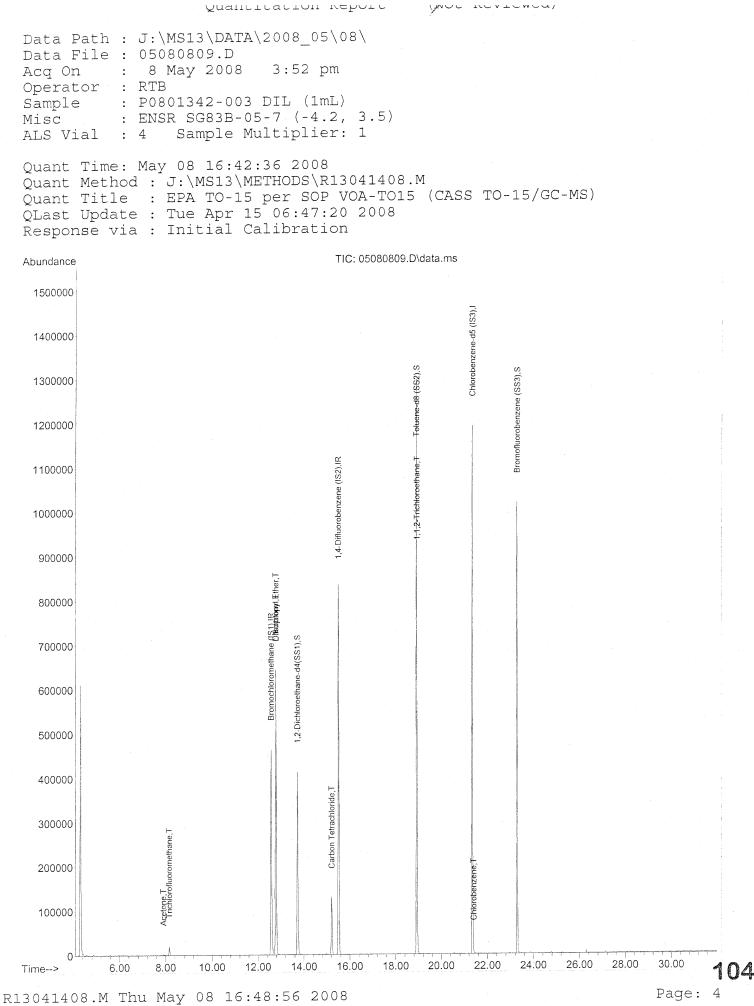
| Abundance | TIC: 05080806.D\data.ms | |
|------------|--|-------------------------|
| 1.15e+07 | | |
| 1.1e+07 | | |
| 1.05e+07 | | |
| 1e+07 | | |
| 9500000 | | |
| 9000000 | | |
| 8500000 | | |
| 8000000 | | |
| 7500000 | | |
| 7000000 | | |
| 6500000 | | |
| 6000000 | | |
| 5500000 | | 1 |
| 5000000 | | |
| 4500000 | | |
| 4000000 | | |
| | | |
| 3500000 | SS SS SS SS SS | |
| 3000000 | ne-d5 ((S2)) a (S22) | |
| 2500000 | diloromethane (IS1),IR Noroethane-d4(SS1),S Toluene-d8 (SS2),S Chlorobenzene-d5 (IS3),I Bromofluorobenzene (SS3),S | |
| 2000000 | Diffuor Chlo Bromot | |
| 1500000 | Biomochloromethane (IS1).IR 2. Dichloroethane-d4(SS1),S 1,4. Difluorobenzene (IS Chlorobenzene-d8 (SS | |
| 1000000 | | - |
| 500000 | | |
| 0 Time> | 6.00 8.00 10.00 12.00 14.00 16.00 18.00 20.00 22.00 24.00 26.00 28.00 | 30.00 |
| | 8.M Thu May 08 15:54:38 2008 | 30.00 Page: 2 |

| Yuuncreacton | | 7 | , , , , , , , , , , | | | | |
|---|---------------|-----------|---|--|--|--|--|
| Data Path : J:\MS13\DATA\2008_05 Data File : 05080806.D Acq On : 8 May 2008 1:45 p Operator : RTB Sample : P0801342-003 (10mL) Misc : ENSR SG83B-05-7 (-4. ALS Vial : 4 Sample Multiplie | m 2, 3.5) | | | | | | |
| Quant Time: May 08 15:48:20 2008 Quant Method : J:\MS13\METHODS\S13041408.M Quant Title : TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD QLast Update : Mon Apr 28 10:06:00 2008 Response via : Initial Calibration | | | | | | | |
| Internal Standards | R.T. | QIon | n Response Conc Units Dev(Min) | | | | |
| 2) 1 / Difluorohonzone (TS2) | 15 51 | 114 | 22619925.000 ng-0.0297171725.000 ng-0.0247034225.000 ng0.00 | | | | |
| Spiked Amount 25.000 5) Toluene-d8 (SS2) Spiked Amount 25.000 | 18.93 | 98 | 414789 22.866 ng -0.02 Recovery = 91.48% ✓ 1080045 25.619 ng 0.00 Recovery = 102.48% ✓ 355347 24.493 ng 0.00 Recovery = 97.96% ✓ | | | | |
| Target Compounds 7) tert-Butylbenzene | 24.79 0.00 | 119 91 | Qvalue 1326 N.D. 🗸 0 N.D. 🖌 | | | | |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

705108/05

Page: 1



Page: 4

| Qualitication | Veborc | (14 | | л, | | | |
|--|--|---|---|---|-----------------------|-----------------------------|-------------------|
| Data Path : J:\MS13\DATA\2008_05 Data File : 05080809.D Acq On : 8 May 2008 3:52 p Operator : RTB Sample : P0801342-003 DIL (1n Misc : ENSR SG83B-05-7 (-4. ALS Vial : 4 Sample Multiplie | om nL) 2, 3.5) | | | | | | |
| Quant Time: May 08 16:42:36 2008 Quant Method : J:\MS13\METHODS\F Quant Title : EPA TO-15 per SOF QLast Update : Tue Apr 15 06:47: Response via : Initial Calibrati | 81304140 > VOA-TO 20 2008 | 15 (C | ASS TO-15/(| GC-MS) | | | |
| Internal Standards | R.T. | QIon | Response | Conc ' | Units | Dev(N | Min) |
| 1) Bromochloromethane (IS1) 37) 1,4-Difluorobenzene (IS2) 56) Chlorobenzene-d5 (IS3) | 12.58 15.51 21.35 | 130 114 82 | | | ng ng ng | - 0 . - 0 . 0 . | .03 .02 .00 |
| System Monitoring Compounds 33) 1,2-Dichloroethane-d4(Spiked Amount 25.000 57) Toluene-d8 (SS2) Spiked Amount 25.000 73) Bromofluorobenzene (SS3) Spiked Amount 25.000 | 18.92 | 98 | Recove 1015105 Recove 328273 | ery = 26.441 ery = | 92 ng 105 ng | .72%✔ -0. .76%✔ 0. | .01 |
| | | | | 2 | | | |
| <pre>Target Compounds 2) Propene 3) Dichlorodifluoromethane 4) Chloromethane 5) Freon 114 6) Vinyl Chloride 7) 1,3-Butadiene 8) Bromomethane 9) Chloroethane 10) Ethanol</pre> | 0.00 0.00 0.00 0.00 | 85 50 135 62 54 | 54 0 0 0 0 0 0 0 0 0 | N.D N.D N.D N.D N.D N.D N.D N.D N.D | | Qval | ue |
| Acetonitrile Acrolein Acetone Trichlorofluoromethane Isopropanol Acrylonitrile 1,1-Dichloroethene tert-Butanol | 7.47 0.00 7.90 8.16 0.00 0.00 0.00 9.37 | 41 56 58 101 45 53 96 59 | 55 0 1561 20143 0 0 0 53 | N.D N.D 0.137 0.827 N.D N.D N.D N.D | ng ng | # | 40 98 |
| 19) Methylene Chloride 20) Allyl Chloride 21) Trichlorotrifluoroethane 22) Carbon Disulfide 23) trans-1,2-Dichloroethene 24) 1,1-Dichloroethane 25) Methyl tert-Butyl Ether 26) Vinyl Acetate | 9.36 0.00 9.77 0.00 0.00 0.00 0.00 | 84 41 151 76 61 63 73 86 | 219 0 384 0 0 0 0 | N.D N.D N.D N.D N.D N.D N.D N.D | • • • • | | |
| 27) 2-Butanone 28) cis-1,2-Dichloroethene 29) Diisopropyl Ether 30) Ethyl Acetate 31) n-Hexane | 0.00 0.00 12.78 0.00 0.00 | 72 61 87 61 57 | 0 | N.D N.D 5.892 N.D N.D | ng | # | 1 |
| 13041408.M Thu May 08 16:48:55 20 | 08 | | Ŧ | 65708/08 |] | Page: | 105 |

(MOL REVIEWED) Quantitation Report Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080809.D Acq On : 8 May 2008 3:52 pm Operator : RTB Sample : P0801342-003 DIL (1mL) Misc : ENSR SG83B-05-7 (-4.2, 3.5) ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 16:42:36 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration

 Response via : Initial Calibration

 Internal Standards
 R.T. Olon Response Conc Units Dev(Min)

 32) Chloroform
 000 72 0
 0 N.D.

 33) Tetrahydrofuran
 000 77 0
 0 N.D.

 35) Ethyl tert-Suyl Ether
 0.00 87 0
 0 N.D.

 36) 1,2-Dichloroethane
 0.00 97 0
 0 N.D.

 39) Isopropyl Acetate
 0.00 56 0
 N.D.

 40) 1-Butanol
 0.00 57 0
 N.D.

 41) Benzene
 15.21 117 93957 5.952 ng 99
 99

 42) Carbon Tetrachloride
 15.21 117 93957 5.952 ng 99
 99

 43) CyCohexane
 0.00 63 0
 N.D.

 44) tert-Amyl Methyl Ether
 0.00 73 0
 N.D.

 45) 1,2-Dichloropropane
 0.00 130 0
 N.D.

 46) Isooctame
 0.00 130 0
 N.D.

 51) dethalacrylate
 0.00 100 0
 N.D.

 52 Methyl Methacrylate
 0.00 100 0
 N.D.

 53) 4-Methyl-2-pentancne
 0.00 75 0
 N.D.

 54) trans-1, 3-Dichloropropene
 0.00 75 0
 N.D.

 55) 1,1,2-Trinchoroethane
 18.94 97 93376
 8.123 ng # 7

 58) Tolene
 19.06 117 0
 N.D.

 <tr R.T. QIon Response Conc Units Dev(Min) Fostoslos Page: 2

R13041408.M Thu May 08 16:48:55 2008

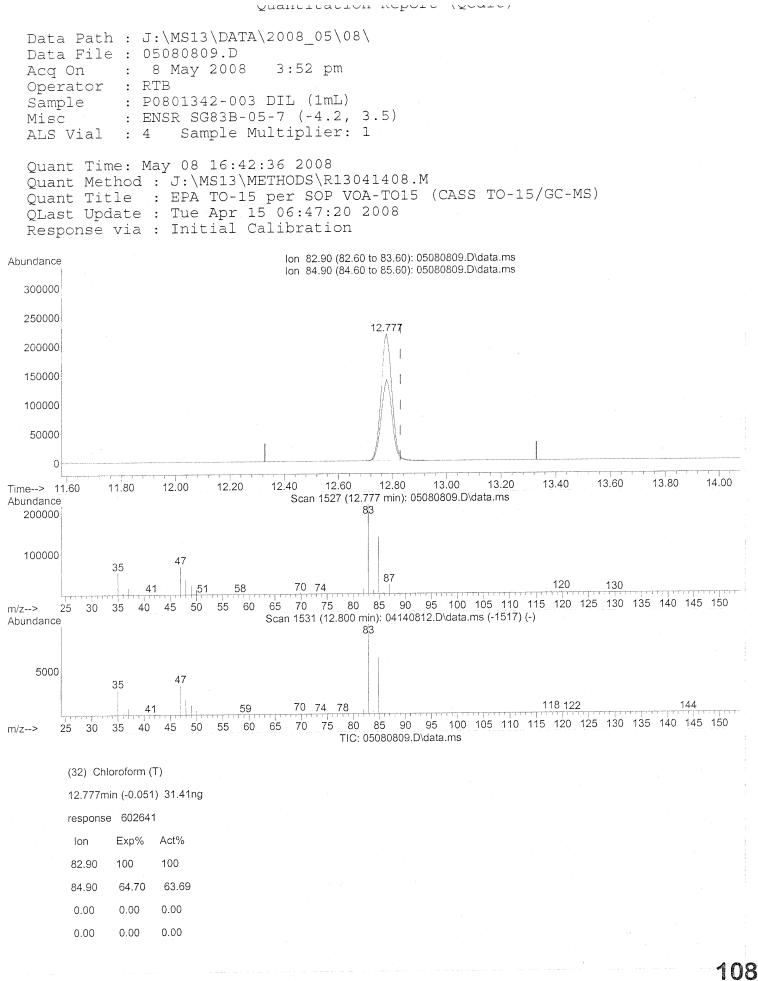
| Quantitation | MCDOT C | (1) | | - / | |
|---|--|--|---|--|------------|
| Data Path : J:\MS13\DATA\2008_05 Data File : 05080809.D Acq On : 8 May 2008 3:52 p Operator : RTB Sample : P0801342-003 DIL (1m Misc : ENSR SG83B-05-7 (-4. ALS Vial : 4 Sample Multiplie | om NL) 2, 3.5) | | | | |
| Quant Time: May 08 16:42:36 2008 Quant Method : J:\MS13\METHODS\R Quant Title : EPA TO-15 per SOB QLast Update : Tue Apr 15 06:47: Response via : Initial Calibrati | 21304140 > VOA-TO 20 2008 .on | 15 (CA | с | | |
| Internal Standards | R.T. | QIon | Response | Conc Units | B Dev(Min) |
| <pre>84) Benzyl Chloride 85) 1,3-Dichlorobenzene 86) 1,4-Dichlorobenzene 87) sec-Butylbenzene 88) p-Isopropyltoluene 89) 1,2,3-Trimethylbenzene 90) 1,2-Dichlorobenzene 91) d-Limonene 92) 1,2-Dibromo-3-Chloropr</pre> | $\begin{array}{c} 25.25 \\ 0.00 \\ 0.00 \\ 25.41 \\ 25.41 \\ 25.41 \\ 0.00 \\ 0.00 \\ 0.00 \\ 26.40 \\ 0.00 \\ 27.81 \\ 27.74 \\ 0.00 \end{array}$ | 91 146 146 105 119 105 146 68 157 57 180 128 57 225 | 0 0 0 61 71 61 0 0 0 79 0 454 173 | N.D. N.D. N.D. N.D. N.D. N.D. N.D. N.D. | |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

205/08/06

107 Page: 3

R13041408.M Thu May 08 16:48:55 2008



RESULTS OF ANALYSIS

Page 1 of 4

ENSR Client: Client Sample ID: Method Blank Client Project ID: Phase B Soil Gas / 04020-023-4311

EPA TO-15

Rusty Bravo

6.0 L Summa Canister

Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13

Test Code:

Analyst:

Instrument ID:

Sampling Media: Test Notes:

CAS Project ID: P0801342 CAS Sample ID: P080508-MB

Date Collected: NA Date Received: NA Date Analyzed: 5/8/08 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

| CAS # | Compound | Result µg∕m³ | MRL µg/m³ | MDL µg/m³ | Result ppbV | MRL ppbV | | Data Qualifier |
|-----------|--|-----------------|--------------|--------------|----------------|-------------|--------|--|
| 75-71-8 | Dichlorodifluoromethane (CFC 12) | ND | 0.50 | 0.050 | ND | 0.10 | 0.010 | |
| 74-87-3 | Chloromethane | ND | 0.10 | 0.050 | ND | 0.048 | 0.024 | |
| 76-14-2 | 1,2-Dichloro-1,1,2,2- tetrafluoroethane (CFC 114) | ND | 0.50 | 0.050 | ND | 0.072 | 0.0072 | |
| 75-01-4 | Vinyl Chloride | ND | 0.10 | 0.050 | ND | 0.039 | 0.020 | |
| 74-83-9 | Bromomethane | ND | 0.10 | 0.050 | ND | 0.026 | 0.013 | |
| 75-00-3 | Chloroethane | ND | 0.10 | 0.050 | ND | 0.038 | 0.019 | |
| 64-17-5 | Ethanol | ND | 5.0 | 0.050 | ND | 2.7 | 0.027 | |
| 67-64-1 | Acetone | 0.42 | 5.0 | 0.073 | 0.18 | 2.1 | 0.031 | J |
| 75-69-4 | Trichlorofluoromethane | ND | 0.10 | 0.050 | ND ND | 0.018 | 0.0089 | |
| 107-13-1 | Acrylonitrile | ND | 0.50 | 0.070 | ND | 0.23 | 0.032 | analysis and the constraint for the second state of the second state |
| 75-35-4 | 1,1-Dichloroethene | ND | 0.10 | 0.050 | ND | 0.025 | 0.013 | |
| 75-65-0 | 2-Methyl-2-Propanol (tert-Butyl Alcohol) | ND | 0.50 | 0.074 | ND | 0.17 | 0.024 | |
| 75-09-2 | Methylene Chloride | ND | 0.50 | 0.050 | ND | 0.14 | 0.014 | |
| 107-05-1 | 3-Chloro-1-propene (Allyl Chloride) | ND | 0.10 | 0.050 | ND | 0.032 | 0.016 | |
| 76-13-1 | Trichlorotrifluoroethane | ND | 0.10 | 0.056 | ND | 0.013 | 0.0073 | |
| 75-15-0 | Carbon Disulfide | ND | 0.50 | 0.12 | ND | 0.16 | 0.039 | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 0.10 | 0.050 | ND | 0.025 | 0.013 | |
| 75-34-3 | 1,1-Dichloroethane | ND | 0.10 | 0.050 | ND | 0.025 | 0.012 | |
| 1634-04-4 | Methyl tert-Butyl Ether | ND | 0.10 | 0.050 | ND | 0.028 | 0.014 | |
| 108-05-4 | Vinyl Acetate | ND | 5.0 | 0.16 | ND | 1.4 | 0.045 | |
| 78-93-3 | 2-Butanone (MEK) | ND | 0.50 | 0.050 | ND | 0.17 | 0.017 | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 0.10 | 0.050 | ND | 0.025 | 0.013 | |
| 108-20-3 | Diisopropyl Ether | ND | 0.50 | 0.059 | ND | 0.12 | 0.014 | |
| 67-66-3 | Chloroform | ND | 0.10 | 0.059 | ND | 0.020 | 0.012 | |

ND = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method. J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

Verified By: R_Cr Date: 5/9/08 TO15SCAN.XLT - Tronox - Henderson - PageNo.

109

RESULTS OF ANALYSIS Page 2 of 4

Client: ENSR Client Sample ID: Method Blank Client Project ID: Phase B Soil Gas / 04020-023-4311

Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13

EPA TO-15

Rusty Bravo

6.0 L Summa Canister

Test Code:

Analyst:

Instrument ID:

Sampling Media: Test Notes:

CAS Project ID: P0801342 CAS Sample ID: P080508-MB

Date Collected: NA Date Received: NA Date Analyzed: 5/8/08 Volume(s) Analyzed: 1.00 Liter(s)

Canister Dilution Factor: 1.00

Date: 5/9/08 TO15SCAN.XLT - Tronox - Henderson - PageNo.

110

| CAS # | Compound | Result μg/m³ | MRL µg/m³ | MDL µg/m³ | Result ppbV | MRL ppbV | MDL ppbV Q | Data Jualifier |
|------------|---------------------------|-----------------|--------------|--------------|----------------|-------------|----------------------|-------------------|
| 637-92-3 | Ethyl tert-Butyl Ether | ND | 0.50 | 0.051 | ND | 0.12 | 0.012 | |
| 107-06-2 | 1,2-Dichloroethane | ND | 0.10 | 0.050 | ND | 0.025 | 0.012 | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 0.10 | 0.050 | ND | 0.018 | 0.0092 | |
| 71-43-2 | Benzene | ND | 0.10 | 0.050 | ND | 0.031 | 0.016 | |
| 56-23-5 | Carbon Tetrachloride | ND | 0.10 | 0.050 | ND | 0.016 | 0.0080 | |
| 994-05-8 | tert-Amyl Methyl Ether | ND | 0.50 | 0.050 | ND | 0.12 | 0.012 | |
| 78-87-5 | 1,2-Dichloropropane | ND | 0.10 | 0.050 | ND | 0.022 | 0.011 | |
| 75-27-4 | Bromodichloromethane | ND | 0.10 | 0.050 | ND | 0.015 | 0.0075 | |
| 79-01-6 | Trichloroethene | ND | 0.10 | 0.050 | ND | 0.019 | 0.0093 | |
| 123-91-1 | 1,4-Dioxane | ND | 0.50 | 0.061 | ND | 0.14 | 0.017 | |
| 80-62-6 | Methyl Methacrylate | ND | 0.50 | 0.075 | ND | 0.12 | 0.018 | |
| 142-82-5 | n-Heptane | ND | 0.50 | 0.064 | ND | 0.12 | 0.016 | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 0.50 | 0.052 | ND | 0.11 | 0.011 | |
| 108-10-1 | 4-Methyl-2-pentanone | ND | 0.50 | 0.056 | ND | 0.12 | 0.014 | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 0.50 | 0.063 | ND | 0.11 | 0.014 | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 0.10 | 0.050 | ND | 0.018 | 0.0092 | |
| 108-88-3 | Toluene | ND | 0.50 | 0.050 | ND | 0.13 | 0.013 | |
| 591-78-6 | 2-Hexanone | ND | 0.50 | 0.076 | ND | 0.12 | 0.019 | |
| 124-48-1 | Dibromochloromethane | ND | 0.10 | 0.068 | ND | 0.012 | 0.0080 | |
| 106-93-4 | 1,2-Dibromoethane | ND | 0.10 | 0.054 | ND | 0.013 | 0.0070 | |
| 111-65-9 | n-Octane | ND | 0.50 | 0.050 | ND | 0.11 | 0.011 | |
| 127-18-4 | Tetrachloroethene | ND | 0.10 | 0.050 | ND | 0.015 | 0.0074 | |
| 108-90-7 | Chlorobenzene | ND | 0.10 | 0.051 | ND | 0.022 | 0.011 | |

ND = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

Verified By:____

RC

RESULTS OF ANALYSIS Page 3 of 4

CAS Project ID: P0801342 CAS Sample ID: P080508-MB

| Test Code: | EPA TO-15 | Date Collected: N. | A |
|-----------------|--|---------------------|---------------|
| Instrument ID: | Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 | Date Received: N. | A |
| Analyst: | Rusty Bravo | Date Analyzed: 5/ | 8/08 |
| Sampling Media: | 6.0 L Summa Canister | Volume(s) Analyzed: | 1.00 Liter(s) |
| Test Notes: | | | |

Canister Dilution Factor: 1.00

| | | Result | MRL | MDL | Result | MRL | MDL | Data |
|-------------|-------------------------------|-------------------|-------------------|-------|--------|--------|--------|-----------|
| CAS # | Compound | μg/m ³ | μg/m ³ | μg/m³ | ppbV | ppbV | ppbV | Qualifier |
| 100-41-4 | Ethylbenzene | ND | 0.50 | 0.062 | ND | 0.12 | 0.014 | |
| 179601-23-1 | m,p-Xylenes | ND | 0.50 | 0.13 | ND | 0.12 | 0.030 | |
| 75-25-2 | Bromoform | ND | 0.50 | 0.076 | ND | 0.048 | 0.0074 | |
| 100-42-5 | Styrene | ND | 0.50 | 0.076 | ND | 0.12 | 0.018 | |
| 95-47-6 | o-Xylene | ND | 0.50 | 0.063 | ND | 0.12 | 0.015 | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 0.10 | 0.064 | ND | 0.015 | 0.0093 | |
| 98-82-8 | Cumene | 0.060 | 0.50 | 0.056 | 0.012 | 0.10 | 0.011 | J |
| 103-65-1 | n-Propylbenzene | ND | 0.50 | 0.052 | ND | 0.10 | 0.011 | |
| 622-96-8 | 4-Ethyltoluene | ND | 0.50 | 0.057 | ND | 0.10 | 0.012 | |
| 108-67-8 | 1,3,5-Trimethylbenzene | ND | 0.50 | 0.060 | ND | 0.10 | 0.012 | |
| 98-83-9 | alpha-Methylstyrene | ND | 0.50 | 0.073 | ND | 0.10 | 0.015 | |
| 95-63-6 | 1,2,4-Trimethylbenzene | ND | 0.50 | 0.069 | ND | 0.10 | 0.014 | |
| 100-44-7 | Benzyl Chloride | ND | 0.10 | 0.086 | ND | 0.019 | 0.017 | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 0.10 | 0.062 | ND | 0.017 | 0.010 | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 0.10 | 0.056 | ND | 0.017 | 0.0093 | |
| 135-98-8 | sec-Butylbenzene | ND | 0.50 | 0.058 | ND | 0.091 | 0.011 | |
| 99-87-6 | 4-Isopropyltoluene (p-Cymene) | ND | 0.50 | 0.065 | ND | 0.091 | 0.012 | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 0.10 | 0.066 | ND | 0.017 | 0.011 | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 0.50 | 0.076 | ND | 0.052 | 0.0079 | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 0.10 | 0.076 | ND | 0.013 | 0.010 | |
| 91-20-3 | Naphthalene | 0.10 | 0.20 | 0.074 | 0.020 | 0.038 | 0.014 | J |
| 87-68-3 | Hexachlorobutadiene | ND | 0.10 | 0.090 | ND | 0.0094 | 0.0084 | |
| 98-06-6 | tert-Butylbenzene | ND | 0.20 | 0.050 | ND | 0.036 | 0.0091 | |
| 104-51-8 | n-Butylbenzene | ND | 0.20 | 0.050 | ND | 0.036 | 0.0091 | |

ND = Compound was analyzed for, but not detected above the laboratory detection limit.

MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method. J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

> Verified By: Re- Date: 5/9/08 TO15SCAN.XLT - Tronox - Henderson - PageNo.:

111

Client:

ENSR

Client Project ID: Phase B Soil Gas / 04020-023-4311

Client Sample ID: Method Blank

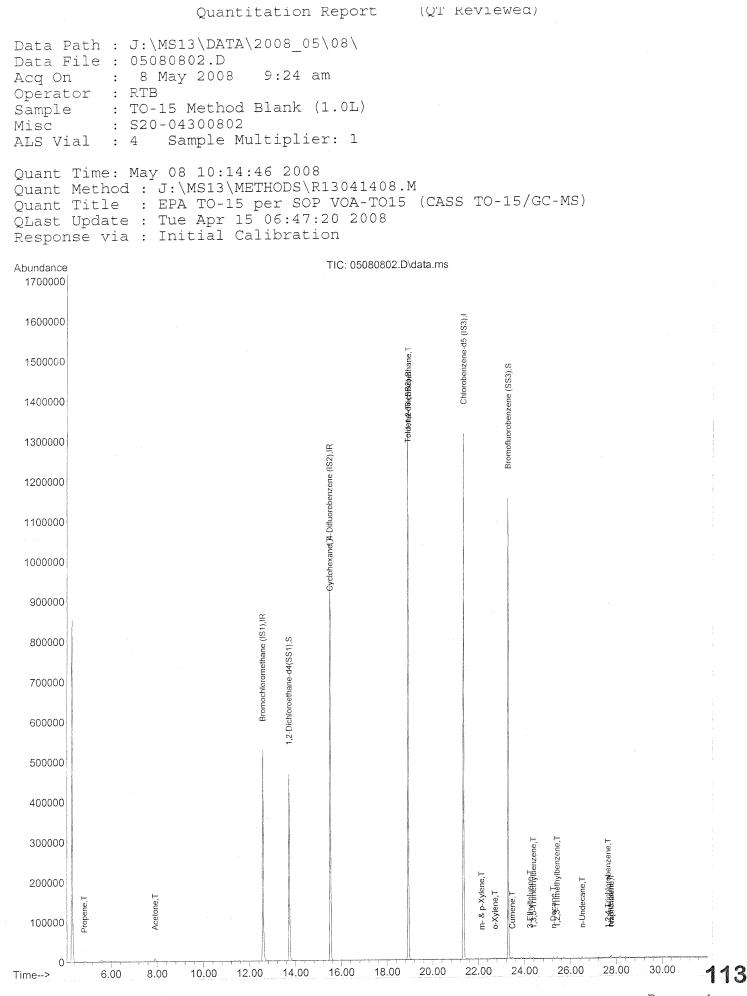
RESULTS OF ANALYSIS Page 4 of 4

| Client: | ENSR | | |
|---------------------------|--|---------------------|---------------|
| Client Sample ID: | Method Blank | CAS Project ID: P | 0801342 |
| Client Project ID: | Phase B Soil Gas / 04020-023-4311 | CAS Sample ID: PO | 080508-MB |
| | Tentatively Identified Comp | ounds | |
| Test Code: | EPA TO-15 | Date Collected: N | А |
| Instrument ID: | Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 | Date Received: N | А |
| Analyst: | Rusty Bravo | Date Analyzed: 5/ | 8/08 |
| Sampling Media: | 6.0 L Summa Canister | Volume(s) Analyzed: | 1.00 Liter(s) |
| Test Notes: | | | |

Canister Dilution Factor: 1.00

| GC/MS | Compound Identification | Concentration | Data |
|----------------|-------------------------|---------------------------------------|-----------|
| Retention Time | | μg/m³ | Qualifier |
| | No Compounds Detected | · · · · · · · · · · · · · · · · · · · | |

Verified By: Rec



R13041408.M Fri May 09 14:44:43 2008

Quantitation Report (QT Keviewed) Data Path : J:\MS13\DATA\2008_05\08\ Data File : 05080802.D Acq On : 8 May 2008 9:24 am Operator : RTB Sample : TO-15 Method Blank (1.0L) Misc : S20-04300802 ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 10:14:46 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards 1) Bromochloromethane (IS1)12.5813024134625.000 ng-0.0337) 1,4-Difluorobenzene (IS2)15.51114103764725.000 ng-0.0256) Chlorobenzene-d5 (IS3)21.358248933425.000 ng0.00 System Monitoring Compounds

 33) 1,2-Dichloroethane-d4(...
 13.72
 65
 432866
 22.365 ng
 -0.03

 Spiked Amount
 25.000
 Recovery
 =
 89.44%

 57) Toluene-d8 (SS2)
 18.93
 98
 1126967
 25.695 ng
 0.00

 Spiked Amount
 25.000
 Recovery
 =
 102.76%

 73) Bromofluorobenzene (SS3)
 23.29
 174
 378759
 25.094 ng
 0.00

 Recovery = 100.36% 🗸 Spiked Amount 25.000

R13041408.M Fri May 09 14:44:43 2008

Quantitation Report (QT Reviewed) Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080802.D Acq On : 8 May 2008 9:24 am Operator : RTB Sample : TO-15 Method Blank (1.0L) Misc : S20-04300802 ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 10:14:46 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) OLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration
 Internal Standards
 R.T. QIon
 Response
 Conc Units Dev

 321
 Chloroform
 0.00
 83
 0
 N.D.

 331
 Tetrahydrofuran
 0.00
 72
 0
 N.D.

 351
 Ethyl tetr-Butyl Ether
 0.00
 72
 0
 N.D.

 351
 J.j.pichloroethane
 12.74
 62
 72
 N.D.

 361
 1.2.Dichloroethane
 0.00
 97
 0
 N.D.

 361
 J.porchloroethane
 0.00
 11
 0
 N.D.

 37
 Isopropyl Acetate
 0.00
 61
 0
 N.D.

 410
 Eenzene
 14.98
 78
 716
 N.D.

 42
 Carbon Tetrachloride
 0.00
 73
 0
 N.D.

 431
 tert-Amyl Methyl Ether
 0.00
 63
 0
 N.D.

 44
 tert-Amyl Methyl Ether
 0.00
 70
 N.D.
 N.D.

 50
 Methyl Methacrylate
 0.00
 70
 N.D.
 N.D.
 </tr R.T. QIon Response Conc Units Dev(Min) Internal Standards 1 8 48 60 93 85 ⁹⁸115 Postoglog Page: 2

R13041408.M Fri May 09 14:44:43 2008

Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080802.D Acg On : 8 May 2008 9:24 am Operator : RTB Sample : TO-15 Method Blank (1.0L) Misc : S20-04300802 ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 10:14:46 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards ------80) alpha-Methylstyrene24.6711854N.D.81) 2-Ethyltoluene24.63105255N.D.82) 1.2.4-Trimethylbenzene24.89105933N.D.

 81)
 2-Ethyltoluene
 24.63
 105
 255
 N.D.

 82)
 1,2,4-Trimethylbenzene
 24.89
 105
 933
 N.D.

 83)
 n-Decane
 25.25
 57
 2264
 0.072 ng
 #

 84)
 Benzyl Chloride
 25.06
 91
 1062
 N.D.

 85)
 1,3-Dichlorobenzene
 25.09
 146
 533
 N.D.

 86)
 1,4-Dichlorobenzene
 25.16
 146
 647
 N.D.

 87)
 sec-Butylbenzene
 25.21
 105
 438
 N.D.

 88)
 p-Isopropyltoluene
 25.41
 119
 1975
 N.D.

 89)
 1,2,3-Trimethylbenzene
 25.60
 146
 52
 N.D.

 90)
 1,2-Dichlorobenzene
 25.60
 146
 52
 N.D.

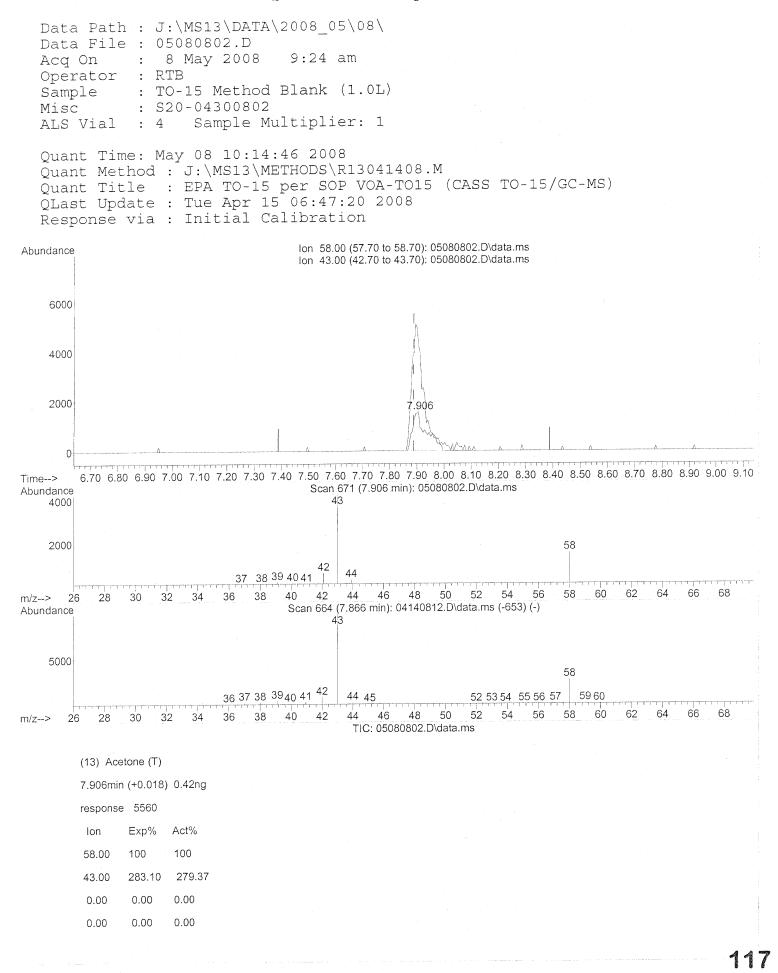
 91)
 d-Limonene
 0.00
 68
 0
 N.D.

 92)
 1,2-Dibromo-3-Chloropr...
 0.00
 157
 0
 N.D.

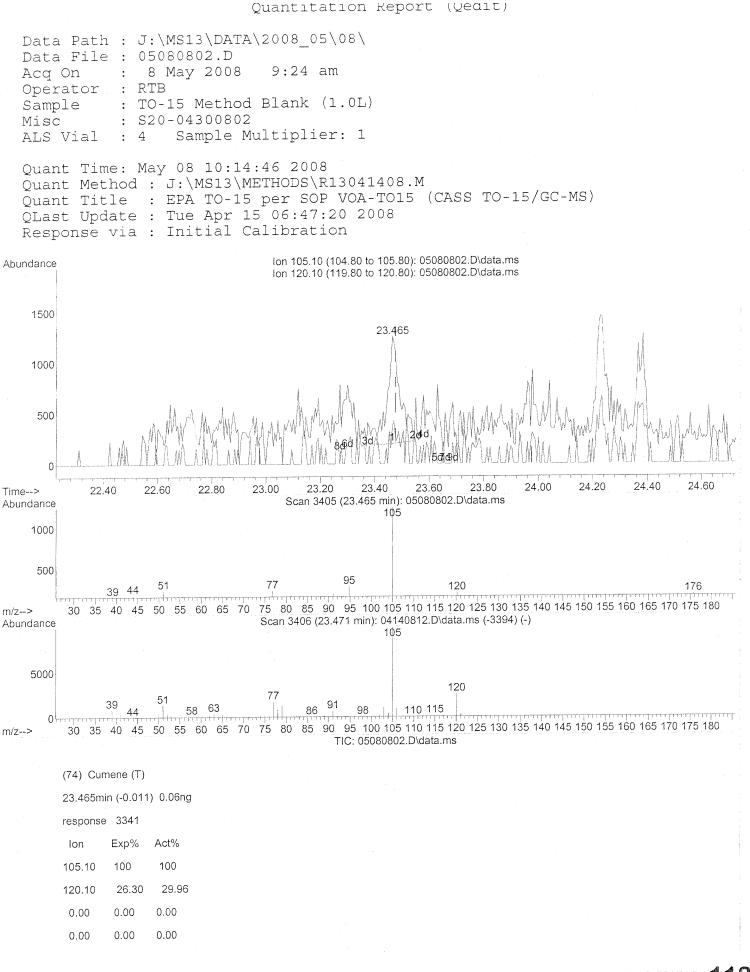
 93)
 n-Undecane
 26.51
 57
 1681
 0.045 ng
 #

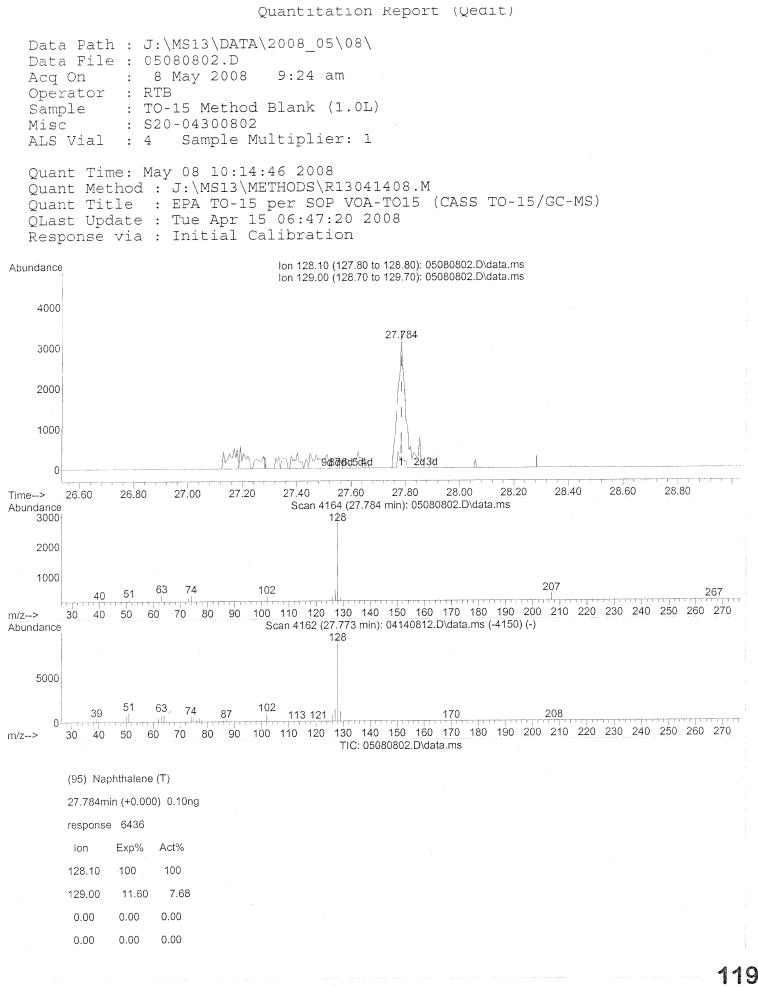
 94)
 1,2,4-Trichlorob 50 90 95 70 90 78 _____

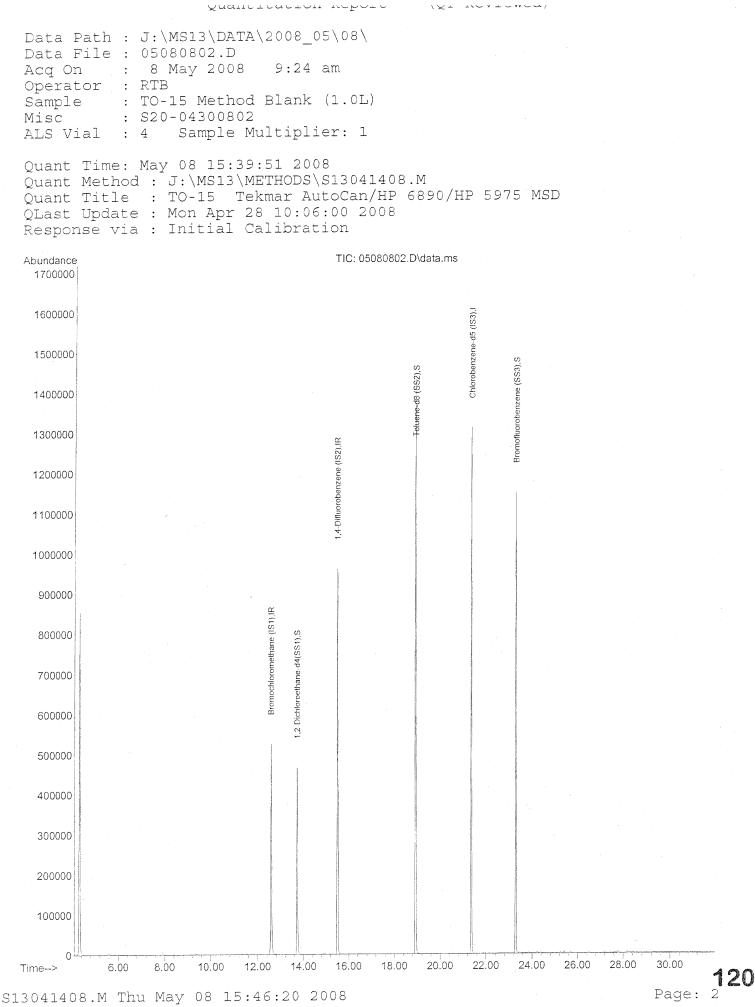
(#) = qualifier out of range (m) = manual integration (+) = signals summed



R13041408.M Fri May 09 10:09:49 2008







Quantituditon Report (VI ICOVICIO) Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080802.D Acq On : 8 May 2008 9:24 am Operator : RTB Sample : TO-15 Method Blank (1.0L) Misc : S20-04300802 ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 15:39:51 2008 Quant Method : J:\MS13\METHODS\S13041408.M Quant Title : TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD QLast Update : Mon Apr 28 10:06:00 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards _____ 1) Bromochloromethane (IS1)12.5813024134625.000 ng-0.033) 1,4-Difluorobenzene (IS2)15.51114103764725.000 ng-0.024) Chlorobenzene-d5 (IS3)21.358248933425.000 ng0.00 System Monitoring Compounds 2) 1,2-Dichloroethane-d4(... 13.72 65 432866 22.365 ng -0.03 Recovery = 89.44% 🗸

 Spiked Amount
 25.000
 Recovery
 =
 89.44%

 5) Toluene-d8 (SS2)
 18.93
 98
 1126967
 25.695 ng
 0.00

 Spiked Amount
 25.000
 Recovery
 =
 102.76%

 6) Bromofluorobenzene (SS3)
 23.29
 174
 378759
 25.094 ng
 0.00

 Recovery = 100.36% 🗸 Spiked Amount 25.000 Ovalue Target Compounds 7)tert-Butylbenzene24.871194258)n-Butylbenzene25.9191147 N.D. N.D.

(#) = qualifier out of range (m) = manual integration (+) = signals summed

205/08/08

121 Page: 1

;13041408.M Thu May 08 15:46:20 2008

QC SUMMARY FORMS

SURROGATE SPIKE RECOVERY RESULTS

Page 1 of 1

Client:ENSRClient Project ID:Phase B Soil Gas / 04020-023-4311

CAS Project ID: P0801342

Test Code: Instrument ID: Analyst: Sampling Media: Test Notes: EPA TO-15 Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 Rusty Bravo 6.0 L Summa Canister(s)

Date(s) Collected: 5/7/08 Date(s) Received: 5/8/08 Date(s) Analyzed: 5/8/08

| | | 1,2-Dichlor | oethane-d4 | Tolue | ene-d8 | Bromofluo | robenzene | |
|--------------------|-----------------|-------------|------------|-----------|------------|-----------|------------|-----------|
| Client Sample ID | CAS Sample ID | % | Acceptance | % | Acceptance | % | Acceptance | Data |
| - | | Recovered | Limits | Recovered | Limits | Recovered | Limits | Qualifier |
| Method Blank | P080508-MB | 89 | 70-130 | 103 | 70-130 | 100 | 70-130 | |
| Lab Control Sample | P080508-LCS | 91 | 70-130 | 103 | 70-130 | 101 | 70-130 | |
| SG83B-05-1 | P0801342-001 | 92 | 70-130 | 102 | 70-130 | 101 | 70-130 | |
| SG83B-05-3 | P0801342-002 | 91 | 70-130 | 106 | 70-130 | 102 | 70-130 | |
| SG83B-05-7 | P0801342-003 | 91 | 70-130 | 102 | 70-130 | 98 | 70-130 | |
| SG83B-05-7 | P0801342-003DUP | 91 | 70-130 | 103 | 70-130 | 99 | 70-130 | |

RE

Date: 5/9/05/ TO15SCAN.XLT - Tronox - Henderson - PageNo.:

LABORATORY CONTROL SAMPLE SUMMARY Page 1 of 3

| Client: | ENSR |
|--------------------|-----------------------------------|
| Client Sample ID: | Lab Control Sample |
| Client Project ID: | Phase B Soil Gas / 04020-023-4311 |

CAS Project ID: P0801342 CAS Sample ID: P080508-LCS

EPA TO-15 Test Code: Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 Instrument ID: Rusty Bravo Analyst: 6.0 L Summa Canister Sampling Media: Test Notes:

Date Collected: NA Date Received: NA Date Analyzed: 5/08/08 Volume(s) Analyzed: NA Liter(s)

| CAS # | Compound | Spike Amount ng | Result ng | % Recovery | CAS Acceptance Limits | Data Qualifier |
|-----------|--|--------------------|--------------|------------|-----------------------------|---|
| 75-71-8 | Dichlorodifluoromethane (CFC 12) | 25.5 | 21.9 | 86 | 69-117 | *************************************** |
| 74-87-3 | Chloromethane | 24.5 | 20.8 | 85 | 53-131 | |
| 76-14-2 | 1,2-Dichloro-1,1,2,2- tetrafluoroethane (CFC 114) | 26.0 | 22.6 | 87 | 58-133 | |
| 75-01-4 | Vinyl Chloride | 24.8 | 20.7 | 83 | 61-127 | |
| 74-83-9 | Bromomethane | 25.0 | 23.9 | 96 | 67-124 | |
| 75-00-3 | Chloroethane | 25.0 | 23.2 | 93 | 69-123 | |
| 64-17-5 | Ethanol | 23.8 | 20.6 | 87 | 56-137 | |
| 67-64-1 | Acetone | 26.8 | 24.9 | 93 | 63-116 | |
| 75-69-4 | Trichlorofluoromethane | 26.3 | 24.4 | 93 | 71-120 | |
| 107-13-1 | Acrylonitrile | 25.5 | 25.5 | 100 | 74-129 | |
| 75-35-4 | 1,1-Dichloroethene | 27.8 | 25.8 | 93 | 77-116 | |
| 75-65-0 | 2-Methyl-2-Propanol (tert-Butyl Alcohol) | 25.8 | 24.4 | 95 | 35-141 | |
| 75-09-2 | Methylene Chloride | 27.8 | 24.0 | 86 | 71-113 | |
| 107-05-1 | 3-Chloro-1-propene (Allyl Chloride) | 26.8 | 29.6 | 110 | 75-127 | |
| 76-13-1 | Trichlorotrifluoroethane | 27.8 | 25.4 | 91 | 63-129 | |
| 75-15-0 | Carbon Disulfide | 25.0 | 23.1 | 92 | 72-122 | |
| 156-60-5 | trans-1,2-Dichloroethene | 26.5 | 24.5 | 92 | 74-118 | |
| 75-34-3 | 1,1-Dichloroethane | 26.8 | 24.6 | 92 | 74-118 | |
| 1634-04-4 | Methyl tert-Butyl Ether | 26.8 | 24.2 | 90 | 72-119 | |
| 108-05-4 | Vinyl Acetate | 25.3 | 28.0 | 111 | 32-163 | |
| 78-93-3 | 2-Butanone (MEK) | 27.0 | 27.0 | 100 | 71-122 | |
| 156-59-2 | cis-1,2-Dichloroethene | 27.0 | 24.4 | 90 | 74-117 | |
| 108-20-3 | Diisopropyl Ether | 26.3 | 22.8 | 87 | 70-131 | |
| 67-66-3 | Chloroform | 29.8 | 28.3 | 95 | 72-113 | |

Verified By: <u>Rc-</u>

Date: 519108 TOISSCAN.XLT - Tronox - Henderson - PageNo.:

124

LABORATORY CONTROL SAMPLE SUMMARY

Page 2 of 3

Client:ENSRClient Sample ID:Lab Control SampleClient Project ID:Phase B Soil Gas / 04020-023-4311

CAS Project ID: P0801342 CAS Sample ID: P080508-LCS

Test Code:EPA TO-15Instrument ID:Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13Analyst:Rusty BravoSampling Media:6.0 L Summa CanisterTest Notes:Test Notes:

Date Collected: NA Date Received: NA Date Analyzed: 5/08/08 Volume(s) Analyzed: NA Liter(s)

| CAS # | Compound | Spike Amount | Result | % Recovery | CAS Acceptance | Data |
|------------|---------------------------|--------------|--------|------------|-------------------|-----------|
| | | ng | ng | | Limits | Qualifier |
| 637-92-3 | Ethyl tert-Butyl Ether | 26.0 | 24.0 | 92 | 74-123 | |
| 107-06-2 | 1,2-Dichloroethane | 26.3 | 23.6 | 90 | 72-117 | |
| 71-55-6 | 1,1,1-Trichloroethane | 26.8 | 25.5 | 95 | 78-114 | |
| 71-43-2 | Benzene | 27.0 | 24.9 | 92 | 73-111 | |
| 56-23-5 | Carbon Tetrachloride | 26.0 | 27.2 | 105 | 78-126 | |
| 994-05-8 | tert-Amyl Methyl Ether | 26.0 | 24.9 | 96 | 81-118 | |
| 78-87-5 | 1,2-Dichloropropane | 26.5 | 24.1 | 91 | 78-117 | |
| 75-27-4 | Bromodichloromethane | 27.8 | 26.8 | 96 | 77-120 | |
| 79-01-6 | Trichloroethene | 27.3 | 26.4 | 97 | 80-116 | |
| 123-91-1 | 1,4-Dioxane | 27.5 | 27.8 | 101 | 79-122 | |
| 80-62-6 | Methyl Methacrylate | 25.8 | 26.3 | 102 | 79-128 | |
| 142-82-5 | n-Heptane | 26.8 | 24.3 | 91 | 77-117 | |
| 10061-01-5 | cis-1,3-Dichloropropene | 25.0 | 25.9 | 104 | 78-112 | |
| 108-10-1 | 4-Methyl-2-pentanone | 27.5 | 24.6 | 89 | 78-128 | |
| 10061-02-6 | trans-1,3-Dichloropropene | 28.0 | 29.8 | 106 | 81-121 | |
| 79-00-5 | 1,1,2-Trichloroethane | 26.3 | 24.7 | 94 | 80-117 | |
| 108-88-3 | Toluene | 26.5 | 27.0 | 102 | 76-116 | |
| 591-78-6 | 2-Hexanone | 26.3 | 25.9 | 98 | 69-131 | |
| 124-48-1 | Dibromochloromethane | 27.0 | 30.2 | 112 | 80-128 | |
| 106-93-4 | 1,2-Dibromoethane | 26.3 | 29.6 | 113 | 79-122 | |
| 111-65-9 | n-Octane | 26.0 | 26.2 | 101 | 78-122 | |
| 127-18-4 | Tetrachloroethene | 26.0 | 27.1 | 104 | 77-118 | |
| 108-90-7 | Chlorobenzene | 26.5 | 27.0 | 102 | 78-117 | |

LABORATORY CONTROL SAMPLE SUMMARY

Page 3 of 3

Client:ENSRClient Sample ID:Lab Control SampleClient Project ID:Phase B Soil Gas / 04020-023-4311

CAS Project ID: P0801342 CAS Sample ID: P080508-LCS

Test Code:EPA TO-15Instrument ID:Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13Analyst:Rusty BravoSampling Media:6.0 L Summa CanisterTest Notes:Test Notes:

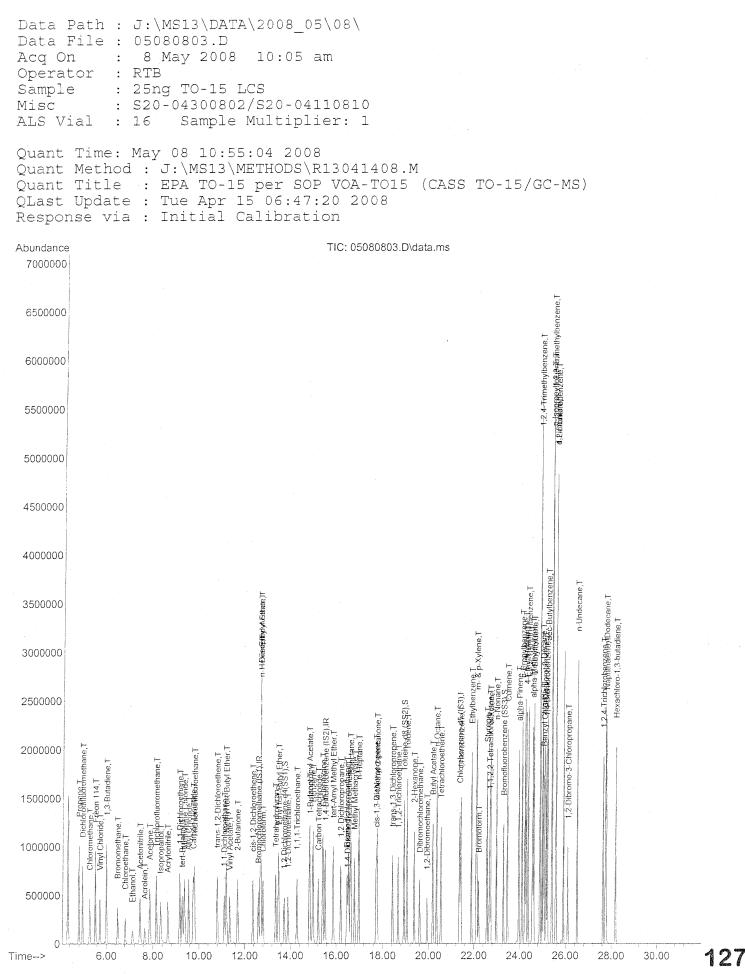
Date Collected: NA Date Received: NA Date Analyzed: 5/08/08 Volume(s) Analyzed: NA Liter(s)

| | | | | | CAS | |
|-------------|-------------------------------|--------------|--------|------------|------------|-----------|
| CAS # | Compound | Spike Amount | Result | % Recovery | Acceptance | Data |
| | | ng | ng | | Limits | Qualifier |
| 100-41-4 | Ethylbenzene | 26.3 | 27.4 | 104 | 79-116 | |
| 179601-23-1 | m,p-Xylenes | 62.5 | 64.8 | 104 | 80-117 | |
| 75-25-2 | Bromoform | 31.3 | 38.5 | 123 | 77-128 | |
| 100-42-5 | Styrene | 26.3 | 27.6 | 105 | 80-124 | |
| 95-47-6 | o-Xylene | 29.8 | 30.4 | 102 | 80-116 | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 29.8 | 31.0 | 104 | 79-120 | |
| 98-82-8 | Cumene | 27.0 | 28.8 | 107 | 81-119 | |
| 103-65-1 | n-Propylbenzene | 26.3 | 28.3 | 108 | 82-120 | |
| 622-96-8 | 4-Ethyltoluene | 26.5 | 28.4 | 107 | 80-119 | |
| 108-67-8 | 1,3,5-Trimethylbenzene | 26.0 | 27.4 | 105 | 80-120 | |
| 98-83-9 | alpha-Methylstyrene | 25.5 | 26.0 | 102 | 54-146 | |
| 95-63-6 | 1,2,4-Trimethylbenzene | 26.0 | 27.3 | 105 | 80-122 | |
| 100-44-7 | Benzyl Chloride | 25.8 | 30.5 | 118 | 85-131 | |
| 541-73-1 | 1,3-Dichlorobenzene | 25.5 | 27.0 | 106 | 81-117 | |
| 106-46-7 | 1,4-Dichlorobenzene | 26.3 | 28.1 | 107 | 81-119 | |
| 135-98-8 | sec-Butylbenzene | 26.8 | 28.7 | 107 | 80-124 | |
| 99-87-6 | 4-Isopropyltoluene (p-Cymene) | 28.8 | 31.6 | 110 | 78-124 | |
| 95-50-1 | 1,2-Dichlorobenzene | 25.8 | 26.1 | 101 | 81-122 | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | 25.8 | 31.7 | 123 | 91-136 | |
| 120-82-1 | 1,2,4-Trichlorobenzene | 26.0 | 28.5 | 110 | 75-138 | |
| 91-20-3 | Naphthalene | 26.3 | 28.6 | 109 | 76-143 | |
| 87-68-3 | Hexachlorobutadiene | 26.3 | 28.8 | 110 | 72-128 | |
| 98-06-6 | tert-Butylbenzene | 26.3 | 22.7 | 86 | 70-130 | |
| 104-51-8 | n-Butylbenzene | 26.8 | 23.5 | 88 | 70-130 | |

Date: 5/9/08

TO15SCAN.XLT - Tronox - Henderson - PageNo.:

26



R13041408.M Thu May 08 10:55:26 2008

| Quantitation | Keport | (Q. | r keviewea |) | | |
|--|---|---|--|--|---|--|
| Data Path : J:\MS13\DATA\2008_05 Data File : 05080803.D Acq On : 8 May 2008 10:05 a Operator : RTB Sample : 25ng TO-15 LCS Misc : S20-04300802/S20-041 ALS Vial : 16 Sample Multipli | m | | | | | |
| Quant Time: May 08 10:55:04 2008 Quant Method : J:\MS13\METHODS\R Quant Title : EPA TO-15 per SOP QLast Update : Tue Apr 15 06:47: Response via : Initial Calibrati | 1304140 VOA-TO 20 2008 | 15 (CA | ASS TO-15/(| GC-MS) | | |
| Internal Standards | | | Response | | Units | Dev(Min) |
| 1) Bromochloromethane (IS1) 37) 1,4-Difluorobenzene (IS2) 56) Chlorobenzene-d5 (IS3) | 12.59 15.52 | 130 114 | 244114 1061640 | 25.000 25.000 | ng | -0.01 |
| System Monitoring Compounds 33) 1,2-Dichloroethane-d4(Spiked Amount 25.000 57) Toluene-d8 (SS2) Spiked Amount 25.000 | 13.73 18.93 | 65 98 | 443758 Recove 1158483 Recove | 22.667 ery = 25.707 ery = | ng 90 ng 102 | -0.02 .68% ⁄ 0.00 .84% ∕ |
| 73) Bromofluorobenzene (SS3) Spiked Amount 25.000 | 23.29 | 174 | 390550 Recove | 25.184 | ng | 0.00 |
| <pre>3) Dichlorodifluoromethane 4) Chloromethane 5) Freon 114 6) Vinyl Chloride 7) 1,3-Butadiene 8) Bromomethane 9) Chloroethane 10) Ethanol 11) Acetonitrile 12) Acrolein 13) Acetone 14) Trichlorofluoromethane 15) Isopropanol 16) Acrylonitrile 17) 1,1-Dichloroethene 18) tert-Butanol 19) Methylene Chloride 20) Allyl Chloride 21) Trichlorotrifluoroethane 22) Carbon Disulfide 23) trans-1,2-Dichloroethene 24) 1,1-Dichloroethane 25) Methyl tert-Butyl Ether 26) Vinyl Acetate 27) 2-Butanone 28) cis-1,2-Dichloroethene 29) Diisopropyl Ether 30) Ethyl Acetate 31) n-Hexane</pre> | 4.95 5.27 5.52 5.72 6.00 6.48 6.82 7.13 7.44 7.64 7.64 7.64 9.16 9.27 9.36 9.55 9.76 10.80 11.19 11.34 12.36 12.69 12.69 12.70 | 85 50 135 64 94 645 16 81 581 536 94 156 63 76 87 61 87 61 | 407641 593092 588675 328456 275127 284504 809787 218995 336805 705880 977864m 534167 347213 924796 371975 614215 314484 1329828 575163 672011 1086599 75223 255480 540182 284170 153450 695486 | 21.912 20.752 22.602 20.701 26.572 23.180 20.648 22.363 22.491 24.924 24.924 25.476 25.757 24.384 25.476 25.476 25.476 25.476 25.476 23.964 29.618 25.430 23.105 24.479 24.575 24.575 24.397 22.839 25.977 22.652 | ng ng ng ng ng ng ng ng ng ng ng ng ng n | 99 97 99 96 77 100 96 95 96 99 # 66 99 # 66 99 # 66 99 # 98 85 94 89 100 95 97 89 95 87 89 95 87 # 90 95 97 89 95 97 89 95 97 89 91 28 |
| 13041408.M Thu May 08 10:55:26 20 | 8 | | 50 | 5/08/08 | Ē | Page: 1 |

Quantitation Report (QT Reviewed) Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080803.D Acg On : 8 May 2008 10:05 am Operator : RTB Sample : 25ng TO-15 LCS Misc : S20-04300802/S20-04110810 ALS Vial : 16 Sample Multiplier: 1 Quant Time: May 08 10:55:04 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration

 Internal Standards
 R.T. Qion Response Conc Units Dev(Min)

 32) Chloroform
 12.80
 83
 646002
 28.343 rg
 99

 34) Tetrahydrofuran
 13.49
 87
 2244651
 25.812 rg
 95

 35) Ethyl tetr-Buryl Ether
 13.49
 87
 2244651
 25.812 rg
 95

 35) 1.2. Jichloroschane
 13.49
 87
 52382
 24.451 rg
 98

 36) 1.2. Jichloroschane
 14.25
 97
 563082
 25.8454 rg
 99

 37) Isopropyl Acetate
 14.85
 56
 324004
 22.8189 rg
 99

 40) 1-Butancl
 14.85
 56
 324004
 22.8189 rg
 99

 41) Carbon Tetrachloride
 15.22
 117
 505854
 27.157 rg
 98

 42) Carbon Tetrachloride
 15.41
 94
 523622
 25.852 rg
 97

 43) Cyclohexane
 16.46
 83
 513269
 26.805 rg
 100

 17 Tothloroethane
 16.46
 83
 513269
 27.61 rg
 92

 44) Lett-Anyl Methyl Ether
 17.77
 58
 36581
 22.812 rg
 92

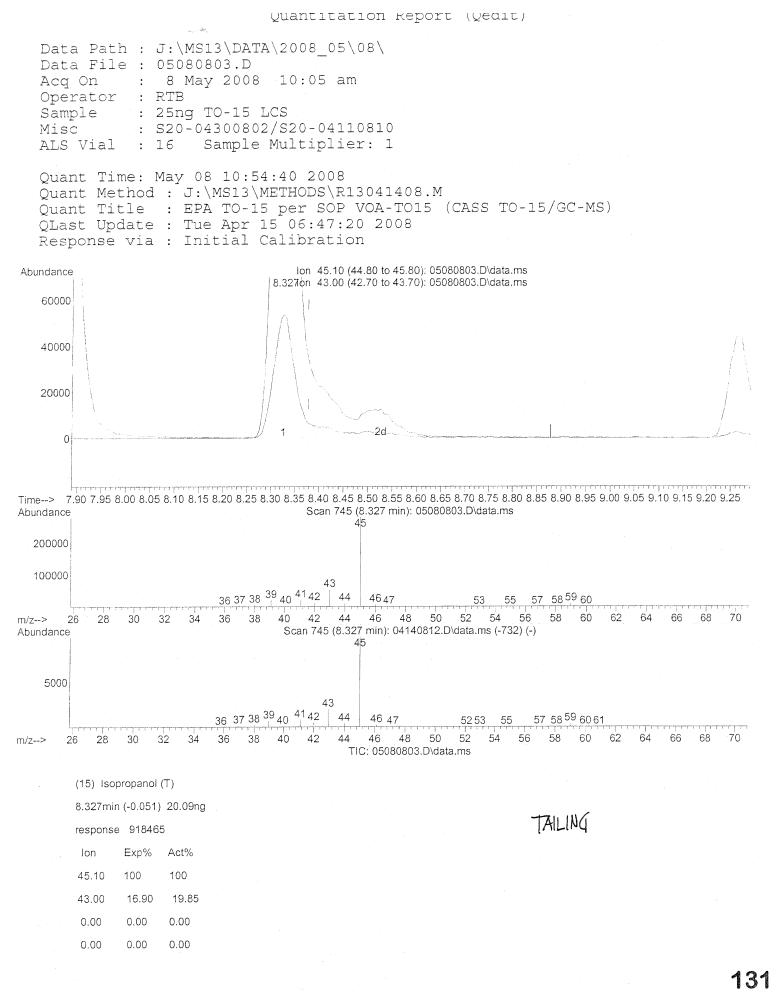
 Internal Standards R.T. QIon Response Conc Units Dev(Min)

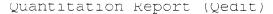
Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080803.D Acq On : 8 May 2008 10:05 am Operator : RTB Sample : 25ng TO-15 LCS Misc : S20-04300802/S20-04110810 ALS Vial : 16 Sample Multiplier: 1 Quant Time: May 08 10:55:04 2008 Ouant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration Internal Standards R.T. QIon Response Conc Units Dev(Min) 80) alpha-Methylstyrene24.5611870747025.970 ng81) 2-Ethyltoluene24.61105168115226.584 ng82) 1,2,4-Trimethylbenzene24.88105158618627.287 ng83) n-Decane24.985787585927.246 ng84) Benzyl Chloride25.0591120731030.469 ng85) 1,3-Dichlorobenzene25.0814684886826.974 ng86) 1,4-Dichlorobenzene25.1514684773428.149 ng87) sec-Butylbenzene25.21105195581728.658 ng88) p-Isopropyltoluene25.40119187886331.568 ng89) 1,2,3-Trimethylbenzene25.5814684125826.084 ng90) 1,2-Dichlorobenzene25.586863628224.141 ng91) d-Limonene26.505792098627.289 ng92) 1,2,4-Trichlorobenzene27.6218057050628.542 ng95) Naphthalene27.745787309225.497 ng 96 98 97 87 95 99 98 96 91 97 100 93 71 87 96 95) Naphthalene27.77128183533328.558ng96) n-Dodecane27.745787309225.497ng 98 85 97) Hexachloro-1,3-butadiene 28.19 225 370854 28.811 ng 99 _____

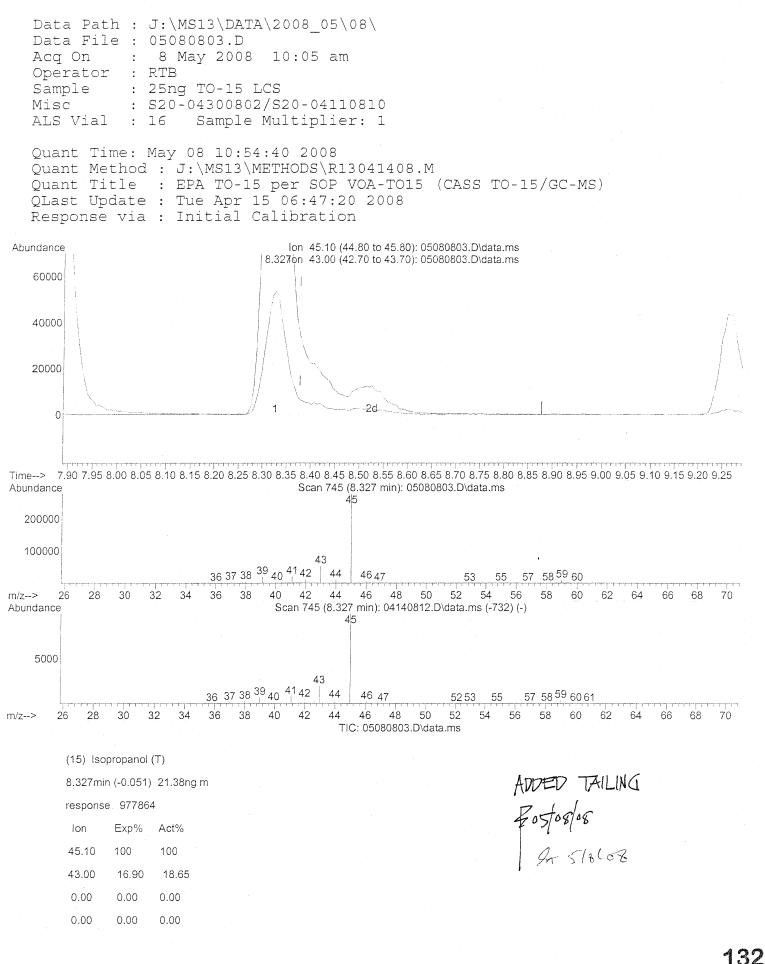
(#) = gualifier out of range (m) = manual integration (+) = signals summed

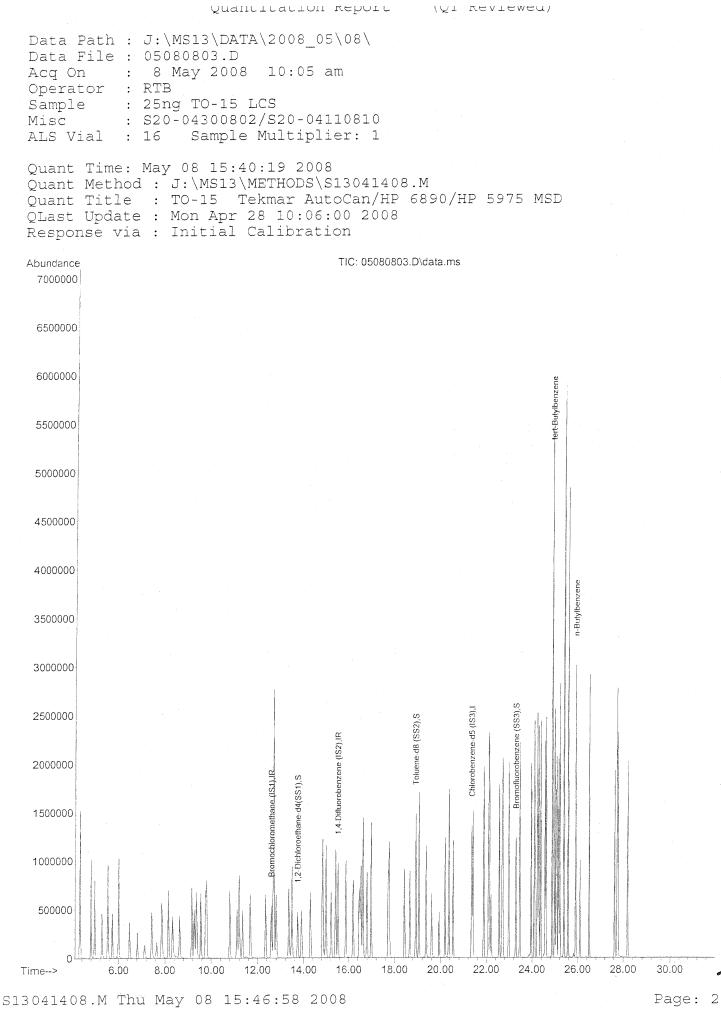
R13041408.M Thu May 08 10:55:26 2008

Fostorlos









Zaarreteneter rectore Data Path : J:\MS13\DATA\2008_05\08\ Data File : 05080803.D Acq On : 8 May 2008 10:05 am Operator : RTB Sample : 25ng TO-15 LCS Misc : S20-04300802/S20-04110810 ALS Vial : 16 Sample Multiplier: 1 Quant Time: May 08 15:40:19 2008 Quant Method : J:\MS13\METHODS\S13041408.M Quant Title : TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD QLast Update : Mon Apr 28 10:06:00 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards _____ 1) Bromochloromethane (IS1)12.5913024411425.000 ng-0.013) 1,4-Difluorobenzene (IS2)15.52114106164025.000 ng-0.014) Chlorobenzene-d5 (IS3)21.358250277225.000 ng0.00 System Monitoring Compounds 2) 1,2-Dichloroethane-d4(... 13.73 65 443758 22.667 ng -0.02 Recovery = 90.68% 🗸

 Spiked Amount
 25.000
 Recovery = 90.68%/

 5) Toluene-d8 (SS2)
 18.93
 98
 1158483
 25.707 ng
 0.00

 Spiked Amount
 25.000
 Recovery = 102.84%/

 6) Bromofluorobenzene (SS3)
 23.29
 174
 390550
 25.184 ng
 0.00

 Recovery = 100
 72%

 Spiked Amount 25.000 Recovery = 100.72% 🖌 Spiked Amount 25.000 Ovalue Target Compounds 7) tert-Butylbenzene24.88119147646226.818ng8) n-Butylbenzene25.9191170294928.670ng - 99 94

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Fistoslas

LABORATORY DUPLICATE SUMMARY RESULTS

Page 1 of 3

| Client: | ENSR | |
|--------------------------|--|-------------|
| Client Sample ID: | SG83B-05-7 | CAS Pro |
| Client Project ID: | Phase B Soil Gas / 04020-023-4311 | CAS Sat |
| Test Code: | EPA TO-15 | Date C |
| Instrument ID: | Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 | Date R |
| Analyst: | Rusty Bravo | Date A |
| Sampling Media: | 6.0 L Summa Canister | Volume(s) A |
| Test Notes: | | |
| Container ID: | SC00791 | |
| | | |

CAS Project ID: P0801342 CAS Sample ID: P0801342-003DUP

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

Initial Pressure (psig): -4.2

TANG

Final Pressure (psig): 3.5

Canister Dilution Factor: 1.73

Date: 5 900 S TOISSCAN.XLT - Tronox - Henderson - PageNo.:

135

| | | Duplicate | | | | | | |
|--|---------------|-----------|---------------|-------|---------|-------|-------|-----------|
| Compound | Sample Result | | Sample Result | | Average | % RPD | RPD | Data |
| | $\mu g/m^3$ | ppbV | μg/m³ | ppbV | μg/m³ | | Limit | Qualifier |
| Dichlorodifluoromethane (CFC 12) | ND | ND | ND | ND | - | - | 25 | |
| Chloromethane | ND | ND | ND | ND | - | - | 25 | |
| 1,2-Dichloro-1,1,2,2-tetrafluoroethane (CFC 114) | ND | ND | ND | ND | - | - | 25 | |
| Vinyl Chloride | ND | ND | ND | ND | · - | · · · | 25 | |
| Bromomethane | ND | ND | ND | ND | - | - | 25 | |
| Chloroethane | ND | ND | ND | ND | - | - | 25 | |
| Ethanol | ND | ND | ND | ND | - | - | 25 | |
| Acetone | 114 | 48.1 | 107 | 45.2 | 110.5 | 6 | 25 | J, B |
| Trichlorofluoromethane | 1,520 | 271 | 1,530 | 273 | 1525 | 0.7 | 25 | |
| Acrylonitrile | ND | ND | ND | ND | - | | 25 | |
| 1,1-Dichloroethene | ND | ND | ND | ND | - | - | 25 | |
| 2-Methyl-2-Propanol (tert-Butyl Alcohol) | ND | ND | ND | ND | - | - | 25 | |
| Methylene Chloride | 9.34 | 2.69 | ND | ND | - | - | 25 | J |
| 3-Chloro-1-propene (Allyl Chloride) | ND | ND | ND | ND | - | - | 25 | |
| Trichlorotrifluoroethane | ND | ND | ND | ND | - | - | 25 | |
| Carbon Disulfide | ND | ND | ND | ND | - | - | 25 | |
| trans-1,2-Dichloroethene | ND | ND | ND | ND | - | - | 25 | |
| 1,1-Dichloroethane | ND | ND | ND | ND | - | - | 25 | |
| Methyl tert-Butyl Ether | ND | ND | ND | ND | - | - | 25 | |
| Vinyl Acetate | ND | ND | ND | ND | - | - | 25 | |
| 2-Butanone (MEK) | 23.0 | 7.80 | 18.7 | 6.34 | 20.85 | 21 | 25 | J |
| cis-1,2-Dichloroethene | ND | ND | ND | ND | - | - | 25 | |
| Diisopropyl Ether | ND | ND | ND | ND | - | - | 25 | |
| Chloroform | 54,300 | 11,100 | 45,700 | 9,370 | 50000 | 17 | 25 | - |

ND = Compound was analyzed for, but not detected above the laboratory detection limit.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

B = Analyte was found in the method blank.

LABORATORY DUPLICATE SUMMARY RESULTS

Page 2 of 3

Client:ENSRClient Sample ID:SG83B-05-7Client Project ID:Phase B Soil Gas / 04020-023-4311

6.0 L Summa Canister

EPA TO-15

Rusty Bravo

SC00791

Test Code:

Analyst:

Test Notes:

Container ID:

Instrument ID:

Sampling Media:

CAS Project ID: P0801342 CAS Sample ID: P0801342-003DUP

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

Initial Pressure (psig):

Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13

-4.2

Final Pressure (psig): 3.5

| | | | | | Canis | ter Dilutio | n Factor: | 1.73 |
|---------------------------|--------|--------|--------|--------|---------|-------------|-----------|-----------|
| | | | Dupli | cate | | | | |
| Compound | Sample | Result | Sample | Result | Average | % RPD | RPD | Data |
| - | μg/m³ | ppbV | µg/m³ | ppbV | μg/m³ | | Limit | Qualifier |
| Ethyl tert-Butyl Ether | ND | ND | ND | ND | - | - | 25 | |
| 1,2-Dichloroethane | ND | ND | ND | ND | - | - | 25 | |
| 1,1,1-Trichloroethane | ND | ND | ND | ND | - | - | 25 | |
| Benzene | 101 | 31.7 | 102 | 32.1 | 101.5 | 1 | 25 | |
| Carbon Tetrachloride | 12,000 | 1,910 | 12,300 | 1,950 | 12150 | 2 | 25 | |
| tert-Amyl Methyl Ether | ND | ND | ND | ND | - | - | 25 | |
| 1,2-Dichloropropane | ND | ND | ND | ND | - | - | 25 | |
| Bromodichloromethane | ND | ND | ND | ND | - | - | 25 | |
| Trichloroethene | 11.4 | 2.13 | 11.9 | 2.22 | 11.65 | 4 | 25 | J |
| 1,4-Dioxane | ND | ND | ND | ND | - | | 25 | |
| Methyl Methacrylate | ND | ND | ND | ND | | - | 25 | |
| n-Heptane | ND | ND | ND | ND | - | - | 25 | |
| cis-1,3-Dichloropropene | ND | ND | ND | ND | - | - | 25 | |
| 4-Methyl-2-pentanone | ND | ND | ND | ND | - | - | 25 | |
| trans-1,3-Dichloropropene | ND | ND | ND | ND | - | | 25 | |
| 1,1,2-Trichloroethane | ND | ND | ND | ND | - | - | 25 | |
| Toluene | ND | ND | ND | ND | - | - | 25 | |
| 2-Hexanone | ND | ND | ND | ND | - | - | 25 | |
| Dibromochloromethane | ND | ND | ND | ND | - | - | 25 | |
| 1,2-Dibromoethane | ND | ND | ND | ND | - | | 25 | |
| n-Octane | ND | ND | ND | ND | - | - | 25 | |
| Tetrachloroethene | 127 | 18.7 | 124 | 18.3 | 125.5 | 2 | 25 | |
| Chlorobenzene | 180 | 39.1 | 197 | 42.7 | 188.5 | 9 | 25 | |

ND = Compound was analyzed for, but not detected above the laboratory detection limit.

J = The analyte was positively identified below the method reporting limit; the associated numerical value is considered estimated.

Verified By: Rc-

Date: 519108 **136** TOISSCAN.XLT - Tronox - Henderson - PageNo.:

LABORATORY DUPLICATE SUMMARY RESULTS

Page 3 of 3

Client:ENSRClient Sample ID:SG83B-05-7Client Project ID:Phase B Soil Gas / 04020-023-4311

6.0 L Summa Canister

EPA TO-15

Rusty Bravo

SC00791

Test Code: Instrument ID:

Analyst:

Test Notes:

Container ID:

Sampling Media:

CAS Project ID: P0801342 CAS Sample ID: P0801342-003DUP

D'1 /

Date: 5/9/05 TO15SCAN.XLT - Tronox - Henderson - PageNo.

1 70

Date Collected: 5/7/08 Date Received: 5/8/08 Date Analyzed: 5/8/08 Volume(s) Analyzed: 0.010 Liter(s) 0.0010 Liter(s)

Initial Pressure (psig):

Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13

-4.2

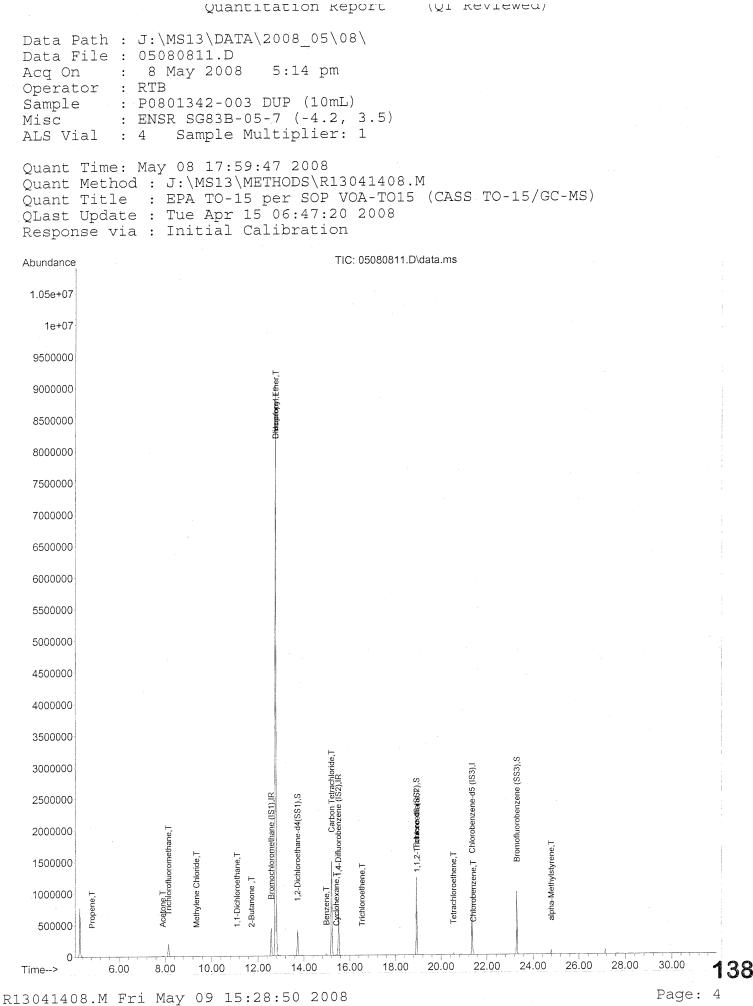
Final Pressure (psig): 3.5

| | Canister Dilution Factor: 1.73 | | | | | 1.73 | | |
|-------------------------------|--------------------------------|---------------|-------|-----------------|-------------|-------|-------|-----------|
| | Duplicate | | | | | | | |
| Compound | Sample | Sample Result | | Sample Result A | | % RPD | RPD | Data |
| | μg/m³ | ppbV | μg/m³ | ppbV | $\mu g/m^3$ | | Limit | Qualifier |
| Ethylbenzene | ND | ND | ND | ND | - | - | 25 | |
| m,p-Xylenes | ND | ND | ND | ND | - | - | 25 | |
| Bromoform | ND | ND | ND | ND | - | - | 25 | |
| Styrene | ND | ND | ND | ND | - | - | 25 | |
| o-Xylene | ND | ND | ND | ND | - | 244 | 25 | |
| 1,1,2,2-Tetrachloroethane | ND | ND | ND | ND | - | - | 25 | |
| Cumene | ND | ND | ND | ND | - | - | 25 | |
| n-Propylbenzene | ND | ND | ND | ND | - | - | 25 | |
| 4-Ethyltoluene | ND | ND | ND | ND | - | - | 25 | |
| 1,3,5-Trimethylbenzene | ND | ND | ND | ND | - | - | 25 | |
| alpha-Methylstyrene | ND | ND | ND | ND | - | - | 25 | |
| 1,2,4-Trimethylbenzene | ND | ND | ND | ND | - | - | 25 | |
| Benzyl Chloride | ND | ND | ND | ND | - | - | 25 | |
| 1,3-Dichlorobenzene | ND | ND | ND | ND | - | - | 25 | |
| 1,4-Dichlorobenzene | ND | ND | ND | ND | - | - | 25 | |
| sec-Butylbenzene | ND | ND | ND | ND | - | - | 25 | |
| 4-Isopropyltoluene (p-Cymene) | ND | ND | ND | ND | - | - | 25 | |
| 1,2-Dichlorobenzene | ND | ND | ND | ND | - | - | 25 | |
| 1,2-Dibromo-3-chloropropane | ND | ND | ND | ND | - | - | 25 | |
| 1,2,4-Trichlorobenzene | ND | ND | ND | ND | - | - | 25 | |
| Naphthalene | ND | ND | ND | ND | - | - | 25 | |
| Hexachlorobutadiene | ND | ND | ND | ND | _ | - | 25 | |
| tert-Butylbenzene | ND | ND | ND | ND | - | - | 25 | |
| n-Butylbenzene | ND | ND | ND | ND | - | - | 25 | |

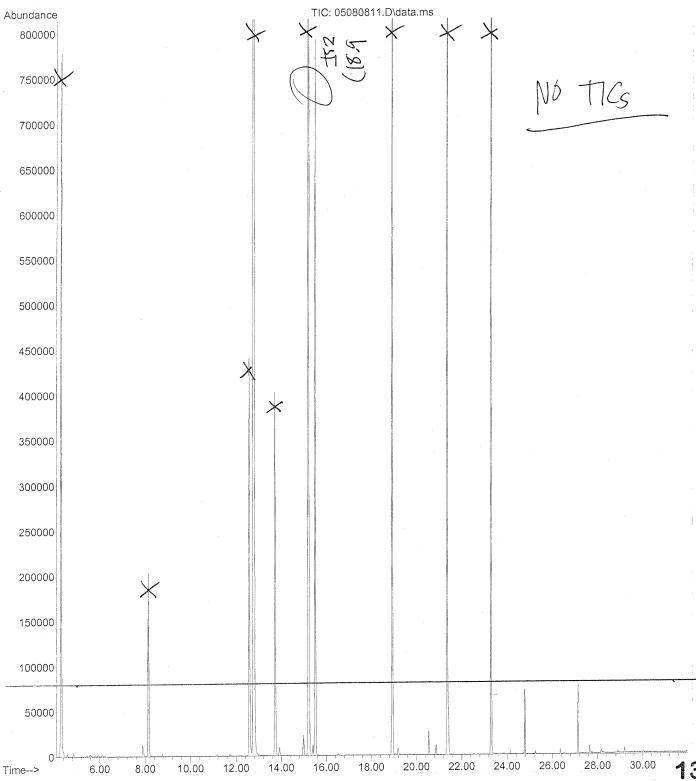
ND = Compound was analyzed for, but not detected above the laboratory detection limit.

Verified By:

Ke-



:J:\MS13\DATA\2008_05\08\05080811.D File : RTB Operator 8 May 2008 5:14 pm using AcqMethod TO15.M Acquired : GCMS13 Instrument : Sample Name: P0801342-003 DUP (10mL) Misc Info : ENSR SG83B-05-7 (-4.2, 3.5) Vial Number: 4



139

Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080811.D Acg On : 8 May 2008 5:14 pm Operator : RTB Sample : P0801342-003 DUP (10mL) Misc : ENSR SG83B-05-7 (-4.2, 3.5) ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 17:59:47 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards 1) Bromochloromethane (IS1)12.5813019955225.000 ng-0.0237) 1,4-Difluorobenzene (IS2)15.5111485324925.000 ng-0.02 56) Chlorobenzene-d5 (IS3) 21.35 82 416006 25.000 ng 0.00 System Monitoring Compounds 33) 1,2-Dichloroethane-d4(... 13.72 65 362203 22.633 ng -0.03 Recovery = 90.52% Spiked Amount 25.000 18.93 98 958033 25.693 ng 0.00 Recoverv = 102.76% / 57) Toluene-d8 (SS2) Recovery = 102.76% 🗸 Spiked Amount 25.000 73) Bromofluorobenzene (SS3) 23.29 174 317099 24.712 ng 0.00 Recovery = 98.84% Spiked Amount 25.000 Qvalue Target Compounds # 72 744 0.045 ng 4.83 42 2) Propene N.D. N.D. 3) Dichlorodifluoromethane 4.99 85 431 0 0.00 50 4) Chloromethane N.D.V 0 0 0 0.00 135 5) Freon 114 N.D. 0.00 62 6) Vinyl Chloride N.D. N.D. N.D. N.D. 0.00 54 0.00 94 7) 1,3-Butadiene 0 0 8) Bromomethane 0.00 64 9) Chloroethane 0.00 45 10) Ethanol 7.47 41 7.68 56 258 N.D. 11) Acetonitrile 57 12) Acrolein N.D. 7.89 58 6861 50 (0.621 ng) # 13) Acetone 8.14 101 100 209697 8.863 ng 14) Trichlorofluoromethane 209 0 N.D.V 0 N.D.V N.D. 8.37 45 15) Isopropanol 0.00 53 16) Acrylonitrile 0.00 96 17) 1,1-Dichloroethene 0.00 59 18) tert-Butanol 517 9.35 84 0.00 41 # 25 0.041 ng 19) Methylene Chloride 0 0 N.D.V 20) Allyl Chloride

 21) Trichlorottili

 22) Carbon Disulfide

 23) trans-1,2-Dichloroethene
 0.00
 61

 24) 1,1-Dichloroethane
 11.10
 63
 898

 25) Methyl tert-Butyl Ether
 0.00
 73
 0

 26) Vinyl Acetate
 0.00
 86
 0

 27) 2-Butanone
 11.72
 72
 840

 22) 78
 87
 826754

 N.D. N.D. N.D. 898 -0.040 ng # 57 N.D. 🗸 N.D. Q.108 ng # 87 N.D.V <u>-81.284 ng</u> M # 1 0.00 61 12.69 57 0 N.D. 30) Ethyl Acetate 212 N.D. 31) n-Hexane

R13041408.M Fri May 09 15:28:49 2008

505/09/08

140

| Data Path : J:\MS13\DATA\2008_05\08\ Data File : 05080811.D Acq On : 8 May 2008 5:14 pm Operator : RTB Sample : P0801342-003 DUP (10mL) Misc : ENSR SG83B-05-7 (-4.2, 3.5) ALS Vial : 4 Sample Multiplier: 1 | | | | | | | | | | |
|--|---|---------------|--------------------|----------------------|---------------------|---------------------------|--|--|--|--|
| Quant Time: May 08 17:59:47 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration | | | | | | | | | | |
| Inter | nal Standards | R.T. | QIon | Respons | e Conc Unit | s Dev(Min) | | | | |
| | | 10 70 | 00 | 7912800 | 121 702 00 | -Su dun 100 | | | | |
| | Chloroform | 12.78 0.00 | 83 72 | ⁷⁹¹²⁸⁰⁰ - | 424.702 lig N.D. | D ² 100 | | | | |
| | Tetrahydrofuran Ethyl tert-Butyl Ether | 0.00 | | 0 | N.D. | | | | | |
| | 1,2-Dichloroethane | 13.73 | | 575 | N.D. | | | | | |
| | 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D.V | | | | | |
| | Isopropyl Acetate | 0.00 | 61 | 0 | N.D. | | | | | |
| | 1-Butanol | 15.02 | 56 | 59 | N.D. | | | | | |
| | Benzene | 14.99 | \geq 78 | 26794 | 0.592 ng | 100 | | | | |
| | Carbon Tetrachloride (| 15.21 | 717 | 1063253 | 71.023 ng | > 99 | | | | |
| | Cyclohexane | 15.41 | 84 | 5267 | 0.315 ng | # 80 | | | | |
| | tert-Amyl Methyl Ether | 0.00 | 73 | 0 | N.D.V | | | | | |
| 45) | 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. 🗸 | | | | | |
| 46) | Bromodichloromethane | 16.45 | 83 | 85 | N.D.K | | | | | |
| 47) | Trichloroethene (| 16.53 | >130 | 764 | <u>(0.069 ng</u> | # 78 | | | | |
| 48) | 1,4-Dioxane | 0.00 | 88 | 0 | N.D. | | | | | |
| | Isooctane | 0.00 | 57 | 0 | N.D. | | | | | |
| | Methyl Methacrylate | 0.00 | 100 | 0 | N.D. | | | | | |
| | n-Heptane | 16.99 | 71 | 72 | N.D. | | | | | |
| | cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. 🗸 | | | | | |
| 53) | 4-Methyl-2-pentanone | 0.00 | 58 | 0 | N.D. | | | | | |
| 54) | trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | M∓ 9 | | | | |
| | 1,1,2-Trichloroethane | 18.94 | 97 | 87069 | <u></u> | 7 4 # 9 | | | | |
| | Toluene | 19.07 | 91 | 1731 | N.D.V | | | | | |
| | 2-Hexanone | 19.18 | 43 | 94 | N.D. | | | | | |
| | Dibromochloromethane | 0.00 | 129 107 | 0 0 | N.D. 🗸 N.D. 🖌 | | | | | |
| | 1,2-Dibromoethane | 0.00 20.36 | 43 | 231 | N.D. | | | | | |
| | Butyl Acetate | 0.00 | | 0 | N.D. | | | | | |
| | n-Octane Tetrachloroethene | 20.54 | \mathbf{R}_{66} | 8417 | 0.718 ng | > 99 | | | | |
| | Chlorobenzene | 21.40 | \sum_{112}^{100} | 32954 | 1.136 ng | > 98 | | | | |
| | Ethylbenzene | 21.89 | 91 | 376 | N.D. | | | | | |
| | m- & p-Xylene | 22.12 | 91 | 797 | N.D. | | | | | |
| | Bromoform | 0.00 | 173 | 0 | N.D./ | | | | | |
| | Styrene | 0.00 | 104 | 0 | N.D. 🗸 | | | | | |
| | o-Xylene | 22.73 | 91 | 164 | N.D. 🗸 | | | | | |
| | n-Nonane | 22.99 | 43 | 222 | N.D. | | | | | |
| | 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | | | | |
| | Cumene | 23.47 | 105 | 115 | N.D. | | | | | |
| | alpha-Pinene | 0.00 | 93 | 0 | N.D. | | | | | |
| | n-Propylbenzene | 0.00 | 91 | 0 | N.D. 🗸 | | | | | |
| | 3-Ethyltoluene | 24.24 | 105 | 239 | N.D. | | | | | |
| | 4-Ethyltoluene | 24.30 | 105 | 102 | N.D. | | | | | |
| 79) | 1,3,5-Trimethylbenzene | 24.38 | 105 | 127 | N.D.V | 141 | | | | |
| | | | | | Kentralie | | | | | |
| 1304140 | 8.M Fri May 09 15:28:49 20 | 08 | | | 100/00 | Page: 2 | | | | |
| | | | | | | | | | | |
| | | | | | 1 | | | | | |

Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080811.D Acq On : 8 May 2008 5:14 pm Operator : RTB Sample : P0801342-003 DUP (10mL) Misc : ENSR SG83B-05-7 (-4.2, 3.5) ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 17:59:47 2008 Ouant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards

 80) alpha-Methylstyrene
 24.79
 118
 925
 -0.041 ng
 #
 4

 81) 2-Ethyltoluene
 24.79
 105
 96
 N.D.

 82) 1,2,4-Trimethylbenzene
 24.88
 105
 432
 N.D.

 83) n-Decane
 24.99
 57
 165
 N.D.

 84) Benzyl Chloride
 0.00
 91
 0
 N.D.

 85) 1,3-Dichlorobenzene
 25.16
 146
 572
 N.D.

 86) 1,4-Dichlorobenzene
 25.41
 105
 114
 N.D.

 87) sec-Butylbenzene
 25.41
 105
 114
 N.D.

 88) p-Isopropyltoluene
 25.40
 119
 223
 N.D.

 90
 1,2-Dichlorobenzene
 25.16
 146
 572
 N.D.

 91) d-Limonene
 0.00
 68
 0
 N.D.

 92) 1,2-Dibromo-3-Chloropr...
 0.00
 157
 0
 N.D.

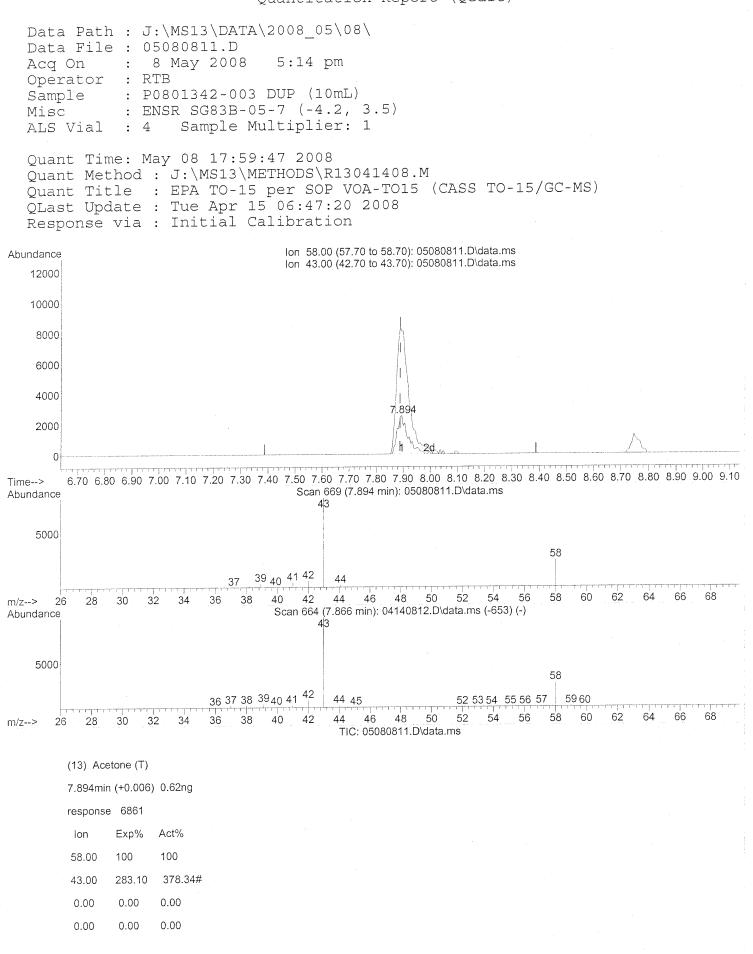
 93) n-Undecane
 26.50
 57
 602
 N.D.

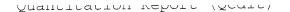
 94) 1,2,4-Trichlorobenzene
 26.50
 57
 602
 N.D.

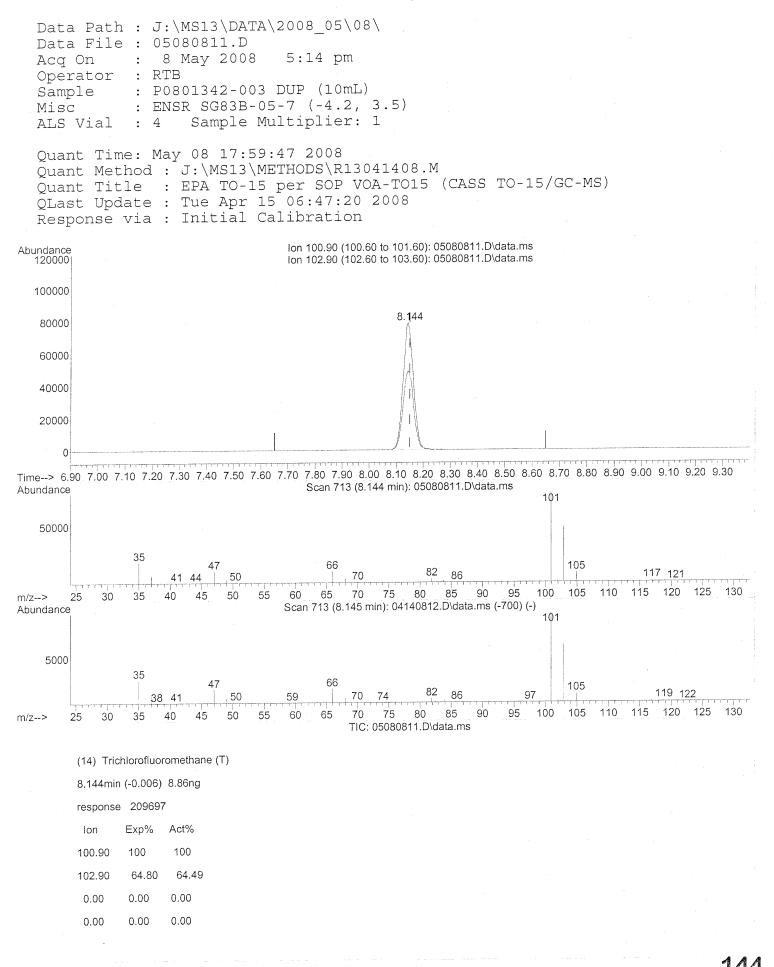
 95) Naphthalene
 27.73
 57
 703

(#) = qualifier out of range (m) = manual integration (+) = signals summed

405/20/108

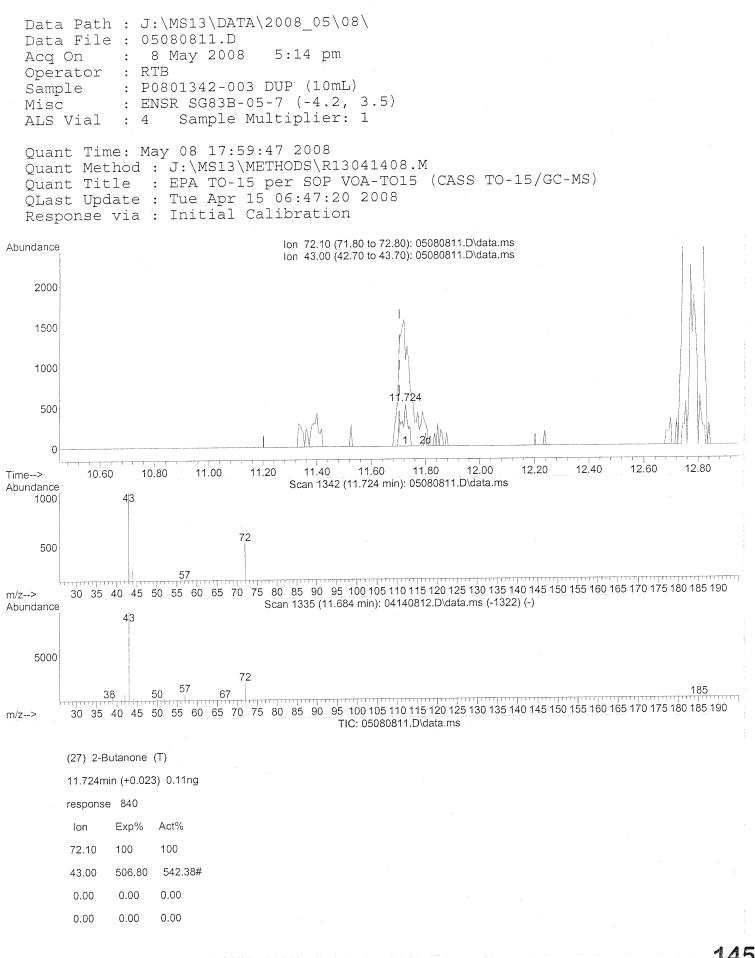


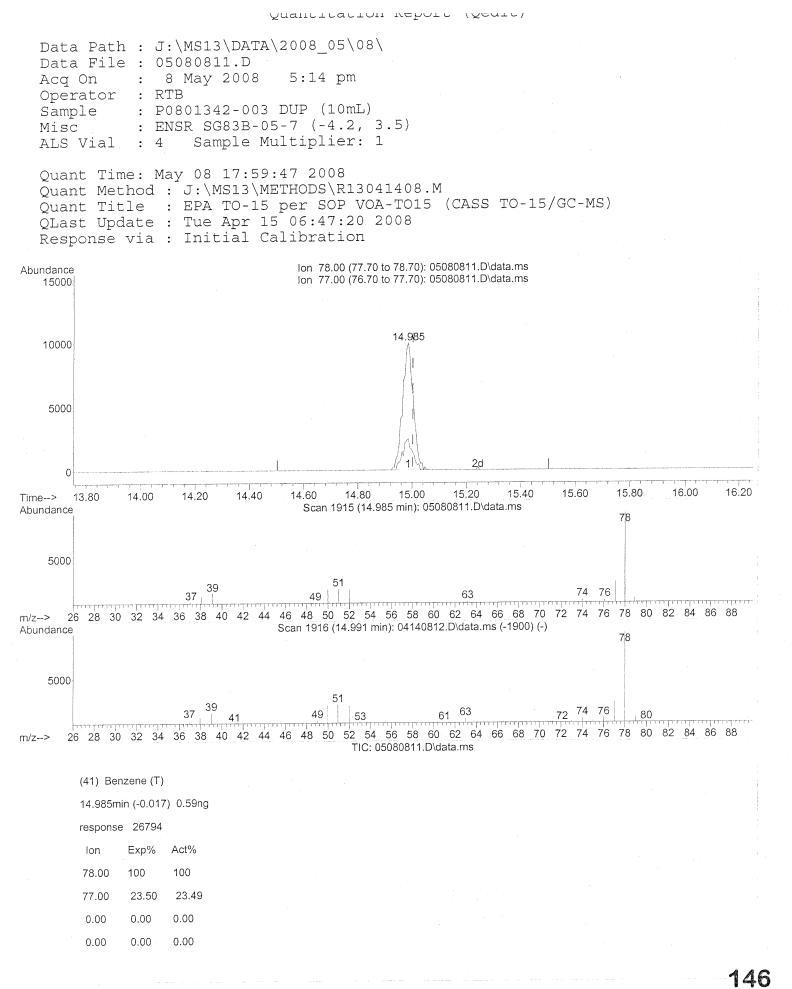




Page:

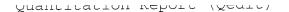
R13041408.M Thu May 08 18:22:43 2008

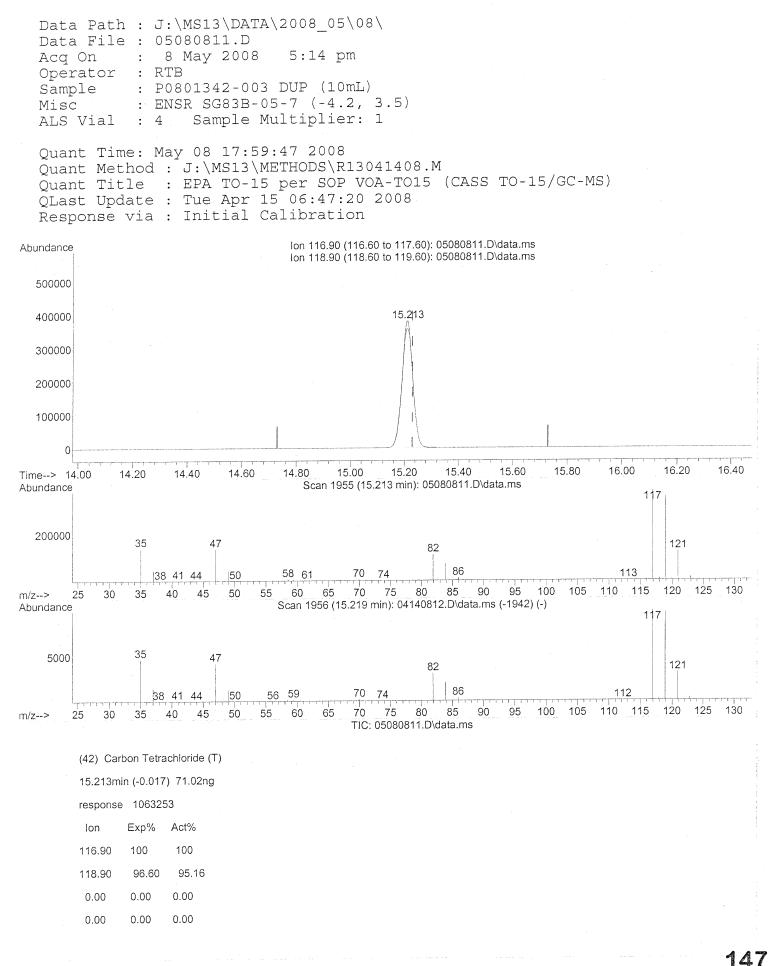




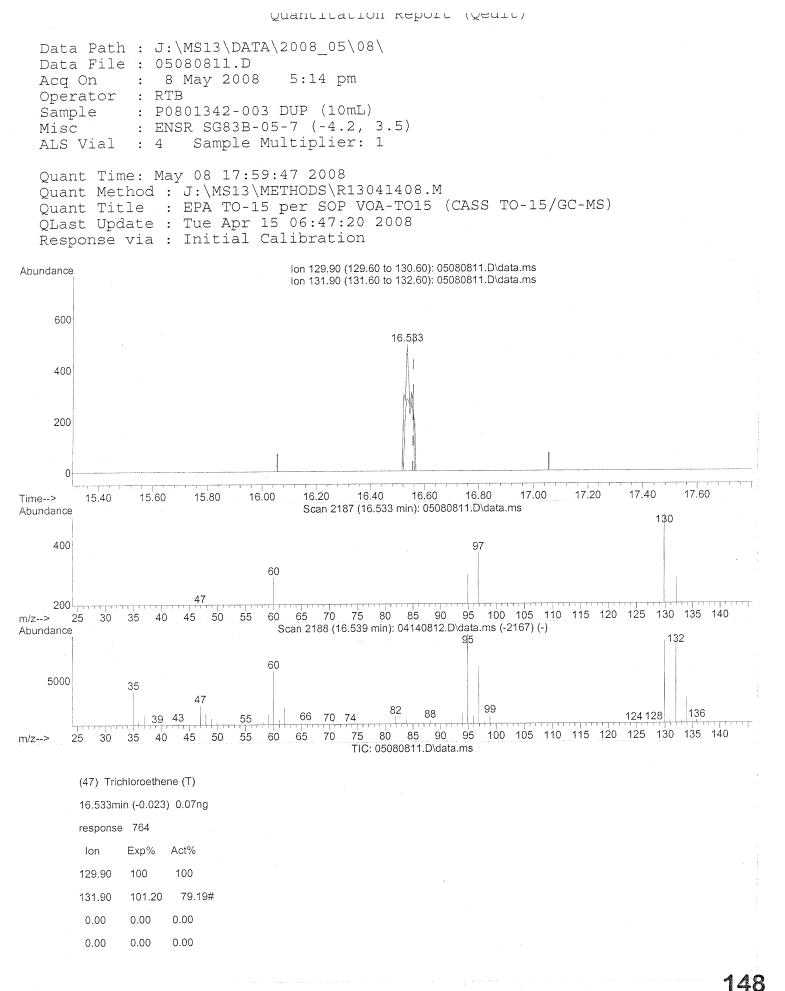
R13041408.M Thu May 08 18:23:39 2008

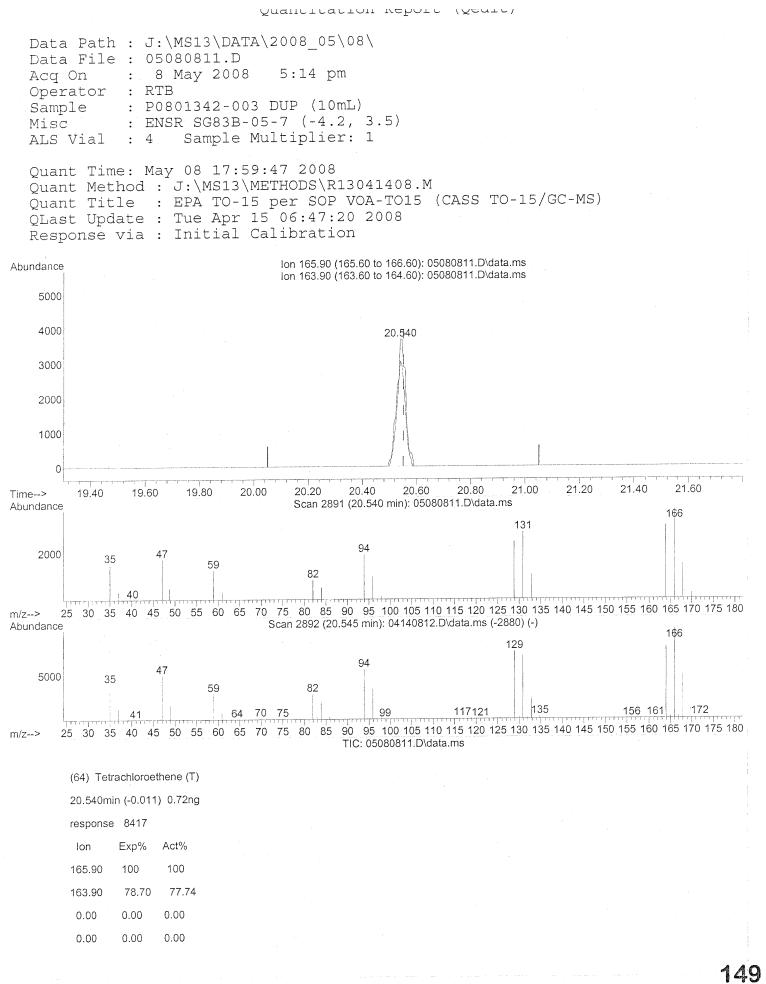
Page: 1

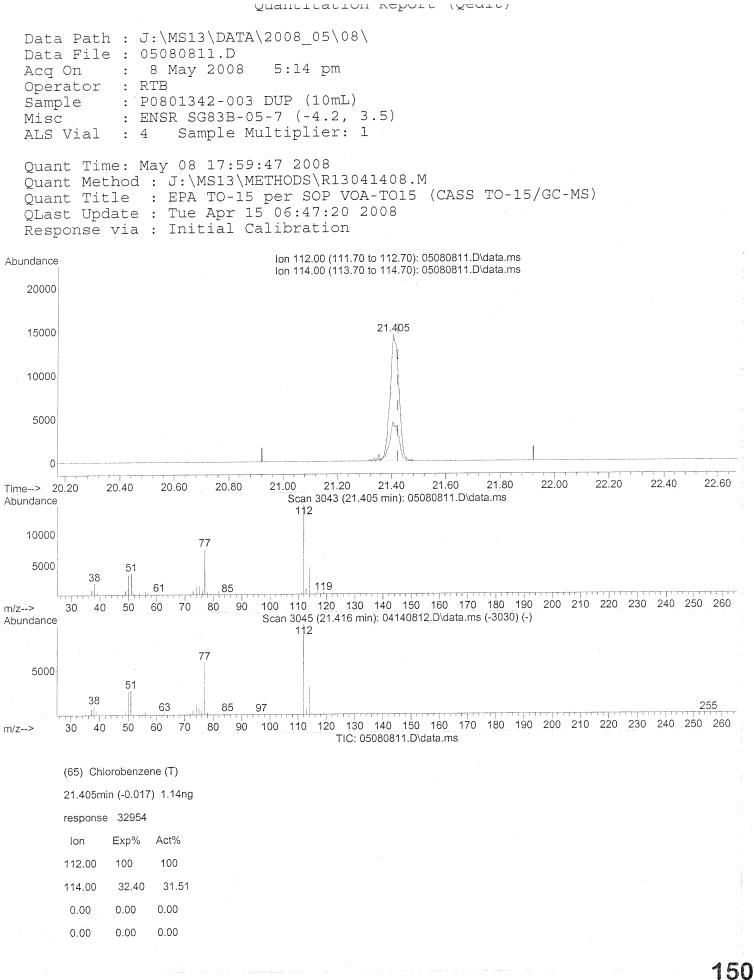




R13041408.M Thu May 08 18:23:44 2008







(NOL REVIEWEU)

Data Path : J:\MS13\DATA\2008_05\08\ Data File : 05080811.D Acq On : 8 May 2008 5:14 pm Operator : RTB Sample : P0801342-003 DUP (10mL) Misc : ENSR SG83B-05-7 (-4.2, 3.5) ALS Vial : 4 Sample Multiplier: 1

Quant Time: May 08 17:59:59 2008 Quant Method : J:\MS13\METHODS\S13041408.M Quant Title : TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD QLast Update : Mon Apr 28 10:06:00 2008 Response via : Initial Calibration

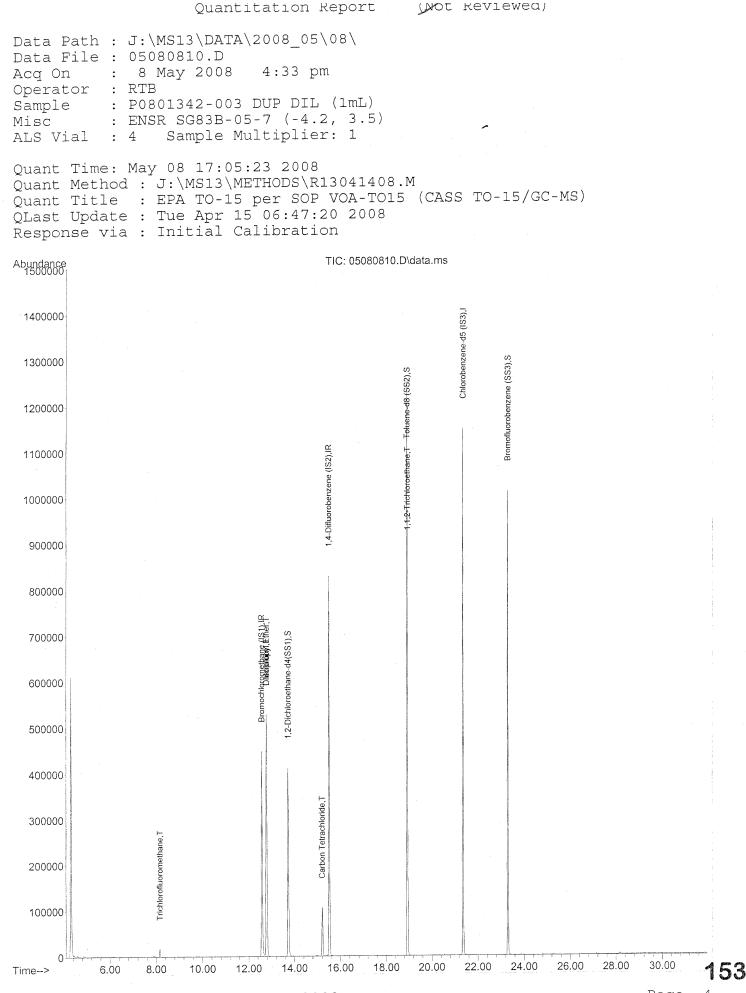
| Abundance TIC: 05080811.D\data.ms | |
|--|---|
| 1.05e+07 | |
| 1e+07 | |
| 950000 | |
| 900000 | |
| 850000 | |
| 800000 | |
| 7500000 | |
| 700000 | |
| 6500000 | |
| 600000 | |
| 5500000 | |
| 500000 | |
| 4500000 | |
| 400000 | |
| 3500000 | |
| 3000000 | |
| 0000001 0000002 0000002 0000002 1 Toluene-d8 (SS2),S 0000005 0000005 Chlorobenzene (IS2),IR Chlorobenzene (IS2),IR 0000005 Bromofluorobenzene (SS3),S SS2),S 0000005 | |
| 500000 0000000 0000000 0000000 0000000 0000000 0000000 0000000 | |
| 000000 000000 1,2-Dichloroethane (1S1),IR 0000005 1,4-Difluorobenzene (1S2),IR 1,4-Difluorobenzene (1S2),IR 1,4-Difluorobenzene (1S2),IR 0000005 Bromofluorobenzene (1S2),IR (132,1),IR | |
| | |
| 500000 | 1 |
| | <u>, , , , , , , , , , , , , , , , , , , </u> |
| Time> 6.00 8.00 10.00 12.00 14.00 16.00 18.00 20.00 22.00 24.00 26.00 | 28.00 30.00 151 Page: 2 |

S13041408.M Thu May 08 18:06:17 2008

QUAIICICACION REPORC (11) C 1/C V L C VV C C/ Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080811.D Acq On : 8 May 2008 5:14 pm Operator : RTB Sample : P0801342-003 DUP (10mL) Misc : ENSR SG83B-05-7 (-4.2, 3.5) ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 17:59:59 2008 Quant Method : J:\MS13\METHODS\S13041408.M Quant Title : TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD QLast Update : Mon Apr 28 10:06:00 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards 1) Bromochloromethane (IS1) 12.58 130 199552 25.000 ng -0.02 3)1,4-Difluorobenzene (IS2)15.5111485324925.000 ng-0.024)Chlorobenzene-d5 (IS3)21.358241600625.000 ng0.00 System Monitoring Compounds 2) 1,2-Dichloroethane-d4(... 13.72 65 362203 22.633 ng -0.03 Recovery = 90.52% 🗸 Spiked Amount 25.000 5) Toluene-d8 (SS2)18.939895803325.693 ng0.00Spiked Amount25.000Recovery = 102.76% / 6) Bromofluorobenzene (SS3) 23.29 174 317099 24.712 ng 0.00 Recovery = 98.84% / Spiked Amount 25.000 Qvalue Target Compounds 24.79 119 1519 N.D. 🗸 7) tert-Butylbenzene 0.00 91 0 N.D. 🗸 8) n-Butylbenzene -

(#) = qualifier out of range (m) = manual integration (+) = signals summed

65/08/08



R13041408.M Thu May 08 17:11:43 2008

| Quantitation | Report | (Ne | St Reviewed | (b | |
|--|--|------------------|----------------------------|--|------------------------------------|
| Data Path : J:\MS13\DATA\2008_05 Data File : 05080810.D Acq On : 8 May 2008 4:33 p Operator : RTB Sample : P0801342-003 DUP DIL Misc : ENSR SG83B-05-7 (-4. ALS Vial : 4 Sample Multiplie | om 」 (1mL) 2, 3.5) | | | | |
| Quant Time: May 08 17:05:23 2008 Quant Method : J:\MS13\METHODS\R Quant Title : EPA TO-15 per SOF QLast Update : Tue Apr 15 06:47: Response via : Initial Calibrati | 21304140 P VOA-TO 20 2008 | 15 (CA | ASS TO-15/0 | GC-MS) | |
| Internal Standards | R.T. | QIon | Response | Conc Units | s Dev(Min) |
| 1) Bromochloromethane (IS1) 37) 1,4-Difluorobenzene (IS2) 56) Chlorobenzene-d5 (IS3) | 12.58 15.51 21.35 | 130 114 82 | 202052 880501 420894 | 25.000 ng 25.000 ng 25.000 ng | -0.03 -0.02 0.00 |
| System Monitoring Compounds 33) 1,2-Dichloroethane-d4(Spiked Amount 25.000 57) Toluene-d8 (SS2) | 18.92 | 98 | Recove 979450 | ery = 92 25.962 ng | -0.01 |
| Spiked Amount 25.000 73) Bromofluorobenzene (SS3) Spiked Amount 25.000 | 23.29 | 174 | Recove 322308 Recove | ery = 103 24.826 ng ery = 99 | 3.84% / 0.00 9.32% / |
| <pre>Target Compounds 2) Propene 3) Dichlorodifluoromethane 4) Chloromethane 5) Freon 114 6) Vinyl Chloride 7) 1,3-Butadiene 8) Bromomethane 9) Chloroethane 10) Ethanol 11) Acetonitrile 12) Acrolein 13) Acetone 14) Trichlorofluoromethane 15) Isopropanol 16) Acrylonitrile 17) 1,1-Dichloroethene 18) tert-Butanol 19) Methylene Chloride 20) Allyl Chloride 21) Trichlorotrifluoroethane 22) Carbon Disulfide 23) trans-1,2-Dichloroethene 24) 1,1-Dichloroethane 25) Methyl tert-Butyl Ether 26) Vinyl Acetate 27) 2-Butanone 28) cis-1,2-Dichloroethene 29) Diisopropyl Ether 30) Ethyl Acetate 31) n-Hexane</pre> | $\begin{array}{c} 4.83\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 7.47\\ 0.00\\ 7.47\\ 0.00\\ 7.47\\ 0.00\\ 7.89\\ 8.16\\ 0.00\\ 0.00\\ 9.39\\ 9.36\\ 0.00\\ 0.00\\ 9.39\\ 9.36\\ 0.00\\ 0.00\\ 9.76\\ 0.00\\$ | 50 135 | 0 0 0 | N.D. N.D. N.D. N.D. N.D. N.D. N.D. N.D. | Qvalue 99 # 1 15 4 |
| 13041408.M Thu May 08 17:11:42 20 | 08 | | | 05/18/18 | Page: 1 |

4

| Quantitation | Report | (Ne | t Reviewed | .) | | |
|---|---|---|---|---|----------|----|
| Data Path : J:\MS13\DATA\2008_0 Data File : 05080810.D Acq On : 8 May 2008 4:33 Operator : RTB Sample : P0801342-003 DUP DI Misc : ENSR SG83B-05-7 (-4 ALS Vial : 4 Sample Multipli | pm L (1mL) .2, 3.5) | | | | | |
| Quant Time: May 08 17:05:23 200 Quant Method : J:\MS13\METHODS\ Quant Title : EPA TO-15 per SO QLast Update : Tue Apr 15 06:47 Response via : Initial Calibrat | R1304140 P VOA-TO :20 2008 ion | 15 (CA | | | | |
| Internal Standards | R.T. | QIon | Response | Conc Units | Dev(Min) | |
| 32) Chloroform 34) Tetrahydrofuran 35) Ethyl tert-Butyl Ether 36) 1,2-Dichloroethane 38) 1,1,1-Trichloroethane 39) Isopropyl Acetate 40) 1-Butanol 41) Benzene 42) Carbon Tetrachloride | 12.78 0.00 0.00 13.72 0.00 0.00 0.00 14.99 15.20 | 83 72 87 62 97 61 56 78 117 | 498718 0 174 0 0 2677 77127 | 26.436 ng N.D. N.D. N.D. N.D. N.D. N.D. N.D. 4.992 ng | 98 | |
| <pre>43) Cyclohexane 44) tert-Amyl Methyl Ether 45) 1,2-Dichloropropane 46) Bromodichloromethane 47) Trichloroethene 48) 1,4-Dioxane 49) Isooctane 50) Methyl Methacrylate 51) n-Heptane 52) cis-1,3-Dichloropropene 53) 4-Methyl-2-pentanone 54) trans-1,3-Dichloropropene</pre> | 15.41 0.00 0.00 0.00 0.00 0.00 0.00 0.00 0 | 84 73 63 130 88 57 100 71 75 58 75 | 56 0 0 0 0 0 0 0 0 0 0 0 0 | N.D. N.D. N.D. N.D. N.D. N.D. N.D. N.D. | | |
| <pre>54) trans-1,3-Dichloropropene 55) 1,1,2-Trichloroethane 58) Toluene 59) 2-Hexanone 60) Dibromochloromethane 61) 1,2-Dibromoethane 62) Butyl Acetate 63) n-Octane 64) Tetrachloroethene 65) Chlorobenzene 66) Ethylbenzene 67) m- & p-Xylene 68) Bromoform 69) Styrene 70) o-Xylene 71) n-Nonane 72) 1,1,2,2-Tetrachloroethane 74) Cumene 75) alpha-Pinene 76) n-Propylbenzene 77) 3-Ethyltoluene 78) 4-Ethyltoluene 79) 1,3,5-Trimethylbenzene</pre> | 18.94 19.08 0.00 0.00 0.00 20.55 21.42 21.89 22.11 0.00 22.73 23.29 0.00 23.48 0.00 24.24 24.24 | 97 91 43 129 107 43 57 166 112 91 173 104 91 43 83 105 93 105 105 | 89424 210 0 0 0 0 485 2552 141 225 0 0 83 215 0 55 0 0 208 208 208 208 | 7.950 ng N.D. N.D. N.D. N.D. N.D. N.D. N.D. N.D | # 7 | 55 |
| 13041408.M Thu May 08 17:11:42 2 | 008 | | Ŧ | 5/15/18 E | Page: 2 | |

(INOL KEVIEWEU) Quantitation keport Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080810.D Acq On : 8 May 2008 4:33 pm Operator : RTB Sample : P0801342-003 DUP DIL (1mL) Misc : ENSR SG83B-05-7 (-4.2, 3.5) ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 17:05:23 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards

 80) alpha-Methylstyrene
 0.00
 118
 0
 N.D.

 81) 2-Ethyltoluene
 24.24
 105
 208
 N.D.

 82) 1,2,4-Trimethylbenzene
 0.00
 105
 0
 N.D.

 83) n-Decane
 25.25
 57
 1615
 N.D.

 84) Benzyl Chloride
 0.00
 146
 0
 N.D.

 85) 1,3-Dichlorobenzene
 0.00
 146
 0
 N.D.

 86) 1,4-Dichlorobenzene
 0.00
 146
 0
 N.D.

 87) sec-Butylbenzene
 25.41
 105
 64
 N.D.

 88) p-Isopropyltoluene
 0.00
 146
 0
 N.D.

 90) 1,2-Dichlorobenzene
 0.00
 146
 0
 N.D.

 91) d-Limonene
 0.00
 68
 0
 N.D.

 92) 1,2-Dibromo-3-Chloropr...
 0.00
 157
 0
 N.D.

 93) n-Undecane
 26.52
 57
 231
 N.D.

 94) 1,2,4-Trichlorobenzene
 0.00
 180
 0
 N.D.

 95) Naphthalene
 27.79
 128
 233
 N.D.

 96) n-Dodecane
 0

(#) = qualifier out of range (m) = manual integration (+) = signals summed

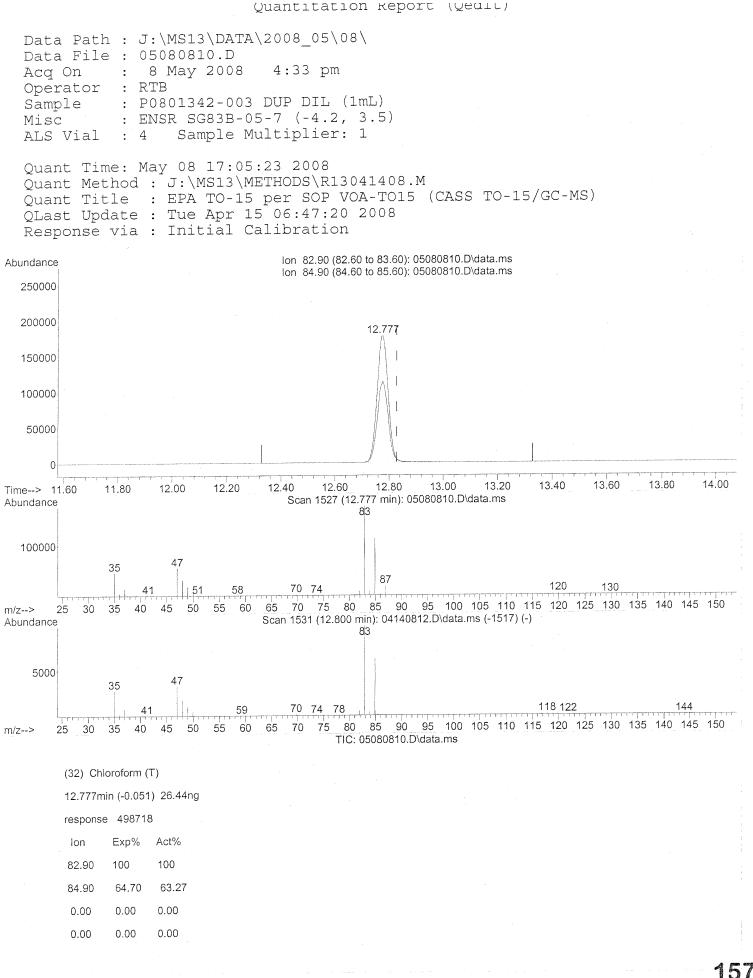
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Prost. 8/08

156 Page: 3

R13041408.M Thu May 08 17:11:42 2008



COLUMBIA ANALYTICAL SERVICES, INC.

RESULTS OF ANALYSIS

Page 1 of 1

Client:ENSRClient Project ID:Phase B Soil Gas / 04020-023-4311

CAS Project ID: P0801342

Internal Standard Area and RT Summary

| Test Code: | EPA TO-15 | |
|-----------------|--|-------------------------|
| Instrument ID: | Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13 | Lab File ID: 05080801.D |
| Analyst: | Rusty Bravo | Date Analyzed: 5/8/08 |
| Sampling Media: | 6.0 L Summa Canister(s) | Time Analyzed: 08:41 |
| Test Notes: | | |

| | IS1 (BCM) | | IS2 (DFB) | , | IS3 (CBZ) | |
|------------------|-----------|-------------|-----------|-------|-----------|-------|
| | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| 24 Hour Standard | 242477 | 12.59 | 1048411 | 15.52 | 500730 | 21.35 |
| Upper Limit | 339468 | 12.92 | 1467775 | 15.85 | 701022 | 21.68 |
| Lower Limit | 145486 | 12.26 | 629047 | 15.19 | 300438 | 21.02 |

| | Client Sample ID | | | | | | |
|----|---------------------------------------|--------|-------|---------|-------|--------|-------|
| 01 | Method Blank | 241346 | 12.58 | 1037647 | 15.51 | 489334 | 21.35 |
| 02 | Lab Control Sample | 244114 | 12.59 | 1061640 | 15.52 | 502772 | 21.35 |
| 03 | SG83B-05-1 | 242853 | 12.58 | 1040889 | 15.51 | 494646 | 21.35 |
| 04 | SG83B-05-3 | 231941 | 12.58 | 1004035 | 15.51 | 467452 | 21.35 |
| 05 | SG83B-05-7 | 226199 | 12.58 | 971717 | 15.51 | 470342 | 21.35 |
| 06 | SG83B-05-1 (Dilution) | 219916 | 12.58 | 959711 | 15.51 | 446346 | 21.35 |
| 07 | SG83B-05-3 (Dilution) | 213718 | 12.58 | 929659 | 15.51 | 442512 | 21.35 |
| 08 | SG83B-05-7 (Dilution) | 205516 | 12.58 | 899772 | 15.51 | 428319 | 21.35 |
| 09 | SG83B-05-7 (Lab Duplicate - Dilution) | 202052 | 12.58 | 880501 | 15.51 | 420894 | 21.35 |
| 10 | SG83B-05-7 (Lab Duplicate) | 199552 | 12.58 | 853249 | 15.51 | 416006 | 21.35 |
| 11 | | | | | | | |

- 12 13 14 15 16 17 18
- 19
- 20

IS1 (BCM) = Bromochloromethane

IS2 (DFB) = 1,4-Difluorobenzene

IS3 (CBZ) = Chlorobenzene-d5

AREA UPPER LIMIT = 140% of internal standard area AREA LOWER LIMIT = 60% of internal standard area RT UPPER LIMIT = 0.33 minutes of internal standard RT RT LOWER LIMIT = 0.33 minutes of internal standard RT

Column used to flag values outside QC limits with an asterisk.

* Values outside of QC limits.

Verified By:____

RG



INITIAL CALIBRATION STANDARDS

Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) Last Update : Tue Apr 15 06:34:47 2008 Response Via : Initial Calibration : J:\MS13\METHODS\ Method File : R13041408.M Method Path

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Calibration Files

5.0 = 04140811.D=04140810.D =04140814.D 1.0 0.5 =04140809.D 50 =04140813.D 0.1 =04140808.D 25 =04140812.D

| | . 1 | Compound | 0.1 | 0.5 | 1.0 | 5.0 | 25 | 20 | | Avg | %RSD |
|-------------|--|----------------------|---------|--|----------|-------|-------|---------|-------------|---------|----------|
| - | Ц Ц | Bromochloromethane | | | : | τ | | | [. | | 80 10 |
| 5) | í [H | | 2.452 | 2.204 | 2.130 | 1.859 | 1.876 | 1.926 | 2.038 | 2.069 | 0.2 |
| 3) | H | Dichlorodifluoromet | .69 | .06 | . 88 | .39 | .44 | 9 С. | .52 | 77. | 12.79 |
| 4) | Ē | шO | .15 | .52 | .36 | .80 | .94 | .87 | .41 | .15 | 8.2 |
| 2) | `. E1 | | .15 | .01 | .78 | .66 | .68 | .72 | .90 | .84 | 9.9 |
| (9) | [→. | 1 | .57 | .14 | 80 00 | .71 | .70 | .67 | . 83 | .93 | 1.0 |
| (_ | [-1 | 1,3-Butadiene | .42 | .39 | .10 | .04 | .21 | .27 | .42 | .26 | 6.8 |
| (8 8 | H | Bromomethane | .46 | .49 | .35 | .31 | .38 | .38 | .47 | .40 | 4.9 |
| 6 | Ē | Chloroethane | .25 | .38 | .16 | .12 | .16 | .13 | .28 | .21 | 7.9 |
| 10) | Εı | | | .81 | .59 | .36 | .22 | .22 | .24 | .41 | 7.1 |
| 11) | Ē | Acetonitrile | 4.507 | .28 | .79 | .57 | .26 | .23 | .29 | .70 | З.8 |
| 12) | [| Acrolein | .16 | .96 | .97 | .95 | .97 | .95 | .98 | .99 | 7.6 |
| | [1 | | | | .67 | .44 | .29 | .24 | .26 | .38 | 3.0 |
| | H | | .36 | .18 | .93 | .71 | .84 | .78 | .91 | .96 | 7.7 |
| | H | Isopropanol | .25 | .49 | .98 | .09 | .61 | .15 | .16 | .68 | 2.2 |
| | [1 | rylonitrile | .60 | .34 | .30 | .11 | .24 | .18 | .23 | .14 | 1.7 |
| | [| , 1- | .79 | .34 | . 35 | .24 | .30 | .28 | .33 | .38 | 3.5 |
| | E; | nol | .76 | .37 | . 19 | .89 | .05 | .93 | .98 | . 88 | 3.4 |
| 19) | [1 | thyl | 2.271 | 1.699 | • 53 | .39 | .42 | .37 | .42 | .59 | 0.1 |
| | | Allyl Chloride | с. 9 | . 08 | .00 | .09 | .38 | .39 | .50 | .12 | 7.6 |
| | E1 | ichlorotriflu | .52 | 30 30 30 30 30 30 30 30 30 30 30 30 30 | .22 | .17 | .17 | .15 | .27 | .26 | 0.3 |
| | H | Ч. | .26 | .21 | .85 | .49 | . 83 | .69 | .91 | .89 | 4.6 |
| | E1 | ans-1,2-Dichlc | .61 | . 53 | .43 | .17 | .32 | .31 | .44 | .40 | 6.0 |
| | H | chloroethane | .22 | .80 | .81 | .62 | .75 | .62 | .75 | .80 | 7.2 |
| | E-1 | l tert-Buty | .66 | .68 | .51 | .18 | .31 | .29 | .53 | .59 | 0.8 |
| | E-1 | | | | .21 | .25 | .29 | .29 | .31 | .27 | 5.0 |
| | Ē | -Butanone | 0.751 | .10 | .03 | .96 | .98 | .96 | .98 | .97 | 1.2 |
| | E | s-1,2-Dich | .56 | .29 | .25 | .13 | .19 | .17 | .24 | .26 | 6.3 |
| | E1 | iisopropyl Ethe | .36 | .32 | .10 | .09 | .18 | .28 | .56 | .27 | 2. 9 |
| | [| يستحم | | 0.517 | .51 | .53 | .60 | .65 | .80 | .60 | 8.7 |
| | E-1 | n-Hexane | 3.505 | .96 | .93 | .65 | .91 | .11 | .92 | .14 | 3.7 |
|) (1 | (| | | | | | | | | | |
| KL304 | 80 5 0 5 0 | .M lue Apr 15 06:46: | 57 200 | ω | | | | | RV. | 3/12/14 | C) |
| |) | | | | | | | | 1.1.1 | | 4 |

Page: 1

EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) Tue Apr 15 06:34:47 2008 Response Via : Initial Calibration : J:\MS13\METHODS' : R13041408.M ••• Path File Last Update Method Method Title

9.80 7.09 5.60 7.10 6.28 2.42 10.05 7.40 5.16 4.84 8.46 7.29 8.34 11.41 6.54 8.07 9.61 1.9¹ %RSD 0.326 0.451 .005 0.521 0.341 .334 1.676 .312 0.227 1.326 0.439 0.944 0.380 0.235 1.570 0.120 0.367 0.971 0.491 Avg \sim \sim \sim .479 1.947 .130 .718 .196 .354 .302 .495 .519 .007 **1.640** .389 0:586 0.535 0.246 .371 0.349 0.944 0.261 2.207 =04140811.D 100 Ч \sim 0 0 0 0 0 0 0 2.176 .949 0.474 .358 .447 0.237 .122 .460 0.310 1.533 .354 1.968 0.958 1.582 0.229 1.271 0.545 2.152 0.501 0.337 50 0 0 0 0 0 0 5.0 .256 0.457 0.470 0.949 0.360 0.235 2.230 2.032 0.977 1.594 0.233 0.335 1.282 0.447 0.314 1.541 0.122 0.354 0.514 Ŋ \sim N - ISTD .467 .147 1.980 0.939 1.207 0.410 0.290 1.506 2.140 0.3333 0.458 0.345 0.408 0.224 1.419 0.110 0.330 0.492 =04140810.D 0.208 0.872 =04140814.D 0 . വ N 0 .642 .366 0.913 0.357 1.233 0.409 0.350 2.045 0.485 0.415 0.332 0.245 1.566 0.116 0.470 2.305 0.914 0.484 0.218 0.341 1.0 \sim 1.0 100 1.790 .590 1.121 .445 1.391 0.419 0.426 0.356 2.037 0.548 0.243 0.348 0.504 0.926 0.398 0.228 1.647 0.122 0.369 0.508 <u>ں</u> 0 \sim \sim 0.941 2.025 1.898 2.604 0.597 1.598 0.420 066.0 0.471 1.645 .425 0.476 2.709 0.524 0.535 0.334 0.214 =04140809.D 0.216 =04140813.D 0.1 0 cis-1, 3-Dichloropro 1,2-Dichloropropane Bromodichloromethan 1,4-Difluorobenzene Carbon Tetrachlorid Methyl Methacrylate 1,1,1-Trichloroetha tert-Amyl Methyl Et Ethyl tert-Butyl Et 1,2-Dichloroethane 1,2-Dichloroethane Isopropyl Acetate Tetrahydrofuran Trichloroethene 0.5 50 1,4-Dioxane Cyclohexane Chloroform n-Heptane Calibration Files sooctane 1-Butano] Compound 0.1 =04140808.D 25 =04140812.D Benzene ЦR HOHHH 34) 35) 36) 32) 33) 37) 38) 39) 40) 41) 42) 43) 50) 51) 52) 44) 45) 46) 46) 48) 49)

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2.254 .923 0.760 713 0 \sim .696 .698 .665 .191 \sim \sim \circ 0 0.689 2.654 2.212 672 0 2.597 2.050 0.648 0.637 2.703 1.885 0.614 9 8 0.5 0.640 2.932 2.056 0.662 2.272 3.203 0.707 2.033 .686 0 Dibromochloromethan R13099408.M Tue Apr 15 06:46 1,2-Dibromoethane 2-Hexanone Toluene ноннн 58) 59) 60) 61)

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6.14 7.26 5.52

1.65 7.61

2.241

2.190

.195

 \sim

2.226

2.254

2.272

2.277

(IS3

Chlorobenzene-d5

56) 57)

(SS2)

Toluene-d8

- ISTD

2.815

2.097

0.679

0.662

5.99

0.363

.402 .526 .328

 \circ 0 0

.372

0 0

0.374

0.338 0.422 0.289

0.348 0.402

0.354 0.426 0.327

0.352 0.378

4-Methyl-2-pentanon trans-1,3-Dichlorop

53)

54) 55)

0.540

0.72

11.96

446319

. 0 0

.490

.481

0 0

.302

0

.301

99

0.2

0.390

1,1,2-Trichloroetha

8.17

517

0

Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) Last Update : Tue Apr 15 06:34:47 2008 Response Via : Initial Calibration Method Path : J:\MS13\METHODS\ Method File : R13041408.M

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Calibration Files

5.0 = 04140811.D1.0 = 04140810.D100 = 04140814.D0.5 =04140809.D 50 =04140813.D 0.1 =04140808.D 25 =04140812.D

| | | Compound | | د | 0. | 10 2 | ы | 0 | 00 | ۲g | QD |
|-----|----------------|-------------------|-------|-------|------|---------|-------|-------|-----------|-------|---------|
| 62) | Ē | Butyl Acetate | 1.716 | 2.099 | 973 | | 2.252 | 2.250 | 2.310 | 2.097 | 9.80 |
| 63) | E⊶ | n-Octane | .66 | .64 | .62 | . 6 | .65 | .67 | .75. | .66 | 4 |
| 64) | Ē | Tetrachloroethene | .84 | .67 | .65 | .62 | .65 | .68 | . 80 | .70 | 1.7 |
| 65) | H | Chlorobenzene | .94 | .81 | .64 | .59 | .64 | .68 | 80 80 | .74 | 7.8 |
| 66) | H | Ethylbenzene | .08 | .25 | .94 | .93 | .14 | .22 | .41 | .14 | 5.5 |
| 67) | Ē | m- & p-Xylene | .94 | .99 | .90 | .93 | .12 | .27 | . 53 | .10 | 6.0 |
| 68) | Ē | romo | .35 | .41 | .44 | .45 | .49 | .51 | .57 | .46 | 5.3 |
| 69) | Ē | Styrene | .87 | .68 | .62 | .66 | .83 | . 8 | .13 | . 81 | 9.7 |
| (02 | H | o-Xylene | .30 | .10 | .00 | .12 | .24 | .35 | .70 | .26 | 0.2 |
| 71) | H | -Nonane | .87 | .78 | .60 | .70 | .81 | .84 | .04 | .81 | 7.6 |
| 72) | Η | ,1,2,2-Tetrac | .12 | .97 | .95 | .01 | .07 | .12 | .29 | .07 | 0.8 |
| 73) | N N | romofluorobenz | .76 | .75 | .76 | .78 | .78 | .76 | .77 | .77 | 1.2 |
| 74) | [1 | umene | .60 | .81 | .67 | .69 | .91 | .04 | .33 | .86 | 8 .9 |
| 75) | E1 | | .40 | .40 | .37 | .41 | .54 | .62 | .91 | .52 | 2.7 |
| 76) | E1 | -Propylbenze | .64 | .83 | .64 | .62 | .96 | .06 | .09 | . 83 | 5.3 |
| (| Η | 3-Ethyltoluene | .95 | .98 | . 88 | .87 | .14 | .34 | .70 | .12 | 9.7 |
| 78) | Εı | -Ethyltoluene | .74 | .62 | .64 | .67 | .95 | .13 | .39 | . 88 | 0.1 |
| 79) | Η | ,3,5-Trimethylben | .35 | .37 | .31 | .32 | .57 | : 78 | .18 | . 55 | 2.7 |
| 80) | Ε | lpha-Methylstyren | .20 | . 10 | .13 | .24 | .40 | .49 | .84 | .35 | 8.6 |
| 81) | E | -Ethyltoluene | .08 | .00 | . 80 | .89 | .18 | .33 | .69 | .14 | 9.6 |
| 82) | Η | l rimeth | .29 | .44 | .30 | .43 | .99 | .67 | .07 | . 89 | 4.9 |
| 83) | [] | -De(| .40 | .46 | .46 | .46 | .61 | .72 | 0 0 | .59 | 4.3 |
| 84) | Ð | zyl Chloride | .26 | .54 | .54 | .90 | .23 | .42 | .87 | .97 | 9.0 |
| 85) | Ξ | ,3-Dichloroben | .46 | .48 | .37 | .36 | .52 | .64 | ο. Ο Ο | .56 | 6.0 |
| 86) | [1 | ,4-Dichlorobenzen | .45 | .38 | .36 | .33 | .46 | .56 | .92 | .49 | 3.6 |
| 87) | H | ec-Butylbenzen | . 28 | .28 | .10 | .13 | .45 | .65 | .83 | .39 | 7.9 |
| 88) | Ξ | -Isopropyltoluen | .43 | .42 | .39 | .50 | .19 | . 88 | .87 | .95 | 3.1 |
| (68 | Η | \geq | .22 | .37 | .29 | .34 | .98 | .64 | .03 | .84 | 5.7 |
| 606 | Ē | ,2-Dichlorobenzen | .19 | .34 | . 29 | .30 | .58 | .93 | .55 | .60 | 0.5 |
| 91) | [| imonene | .03 | .05 | .99 | .06 | ЭЭ | .57 | | . 31 | 1.3 |
| 92) | [1 | 2 | .23 | .38 | .42 | .41 | .46 | .48 | .55 | .42 | 4.1 |
| 63) | E- | n-Undecane | .49 | .50 | .43 | .55 | .69 | . 83 | .23 | .67 | 6.7 |
| | | | | | | | | | | | |

Page: 3

RA 4/15/08

R13(9,408.M Tue Apr 15 06:46:57 2008

: EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) te : Tue Apr 15 06:34:47 2008 Response Via : Initial Calibration Method Path : J:\MS13\METHODS\ Method File : R13041408.M Last Update Title

Ava 5.0 =04140811.D 100 С С 5 1.0 =04140810.D 100 =04140814.D С . س л. О.Т 0. 5 0.5 =04140809.D 50 =04140813.D 0.1 Calibration Files Compound 0.1 =04140808.D 25 =04140812.D

| | | [mon | 0.1 | 0.5 | 0.1 0.5 1.0 5.0 | 5.0 | 25 | 50 | 100 | Avg | %RSD |
|-----------------|----------------|---------------------|--------------------|-------|-------------------|-------------|-------|-------|---|-------------------------------------|--------------|
| : { | 1 | | | | 1 1 1 1 | | | | | | |
| 94) | [1 | 1,2,4-Trichlorobenz | 0.796 | 0.875 | 0.876 | 0.899 | 0.992 | 1.100 | 0.796 0.875 0.876 0.899 0.992 1.100 1.419 0.994 | 0.994 | 90 LC |
| 95) | Εı | Naphthalene | 2.304 | 3.074 | 2.304 3.074 2.991 | 3.156 | 3.394 | 3.589 | 3.861 | 3.156 3.394 3.589 3.861 3.196 | 10.9 . L |
| 96) | [| n-Dodecane | 1.646 | 1.521 | 1.646 1.521 1.441 | 1.519 1.671 | 1.671 | 1.830 | 1.830 2.290 1.703 | 1 703 | 90.01 1.0 |
| 97) | [1 | Hexachloro-1,3-buta | 3-buta 0.415 0.578 | 0.578 | 0.567 | 0.545 | 0.638 | 0.724 | 1.012 | 0.567 0.545 0.638 0.724 1.012 0.640 | 29.53 |
| 1 | | | | | | | | | | | |
| | Ċ | | | | | | | | | | |

(#) = Out of Range

R130**9**408.M Tue Apr 15 06:46:57 2008

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Page:

Primary Source Standards Concentrations (Working & Initial Calibration)

4ng/L Std. ID: <u>S20-04030801</u> 20ng/L Std. ID: <u>S20-03210809</u> 200ng/L Std. ID: <u>S20-04020808</u>

| 200ng/L Std. ID: s | 520-04020808 | | | | | | ICAL C | oncentra | tions (Pi | rim <mark>ary So</mark> | urce) | |
|--------------------------|-------------------|-----------|-----------|----------|---|--------------|--------|----------|-----------|-------------------------|-------|-------|
| Dilution Factors: | | 5 | 50 | 250 | Working STD Conc.(ng/L): | NOAKS STORES | 20 | 20 | 20 | 200 | 200 | 200 |
| | Source Std. | Primary V | Vorking S | tandards | Injection (L): | 0.025 | 0.025 | 0.050 | 0.25 | 0.125 | 0.25 | 0.50 |
| Compounds | mg/m ³ | 200ng/L | 20ng/L | 4ng/L | ICAL Points: | 0.1ng | 0.5ng | 1ng | 5ng | 25ng | 50ng | 100ng |
| Propene | 1.08 | 216 | 21.6 | 4.32 | | NA | 0.540 | 1.08 | 5.40 | 27.0 | 54.0 | 108 |
| Dichlorodifluoromethane | 1.04 | 208 | 20.8 | 4.16 | 2////////// | NA | 0.520 | 1.04 | 5.20 | 26.0 | 52.0 | 104 |
| Chloromethane | 1.02 | 204 | 20.4 | 4.08 | Z////////// | NA | 0.510 | 1.02 | 5.10 | 25.5 | 51.0 | 102 |
| Freon-114 | 1.07 | 214 | 21.4 | 4.28 | AIIIIIIIIIIIIIIII | NA | 0.535 | 1.07 | 5.35 | 26.8 | 53.5 | 107 |
| Vinyl Chloride | 1.03 | 206 | 20.6 | 4.12 | ZIIIIIIIIIIII | NA | 0.515 | 1.03 | 5.15 | 25.8 | 51.5 | 103 |
| 1,3-Butadiene | 1.09 | 218 | 21.8 | 4.36 | Z///////////////////////////////////// | NA | 0.545 | 1.09 | 5.45 | 27.3 | 54.5 | 109 |
| Bromomethane | 1.05 | 210 | 21.0 | 4.20 | VIIIIIIIV | NA | 0.525 | 1.05 | 5.25 | 26.3 | 52.5 | 105 |
| Chloroethane | 1.05 | 210 | 21.0 | 4.20 | AIIIIIIIIIII | NA | 0.525 | 1.05 | 5.25 | 26.3 | 52.5 | 105 |
| Ethanol | 0.91 | 182 | 18.2 | 3.64 | Z///////////////////////////////////// | NA | 0.455 | 0.910 | 4.55 | 22.8 | 45.5 | 91.0 |
| Acetonitrile | 0.980 | 196 | 19.6 | 3.92 | | NA | 0.490 | 0.980 | 4.90 | 24.5 | 49.0 | 98.0 |
| Acrolein | 0.960 | 192 | 19.2 | 3.84 | ZIIIIIIIIIIII | NA | 0.480 | 0.960 | 4.80 | 24.0 | 48.0 | 96.0 |
| Acetone | 1.11 | 222 | 22.2 | 4.44 | ZIIIIIIIIIIIIIIIII | NA | 0.555 | 1.11 | 5.55 | 27.8 | 55.5 | 111 |
| Trichlorofluoromethane | 1.04 | 208 | 20.8 | 4.16 | ZHHHHHH | NA | 0.520 | 1.04 | 5.20 | 26.0 | 52.0 | 104 |
| Isopropanol | 1.03 | 206 | 20.6 | 4.12 | | NA | 0.515 | 1.03 | 5.15 | 25.8 | 51.5 | 103 |
| Acrylonitrile | 1.010 | 202 | 20.2 | 4.04 | YIIIIIIIIII | NA | 0.505 | 1.01 | 5.05 | 25.3 | 50.5 | 101 |
| 1,1-Dichloroethene | 1.13 | 226 | 22.6 | 4.52 | VIIIIIIIIII | NA | 0.565 | 1.13 | 5.65 | 28.3 | 56.5 | 113 |
| tert-Butanol | 1.020 | 204 | 20.4 | 4.08 | VIIIIIIIV | NA | 0.510 | 1.02 | 5.10 | 25.5 | 51.0 | 102 |
| Methylene Chloride | 1.12 | 224 | 22.4 | 4.48 | AIIIIIIIIIA | NA | 0.560 | 1.12 | 5.60 | 28.0 | 56.0 | 112 |
| Allyl Chloride | 1.05 | 210 | 21.0 | 4.20 | | NA | 0.525 | 1.05 | 5.25 | 26.3 | 52.5 | 105 |
| Trichlorotrifluoroethane | 1.14 | 228 | 22.8 | 4.56 | (////////// | NA | 0.570 | 1.14 | 5.70 | 28.5 | 57.0 | 114 |
| Carbon Disulfide | 1.00 | 200 | 20.0 | 4.00 | $\Delta $ | NA | 0.500 | 1.00 | 5.00 | 25.0 | 50.0 | 100 |
| trans-1,2-Dichloroethene | 1.10 | 220 | 22.0 | 4.40 | | NA | 0.550 | 1.10 | 5.50 | 27.5 | 55.0 | 110 |
| 1,1-Dichloroethane | 1.11 | 222 | 22.2 | 4.44 | | NA | 0.555 | 1.11 | 5.55 | 27.8 | 55.5 | 111 |
| Methyl tert-Butyl Ether | 1.11 | 222 | 22.2 | 4.44 | | NA | 0.555 | 1.11 | 5.55 | 27.8 | 55.5 | 111 |
| Vinyl Acetate | 0.98 | 196 | 19.6 | 3.92 | | NA | 0.490 | 0.980 | 4.90 | 24.5 | 49.0 | 98.0 |
| 2-Butanone | 1.12 | 224 | 22.4 | 4.48 | <i>2///////////</i> | NA | 0.560 | 1.12 | 5.60 | 28.0 | 56.0 | 112 |
| cis-1,2-Dichloroethene | 1.11 | 222 | 22.2 | 4.44 | | NA | 0.555 | 1.11 | 5.55 | 27.8 | 55.5 | 111 |
| Diisopropyl Ether | 1.03 | 206 | 20.6 | 4.12 | <i>111111111111</i> | NA | 0.515 | 1.03 | 5.15 | 25.8 | 51.5 | 103 |
| Ethyl Acetate | 1.27 | 254 | 25.4 | 5.08 | | NA | 0.635 | 1.27 | 6.35 | 31.8 | 63.5 | 127 |
| n-Hexane | 1.12 | 224 | 22.4 | 4.48 | | NA | 0.560 | 1.12 | 5.60 | 28.0 | 56.0 | 112 |
| Chloroform | 1.29 | 258 | 25.8 | 5.16 | | NA | 0.645 | 1.29 | 6.45 | 32.3 | 64.5 | 129 |
| Tetrahydrofuran | 1.11 | 222 | 22.2 | 4.44 | | NA | 0.555 | 1.11 | 5.55 | 27.8 | 55.5 | 111 |
| Ethyl tert-Butyl Ether | 1.05 | 210 | 21.0 | 4.20 | X//////////A | NA | 0.525 | 1.05 | 5.25 | 26.3 | 52.5 | 105 |
| 1,2-Dichloroethane | 1.10 | 220 | 22.0 | 4.40 | | NA | 0.550 | 1.10 | 5.50 | 27.5 | 55.0 | 110 |
| 1,1,1-Trichloroethane | 1.10 | 220 | 22.0 | 4.40 | <i>SHHHHHH</i> | NA | 0.550 | 1.10 | 5.50 | 27.5 | 55.0 | 110 |
| Isopropyl Acetate | 1.010 | 202 | 20.2 | 4.04 | | NA | 0.505 | 1.01 | 5.05 | 25.3 | 50.5 | 101 |
| 1-Butanol | 0.910 | 182 | 18.2 | 3.64 | S////////// | NA | 0.455 | 0.910 | 4.55 | 22.8 | 45.5 | 91.0 |
| Benzene | 1.10 | 220 | 22.0 | 4.40 | | NA | 0.550 | 1.10 | 5.50 | 27.5 | 55.0 | 110 |
| Carbon Tetrachloride | 1.07 | 214 | 21.4 | 4.28 | | NA | 0.535 | 1.07 | 5.35 | 26.8 | 53.5 | 107 |
| Cyclohexane | 1.11 | 222 | 22.2 | 4.44 | AIIIIIIIIIIIA | NA | 0.555 | 1.11 | 5.55 | 27.8 | 55.5 | 111 |
| ert-Amyl Methyl Ether | 1.04 | 208 | 20.8 | 4.16 | | NA | 0.520 | 1.04 | 5.20 | 26.0 | 52.0 | 104 |
| 1,2-Dichloropropane | 1.09 | 218 | 21.8 | 4.36 | | NA | 0.545 | 1.09 | 5.45 | 27.3 | 54.5 | 109 |
| Bromodichloromethane | 1.15 | 230 | 23.0 | 4.60 | AIIIIIIIIIIA | NA | 0.575 | 1.15 | 5.75 | 28.8 | 57.5 | 115 |
| Trichloroethene | 1.14 | 228 | 22.8 | 4.56 | AIIIIIIIIIIII | NA | 0.570 | 1.14 | 5.70 | 28.5 | 57.0 | 114 |
| 1,4-Dioxane | 1.15 | 230 | 23.0 | 4.60 | | NA | 0.575 | 1.15 | 5.75 | 28.8 | 57.5 | 115 |
| sooctane | 1.04 | 208 | 20.8 | 4.16 | | NA | 0.520 | 1.04 | 5.20 | 26.0 | 52.0 | 104 |
| Methyl Methacrylate | 1.06 | 212 | 21.2 | 4.24 | | NA | 0.530 | 1.06 | 5.30 | 26.5 | 53.0 | 106 |
| n-Heptane | 1.11 | 222 | 22.2 | 4.44 | | NA | 0.555 | 1.11 | 5.55 | 27.8 | 55.5 | 111 |
| cis-1,3-Dichloropropene | 1.04 | 208 | 20.8 | 4.16 | | NA | 0.520 | 1.04 | 5.20 | 26.0 | 52.0 | 104 |
| 1-Methyl-2-pentanone | 1.05 | 210 | 21.0 | 4.20 | | NA | 0.525 | 1.05 | 5.25 | 26.3 | 52.5 | 105 |
| rans-1,3-Dichloropropene | 1.16 | 232 | 23.2 | 4.64 | 4////////// | NA | 0.580 | 1.16 | 5.80 | 29.0 | 58.0 | 116 |
| 1,1,2-Trichloroethane | 1.09 | 218 | 21.8 | 4.36 | | NA | 0.545 | 1.09 | 5.45 | 27.3 | 54.5 | 109 |
| Toluene | 1.10 | 220 | 22.0 | 4.40 | | NA | 0.550 | 1.10 | 5.50 | 27.5 | 55.0 | 110 |
| 2-Hexanone | 1.02 | 204 | 20.4 | 4.08 | 4////////// | NA | 0.510 | 1.02 | 5.10 | 25.5 | 51.0 | 102 |
| Dibromochloromethane | 1.11 | 222 | 22.2 | 4.44 | | NA | 0.555 | 1.11 | 5.55 | 27.8 | 55.5 | 111 |
| ,2-Dibromoethane | 1.09 | 218 | 21.8 | 4.36 | | NA | 0.545 | 1.09 | 5.45 | 27.3 | 54.5 | 109 |
| n-Butyl Acetate | 1.05 | 210 | 21.0 | 4.20 | | NA | 0.525 | 1.05 | 5.25 | 26.3 | 52.5 | 105 |
| n-Octane | 1.04 | 208 | 20.8 | 4.16 | | NA | 0.520 | 1.04 | 5.20 | 26.0 | 52.0 | 104 |
| etrachloroethene | 1.09 | 218 | 21.8 | 4.36 | | NA | 0.545 | 1.09 | 5.45 | 27.3 | 54.5 | 109 |
| Chlorobenzene | 1.10 | 220 | 22.0 | 4.40 | | NA | 0.550 | 1.10 | 5.50 | 27.5 | 55.0 | 110 |
| Ethylbenzene | 1.08 | 216 | 21.6 | 4.32 | XIIIIIIIIIIIIX | NA | 0.540 | 1.08 | 5.40 | 27.0 | 54.0 | 108 |
| n-&p-Xylene | 2.58 | 516 | 51.6 | 10.32 | | NA | 1.29 | 2.58 | 12.9 | 64.5 | 129 | 258 |

DA 4/15/08

Primary Source Standards Concentrations (Working & Initial Calibration)

4ng/L Std. ID: S20-04030801 20ng/L Std. ID: S20-03210809 200ng/L Std. ID: S20-04020808

| 200ng/L Std. ID: | S20-04020808 | | | | | | ICAL C | oncentra | tions (Pr | imary So | urce) | |
|-----------------------------|-------------------|-----------|---------------|----------|--|--------------|--------------|------------|------------|-------------|-------------|-------|
| | | | | | Working STD | | | | | | | |
| Dilution Factors: | : | 5 | 50 | 250 | Conc.(ng/L): | 0 | 20 | 20 | 20 | 200 | 200 | 200 |
| | Source Std. | Primary V | Vorking S | tandards | Injection (L): | 0.025 | 0.025 | 0.05 | 0.25 | 0.125 | 0.25 | 0.50 |
| Compounds | mg/m ³ | 200ng/L | <u>20ng/L</u> | 4ng/L | ICAL Points: | <u>0.1ng</u> | <u>0.5ng</u> | <u>1ng</u> | <u>5ng</u> | <u>25ng</u> | <u>50ng</u> | 100ng |
| Bromoform | 1.31 | 262 | 26.2 | 5.24 | | NA | 0.655 | 1.31 | 6.55 | 32.8 | 65.5 | 131 |
| Styrene | 1.08 | 216 | 21.6 | 4.32 | χ | NA | 0.540 | 1.08 | 5.40 | 27.0 | 54.0 | 108 |
| o-Xylene | 1.22 | 244 | 24.4 | 4.88 | | NA | 0.610 | 1.22 | 6.10 | 30.5 | 61.0 | 122 |
| n-Nonane | 1.03 | 206 | 20.6 | 4.12 | | NA | 0.515 | 1.03 | 5.15 | 25.8 | 51.5 | 103 |
| 1,1,2,2-Tetrachloroethane | 1.23 | 246 | 24.6 | 4.92 | ΔΗΠΗΠΗΙΑ | NA | 0.615 | 1.23 | 6.15 | 30.8 | 61.5 | 123 |
| Cumene | 1.08 | 216 | 21.6 | 4.32 | | NA | 0.540 | 1.08 | 5.40 | 27.0 | 54.0 | 108 |
| alpha-Pinene | 1.06 | 212 | 21.2 | 4.24 | | NA | 0.530 | 1.06 | 5.30 | 26.5 | 53.0 | 106 |
| n-Propylbenzene | 1.05 | 210 | 21.0 | 4.20 | | NA | 0.525 | 1.05 | 5.25 | 26.3 | 52.5 | 105 |
| 3-Ethyltoluene | 1.02 | 204 | 20.4 | 4.08 | | NA | 0.510 | 1.02 | 5.10 | 25.5 | 51.0 | 102 |
| 4-Ethyltoluene | 1.11 | 222 | 22.2 | 4.44 | | NA | 0.555 | 1.11 | 5.55 | 27.8 | 55.5 | 111 |
| 1,3,5-Trimethylbenzene | 1.08 | 216 | 21.6 | 4.32 | | NA | 0.540 | 1.08 | 5.40 | 27.0 | 54.0 | 108 |
| alpha-Methylstyrene | 1.02 | 204 | 20.4 | 4.08 | AIIIIIIIIIIIA | NA | 0.510 | 1.02 | 5.10 | 25.5 | 51.0 | 102 |
| 2-Ethyltoluene | 0.990 | 198 | 19.8 | 3.96 | /////////////////////////////////////// | NA | 0.495 | 0.990 | 4.95 | 24.8 | 49.5 | 99.0 |
| 1,2,4-Trimethylbenzene | 1.10 | 220 | 22.0 | 4.40 | <i>())))))))))))))))))))))))))))))))))))</i> | NA | 0.550 | 1.10 | 5.50 | 27.5 | 55.0 | 110 |
| n-Decane | 1.04 | 208 | 20.8 | 4.16 | | NA | 0.520 | 1.04 | 5.20 | 26.0 | 52.0 | 104 |
| Benzyl Chloride | 1.07 | 214 | 21.4 | 4.28 | <i>11111111111</i> | NA | 0.535 | 1.07 | 5.35 | 26.8 | 53.5 | 107 |
| 1,3-Dichlorobenzene | 1.06 | 212 | 21.2 | 4.24 | | NA | 0.530 | 1.06 | 5.30 | 26.5 | 53.0 | 106 |
| 1,4-Dichlorobenzene | 1.10 | 220 | 22.0 | 4.40 | X/////////// | NA | 0.550 | 1.10 | 5.50 | 27.5 | 55.0 | 110 |
| sec-Butylbenzene | 1.07 | 214 | 21.4 | 4.28 | AIIIIIIIIIIIIA | NA | 0.535 | 1.07 | 5.35 | 26.8 | 53.5 | 107 |
| p-Isopropyltoluene | 1.180 | 236 | 23.6 | 4.72 | 811111111111 | NA | 0.590 | 1.18 | 5.90 | 29.5 | 59.0 | 118 |
| 1,2,3-Trimethylbenzene | 1.10 | 220 | 22.0 | 4.40 | <i>XIIIIIIIIIIIX</i> | NA | 0.550 | 1.10 | 5.50 | 27.5 | 55.0 | 110 |
| 1,2-Dichlorobenzene | 1.08 | 216 | 21.6 | 4.32 | AIIIIIIIIIII | NA | 0.540 | 1.08 | 5.40 | 27.0 | 54.0 | 108 |
| d-Limonene | 1.06 | 212 | 21.2 | 4.24 | | NA | 0.530 | 1.06 | 5.30 | 26.5 | 53.0 | 106 |
| 1,2-Dibromo-3-chloropropane | 1.04 | 208 | 20.8 | 4.16 | | NA | 0.520 | 1.04 | 5.20 | 26.0 | 52.0 | 104 |
| n-Undecane | 1.05 | 210 | 21.0 | 4.20 | | NA | 0.525 | 1.05 | 5.25 | 26.3 | 52.5 | 105 |
| 1,2,4-Trichlorobenzene | 1.12 | 224 | 22.4 | 4.48 | 2///////////// | NA | 0.560 | 1.12 | 5.60 | 28.0 | 56.0 | 112 |
| Naphthalene | 1.05 | 210 | 21.0 | 4.20 | | NA | 0.525 | 1.05 | 5.25 | 26.3 | 52.5 | 105 |
| n-Dodecane | 1.06 | 212 | 21.2 | 4.24 | /////////////////////////////////////// | NA | 0.530 | 1.06 | 5.30 | 26.5 | 53.0 | 106 |
| Hexachloro-1,3-butadiene | 1.11 | 222 | 22.2 | 4.44 | | NA | 0.555 | 1.11 | 5.55 | 27.8 | 55.5 | 111 |

*Enter Information in the Solid Shaded Areas ONLY.

Q:\T015 Std. Concentrations\MS13 Std. Conc\R13041408\ICAL Conc. (Primary Source)

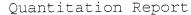
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| 2 | 0.5 | 1 | 25 | J:\MS13\DATA\2008_04\14\04140809.D |
| 3 | 1.0 | 1 | 25 | J:\MS13\DATA\2008_04\14\04140810.D |
| 4 | 5.0 | 5 | 25 | J:\MS13\DATA\2008_04\14\04140811.D |
| 5 | 25 | 27 | 25 | J:\MS13\DATA\2008_04\14\04140812.D |
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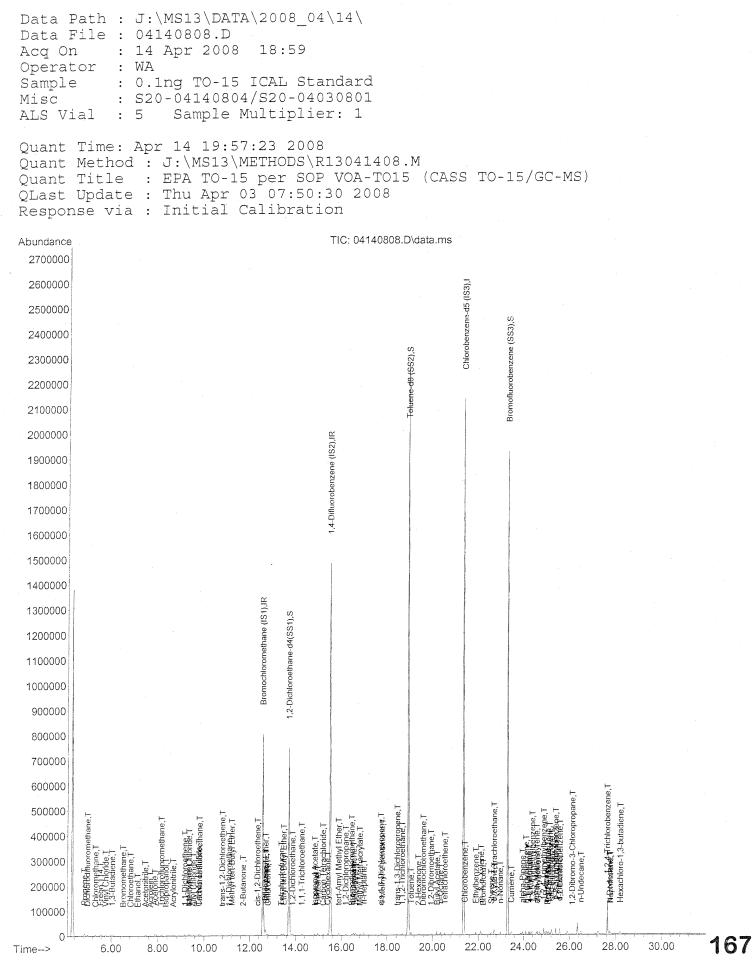
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|---|-----|-------------------|-------------------|-------------------|--|--|--|
| | | | | | | | |
| 1 | 0.1 | Apr 15 06:33 2008 | Apr 14 19:57 2008 | 14 Apr 2008 18:59 | | | |
| 2 | 0.5 | Apr 15 06:33 2008 | Apr 14 20:24 2008 | 14 Apr 2008 19:40 | | | |
| 3 | 1.0 | Apr 15 06:33 2008 | Apr 15 06:20 2008 | 14 Apr 2008 20:21 | | | |
| 4 | 5.0 | Apr 15 06:34 2008 | Apr 15 06:22 2008 | 14 Apr 2008 21:01 | | | |
| 5 | 25 | Apr 15 06:34 2008 | Apr 15 06:24 2008 | 14 Apr 2008 21:43 | | | |
| 6 | 50 | Apr 15 06:34 2008 | Apr 15 06:26 2008 | 14 Apr 2008 22:24 | | | |
| 7 | 100 | Apr 15 06:34 2008 | Apr 15 06:28 2008 | 14 Apr 2008 23:04 | | | |

R13041408.M Tue Apr 15 15:34:11 2008

18# 4/15/08



(QT Reviewed)



R13041408.M Mon Apr 14 20:02:24 2008

| Data Path : J:\MS13\DATA\2008_04 Data File : 04140808.D Acq On : 14 Apr 2008 18:59 Operator : WA Sample : 0.1ng TO-15 ICAL Sta Misc : S20-04140804/S20-040 ALS Vial : 5 Sample Multiplie Quant Time: Apr 14 19:57:23 2008 | indard 30801 er: 1 | 0 | | | | |
|---|---|--|---|--|----------------------------------|--|
| Quant Method : J:\MS13\METHODS\R Quant Title : EPA TO-15 per SOP QLast Update : Thu Apr 03 07:50: Response via : Initial Calibrati | • VOA-TO 30 2008 | 15 (C | ASS TO-15/ | GC-MS) | | |
| Internal Standards | R.T. | QIon | Response | Conc Un: | its Dev | (Min) |
| 1) Bromochloromethane (IS1) 37) 1,4-Difluorobenzene (IS2) 56) Chlorobenzene-d5 (IS3) | 15.51 | 114 | 1516799 | 25.000 ng | g - C | 0.02 |
| System Monitoring Compounds 33) 1,2-Dichloroethane-d4(Spiked Amount 25.000 57) Toluene-d8 (SS2) Spiked Amount 25.000 73) Bromofluorobenzene (SS3) Spiked Amount 25.000 | 18.93 | 98 | Recov 583069 | ery = 28.176 ng ery = | 86.76% g -C L12.72% g C | 0.01 |
| 6) Vinyl Chloride 7) 1,3-Butadiene 8) Bromomethane 9) Chloroethane 10) Ethanol 11) Acetonitrile 12) Acrolein 13) Acetone 14) Trichlorofluoromethane 15) Isopropanol 16) Acrylonitrile 17) 1,1-Dichloroethene 18) tert-Butanol 19) Methylene Chloride 20) Allyl Chloride 21) Trichlorotrifluoroethane | 5.33 5.55 5.77 6.04 6.52 6.84 7.14 7.68 7.89 8.16 8.35 8.67 9.34 9.36 9.36 9.561 9.78 10.80 11.22 0.00 11.71 12.34 12.70 12.72 | 85 50 135 62 54 94 45 46 58 105 58 41 56 84 11 61 33 62 17 61 87 61 87 61 87 61 | 6457 5608 3051 4872 3496 2038 1736 905 5842 1483 4681 4631 7164m 2141m 2684 3724 3365 1936 2299 8282 3798 4737 8322 0 1112 3769 1863 223 | 0.101 ng 0.113 ng 0.096 ng 0.086 ng 0.079 ng 0.070 ng 0.070 ng 0.094 ng 0.087 ng 0.083 ng 0.083 ng 0.083 ng 0.083 ng 0.083 ng 0.121 ng 0.049 ng 0.121 ng 0.047 ng 0.121 ng 0.047 ng 0.121 ng 0.047 ng 0.054 ng | | 93 85 89 94 61 91 85 44 83 73 56 96 72 84 95 88 95 88 96 93 |

213041408.M Mon Apr 14 20:02:23 2008

*j*84 4/15/08

| | Quantitation | Report | (QT | Reviewed) | | |
|--|--|--|--|---|--|---|
| Data F Acq Or Operat Sample Misc | Path : J:\MS13\DATA\2008_04 File : 04140808.D n : 14 Apr 2008 18:59 for : WA e : 0.1ng TO-15 ICAL Sta : S20-04140804/S20-040 Lal : 5 Sample Multiplie | ndard 30801 | | | | |
| Quant Quant QLast | Time: Apr 14 19:57:23 2008 Method : J:\MS13\METHODS\R Title : EPA TO-15 per SOF Update : Thu Apr 03 07:50: nse via : Initial Calibrati | 21304140 > VOA-TC 30 2008 |)15 (CA | SS TO-15/G | C-MS) | |
| Inter | nal Standards | | | | | |
| 34) 35) 36) 38) 39) 40) 41) 42) 43) 42) 42) 43) 42) 43) 42) 43) 42) 43) 42) 42) 43) 42) 43) 42) 52) 53) 53) 53) 54) 55) 52) 53) 52) 52) 53) 52) 53) 52) 52) 52) 52) 52) 52) 52) 52) 52) 52 | Chloroform Tetrahydrofuran Ethyl tert-Butyl Ether 1,2-Dichloroethane 1,1,1-Trichloroethane Isopropyl Acetate 1-Butanol Benzene Carbon Tetrachloride Cyclohexane tert-Amyl Methyl Ether 1,2-Dichloropropane Bromodichloromethane Trichloroethene 1,4-Dioxane Isooctane Methyl Methacrylate n-Heptane cis-1,3-Dichloropropene 4-Methyl-2-pentanone trans-1,3-Dichloropropene 1,1,2-Trichloroethane Toluene 2-Hexanone Dibromochloromethane 1,2-Dibromoethane Butyl Acetate n-Octane Tetrachloroethene Chlorobenzene Ethylbenzene m- & p-Xylene Bromoform Styrene o-Xylene n-Nonane 1,1,2,2-Tetrachloroethane Cumene | 12.79 13.39 13.49 13.88 14.27 14.85 14.92 14.92 14.92 15.40 15.40 15.40 15.40 16.45 16.45 16.53 16.53 16.62 16.81 16.98 17.79 18.45 18.67 19.07 19.39 19.60 19.39 19.60 19.94 20.21 20.35 20.54 21.41 21.89 22.58 22.70 23.46 23.96 | $\begin{array}{c} 83\\72\\87\\62\\97\\61\\56\\78\\117\\84\\73\\83\\130\\85\\70\\71\\58\\75\\97\\129\\143\\129\\107\\43\\57\\62\\112\\91\\173\\104\\43\\83\\105\\93\end{array}$ | 4622 1382 2636 3788 3981 1323m 786 10664 2729 3527 6246 3114 3734 2307 1490 10380 424m 2860 3001 2240 2661 2577 10684 6288 2380 2269 5465 2088 2776 6476 10115 15249 1417 6125 8518 5870 4182 8543 4522 | 0.106 ng 0.071 ng 0.082 ng 0.087 ng 0.102 ng 0.102 ng 0.075 ng 0.030 ng 0.082 ng 0.090 ng 0.081 ng 0.105 ng 0.105 ng 0.107 ng 0.102 ng 0.102 ng 0.084 ng 0.084 ng 0.045 ng 0.045 ng 0.068 ng 0.107 ng 0.068 ng 0.111 ng 0.119 ng 0.069 ng 0.111 ng 0.119 ng 0.089 ng 0.115 ng 0.098 ng 0.092 ng 0.135 ng 0.092 ng 0.135 ng 0.097 ng 0.122 ng 0.099 ng 0.125 ng 0.125 ng 0.127 ng 0.099 ng 0.127 ng 0.099 ng 0.127 ng 0.099 ng 0.127 ng 0.097 ng 0.106 ng 0.117 ng 0.095 ng 0.091 ng | <pre># 83 ## 84 # 66 # 79 95 88 # 84 90 95 88 84 90 95 88 84 90 95 84 # 75 90 94 91 94 95 88 87 97 88 50 97 98 85 67 97 80 85 67 97 80 85 67 97 82 87 87 87 87 87 87 87 87 87 87 87 87 87</pre> |
| 78) 4 | 3-Ethyltoluene 4-Ethyltoluene 1,3,5-Trimethylbenzene | 24.28 | 105 | 9132 9233 7705 | 0.094 ng 0.101 ng 0.094 ng | 95 96 92 169 |

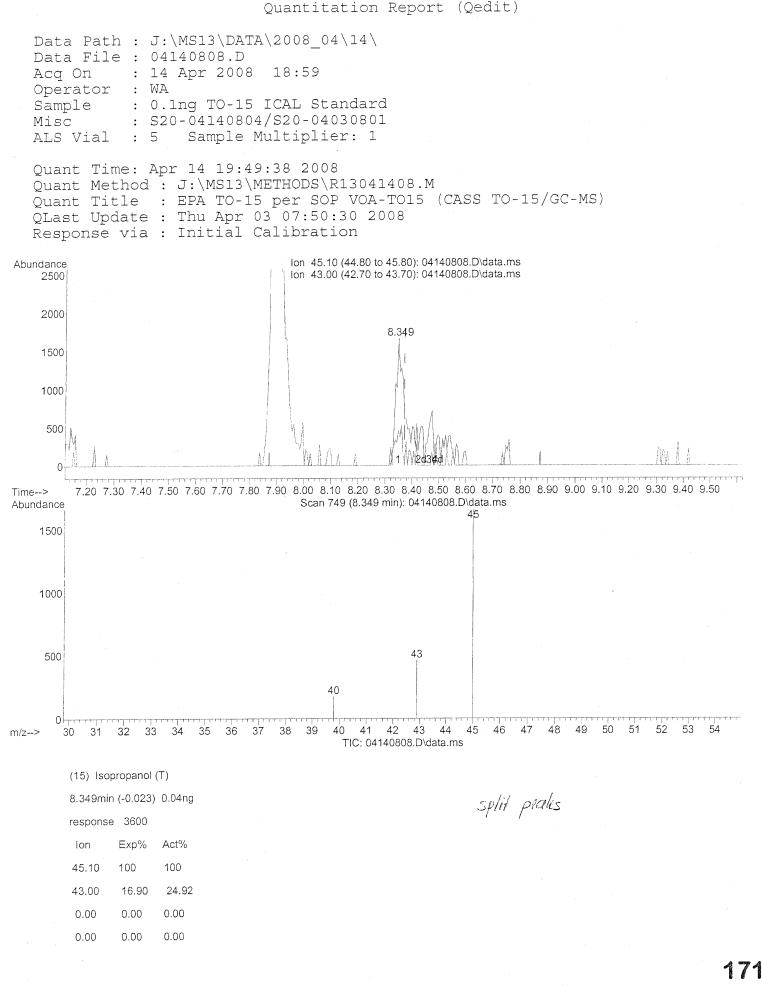
213041408.M Mon Apr 14 20:02:23 2008

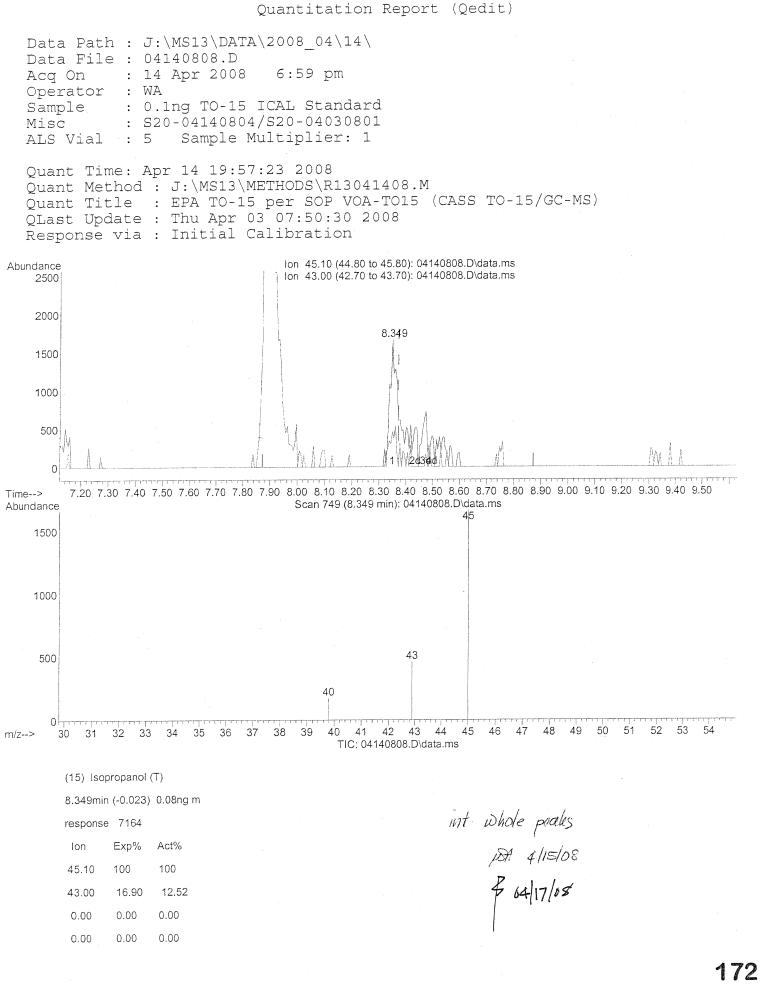
jor 4/15/08

Page: 2

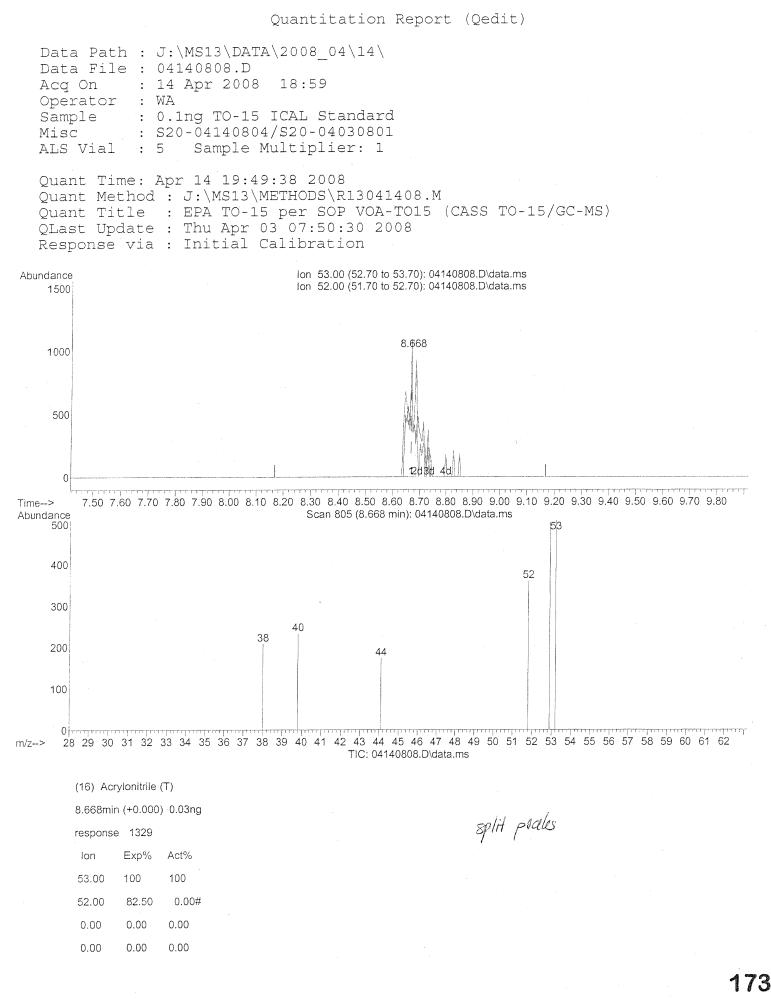
| Data Path : J:\MS13\DATA\2008_04 Data File : 04140808.D Acq On : 14 Apr 2008 18:59 Operator : WA Sample : 0.1ng TO-15 ICAL Stat Misc : S20-04140804/S20-040 ALS Vial : 5 Sample Multiplie | ndard 30801 | | | | | | |
|---|---|--|--|--|------|--|--|
| Quant Time: Apr 14 19:57:23 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Thu Apr 03 07:50:30 2008 Response via : Initial Calibration | | | | | | | |
| Internal Standards | R.T. | QIon | Response | Conc Units | Dev(| Min) | |
| <pre>82) 1,2,4-Trimethylbenzene 83) n-Decane 84) Benzyl Chloride 85) 1,3-Dichlorobenzene 86) 1,4-Dichlorobenzene 87) sec-Butylbenzene 88) p-Isopropyltoluene 89) 1,2,3-Trimethylbenzene 90) 1,2-Dichlorobenzene 91) d-Limonene 92) 1,2-Dibromo-3-Chloropr 93) n-Undecane 94) 1,2,4-Trichlorobenzene 95) Naphthalene</pre> | 24.61 24.88 24.99 25.04 25.16 25.22 25.40 25.58 25.57 26.12 26.50 27.64 27.74 | 105 57 91 146 146 105 119 105 146 68 157 57 180 128 57 | 3727 9246 7668 4427 4100 4705 4842 10671 8726 7415 3907 3330 728 4750 2704 7336 5292 1398 | 0.088 ng 0.082 ng 0.060 ng 0.110 ng 0.115 ng | | 97 98 99 88 87 94 84 93 | |

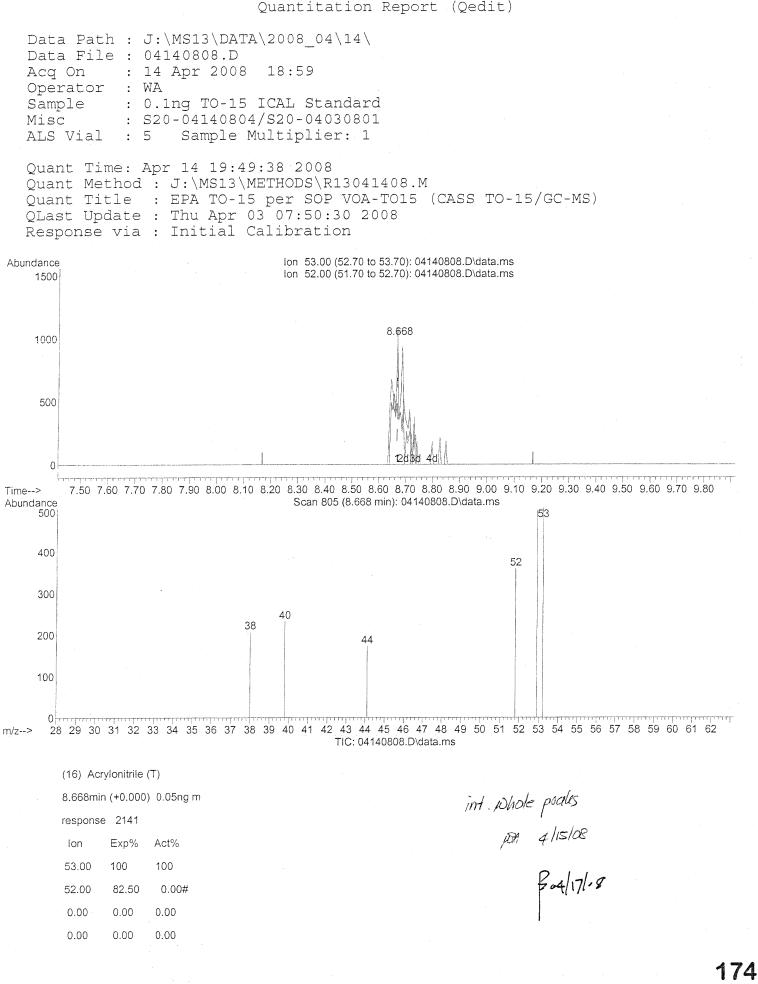
(#) = qualifier out of range (m) = manual integration (+) = signals summed



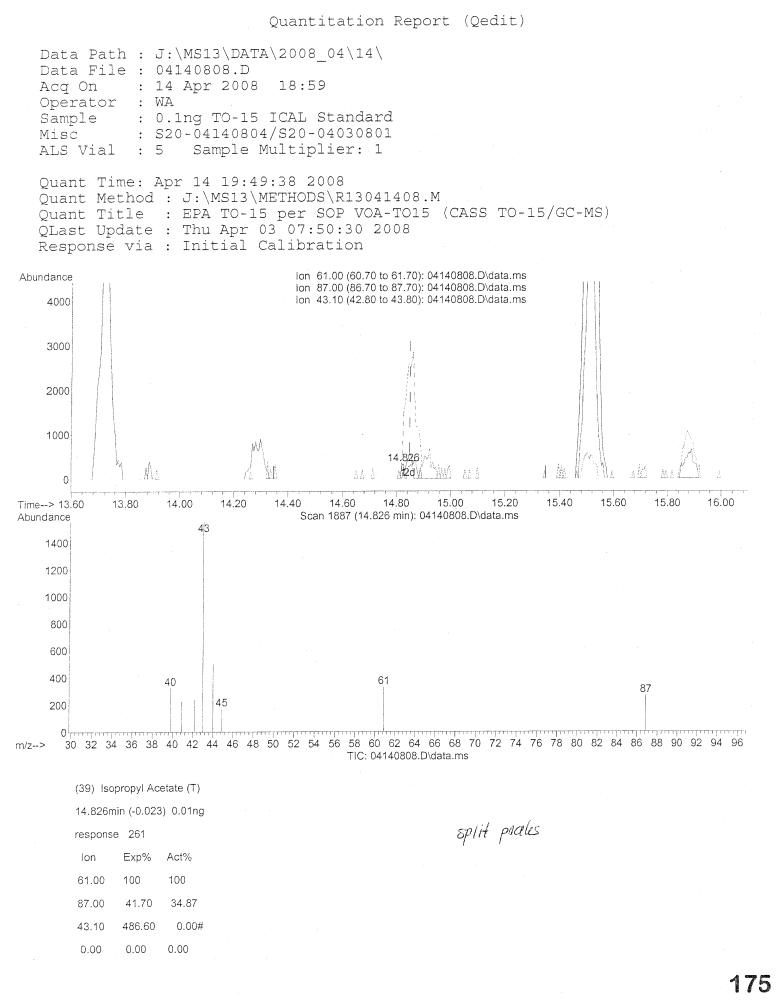


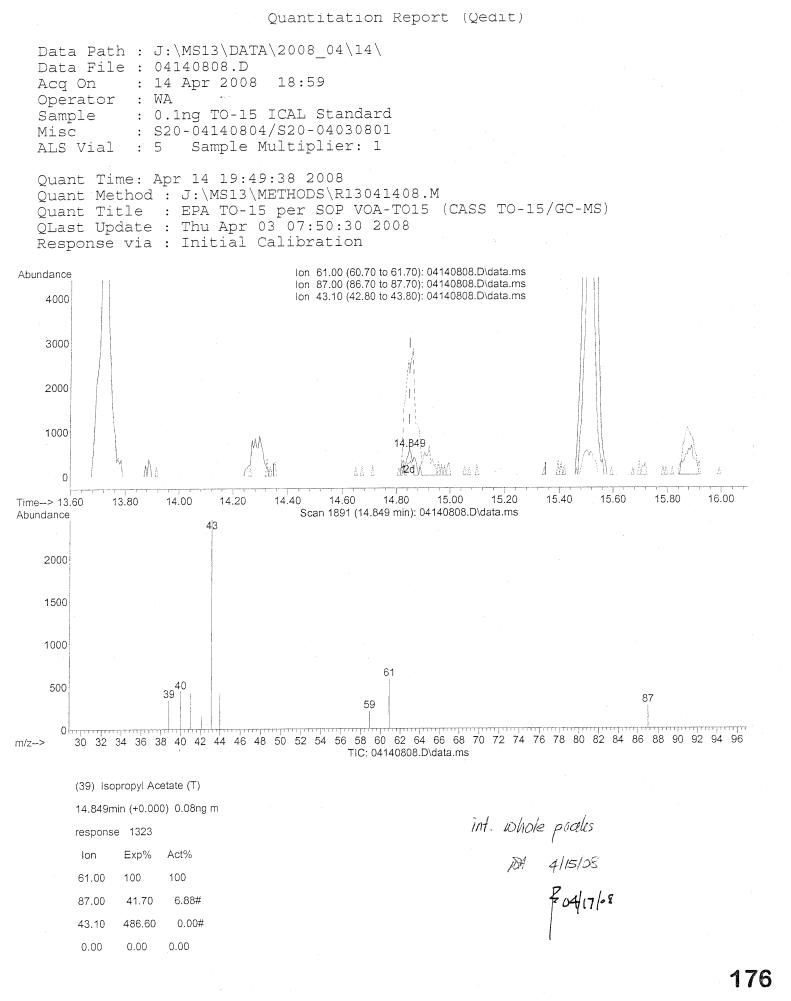
R13041408.M Tue Apr 15 15:33:31 2008

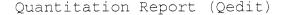


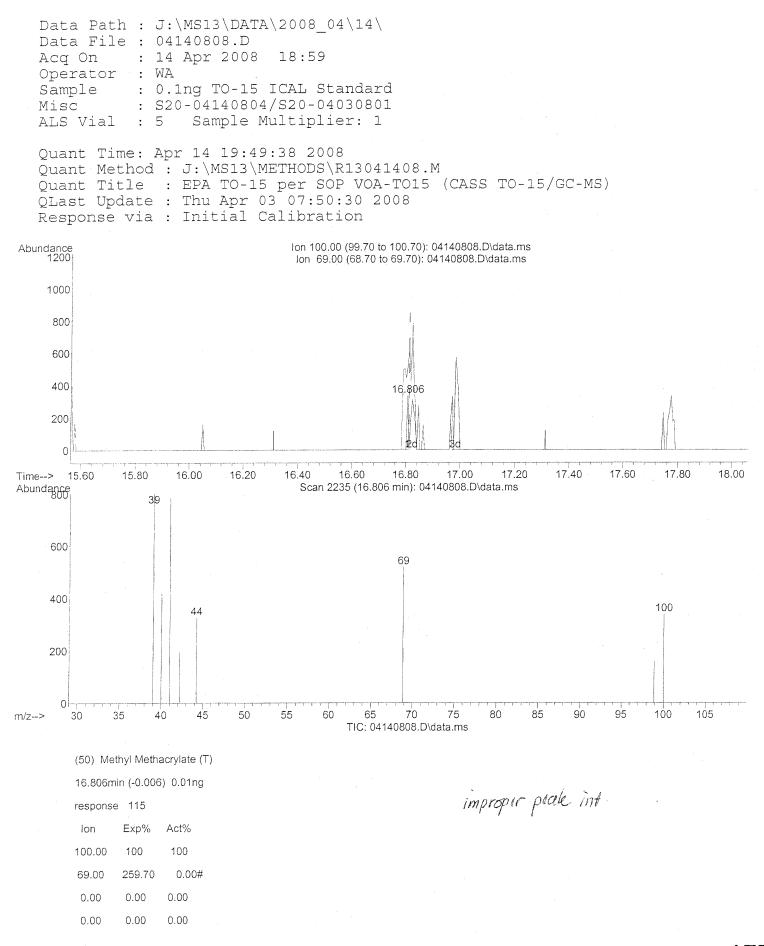


Page: 1



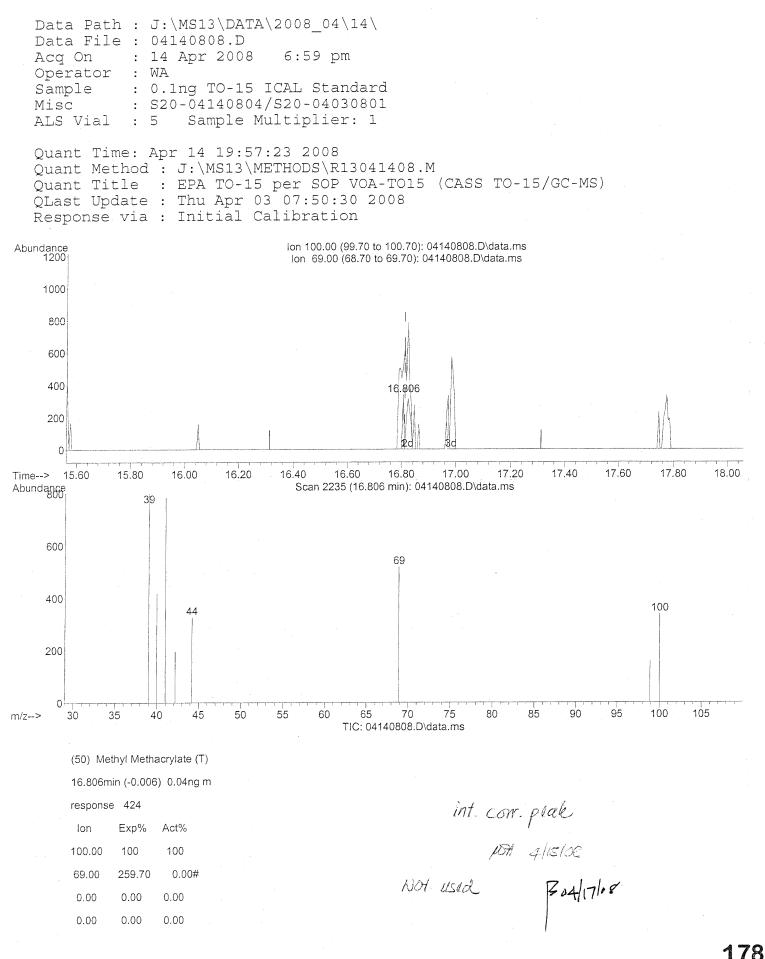




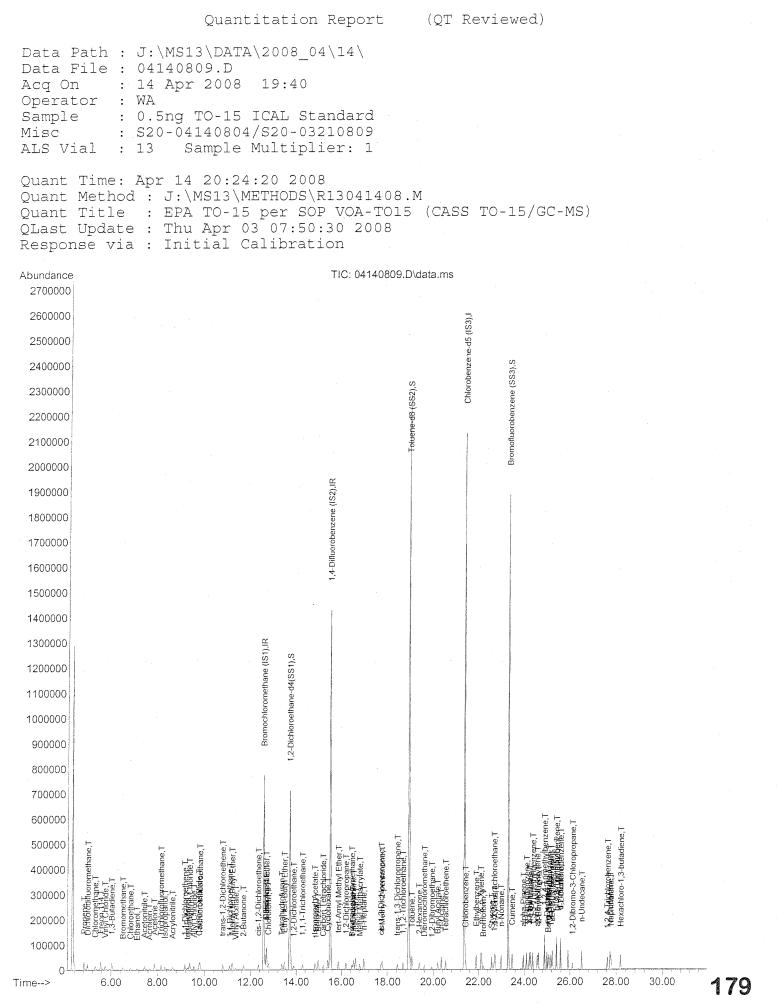


Page: 1





R13041408.M Tue Apr 15 15:32:53 2008



R13041408.M Mon Apr 14 20:27:49 2008

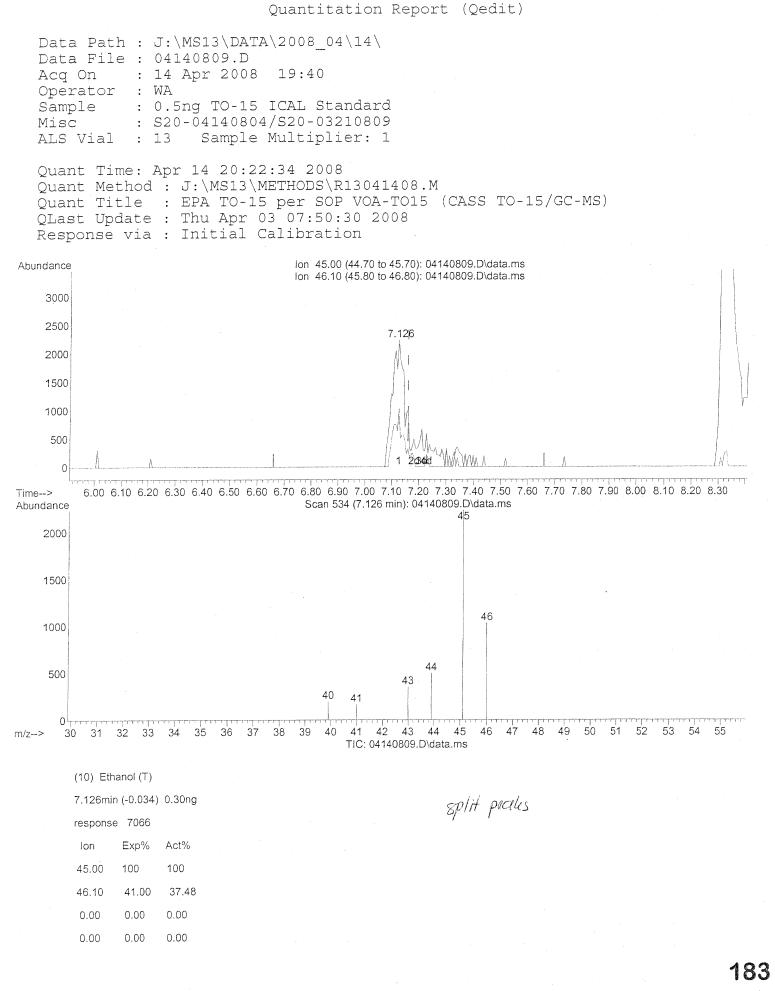
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|---|---|---|---|---|---|------------------------------------|
| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
| 1) Bromochloromethane (IS1) 37) 1,4-Difluorobenzene (IS2) 56) Chlorobenzene-d5 (IS3) | 12.58 15.51 21.35 | 130 114 82 | 321232 1468142 744311 | 25.000 25.000 25.000 | ng ng ng | -0.03 -0.02 0.00 |
| System Monitoring Compounds 33) 1,2-Dichloroethane-d4(Spiked Amount 25.000 57) Toluene-d8 (SS2) | | | 654334 Recove 1694424 | ery = | 87 | .24% |
| Spiked Amount 25.000 73) Bromofluorobenzene (SS3) Spiked Amount 25.000 | | | Recove 565263 | ery = | 112 ng | .96% 0.00 |
| <pre>3) Dichlorodifluoromethane 4) Chloromethane 5) Freon 114 6) Vinyl Chloride 7) 1,3-Butadiene 8) Bromomethane 9) Chloroethane 10) Ethanol 11) Acetonitrile 12) Acrolein 13) Acetone 14) Trichlorofluoromethane 15) Isopropanol 16) Acrylonitrile 17) 1,1-Dichloroethene 18) tert-Butanol 19) Methylene Chloride 20) Allyl Chloride 21) Trichlorotrifluoroethane 22) Carbon Disulfide 23) trans-1,2-Dichloroethene 24) 1,1-Dichloroethane 25) Methyl tert-Butyl Ether 26) Vinyl Acetate 27) 2-Butanone 28) cis-1,2-Dichloroethene</pre> | 4.99 5.31 5.56 5.75 6.03 6.52 6.85 7.13 7.44 7.66 7.88 8.16 8.33 8.65 9.17 9.28 9.37 9.56 9.37 9.56 9.78 10.80 11.21 11.21 11.36 11.69 | 85 132 94 45 145 156 145 156 156 156 156 156 156 156 15 | 23085 13818 20819 16796 10055 9331 10603m 26981 5948 13071 21260 36380m 15227 9768 | 0.422 0.427 0.528 0.420 0.427 0.402 0.385 0.445 0.445 0.447 0.321 0.533 0.412 0.434 0.389 0.401 0.386 0.454 0.350 0.493 0.353 0.397 0.385 0.385 0.385 0.387 0.385 0.396 | ng n | 99 97 96 98 # 71 98 |

De 4/16/08

| Quantitation | Report | (Q1 | Reviewed) | | |
|---|---|---|---|--|--|
| Data Path : J:\MS13\DATA\2008_04 Data File : 04140809.D Acq On : 14 Apr 2008 19:40 Operator : WA Sample : 0.5ng TO-15 ICAL Sta | | | | | |
| Sample : 0.5ng TO-15 ICAL Sta Misc : S20-04140804/S20-032 ALS Vial : 13 Sample Multipli | ler: 1 | | | | |
| Quant Time: Apr 14 20:24:20 2008 Quant Method : J:\MS13\METHODS\F Quant Title : EPA TO-15 per SOF QLast Update : Thu Apr 03 07:50: Response via : Initial Calibrati | 81304140 9 VOA-TO 30 2008 .on |)15 (CA } | | | |
| Internal Standards | R.T. | QIon | Response | Conc Units | B Dev(Min) |
| <pre>34) Tetrahydrofuran 35) Ethyl tert-Butyl Ether 36) 1,2-Dichloroethane 38) 1,1,1-Trichloroethane 39) Isopropyl Acetate 40) 1-Butanol 41) Benzene 42) Carbon Tetrachloride 43) Cyclohexane 44) tert-Amyl Methyl Ether 45) 1,2-Dichloropropane 46) Bromodichloromethane 47) Trichloroethene 48) 1,4-Dioxane 49) Isooctane 50) Methyl Methacrylate 51) n-Heptane 52) cis-1,3-Dichloropropene 53) 4-Methyl-2-pentanone 54) trans-1,3-Dichloropropene 55) 1,1,2-Trichloroethane 58) Toluene 59) 2-Hexanone 60) Dibromochloromethane 61) 1,2-Dibromoethane 62) Butyl Acetate 63) n-Octane 64) Tetrachloroethene 65) Chlorobenzene 66) Ethylbenzene 67) m- & p-Xylene 68) Bromoform 69) Styrene 70) o-Xylene 71) n-Nonane 72) 1,1,2,2-Tetrachloroethane 74) Cumene</pre> | 13.37 13.49 13.89 14.29 14.84 14.99 15.21 15.41 15.41 15.40 16.20 16.46 16.53 16.50 16.62 16.81 16.98 17.73 18.43 18.67 19.06 19.38 19.06 19.38 19.94 20.19 20.35 20.54 21.41 21.89 20.57 22.21 22.57 22.69 23.46 | $\begin{array}{c} 72\\ 87\\ 62\\ 97\\ 61\\ 56\\ 78\\ 117\\ 84\\ 73\\ 63\\ 83\\ 130\\ 857\\ 100\\ 71\\ 75\\ 58\\ 75\\ 97\\ 43\\ 129\\ 107\\ 43\\ 57\\ 166\\ 291\\ 173\\ 104\\ 43\\ 83\\ 105 \end{array}$ | 7992 12073 17281 17708 7193 9296 44943 13169 16422 28263 12735 14378 11916 7698 50293 3798 12027 15511 10899 14502 10473 48005 31211 10573 10738 32810 10003 11015 29701 52327 76463 8077 27145 38188 27324 17907 45202 | 0.436 Hg 0.410 ng 0.434 ng 0.381 ng 0.427 ng 0.427 ng 0.426 ng 0.426 ng 0.426 ng 0.421 ng 0.421 ng 0.423 ng 0.436 ng 0.366 ng 0.391 ng 0.388 ng 0.366 ng 0.391 ng 0.388 ng 0.464 ng 0.544 ng 0.547 ng 0.469 ng 0.449 ng 0.547 ng 0.547 ng 0.547 ng 0.568 ng 0.521 ng 1.148 ng 0.565 ng 0.478 ng 0.536 ng 0.464 ng 0.536 ng 0.464 ng 0.536 ng 0.464 ng 0.536 ng 0.464 ng 0.536 ng 0.464 ng 0.523 ng 0.511 ng | <pre>94 78 91 95 43 # 32 98 99 99 99 98 97 97 67 7 97 67 7 97 67 7 97 67 7 97 88 99 99 98 97 99 98 97 97 67 4 99 99 99 98 97 97 67 4 99 99 88 97 97 67 67 67 88 93 76 96 88 94 93 76 95 88 94 94 93 76 95 88 94 94 95 88 97 97 67 67 88 99 88 99 99 88 97 97 67 67 88 99 88 99 99 88 97 97 67 67 88 99 99 88 99 99 88 97 97 67 67 67 67 88 99 88 99 87 99 99 88 97 97 67 67 88 99 88 99 83 99 99 88 99 99 88 99 99 88 99 99 88 99 99</pre> |
| 76) n-Propylbenzene | 24.10 24.23 24.28 | 91 105 105 | 22191 59886 45370 43433 38128 | 0.456 ng 0.509 ng 0.475 ng 0.485 ng 0.476 ng | 95 |

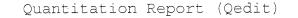
Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140809.D Acq On : 14 Apr 2008 19:40 Operator : WA Sample : 0.5ng TO-15 ICAL Standard Misc : S20-04140804/S20-03210809 ALS Vial : 13 Sample Multiplier: 1 Quant Time: Apr 14 20:24:20 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Thu Apr 03 07:50:30 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards 11111111111111111111111111180) alpha-Methylstyrene24.56118175750.432 ng8781) 2-Ethyltoluene24.61105443410.468 ng9882) 1,2,4-Trimethylbenzene24.88105400980.467 ng9383) n-Decane24.9857226370.425 ng8284) Benzyl Chloride25.0491246110.367 ng10085) 1,3-Dichlorobenzene25.08146226080.547 ng9986) 1,4-Dichlorobenzene25.16146226080.547 ng9987) sec-Butylbenzene25.21105522470.508 ng9588) p-Isopropyltoluene25.39119426060.518 ng8991,2-Dichlorobenzene25.58146216740.504 ng9991) d-Limonene25.5868166670.411 ng9192) 1,2-Dibromo-3-Chloropr...26.1115759140.499 ng4793) n-Undecane27.63180145910.504 ng9995) Naphthalene27.7457239960.436 ng9097) Hexachloro-1,3-butadiene28.1922595560.556 ng95

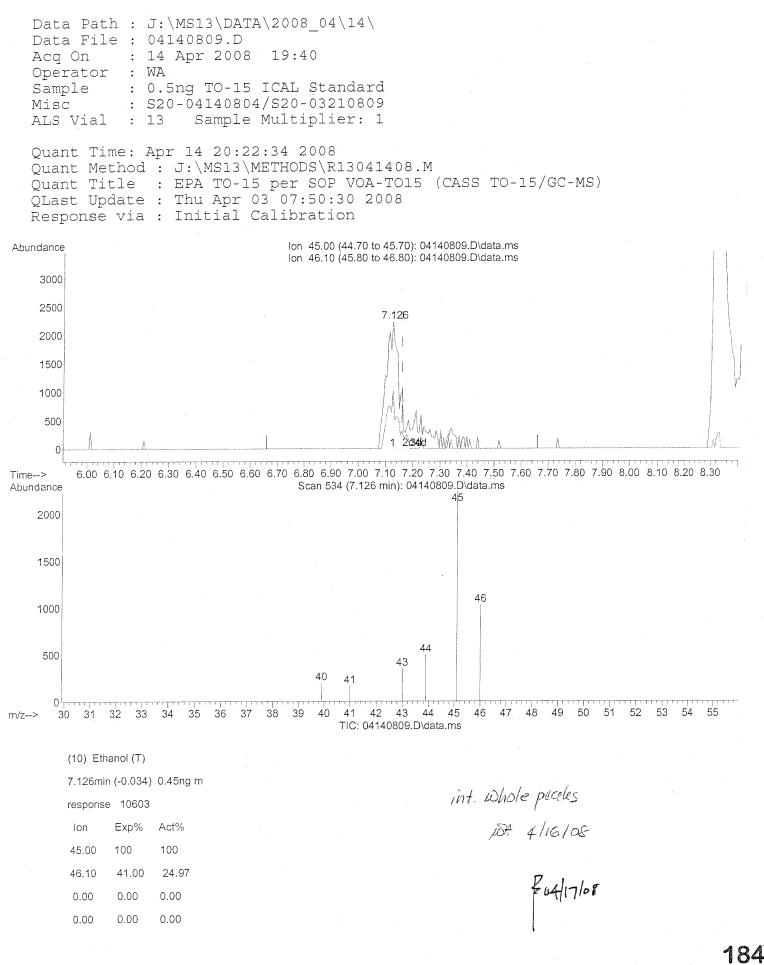
(#) = qualifier out of range (m) = manual integration (+) = signals summed

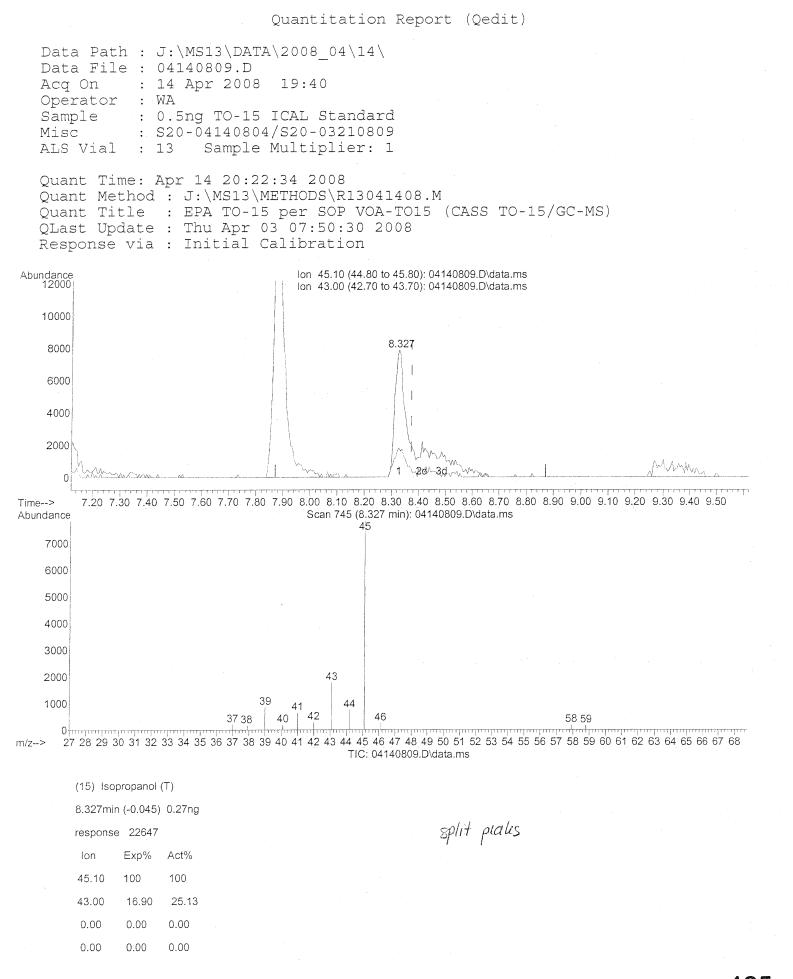


R13041408.M Mon Apr 14 20:23:25 2008

Page: 1

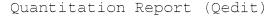


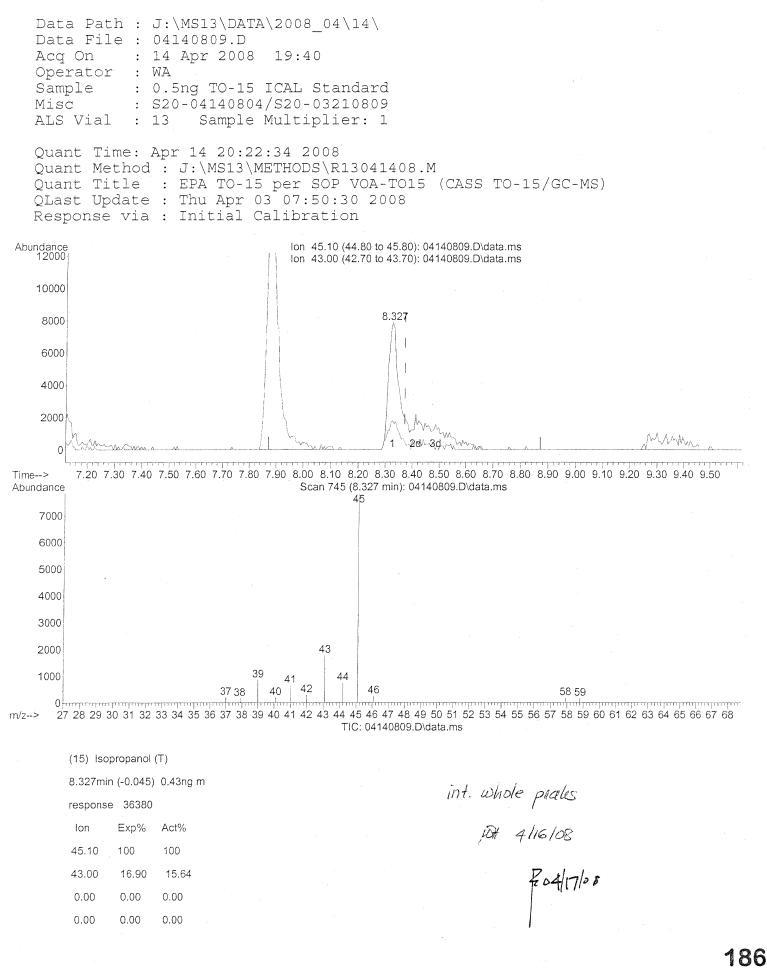




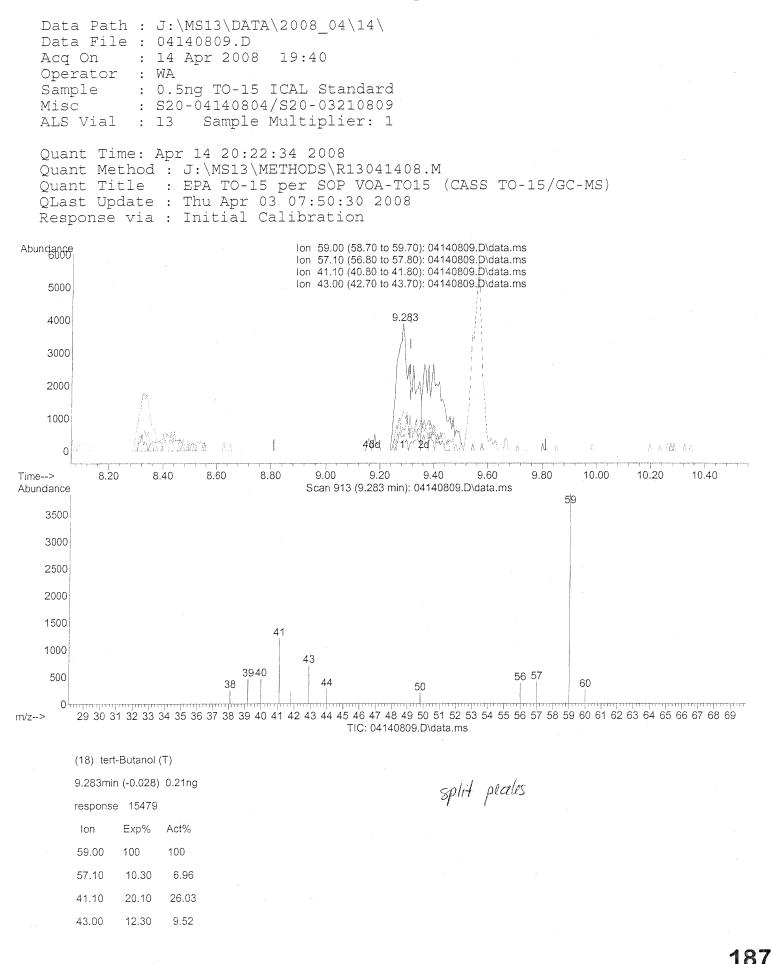
R13041408.M Mon Apr 14 20:24:03 2008

Page: 1

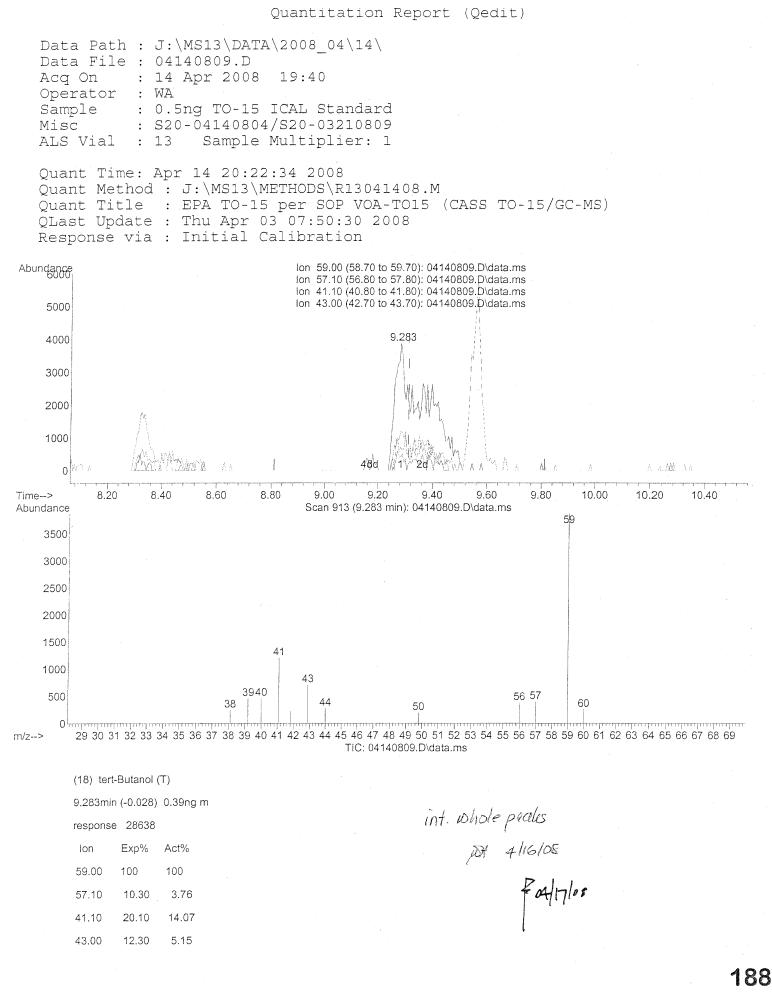


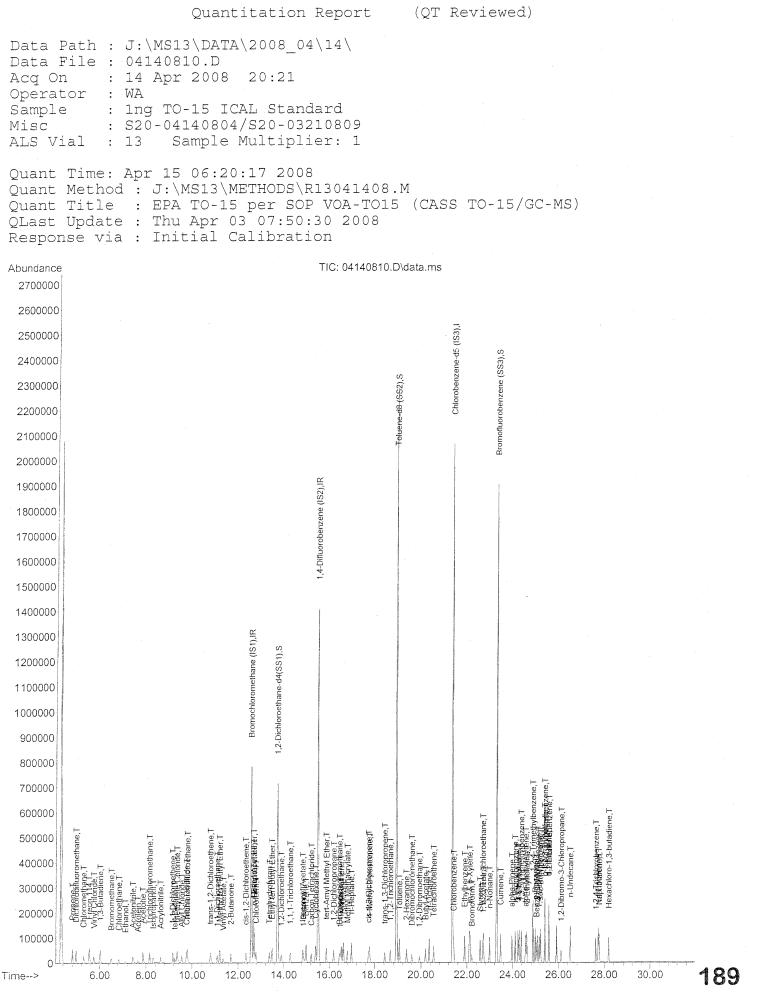


R13041408.M Mon Apr 14 20:24:11 2008



R13041408.M Mon Apr 14 20:24:25 2008





R13041408.M Tue Apr 15 06:21:26 2008

Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140810.D Acg On : 14 Apr 2008 20:21 Operator : WA Sample : 1ng TO-15 ICAL Standard Misc : S20-04140804/S20-03210809 ALS Vial : 13 Sample Multiplier: 1 Ouant Time: Apr 15 06:20:17 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Thu Apr 03 07:50:30 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards _____ 1) Bromochloromethane (IS1)12.5813031446125.000 ng-0.0337) 1,4-Difluorobenzene (IS2)15.51114145464725.000 ng-0.0256) Chlorobenzene-d5 (IS3)21.358273508325.000 ng0.00 System Monitoring Compounds

 Spiked Amount
 25.000
 Recovery
 = 100.000

 Target Compounds
 Qvalue

 2) Propene
 4.81
 42
 28935
 0.777 ng
 96

 3) Dichlorodifluoromethane
 4.97
 85
 50859
 0.808 ng
 99

 4) Chloromethane
 5.29
 50
 43114
 0.814 ng
 92

 5) Freon 114
 5.54
 135
 23952
 0.936 ng
 99

 6) Vinyl Chloride
 5.74
 62
 3763
 0.770 ng
 94

 7) 1, 3-Butadiene
 6.02
 54
 23952
 0.936 ng
 99

 6) Vinyl Chloride
 5.74
 62
 3763
 0.770 ng
 94

 9) Chloromethane
 6.151
 94
 17846
 0.729 ng
 94

 9) Chloromethane
 6.81
 61
 1800
 0.651 ng
 95

 10) Ethanol
 7.11
 45
 18205m
 0.761 ng
 96

 12) Accolein
 7.66
 56
 11810
 0.651 ng
 96

 13) Acetone
 7.87
 58
 23380
 0.7761 ng
 96

 17) 1,1-⁹³190

| Data Path : J:\MS13\DATA\2008_04 Data File : 04140810.D Acq On : 14 Apr 2008 20:21 Operator : WA Sample : 1ng TO-15 ICAL Stand Misc : S20-04140804/S20-032 ALS Vial : 13 Sample Multipl: Quant Time: Apr 15 06:20:17 2008 | dard 210809 ier: 1 | | | | |
|---|---|--|--|--|---|
| Quant Method : J:\MS13\METHODS\F Quant Title : EPA TO-15 per SOF QLast Update : Thu Apr 03 07:50: Response via : Initial Calibrati | R1304140 P VOA-TC :30 2008 |)15 (CA | SS TO-15/0 | BC-MS) | |
| Internal Standards | | | | | |
| <pre>34) Tetrahydrofuran 35) Ethyl tert-Butyl Ether 36) 1,2-Dichloroethane 38) 1,1,1-Trichloroethane 39) Isopropyl Acetate 40) 1-Butanol 41) Benzene 42) Carbon Tetrachloride 43) Cyclohexane 44) tert-Amyl Methyl Ether 45) 1,2-Dichloropropane 46) Bromodichloromethane 47) Trichloroethene 48) 1,4-Dioxane 49) Isooctane 50) Methyl Methacrylate 51) n-Heptane 52) cis-1,3-Dichloropropene 53) 4-Methyl-2-pentanone 54) trans-1,3-Dichloropropene 53) 4-Methyl-2-pentanone 54) trans-1,3-Dichloropropene 55) 1,1,2-Trichloroethane 59) 2-Hexanone 60) Dibromochloromethane 61) 1,2-Dibromoethane 62) Butyl Acetate 63) n-Octane 64) Tetrachloroethene 65) Chlorobenzene 66) Ethylbenzene 67) m- & p-Xylene 68) Bromoform</pre> | 18.67 19.07 19.37 19.60 19.93 20.19 20.35 20.54 21.41 21.89 22.12 22.21 22.57 22.71 22.98 | 97 91 43 129 107 43 57 166 112 91 173 104 91 43 | 18987 87414 56543 20024 19169 60902 18987 20868 53156 93406 144511 16953 51489 71744 48600 | 0.850 ng 1.002 ng 0.822 ng 0.998 ng 0.851 ng 0.863 ng 1.049 ng 1.029 ng 0.942 ng 2.196 ng 1.201 ng 0.919 ng 1.020 ng 0.835 ng | 96 94 81 98 94 83 93 97 95 94 89 98 95 93 # |
| 74) Cumene 75) alpha-Pinene 76) n-Propylbenzene 77) 3-Ethyltoluene | 23.46 23.96 24.10 24.23 24.28 | 105 93 91 105 105 | 34347 85006 42697 112564 86580 86229 73399 | 1.015 ng 0.973 ng 0.888 ng 0.969 ng 0.917 ng 0.975 ng 0.928 ng | 97 98 88 93 92 97 ⁹⁷ 191 |

13041408.M Tue Apr 15 06:21:26 2008

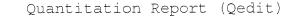
101 4/16/08

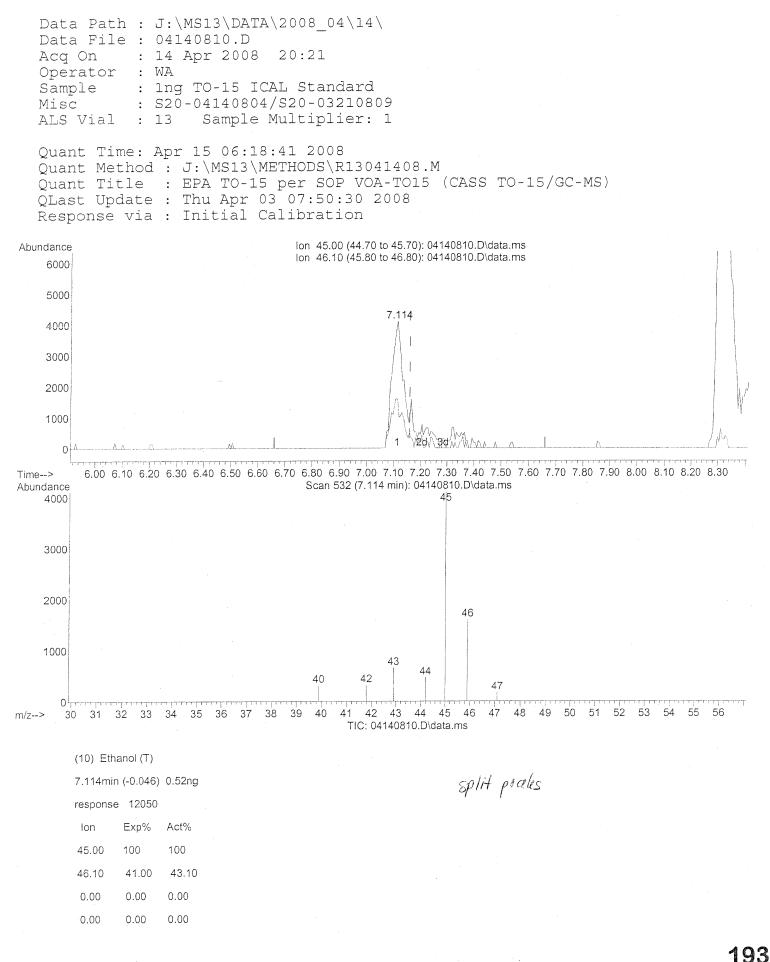
Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140810.D Acq On : 14 Apr 2008 20:21 Operator : WA Sample : 1ng TO-15 ICAL Standard Misc : S20-04140804/S20-03210800 : S20-04140804/S20-03210809 ALS Vial : 13 Sample Multiplier: 1 Quant Time: Apr 15 06:20:17 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Thu Apr 03 07:50:30 2008 Response via : Initial Calibration Internal Standards R.T. QIon Response Conc Units Dev(Min) And the standardsA.1. Qion ResponseConc ontes Dev(Min80) alpha-Methylstyrene24.56118341560.850 ng9381) 2-Ethyltoluene24.61105815530.872 ng9682) 1,2,4-Trimethylbenzene24.88105745380.879 ng9683) n-Decane24.9857447550.851 ng8984) Benzyl Chloride25.0491484450.731 ng9485) 1,3-Dichlorobenzene25.08146429501.039 ng9986) 1,4-Dichlorobenzene25.16146441991.083 ng9487) sec-Butylbenzene25.21105977170.961 ng9588) p-Isopropyltoluene25.39119830691.023 ng9089) 1,2,3-Trimethylbenzene25.58146410870.967 ng9690) 1,2-Dichlorobenzene25.5868309430.772 ng9492) 1,2-Dibromo-3-Chloropr...26.11157131301.122 ng#94) 1,2,4-Trichlorobenzene27.63180288411.008 ng9995) Naphthalene27.77128923401.081 ng9596) n-Dodecane27.7457449230.827 ng9097) Hexachloro-1,3-butadiene28.19225185131.090 ng95 _____ 93 _____

(#) = qualifier out of range (m) = manual integration (+) = signals summed

13041408.M Tue Apr 15 06:21:26 2008

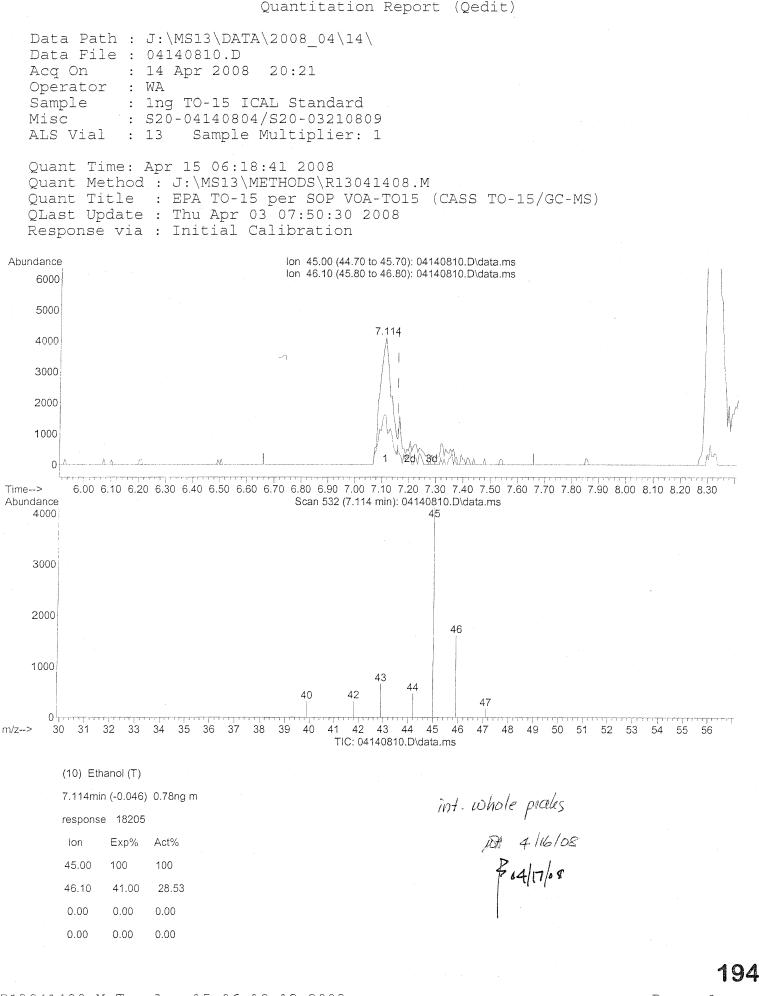
181 4/16/08



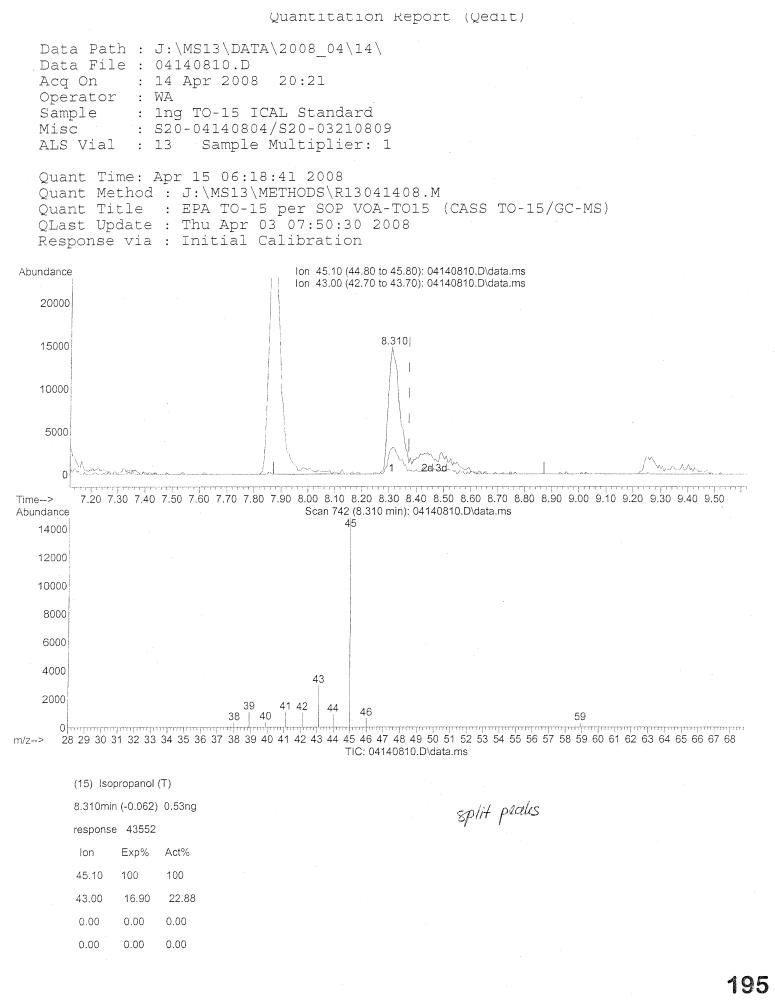


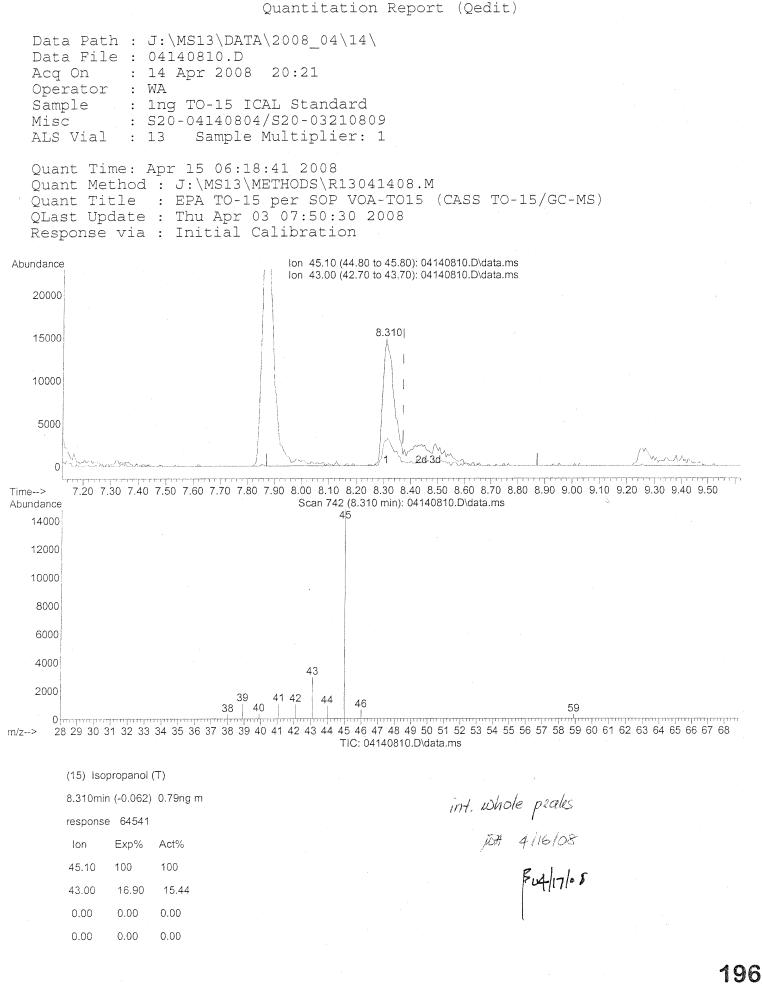
R13041408.M Tue Apr 15 06:19:06 2008

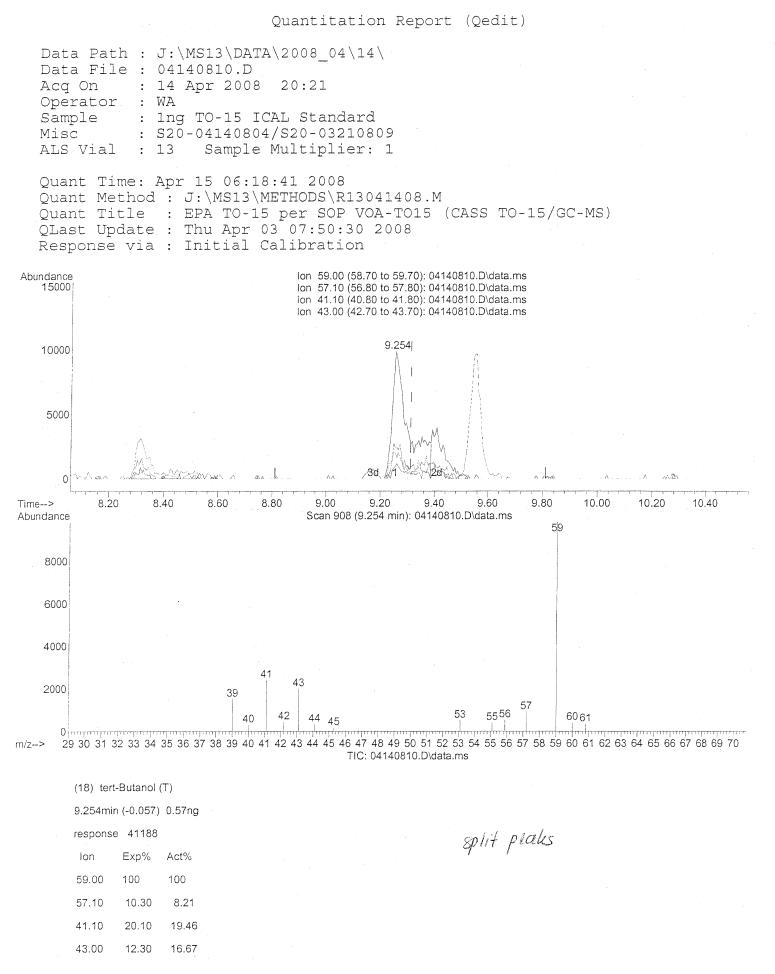
Page: 1



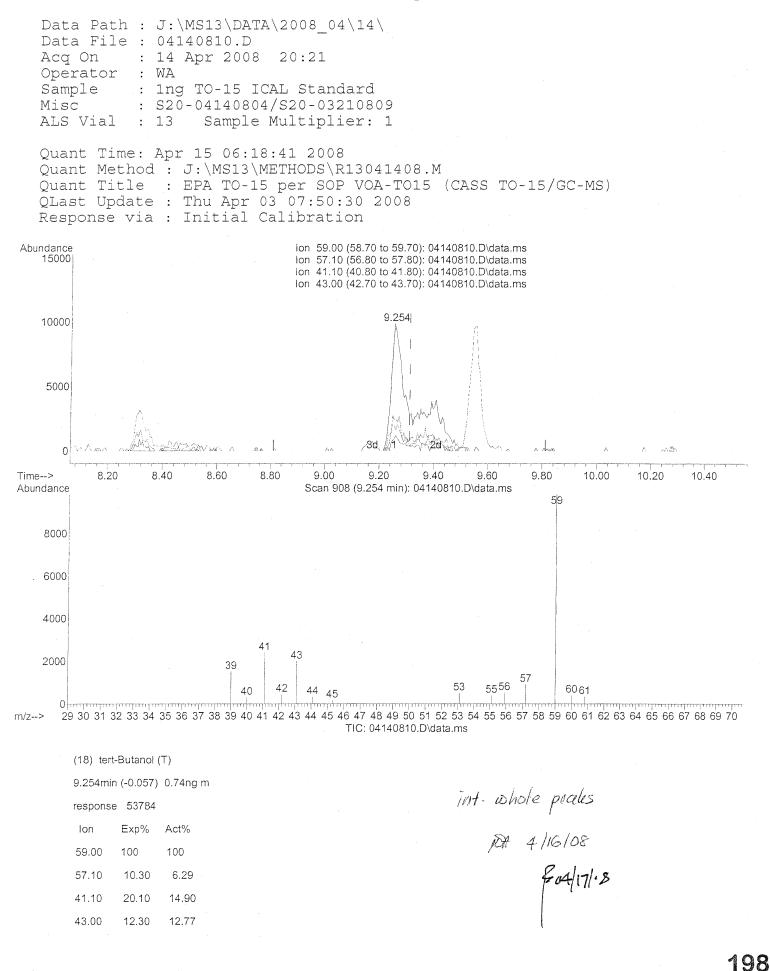
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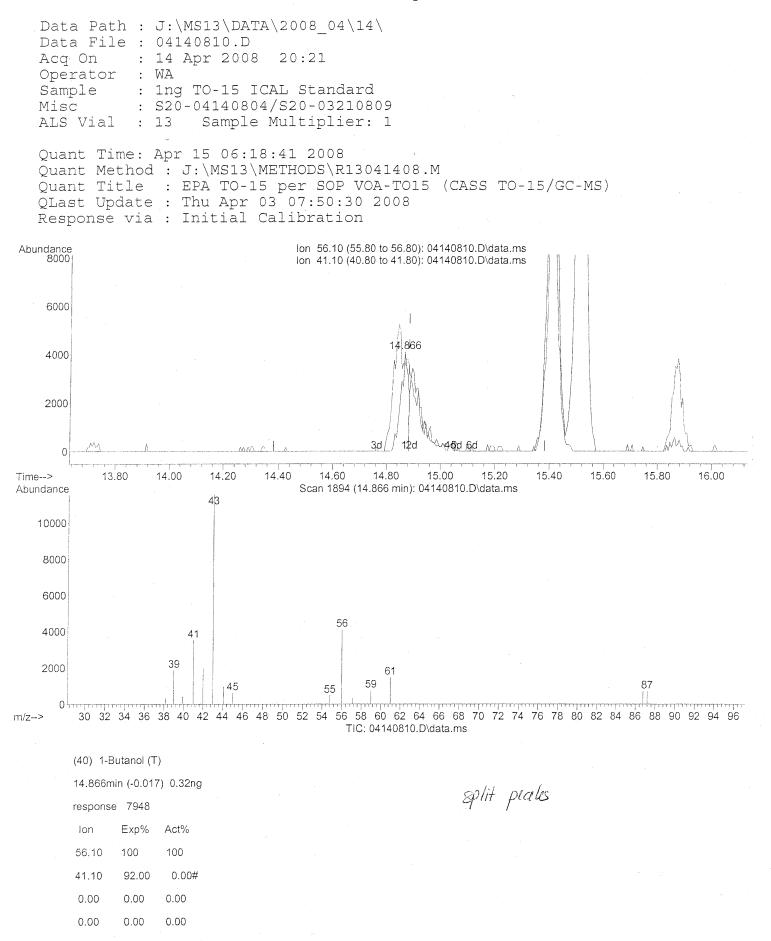


R13041408.M Tue Apr 15 06:19:36 2008

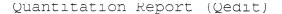


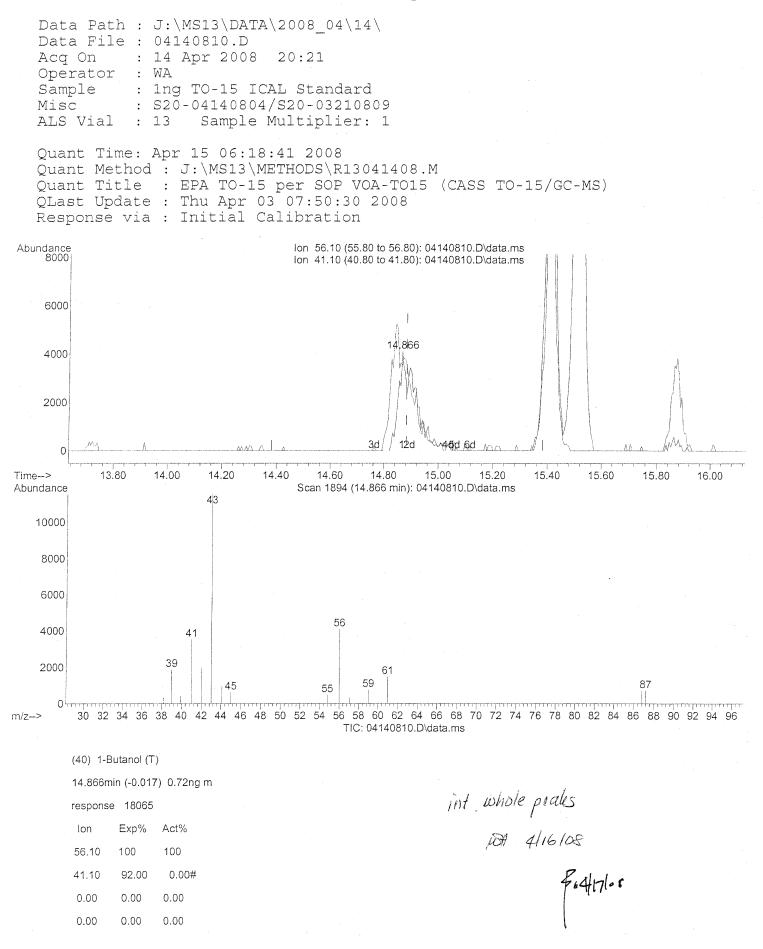
Page: 1



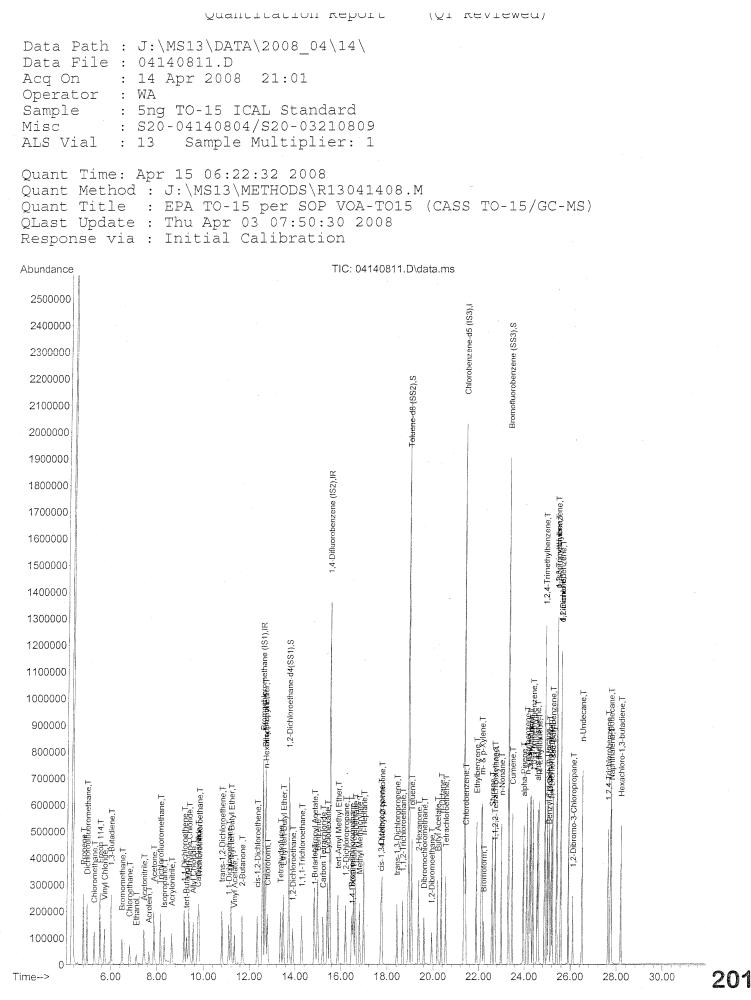


R13041408.M Tue Apr 15 06:20:15 2008





R13041408.M Tue Apr 15 06:20:20 2008



R13041408.M Tue Apr 15 06:24:08 2008

Page: 4

| Data Path : J:\MS13\DATA\2008_04 Data File : 04140811.D Acq On : 14 Apr 2008 21:01 Operator : WA Sample : 5ng TO-15 ICAL Stand Misc : S20-04140804/S20-032 ALS Vial : 13 Sample Multipli | ard 10809 er: 1 | | | diφ, | | |
|--|--|---|---|---|--|--|
| Quant Time: Apr 15 06:22:32 2008 Quant Method : J:\MS13\METHODS\R Quant Title : EPA TO-15 per SOP QLast Update : Thu Apr 03 07:50: Response via : Initial Calibrati | 1304140 VOA-TO 30 2008 | 15 (C | ASS TO-15/0 | GC-MS) | | |
| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
| 1) Bromochloromethane (IS1) 37) 1,4-Difluorobenzene (IS2) 56) Chlorobenzene-d5 (IS3) | 12.58 15.51 21.35 | 130 114 82 | 313584 1406515 715799 | 25.000 25.000 25.000 | ng ng ng | -0.02 -0.02 0.00 |
| System Monitoring Compounds 33) 1,2-Dichloroethane-d4(Spiked Amount 25.000 57) Toluene-d8 (SS2) | 13.73 18.93 | 65 98 | 620740 Recove 1613556 | 21.197 ery = 27.961 | ng 84 ng | -0.02 .80% -0.01 |
| Spiked Amount 25.000 73) Bromofluorobenzene (SS3) Spiked Amount 25.000 | | | 563169 | | ng | 0.00 |
| <pre>Target Compounds 2) Propene 3) Dichlorodifluoromethane 4) Chloromethane 5) Freon 114 6) Vinyl Chloride 7) 1,3-Butadiene 8) Bromomethane 9) Chloroethane 10) Ethanol 11) Acetonitrile 12) Acrolein 13) Acetone 14) Trichlorofluoromethane 15) Isopropanol 16) Acrylonitrile 17) 1,1-Dichloroethene 18) tert-Butanol 19) Methylene Chloride 20) Allyl Chloride 21) Trichlorotrifluoroethane 22) Carbon Disulfide 23) trans-1,2-Dichloroethene 24) 1,1-Dichloroethane 25) Methyl tert-Butyl Ether 26) Vinyl Acetate 27) 2-Butanone 28) cis-1,2-Dichloroethene 29) Diisopropyl Ether 30) Ethyl Acetate 31) n-Hexane</pre> | 4.96 5.28 5.53 6.00 6.49 6.82 7.11 7.43 7.65 7.87 | 85 50 135 62 54 94 64 45 41 56 58 | 175637 139777 86439 73739 78143m 219546 57551 100677 | 3.532 3.391 4.380 3.632 3.641 3.541 3.114 3.361 3.729 3.184 4.209 | ng ng ng ng ng ng ng ng ng ng | 96 # 74 99 92 96 99 # 59 |

DA 4/16/08

Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140811.D Acq On : 14 Apr 2008 21:01 Operator : WA Sample : 5ng TO-15 ICAL Standard Misc : S20-04140804/S20-03210809 ALS Vial : 13 Sample Multiplier: 1 Quant Time: Apr 15 06:22:32 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Thu Apr 03 07:50:30 2008 Response via : Initial Calibration

 Internal Scandards
 R.T. Qion Response Conc Units Dev(Min)

 32) Chloroform
 12.79
 B3
 173682
 4.201 ng
 99

 34) Tetrahydrofuran
 13.36
 72
 65400
 3.533 ng
 96

 35) Ethyl Tetr-Butyl Ether
 13.48
 87
 99161
 3.236 ng
 # 84

 35) I.p.Dichlorothane
 14.29
 97
 144607
 4.003 ng
 97

 38) I.j.l.-Trichlorothane
 14.29
 97
 144607
 4.003 ng
 # 22

 40) 1-Butanol
 14.85
 56
 65220
 3.530 ng
 # 49

 41.Benzane
 14.85
 78
 372533
 3.783 ng
 99

 42) Cerbon Tetrachloride
 15.21
 117
 12340
 4.010 ng
 # 78

 43) Cyclobexane
 16.42
 63
 105707
 3.828 ng
 92

 43) Locotane
 16.42
 57
 41519
 3.628 ng
 92

 45) 1.2-Dichloropropene
 17.73
 75
 144058
 3.648 ng
 100

 50) Methyl Methacrylate
 16.79
 100
 32918
 3.732 ng
 # 69

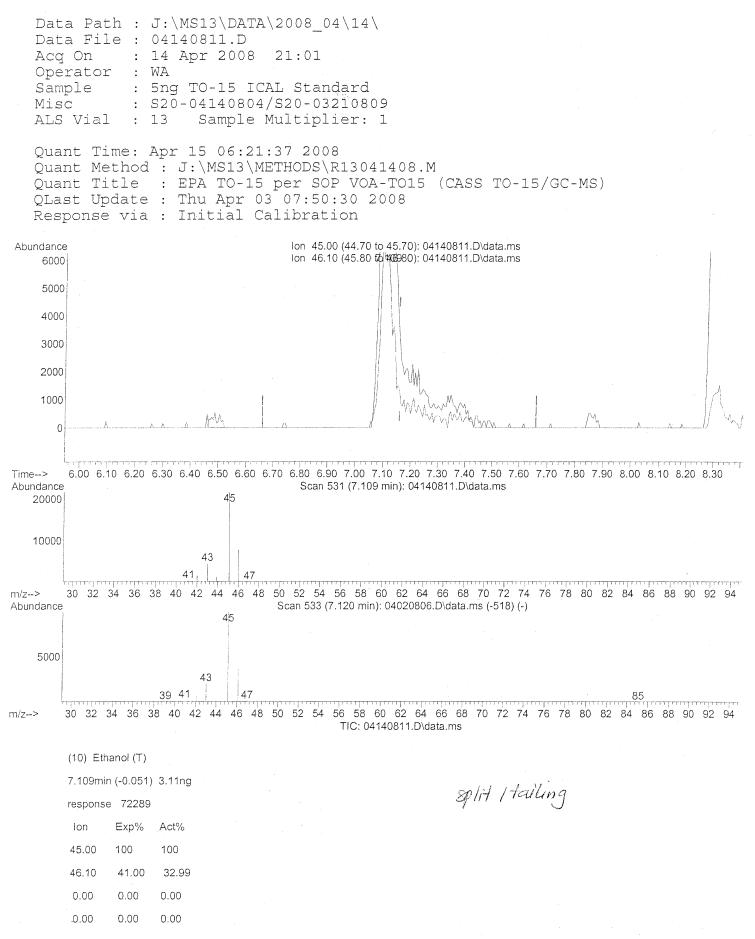
 51) n-Heptane
 Internal Standards R.T. QIon Response Conc Units Dev(Min)

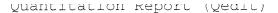
13041408.M Tue Apr 15 06:24:08 2008

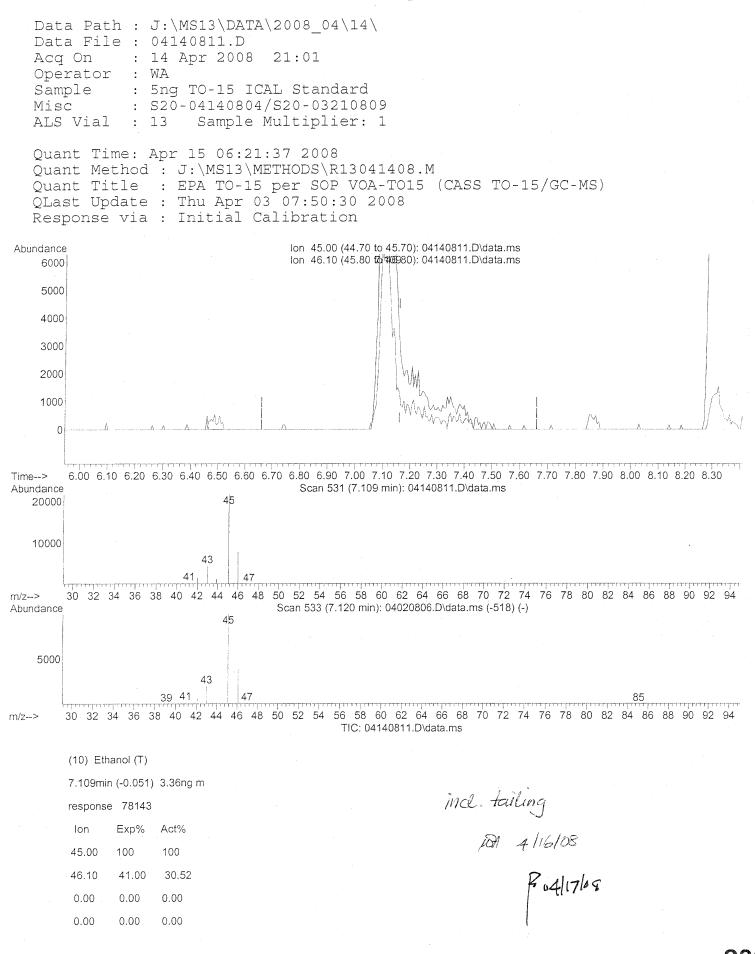
| Data Path : J:\MS13\DATA\2008_04 Data File : 04140811.D Acq On : 14 Apr 2008 21:01 Operator : WA Sample : 5ng TO-15 ICAL Standa Misc : S20-04140804/S20-0323 ALS Vial : 13 Sample Multiplie | ard 10809 | | | | |
|---|--|--|---|--|---|
| Quant Time: Apr 15 06:22:32 2008 Quant Method : J:\MS13\METHODS\R Quant Title : EPA TO-15 per SOP QLast Update : Thu Apr 03 07:50:3 Response via : Initial Calibratic | VOA-TO 30 2008 | | SS TO-15/G | C-MS) | |
| Internal Standards | R.T. | QIon | Response | Conc Units | Dev(Min) |
| <pre>82) 1,2,4-Trimethylbenzene 83) n-Decane 84) Benzyl Chloride 85) 1,3-Dichlorobenzene 86) 1,4-Dichlorobenzene 87) sec-Butylbenzene 88) p-Isopropyltoluene 89) 1,2,3-Trimethylbenzene 90) 1,2-Dichlorobenzene 91) d-Limonene 92) 1,2-Dibromo-3-Chloropr 93) n-Undecane 94) 1 2 4-Trichlorobenzene</pre> | 24.61 24.88 24.98 25.05 25.08 25.16 25.21 25.40 25.58 25.58 26.11 26.50 27.62 27.74 | 105 57 91 146 146 105 119 105 146 68 157 57 180 128 57 | 410666 384094 218596 292347 207129 210036 480271 423486 369517 202226 162016 62082 233672 144138 474438 230491 | 4.654 ng 4.271 ng 4.529 ng 5.144 ng 5.286 ng 4.852 ng 4.852 ng 4.630 ng 4.630 ng 4.888 ng 4.151 ng 5.448 ng 4.303 ng 5.172 ng 5.706 ng | 97 94 87 92 99 100 95 91 95 |

(#) = qualifier out of range (m) = manual integration (+) = signals summed



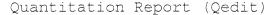


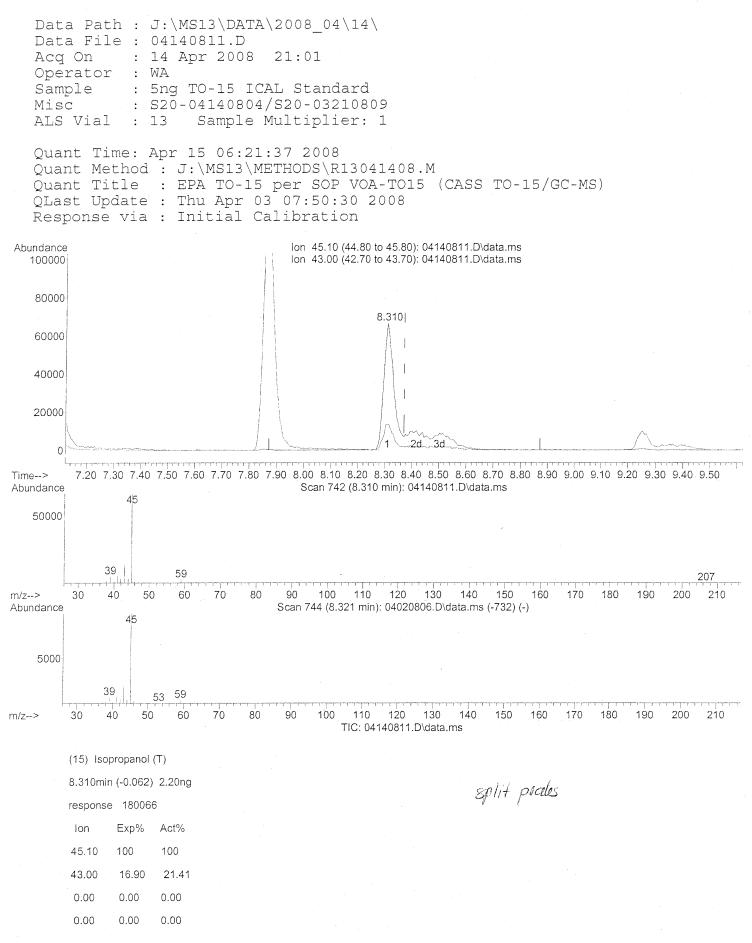




R13041408.M Tue Apr 15 06:22:07 2008

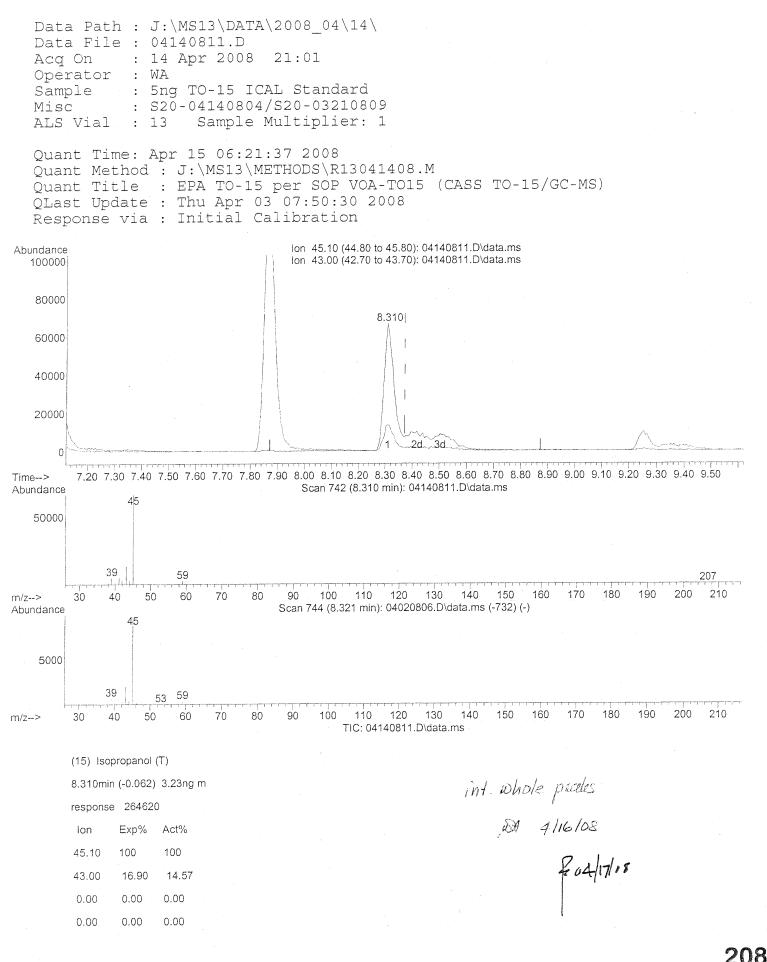
Page: 1

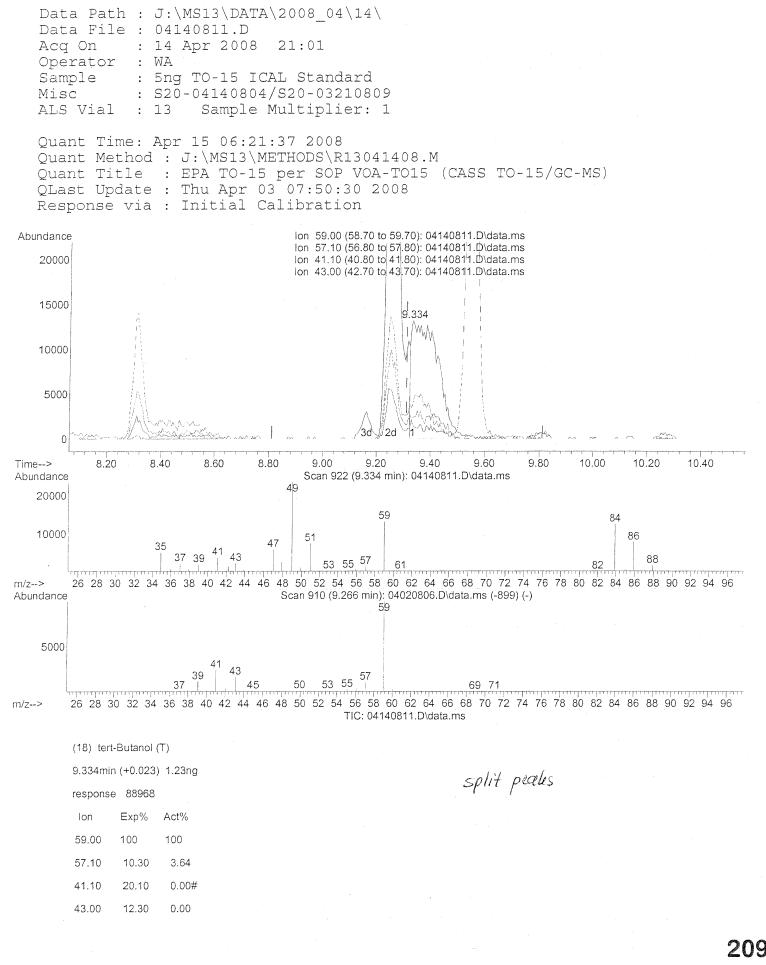


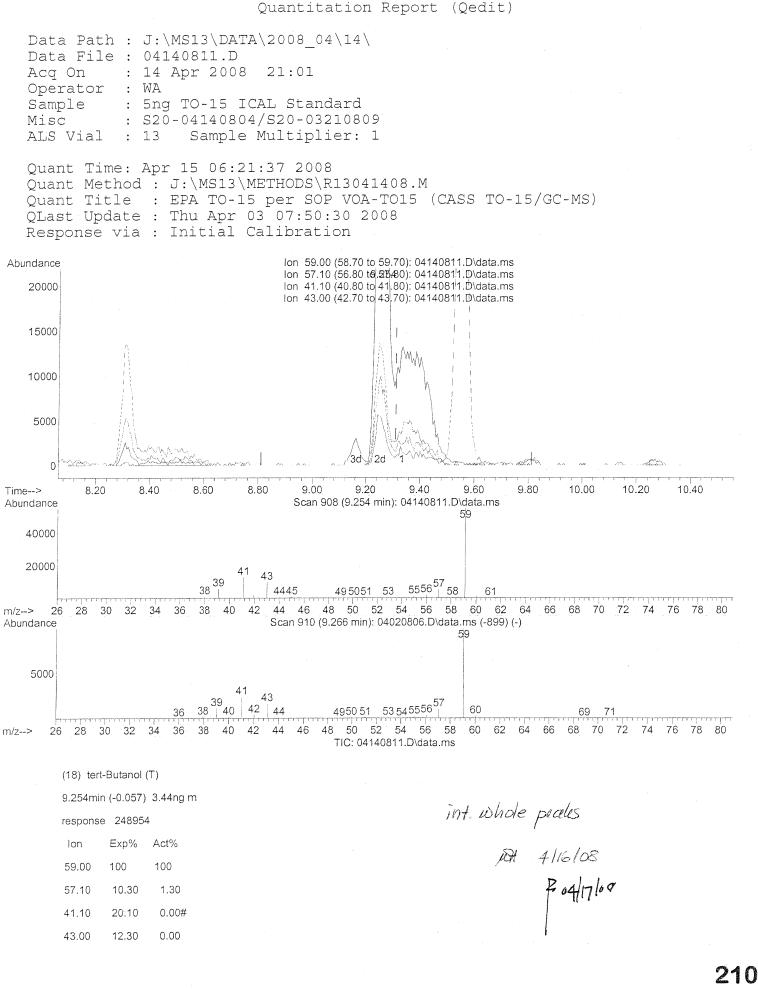


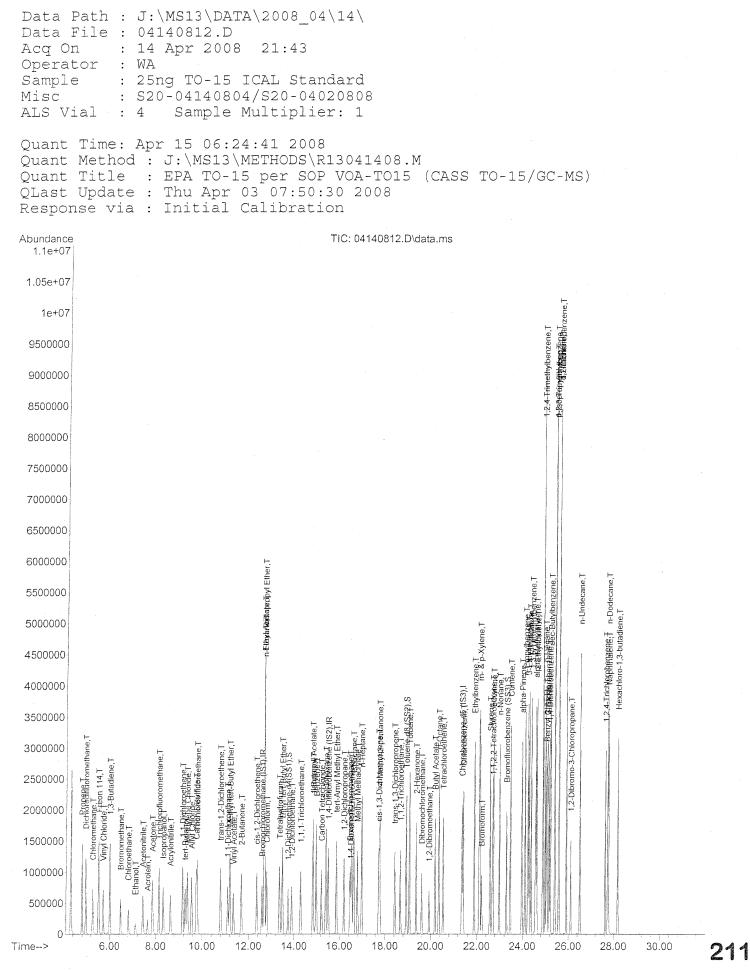
R13041408.M Tue Apr 15 06:22:16 2008











R13041408.M Tue Apr 15 06:25:47 2008

Quantitation keport (QT keviewed) Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140812.D Acq On : 14 Apr 2008 21:43 Operator : WA Sample : 25ng TO-15 ICAL Standard Misc : S20-04140804/S20-04020808 ALS Vial : 4 Sample Multiplier: 1 Quant Time: Apr 15 06:24:41 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Thu Apr 03 07:50:30 2008 Response via : Initial Calibration Internal Standards R.T. QIon Response Conc Units Dev(Min) 1) Bromochloromethane (IS1)12.6013033207025.000 ng-0.0137) 1,4-Difluorobenzene (IS2)15.52114146703225.000 ng-0.0156) Chlorobenzene-d5 (IS3)21.358276215225.000 ng0.00

 System Monitoring Compounds

 33) 1,2-Dichloroethane-d4(...
 13.73
 65
 674919
 21.764 ng
 -0.02

 Spiked Amount
 25.000
 Recovery
 =
 87.04%

 57) Toluene-d8 (SS2)
 18.93
 98
 1696875
 27.616 ng
 -0.01

 Spiked Amount
 25.000
 Recovery
 =
 110.48%

 73) Bromofluorobenzene (SS3)
 23.29
 174
 594489
 30.286 ng
 0.00

 Spiked Amount
 25.000
 Recovery
 =
 121.16%

 System Monitoring Compounds

 Spiked Amount
 25.000
 Recovery - Harmet

 Target Compounds
 Qvalue

 2) Propene
 4.79
 42
 672960
 17.105
 ng
 90

 3) Dichlorodifluoromethane
 5.27
 50
 996447
 17.815
 ng
 97

 5) Freen 114
 5.52
 135
 601060
 22.233
 ng
 100

 6) Vinyl Chloride
 5.72
 62
 926947
 18.099
 ng
 97

 7) 1,3-Butadiene
 6.00
 54
 801354
 19.711
 17
 75

 8) Bromomethane
 6.48
 94
 482161
 18.653
 ng
 98

 9) Chloroethane
 6.82
 64
 407498
 16.251
 ng
 95

 10) Ethanol
 7.13
 45
 369567m
 15.011
 ng
 91

 11) Acetonitrile
 7.64
 41
 1063940
 17.066
 ng
 93

 13) Acetone
 7.87
 58
 476809
 18.825
 ng
 93

 14) Trichlorofluoromethane
 8.14
 101
 9819231
 18.268
 93
 93
 Recovery = 121.16% Spiked Amount 25.000 ⁹³212

13041408.M Tue Apr 15 06:25:46 2008

Jost 4/16/08

| Quantitation | keport | (| UT KEVLEWEC | l) | | | |
|--|--|---|---|--|---|--|---|
| Data Path : J:\MS13\DATA\2008_04 Data File : 04140812.D Acq On : 14 Apr 2008 21:43 Operator : WA Sample : 25ng TO-15 ICAL Star Misc : S20-04140804/S20-040 ALS Vial : 4 Sample Multiplie | ndard 020808 | | | | | | |
| Quant Time: Apr 15 06:24:41 2008 Quant Method : J:\MS13\METHODS\F Quant Title : EPA TO-15 per SOF QLast Update : Thu Apr 03 07:50: Response via : Initial Calibrati | R1304140 P VOA-TO :30 2008 | 15 (C | CASS TO-15/ | GC-MS) | | | |
| Internal Standards | R.T. | QION | Response | Conc | Units | Dev | (Min) |
| <pre>32) Chloroform 34) Tetrahydrofuran 35) Ethyl tert-Butyl Ether 36) 1,2-Dichloroethane 38) 1,1,1-Trichloroethane 39) Isopropyl Acetate 40) 1-Butanol 41) Benzene 42) Carbon Tetrachloride 43) Cyclohexane 44) tert-Amyl Methyl Ether 45) 1,2-Dichloropropane 46) Bromodichloromethane 47) Trichloroethene 48) 1,4-Dioxane 49) Isooctane 50) Methyl Methacrylate 51) n-Heptane 52) cis-1,3-Dichloropropene</pre> | 12.80 13.36 13.49 13.90 14.30 14.84 14.84 14.99 15.22 15.41 15.87 16.20 16.46 16.54 16.49 16.62 16.80 16.98 17.73 17.77 | 83 72 87 62 97 61 56 78 117 84 73 63 83 130 88 57 100 71 75 58 | 956625 360780 556742 824015 830170 345329 447826 2068389 718394 767539 1448190 576719 756240 525239 396979 2351708 190312 578179 823511 577342 | 21.851 18.405 17.155 18.894 22.032 20.309 17.785 20.085 22.376 20.298 19.521 20.028 22.466 24.045 22.001 19.710 20.691 20.955 19.426 20.749 | ng ng ng ng ng ng ng ng ng ng ng ng ng n | ++++++++++++++++++++++++++++++++++++++ | 98 98 99 98 30 47 98 99 81 91 98 100 98 84 70 68 82 98 87 |
| <pre>55) 1,1,2-Trichloroethane 58) Toluene 59) 2-Hexanone 60) Dibromochloromethane 61) 1,2-Dibromoethane 62) Butyl Acetate 63) n-Octane 64) Tetrachloroethene 65) Chlorobenzene 66) Ethylbenzene 67) m- & p-Xylene 68) Bromoform 69) Styrene 70) o-Xylene 71) n-Nonane 72) 1,1,2,2-Tetrachloroethane</pre> | 18.67 19.07 19.37 19.61 19.94 20.19 20.36 20.55 21.42 21.89 22.13 22.21 22.57 22.57 22.57 22.98 22.69 23.47 23.97 24.10 24.23 24.28 | 97 91 43 129 107 43 57 166 112 91 173 104 91 173 104 93 105 105 | 481442 2224726 1719849 583893 559507 1805701 515467 547110 1376096 2587658 4184617 497365 1508280 2083003 1425405 1013401 2395512 1244786 3178462 2444332 2502986 | 21.367 24.602 24.112 28.073 23.950 25.220 22.585 26.534 25.704 25.169 61.330 33.995 25.951 28.561 23.621 28.884 26.455 24.972 26.384 24.967 | ng ng ng ng ng ng ng ng ng ng ng ng ng n | # | 92 96 79 98 99 83 97 99 100 91 89 98 95 90 78 97 98 97 98 90 |

13041408.M Tue Apr 15 06:25:46 2008

101 4/16/08

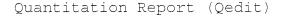
Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140812.D Acg On : 14 Apr 2008 21:43 Operator : WA
 Sample
 : 25ng TO-15 ICAL Standard

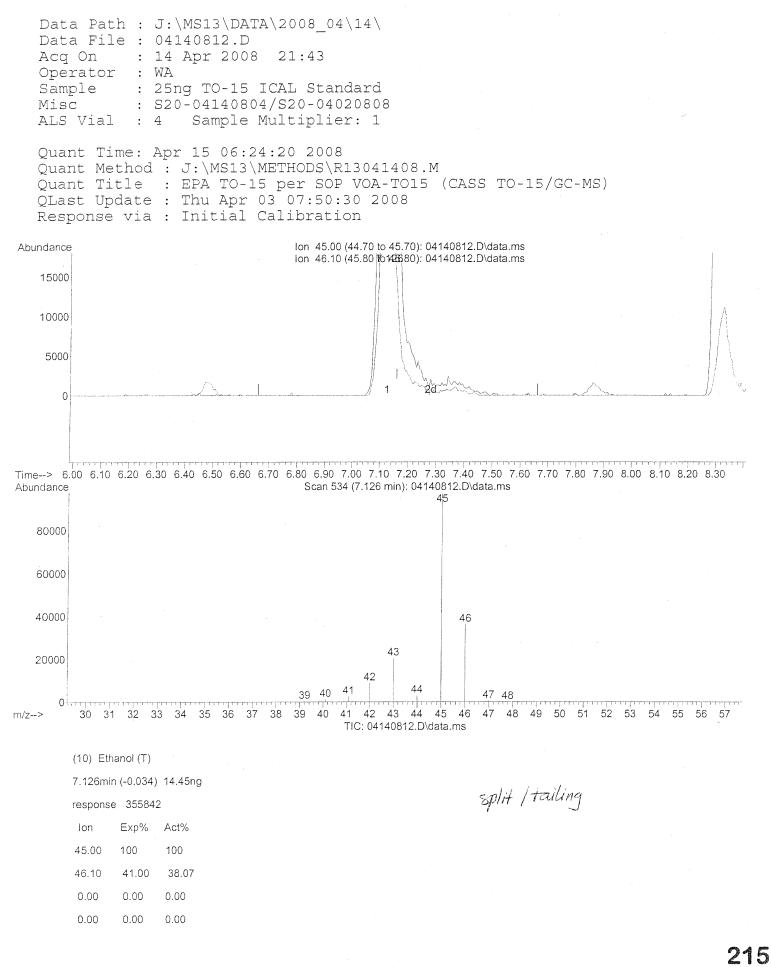
 Misc
 : S20-04140804/S20-04020808
 ALS Vial : 4 Sample Multiplier: 1 Quant Time: Apr 15 06:24:41 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Thu Apr 03 07:50:30 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards 11111111111111111111111180) alpha-Methylstyrene24.56118108797326.108 ng9681) 2-Ethyltoluene24.61105241090124.872 ng9782) 1,2,4-Trimethylbenzene24.88105250886928.548 ng9683) n-Decane24.9957127778523.445 ng8884) Benzyl Chloride25.0591182564226.564 ng9485) 1,3-Dichlorobenzene25.08146122970028.682 ng10086) 1,4-Dichlorobenzene25.16146122531328.963 ng9987) sec-Butylbenzene25.21105282452926.797 ng9588) p-Isopropyltoluene25.40119287209534.115 ng9189) 1,2,3-Trimethylbenzene25.58146130668829.663 ng10091) d-Limonene25.5868107749025.929 ng9492) 1,2-Dibromo-3-Chloropr...26.1115737161630.629 ng#94) 1,2,4-Trichlorobenzene27.6318084660128.533 ng9595) Naphthalene27.77128272139930.741 ng9896) n-Dodecane27.7457135034823.984 ng8797) Hexachloro-1,3-butadiene28.1922554095530.728 ng100 _____ _____

(#) = qualifier out of range (m) = manual integration (+) = signals summed

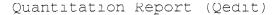
13041408.M Tue Apr 15 06:25:46 2008

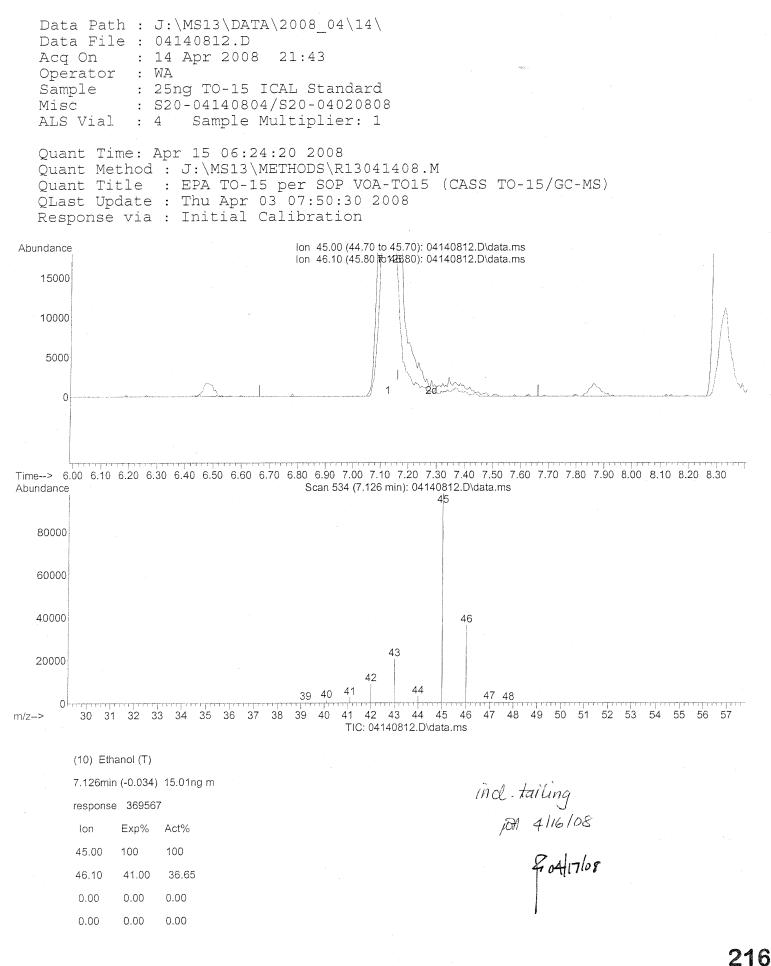
PA 4/16/08

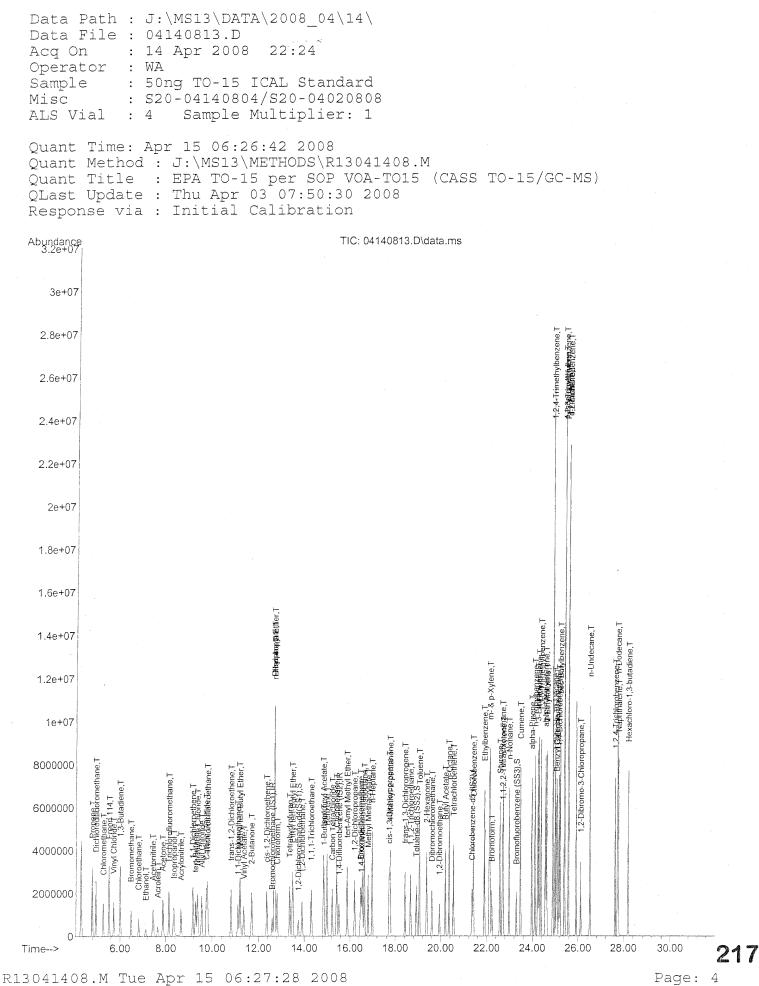




R13041408.M Tue Apr 15 06:24:38 2008







| Data Path : J:\MS13\DATA\2008_04\14\ Data File : 04140813.D Acq On : 14 Apr 2008 22:24 Operator : WA Sample : 50ng TO-15 ICAL Standard Misc : S20-04140804/S20-04020808 ALS Vial : 4 Sample Multiplier: 1 Quant Time: Apr 15 06:26:42 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Thu Apr 03 07:50:30 2008 Response via : Initial Calibration | | | | | | | | | |
|---|---|--|---|--|---|---|--|--|--|
| Internal Standards | | QIon | Response | Conc | Units | Dev(Min) | | | |
| 1) Bromochloromethane (IS1) 37) 1,4-Difluorobenzene (IS2) 56) Chlorobenzene-d5 (IS3) | 12.60 | 130 | 359135 | 25.000 |) na | 0.00 | | | |
| System Monitoring Compounds 33) 1,2-Dichloroethane-d4(Spiked Amount 25.000 57) Toluene-d8 (SS2) Spiked Amount 25.000 73) Bromofluorobenzene (SS3) Spiked Amount 25.000 | 18.93 | 98 | Recov 1797014 Recov | ery = 27.224 ery = 29.723 | 84 ng 108 ng | .28% 0.00 .88% 0.00 | | | |
| 13) Acetone 14) Trichlorofluoromethane 15) Isopropanol 16) Acrylonitrile 17) 1,1-Dichloroethene 18) tert-Butanol 19) Methylene Chloride | 4.95 5.28 5.53 5.72 6.00 6.48 6.82 7.14 7.45 7.65 7.65 7.87 8.14 8.34 9.28 9.28 9.28 9.281 9.281 9.76 10.81 11.20 11.20 11.68 12.36 12.70 12.70 | 85 535 54 44 58 153 59 41 156 53 62 17 63 76 217 61 51 | 2106258 1321916 1980229 1778388 1041693 855335 802216m 2274356 658477 989790 2079897 3077080m 1587115 1039085 2885669 1106480 1804544 947778 4089706 1829530 2089949 3425202 208979 772056 | 35.317 34.819 45.213 35.751 40.447 37.263 31.541 30.129 33.733 31.805 36.133 36.056 32.835 36.268 38.154 34.776 36.752 40.079 42.638 32.370 36.213 35.955 35.500 31.862 34.783 36.226 37.093 49.781 | ng ng ng ng ng ng ng ng ng ng ng ng ng n | 99 97 99 96 76 100 96 99 99 99 99 99 99 99 97 98 97 98 97 98 97 98 91 96 92 95 88 91 96 92 95 88 91 | | | |

.13041408.M Tue Apr 15 06:27:27 2008

jan 4/16/08

²218

Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140813.D Acq On : 14 Apr 2008 22:24 Operator : WA Sample : 50ng TO-15 ICAL Standard Misc : S20-04140804/S20-04020808 ALS Vial : 4 Sample Multiplier: 1 Quant Time: Apr 15 06:26:42 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Thu Apr 03 07:50:30 2008 Response via : Initial Calibration

 Internal Standards
 R.T. Qion Response Conc Units Dev(Min)

 32) Chioroform
 12.81
 83
 1994360
 42.122 ng
 98

 33) Ethyl tert-Butyl Ether
 13.36
 72
 764002
 36.038 ng
 98

 34) Tetrahydrofuran
 13.36
 72
 764002
 36.038 ng
 98

 35) Ethyl tert-Butyl Ether
 13.49
 87
 1193455
 34.004 ng
 #
 55

 36) 1.1.1-Trichloroethane
 14.30
 97
 1742008
 42.924 ng
 97

 38) 1.1.1-Trichloroethane
 14.84
 61
 731723
 39.953 ng
 #
 28

 40) 1-Butanol
 14.86
 56
 968057
 35.694 ng
 #
 62

 41 benzene
 15.67
 73
 3119689
 39.043 ng
 99
 91

 42) Carbon Tetrachloride
 15.67
 73
 3119689
 39.043 ng
 99
 91

 43) Locothane
 16.63
 57
 5039242
 39.214 ng
 70
 50

 50) Methyl Methatrylate
 16.61
 100
 403557
 41.333 ng
 #
 82

 51) Locothane
 19. Internal Standards R.T. QIon Response Conc Units Dev(Min)

13041408.M Tue Apr 15 06:27:27 2008 Page: 2

Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140813.D Acq On : 14 Apr 2008 22:24 Operator : WA Sample : 50ng TO-15 ICAL Standard Misc : S20-04140804/S20-04020808 ALS Vial : 4 Sample Multiplier: 1 Quant Time: Apr 15 06:26:42 2008 Ouant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Thu Apr 03 07:50:30 2008 Response via : Initial Calibration Internal Standards R.T. QIon Response Conc Units Dev(Min) 80) alpha-Methylstyrene24.57118249436855.717ng9681) 2-Ethyltoluene24.62105541081251.961ng9882) 1,2,4-Trimethylbenzene24.89105661344070.050ng9883) n-Decane24.9957293649950.154ng8884) Benzyl Chloride25.0591424641857.516ng9585) 1,3-Dichlorobenzene25.08146285913562.075ng10086) 1,4-Dichlorobenzene25.16146281432061.922ng9987) sec-Butylbenzene25.22105639840256.506ng9688) p-Isopropyltoluene25.41119750508582.980ng9189) 1,2,3-Trimethylbenzene25.58146342752672.428ng9990) 1,2-Dichlorobenzene25.5868273862061.346ng9592) 1,2-Dibromo-3-Chloropr...26.1115783337663.938ng#62

 90) 1,2-Dichlorobenzene
 25.58
 146
 3427526
 72.428
 1g
 99

 91) d-Limonene
 25.58
 68
 2738620
 61.346
 ng
 95

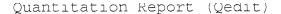
 92) 1,2-Dibromo-3-Chloropr...
 26.11
 157
 833376
 63.938
 ng
 #
 62

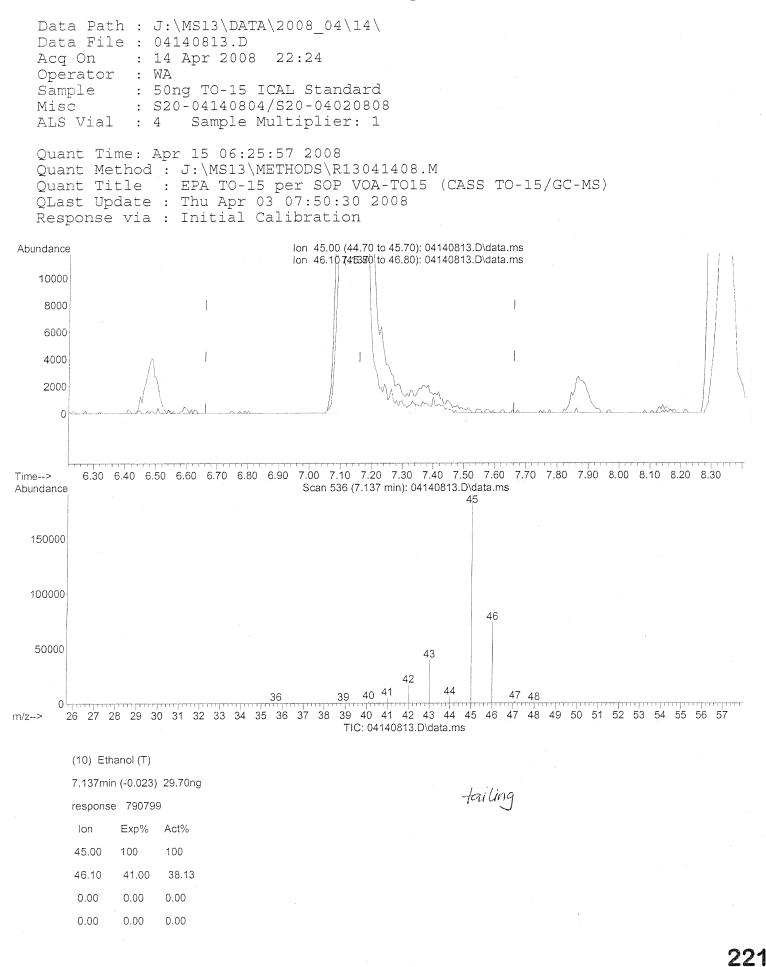
 93) n-Undecane26.5057314646550.655 ng94) 1,2,4-Trichlorobenzene27.63180201784663.304 ng95) Naphthalene27.78128617126064.890 ng96) n-Dodecane27.7457317722552.528 ng97) Hexachloro-1,3-butadiene28.19225131524169.543 ng 87 96 99 86 100 _____

(#) = gualifier out of range (m) = manual integration (+) = signals summed

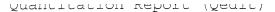
13041408.M Tue Apr 15 06:27:27 2008

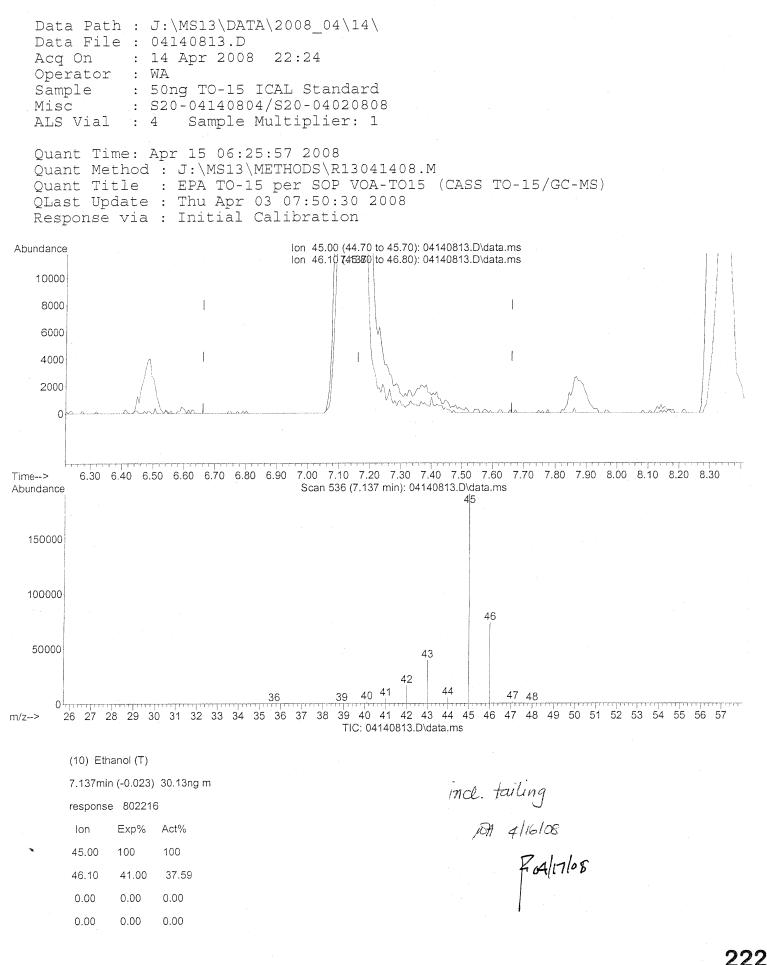
184 4/16/08

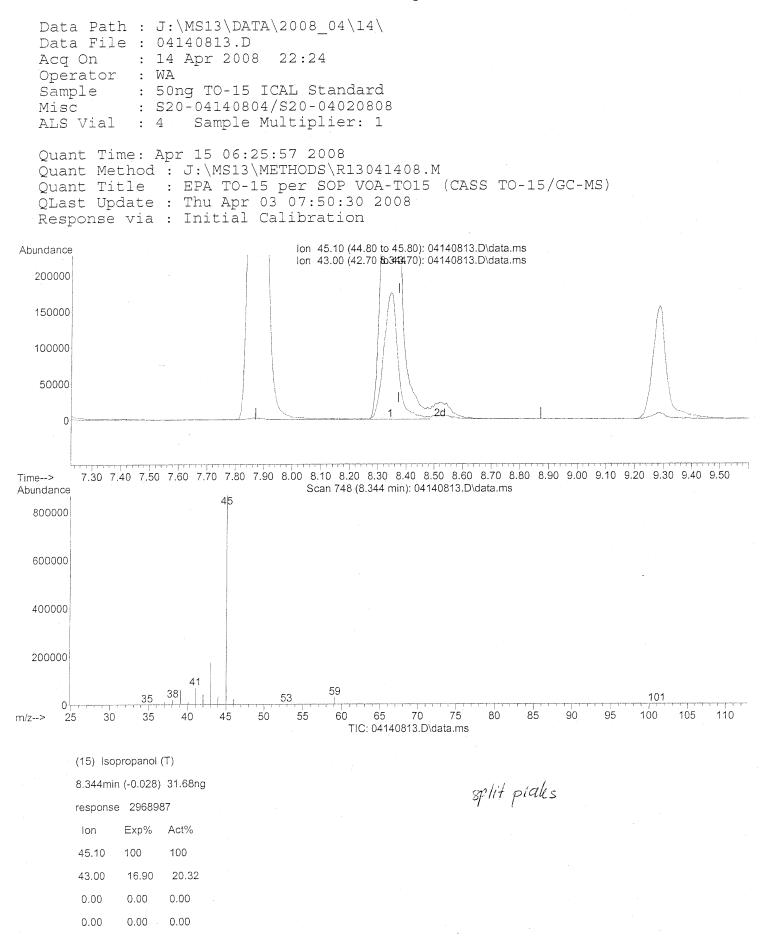


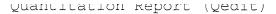


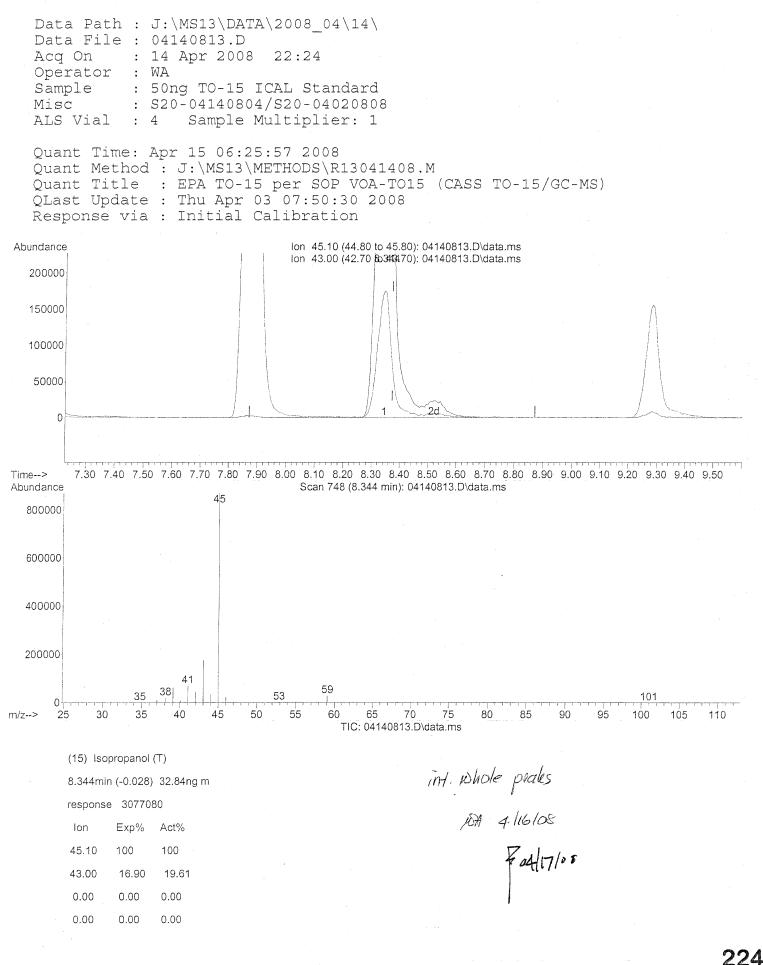
R13041408.M Tue Apr 15 06:26:20 2008

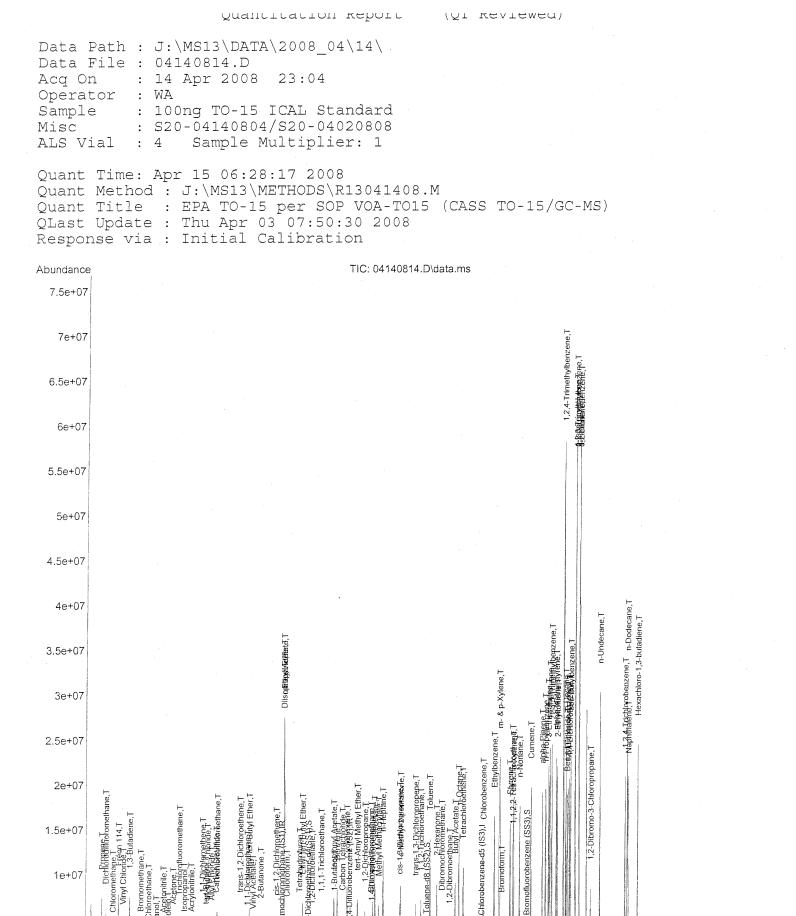


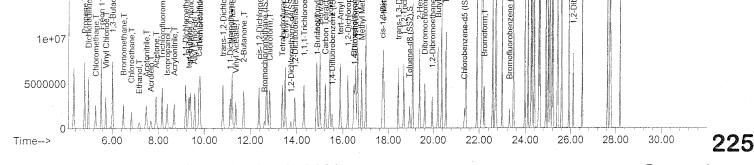












pon

R13041408.M Tue Apr 15 06:28:52 2008

1e+07

| Quantitation | Report | (Ç |)T Reviewed | d) | | | |
|--|--|---|---|--|---|-------------|--|
| Data Path : J:\MS13\DATA\2008_04 Data File : 04140814.D Acq On : 14 Apr 2008 23:04 Operator : WA Sample : 100ng TO-15 ICAL Sta Misc : S20-04140804/S20-040 ALS Vial : 4 Sample Multiplie | · · · · | | | | | | |
| Quant Time: Apr 15 06:28:17 2008 Quant Method : J:\MS13\METHODS\R Quant Title : EPA TO-15 per SOP QLast Update : Thu Apr 03 07:50: Response via : Initial Calibrati | 1304140 VOA-TO 30 2008 | 15 (C | ASS TO-15/ | (GC-MS) | | | |
| Internal Standards | R.T. | QIon | Response | e Conc | Units | Dev () | Min) |
| 1) Bromochloromethane (IS1) 37) 1,4-Difluorobenzene (IS2) 56) Chlorobenzene-d5 (IS3) | 15.53 | 114 | 1673737 | 25.000 | ng | 0 | .00 |
| System Monitoring Compounds 33) 1,2-Dichloroethane-d4(Spiked Amount 25.000 57) Toluene-d8 (SS2) Spiked Amount 25.000 73) Bromofluorobenzene (SS3) Spiked Amount 25.000 | | | Recov 674346 | ery = | ng | .68% 0 | |
| <pre>7) 1,3-Butadiene 8) Bromomethane 9) Chloroethane 10) Ethanol 11) Acetonitrile 12) Acrolein 13) Acetone 14) Trichlorofluoromethane 15) Isopropanol 16) Acrylonitrile 17) 1,1-Dichloroethene 18) tert-Butanol 19) Methylene Chloride 20) Allyl Chloride 21) Trichlorotrifluoroethane 22) Carbon Disulfide 23) trans-1,2-Dichloroethene 24) 1,1-Dichloroethane 25) Methyl tert-Butyl Ether 26) Vinyl Acetate 27) 2-Butanone 28) cis-1,2-Dichloroethene 29) Diisopropyl Ether 30) Ethyl Acetate</pre> | $\begin{array}{c} 4.96\\ 5.28\\ 5.53\\ 5.73\\ 6.00\\ 6.49\\ 6.82\\ 7.17\\ 7.67\\ 7.89\\ 8.15\\ 8.38\\ 8.67\\ 9.31\\ 9.38\\ 9.57\\ 9.31\\ 9.38\\ 9.57\\ 9.81\\ 9.38\\ 9.57\\ 10.81\\ 11.12\\ 11.21\\ 11.37\\ 11.70\\ 12.37\\ 12.70\\ 12.71\\ \end{array}$ | 850 1362 544 4550 5524 445 50453 59411 613362 171 6787 6787 671 671 671 671 671 671 671 671 671 67 | 3729310 3097182 4423044 4005951 2351722 2051767 1718886m 4896836 1427932 2130633 4597540 6510791m 3418887 2288970 6157599 2423001 3992555 2198878 8961871 | 73.339 58.412 100.368 75.661 86.326 79.707 71.687 61.166 68.815 65.349 73.695 75.515 65.827 74.024 79.635 70.310 76.255 84.017 93.726 67.208 76.545 75.490 74.911 67.319 74.867 90.272 123.759 | ng ng ng ng ng ng ng ng ng ng ng ng ng n | # # # | 91 99 97 99 95 75 99 96 96 99 96 99 99 97 92 89 97 92 89 95 |

jBi 4/16/08

Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140814.D Acg On : 14 Apr 2008 23:04 Operator : WA Sample : 100ng TO-15 ICAL Standard Misc : S20-04140804/S20-04020808 ALS Vial : 4 Sample Multiplier: 1 Quant Time: Apr 15 06:28:17 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Thu Apr 03 07:50:30 2008 Response via : Initial Calibration

 Internal Standards
 R.T. Qion Response Conc Units Dev(Min)

 22: Chloroform
 12.63
 83
 4315526
 86.380 ng
 99

 36: Echyl tert-Butyl Ether
 13.57
 72
 1589023
 71.018 ng
 99

 36: 1.2-Dichloroethane
 14.31
 97
 3932525
 91.634 ng
 98

 37: 1.1.1-Trichloroethane
 14.31
 97
 3932525
 91.634 ng
 98

 38: 1.1.1-Trichloroethane
 14.31
 97
 393255
 91.634 ng
 97

 40: 1-Sutanol
 14.88
 56
 1562836
 85.713 ng
 # 39

 41: Beraree
 15.03
 17
 3543916
 96.75.0 ng
 98

 42: Cathon Tetrachloride
 15.23
 117
 3543916
 96.75.0 ng
 98

 43: Lychekare
 16.21
 63
 2710579
 82.504 ng
 96

 43: Lychekare
 16.21
 83
 3667281
 90.601 ng
 90

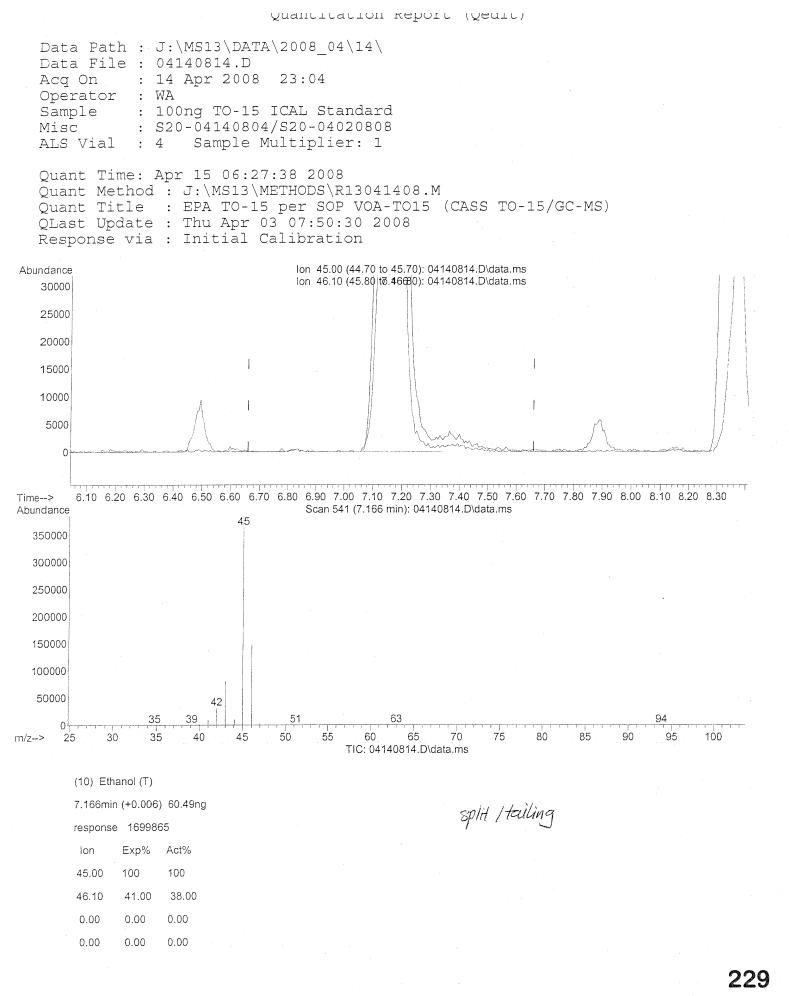
 44: tert-Amyl Methyl Ether
 15.88
 73<700525</td>
 82.635 ng
 91

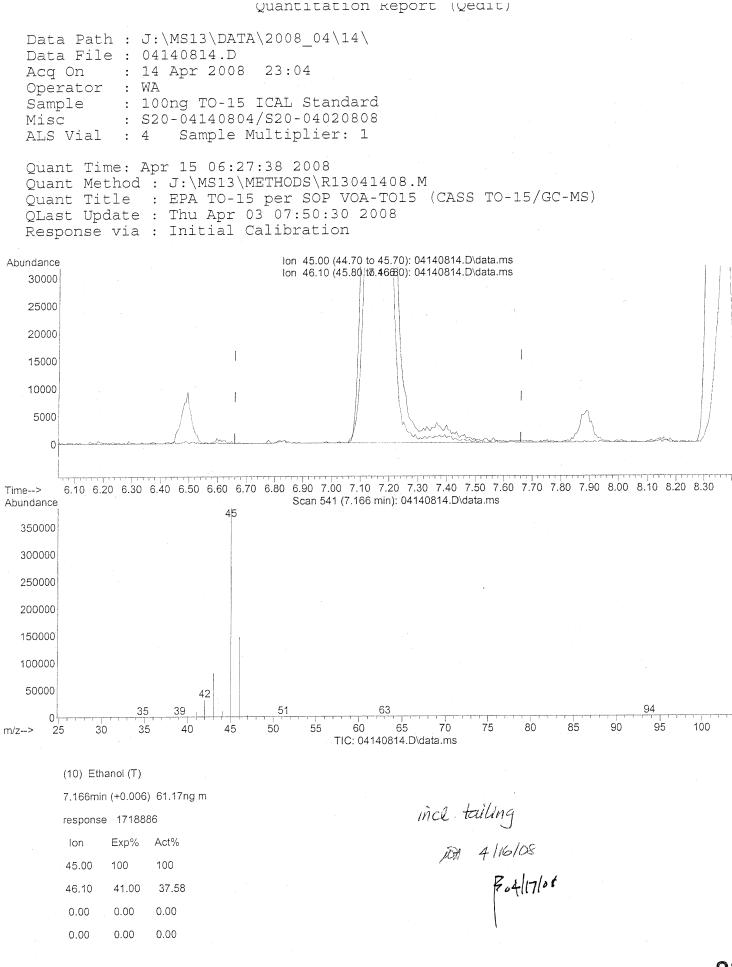
 45: Isooctane
 16.51
 82
 011 ng
 100

 47
 74
 75 4 Internal Standards R.T. QIon Response Conc Units Dev(Min)

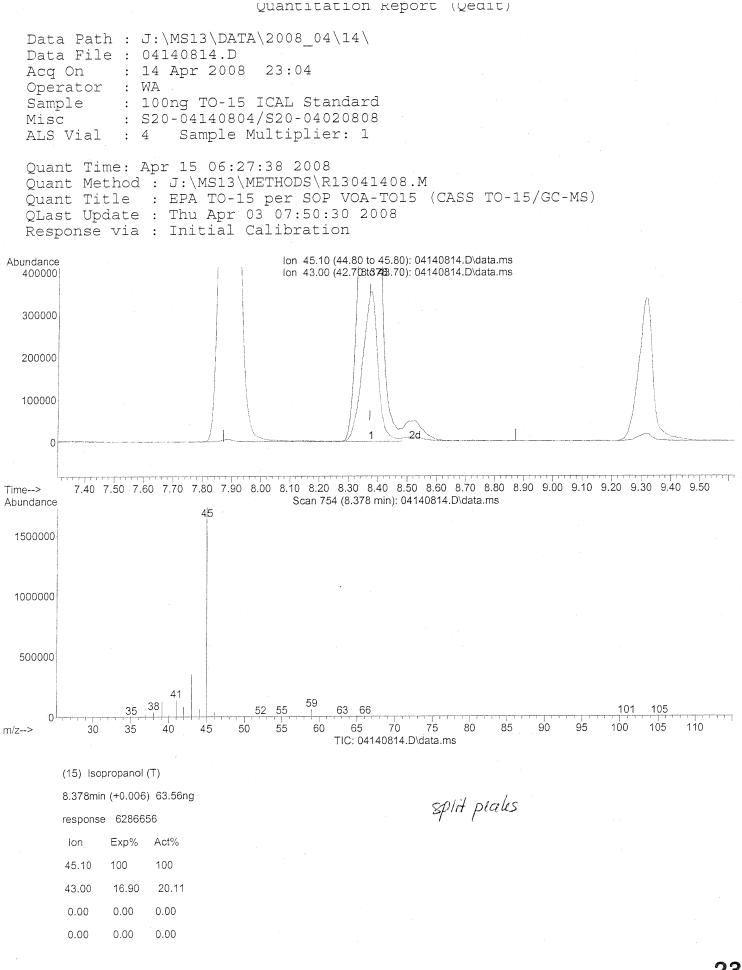
Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140814.D Acq On : 14 Apr 2008 23:04 Operator : WA Sample : 100ng TO-15 ICAL Standard Misc : S20-04140804/S20-04020808 ALS Vial : 4 Sample Multiplier: 1 Quant Time: Apr 15 06:28:17 2008 Ouant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Thu Apr 03 07:50:30 2008 Response via : Initial Calibration Internal Standards R.T. QIon Response Conc Units Dev(Min) 80) alpha-Methylstyrene24.571186563154137.807 ng9581) 2-Ethyltoluene24.6310512754976115.139 ng9682) 1,2,4-Trimethylbenzene24.9010515629562155.615 ng8783) n-Decane25.00577447447119.568 ng8584) Benzyl Chloride25.069110726388136.566 ng9985) 1,3-Dichlorobenzene25.091467724653157.649 ng9986) 1,4-Dichlorobenzene25.171467381200152.660 ng9987) sec-Butylbenzene25.2210514301505118.722 ng8988) p-Isopropyltoluene25.4111915925026165.511 ng#7089) 1,2,3-Trimethylbenzene25.591469624011191.165 ng9491) d-Limonene25.59687792659164.085 ng10092) 1,2-Dibromo-3-Chloropr...26.111572017145145.472 ng#6593) n-Undecane26.51578173344123.688 ng82 _____ 93) n-Undecane26.51578173344123.688 ng94) 1,2,4-Trichlorobenzene27.641805538389163.325 ng95) Naphthalene27.7812814124345139.605 ng96) n-Dodecane27.74578456134131.415 ng 82 95 95 80 97) Hexachloro-1,3-butadiene 28.19 225 3915632 194.616 ng 99 _____

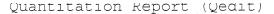
(#) = qualifier out of range (m) = manual integration (+) = signals summed

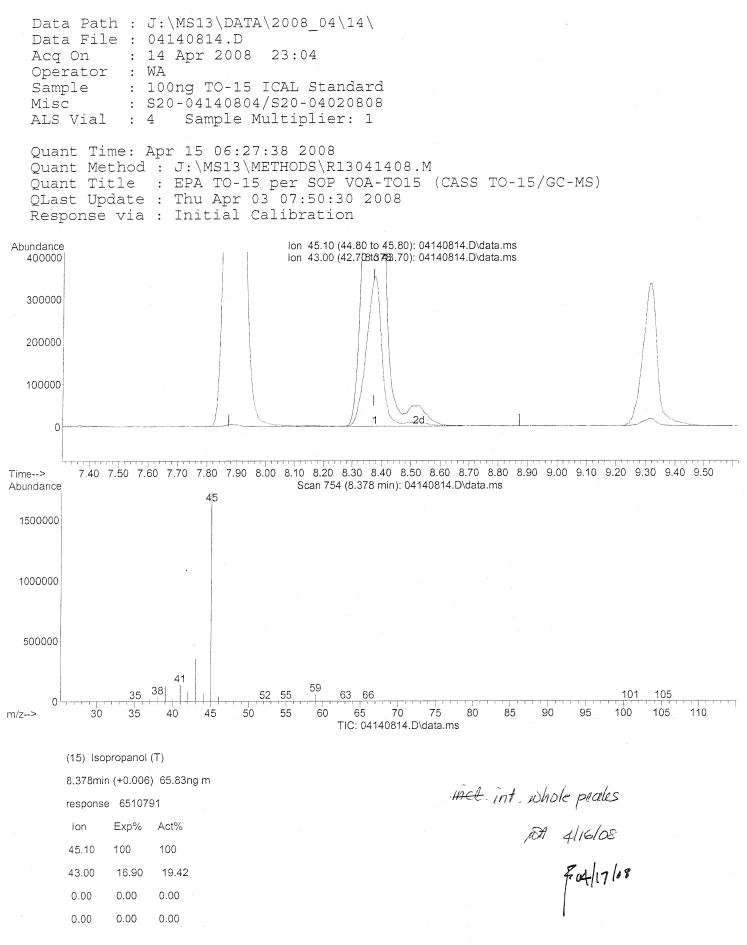


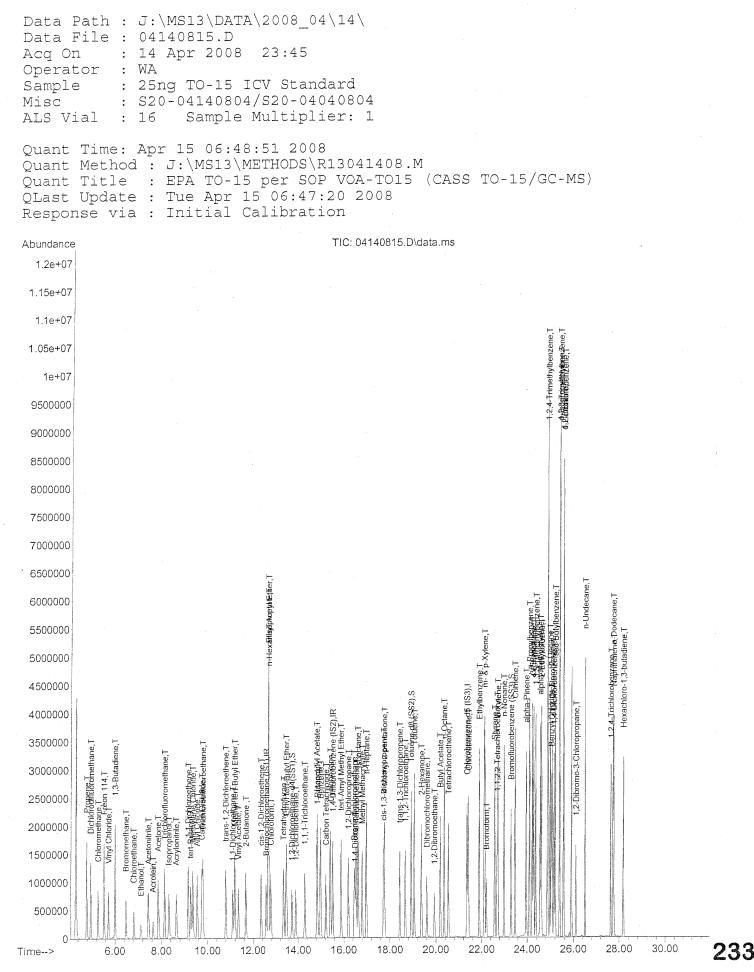


R13041408.M Tue Apr 15 06:28:05 2008









R13041408.M Tue Apr 15 06:49:40 2008

| Data Path : J:\MS13\DATA\2008_04 Data File : 04140815.D Acq On : 14 Apr 2008 23:45 Operator : WA Sample : 25ng TO-15 ICV Stand Misc : S20-04140804/S20-040 ALS Vial : 16 Sample Multiplie | | | | | | |
|--|--|---|--|--|--|--|
| Quant Time: Apr 15 06:48:51 2008 Quant Method : J:\MS13\METHODS\R Quant Title : EPA TO-15 per SOP QLast Update : Tue Apr 15 06:47:2 Response via : Initial Calibratio | 1304140 VOA-TO 20 2008 | 8.M 15 (C | ASS TO-15/ | GC-MS) | | |
| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
| 1) Bromochloromethane (IS1) 37) 1,4-Difluorobenzene (IS2) 56) Chlorobenzene-d5 (IS3) | | | | | | |
| System Monitoring Compounds 33) 1,2-Dichloroethane-d4(Spiked Amount 25.000 57) Toluene-d8 (SS2) Spiked Amount 25.000 73) Bromofluorobenzene (SS3) Spiked Amount 25.000 | 13.73 18.93 23.29 | 65 98 174 | 678298 | 24.687 ery = 24.904 ery = 24.454 ery = | ng | 0.00 |
| <pre>Target Compounds 2) Propene 3) Dichlorodifluoromethane 4) Chloromethane 5) Freon 114 6) Vinyl Chloride 7) 1,3-Butadiene 8) Bromomethane 9) Chloroethane 10) Ethanol 11) Acetonitrile 12) Acrolein 13) Acetone 14) Trichlorofluoromethane 15) Isopropanol 16) Acrylonitrile 17) 1,1-Dichloroethene 18) tert-Butanol 19) Methylene Chloride 20) Allyl Chloride 21) Trichlorotrifluoroethane 22) Carbon Disulfide 23) trans-1,2-Dichloroethene 24) 1,1-Dichloroethane 25) Methyl tert-Butyl Ether 26) Vinyl Acetate 27) 2-Butanone 28) cis-1,2-Dichloroethene 29) Diisopropyl Ether 30) Ethyl Acetate 31) n-Hexane</pre> | 4.95 5.27 5.52 5.71 5.99 6.48 6.81 7.12 7.44 7.64 7.64 7.64 7.86 8.14 8.32 8.64 9.16 9.26 | 85 50 135 62 54 94 64 45 41 58 1015 53 96 59 | 718887 1089924 1100149 587100 501238 514214m 1439658 393970 573236 1195874 1795159m 929763 593272 1706507 | 22.673 24.185 23.083 30.131 25.899 25.623 22.644 24.123 24.550 25.739 25.071 23.820 26.906 26.703 27.302 | ng ng ng ng ng ng ng ng ng ng ng ng | 97 98 63 99 98 90 90 |

DH 4/16/08

Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140815.D Acq On : 14 Apr 2008 23:45 Operator : WA Sample : 25ng TO-15 ICV Standard Misc : S20-04140804/S20-04040804 ALS Vial : 16 Sample Multiplier: 1 Quant Time: Apr 15 06:48:51 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration
 Internal Standards
 R.T. Qion Response Conc Units Dev (Min)

 22) Chloroform
 12.60
 63
 1101851
 29.333 ng
 99

 33) Tetrahydrofuran
 13.35
 72
 414917
 26.561 ng
 96

 33) Tetrahydrofuran
 13.48
 87
 680275
 25.226 ng
 #
 83

 34) Tetrahydrofuran
 14.29
 97
 953172
 25.625 ng
 99

 33) Isopropyl Acetace
 14.85
 56
 615240
 25.675 ng
 #
 35

 40) 1-Butanol
 14.85
 56
 615240
 25.686 ng
 #
 80

 41) Eperane
 14.85
 56
 615240
 25.686 ng
 #
 80

 42) Carbon Tetrachloride
 15.22
 117
 820346
 25.987 ng
 100

 43) Epoctane
 16.46
 83
 800346
 25.986 ng
 #
 99

 45) Isooctane
 16.62
 57
 2417594
 24.933 ng
 73

 46) Epromolichioromethane
 16.62
 57
 2417594
 24.933 ng
 73

 51) n-Heptane
 16.62
 57
 2417594
 24.933 ng
 73

 52) Metyl Methacrylate
 16.867
 97
 564173
 30.232 ng
 100

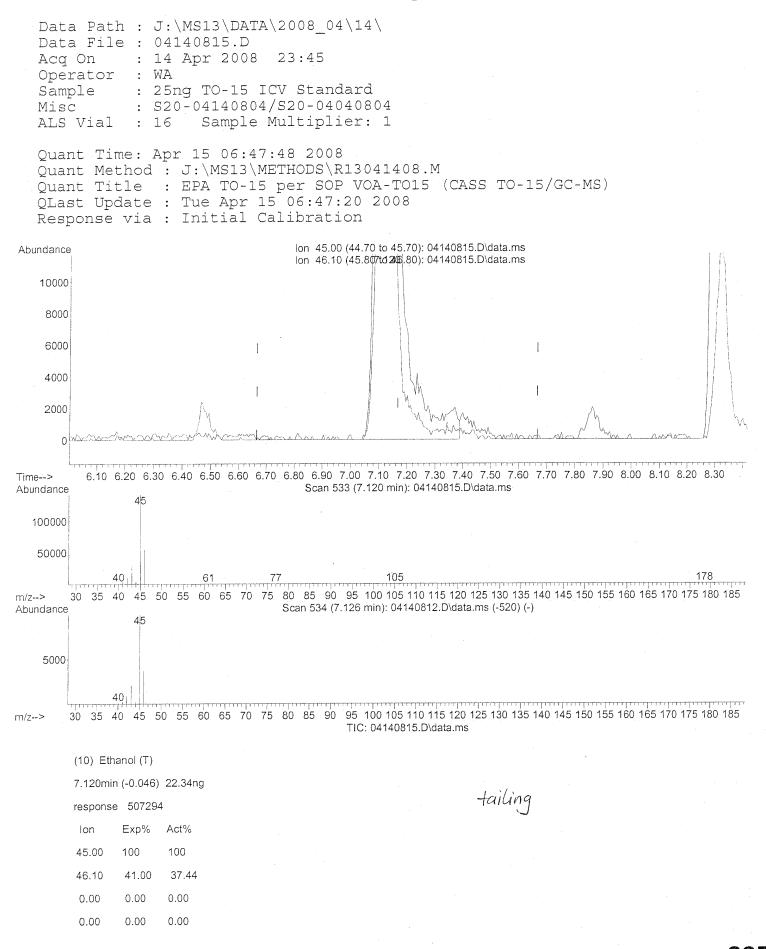
 53) Metyl Methacrylate Internal Standards R.T. QIon Response Conc Units Dev(Min)

Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140815.D Acq On : 14 Apr 2008 23:45 Operator : WA Sample : 25ng TO-15 ICV Standard Misc : S20-04140804/S20-04040804 ALS Vial : 16 Sample Multiplier: 1 Quant Time: Apr 15 06:48:51 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards 80) alpha-Methylstyrene24.56118114766223.553ng81) 2-Ethyltoluene24.61105267697723.667ng82) 1,2,4-Trimethylbenzene24.88105265241425.511ng83) n-Decane24.9957143840125.017ng 96 97 82)1,2,4-Trimethylbenzene24.0010183)n-Decane24.9957143840125.017 ng84)Benzyl Chloride25.0591199656928.171 ng85)1,3-Dichlorobenzene25.08146131813023.418 ng86)1,4-Dichlorobenzene25.16146129946024.124 ng87)sec-Butylbenzene25.21105311691125.534 ng88)p-Isopropyltoluene25.40119314216329.517 ng89)1,2,3-Trimethylbenzene25.41105280303927.414 ng90)1,2-Dichlorobenzene25.58146134185723.261 ng91)d-Limonene26.1115740131926.260 ng 97 88 94 99 99 95 91 96 #) # #) = qualifier out of #

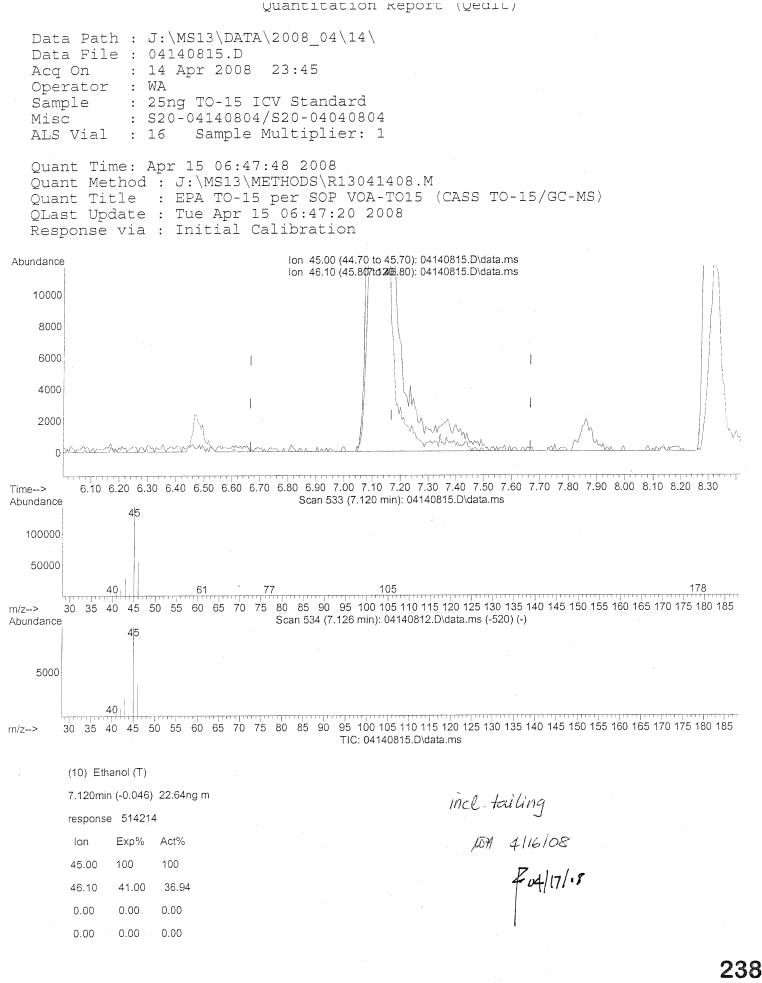
(#) = qualifier out of range (m) = manual integration (+) = signals summed

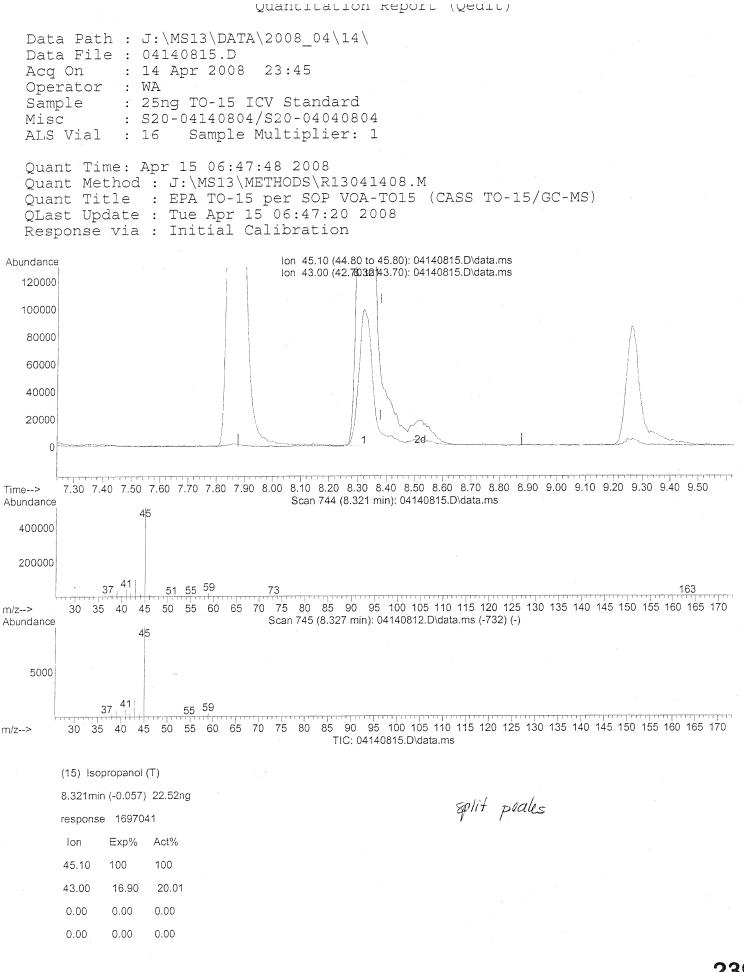
13041408.M Tue Apr 15 06:49:39 2008

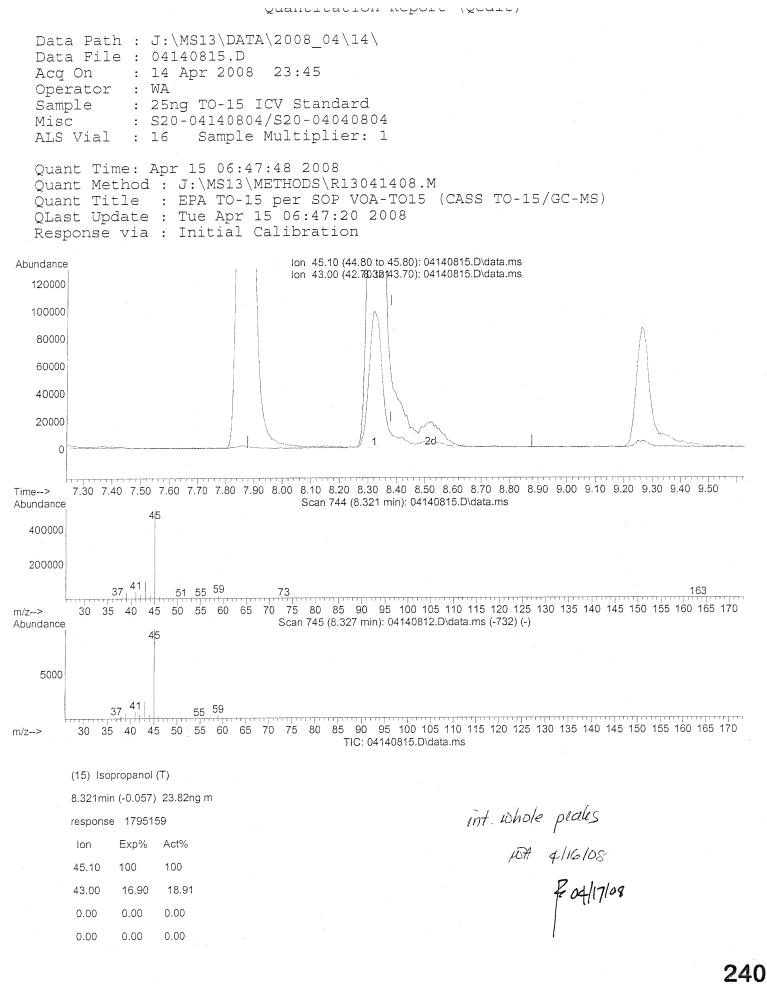
A116108



R13041408.M Tue Apr 15 06:48:34 2008







INITIAL CALIBRATION VERIFICATION CHECK SHEET

Data File Name: 04140815.D Data File Path: J:\MS13\DATA\2008_04\14\ Operator: WA Date Acquired: 4/14/08 23:45 Acq. Method File: TO15.M Sample Name: 25ng TO-15 ICV Standard Misc Info: S20-04140804/S20-04040804 Instrument Name: GCMS13

| | Name | Ret. | Amt. | Spike | % | Lower | Upper | * OR |
|----------|--------------------------|-------|-------------|-----------------|------------------|--------------|-------|-------------|
| <u>#</u> | Compound | Time | <u>(ng)</u> | <u>Amt.(ng)</u> | <u>Rec.</u> | <u>Limit</u> | Limit | <u>Fail</u> |
| 2) | Propene | 4.79 | 24,40 | 26.3 | 92.8 | 70 | 130 | * |
| 3) | Dichlorodifluoromethane | 4.95 | 22.67 | 25.5 | 88.9 | 70 | 130 | * |
| 4) | Chloromethane | 5.27 | 22.68 | 24.5 | 92.6 | 70 | 130 | * |
| 5) | Freon 114 | 5.52 | 24.19 | 26.0 | 93.0 | 70 | 130 | * |
| 6) | Vinyl Chloride | 5.71 | 23.08 | 24.8 | 93.1 | 70 | 130 | * |
| 7) | 1,3-Butadiene | 5.99 | 30.13 | 30.0 | 100.4 | 70 | 130 | * |
| 8) | Bromomethane | 6.48 | 25.90 | 25.0 | 103.6 | 70 | 130 | * |
| 9) | Chloroethane | 6.81 | 25.62 | 25.0 | 102.5 | 70 | 130 | * |
| 10) | Ethanol | 7.12 | 22.64 | 23.8 | 95.1 | 70 | 130 | * |
| 11) | Acetonitrile | 7.44 | 24.12 | 25.3 | 95.3 | 70 | 130 | * |
| 12) | Acrolein | 7.64 | 24.55 | 24.8 | 99.0 | 70 | 130 | * |
| 13) | Acetone | 7.86 | 25.74 | 26.8 | 96.0 | 70 | 130 | * |
| 14) | Trichlorofluoromethane | 8.14 | 25.07 | 26.3 | 95.3 | 70 | 130 | * |
| 15) | Isopropanol | 8.32 | 23.82 | 25.8 | 92.3 | 70 | 130 | * |
| 16) | Acrylonitrile | 8.64 | 26.91 | 25.5 | 105.5 | 70 | 130 | * |
| 17) | 1,1-Dichloroethene | 9.16 | 26.70 | 27.8 | 96.1 | 70 | 130 | * |
| 18) | tert-Butanol | 9.26 | 27.30 | 25.8 | 105.8 | 70 | 130 | * |
| 19) | Methylene Chloride | 9.36 | 24.39 | 27.8 | 87.7 | 70 | 130 | * |
| 20) | Allyl Chloride | 9.54 | 31.46 | 26.8 | 117.4 | 70 | 130 | * |
| 21) | Trichlorotrifluoroethane | 9.80 | 25.40 | 27.8 | 91.4 | 70 | 130 | * |
| 22) | Carbon Disulfide | 9.76 | 24.92 | 25.0 | 99.7 | 70 | 130 | * |
| 23) | trans-1,2-Dichloroethene | 10.80 | 25.91 | 26.5 | 97.8 | 70 | 130 | * |
| 24) | 1,1-Dichloroethane | 11.10 | 25.83 | 26.8 | 96.4 | 70 | 130 | * |
| 25) | Methyl tert-Butyl Ether | 11.19 | 25.62 | 26.8 | 95.6 | 70 | 130 | * |
| 26) | Vinyl Acetate | 11.35 | 29.96 | 25.3 | 118.4 | 70 | 130 | * |
| 27) | 2-Butanone | 11.68 | 27.93 | 27.0 | 103.4 | 70 | 130 | * |
| 28) | cis-1,2-Dichloroethene | 12.36 | 25.85 | 27.0 | 9 5.7 | 70 | 130 | * |
| 29) | Diisopropyl Ether | 12.69 | 23.70 | 26.3 | 90.1 | 70 | 130 | * |
| 30) | Ethyl Acetate | 12.69 | 28.00 | 29.3 | 95.6 | 70 | 130 | * |
| 31) | n-Hexane | 12.70 | 24.38 | 27.0 | 90.3 | 70 | 130 | * |
| 32) | Chloroform | 12.80 | 29.33 | 29.8 | 98.4 | 70 | 130 | * |
| 34) | Tetrahydrofuran | 13.35 | 26.56 | 26.8 | 99.1 | 70 | 130 | * |
| 35) | Ethyl tert-Butyl Ether | 13.48 | 25.23 | 26.0 | 97.0 | 70 | 130 | * |
| 36) | 1,2-Dichloroethane | 13.89 | 24.96 | 26.3 | 94.9 | 70 | 130 | * |
| 38) | 1,1,1-Trichloroethane | 14.29 | 25.42 | 26.8 | 94.9 | 70 | 130 | * |
| 39) | Isopropyl Acetate | 14.83 | 25.87 | 25.5 | 101.5 | 70 | 130 | * |

J:\MS13\ICV_0907.CRT

pt 4/16/08

4/15/08 6:50 AM **241**

INITIAL CALIBRATION VERIFICATION CHECK SHEET

Data File Name: 04140815.D Data File Path: J:\MS13\DATA\2008_04\14\ Operator: WA Date Acquired: 4/14/08 23:45 Acq. Method File: TO15.M Sample Name: 25ng TO-15 ICV Standard Misc Info: S20-04140804/S20-04040804 Instrument Name: GCMS13

| | Name | Ret. | Amt. | Spike | % | Lower | Upper | * OR |
|----------|---------------------------|-------|-------------|-----------------|-------|--------------|--------------|-------|
| <u>#</u> | Compound | Time | <u>(ng)</u> | <u>Amt.(ng)</u> | Rec. | <u>Limit</u> | <u>Limit</u> | Fail |
| 40) | 1-Butanol | 14.85 | 25.05 | 24.8 | 101.0 | 70 | 130 | * |
| 41) | Benzene | 14.99 | 24.72 | 27.0 | 91.6 | 70 | 130 | * |
| 42) | Carbon Tetrachloride | 15.22 | 25.99 | 26.0 | 99.9 | 70 | 130 | * |
| 43) | Cyclohexane | 15.41 | 25.59 | 26.8 | 95.5 | 70 | 130 | * |
| 44) | tert-Amyl Methyl Ether | 15.87 | 25.75 | 26.0 | 99.0 | 70 | 130 | * |
| 45) | 1,2-Dichloropropane | 16.20 | 24.80 | 26.5 | 93.6 | 70 | 130 | * |
| 46) | Bromodichloromethane | 16.46 | 27.14 | 27.8 | 97.6 | 70 | 130 | * |
| 47) | Trichloroethene | 16.54 | 25.20 | 27.3 | 92.3 | 70 | 130 | * |
| 48) | 1,4-Dioxane | 16.49 | 27.93 | 27.5 | 101.6 | 70 | 130 | * |
| 49) | Isooctane | 16.62 | 24.93 | 26.3 | 94.8 | 70 | 130 | * |
| 50) | Methyl Methacrylate | 16.80 | 25.95 | 25.8 | 100.6 | 70 | 130 | . · * |
| 51) | n-Heptane | 16.98 | 24.24 | 26.8 | 90.5 | 70 | 130 | * |
| 52) | cis-1,3-Dichloropropene | 17.73 | 26.16 | 25.0 | 104.7 | 70 | 130 | * |
| 53) | 4-Methyl-2-pentanone | 17.77 | 25.46 | 27.5 | 92.6 | 70 | 130 | * |
| 54) | trans-1,3-Dichloropropene | 18.43 | 30.23 | 28.0 | 108.0 | 70 | 130 | * |
| 55) | 1,1,2-Trichloroethane | 18.67 | 24.56 | 26.3 | 93.4 | 70 | 130 | * |
| 58) | Toluene | 19.07 | 24.74 | 26.5 | 93.3 | 70 | 130 | * |
| 59) | 2-Hexanone | 19.37 | 25.37 | 26.3 | 96.5 | 70 | 130 | * |
| 60) | Dibromochloromethane | 19.61 | 27.20 | 27.0 | 100.8 | 70 | 130 | * |
| 61) | 1,2-Dibromoethane | 19.94 | 27.00 | 26.3 | 102.7 | 70 | 130 | * |
| 62) | Butyl Acetate | 20.19 | 28.01 | 26.3 | 106.5 | 70 | 130 | * |
| 63) | n-Octane | 20.35 | 25.18 | 26.0 | 96.8 | 70 | 130 | * |
| 64) | Tetrachloroethene | 20.55 | 24.00 | 26.0 | 92.3 | 70 | 130 | * |
| 65) | Chlorobenzene | 21.42 | 24.37 | 26.5 | 91.9 | 70 | 130 | * |
| 66) | Ethylbenzene | 21.89 | 25.30 | 26.3 | 96.2 | 70 | 130 | * |
| 67) | m- & p-Xylene | 22.13 | 59.91 | 62.5 | 95.9 | 70 | 130 | * |
| 68) | Bromoform | 22.21 | 33.79 | 31.3 | 108.0 | 70 | 130 | * |
| 69) | Styrene | 22.57 | 25.05 | 26.3 | 95.3 | 70 | 130 | * |
| 70) | o-Xylene | 22.72 | 28.25 | 29.8 | 94.8 | 70 | 130 | * |
| 71) | n-Nonane | 22.98 | 24.43 | 26.0 | 94.0 | 70 | 130 | * |
| 72) | 1,1,2,2-Tetrachloroethane | 22.69 | 29.03 | 29.8 | 97.4 | 70 | 130 | * |
| 74) | Cumene | 23.47 | 26.32 | 27.0 | 97.5 | 70 | 130 | * |
| 75) | alpha-Pinene | 23.97 | 25.04 | 26.3 | 95.2 | 70 | 130 | * |
| , 76) | n-Propylbenzene | 24.10 | 25.93 | 26.3 | 98.6 | 70 | 130 | * |
| 77). | 3-Ethyltoluene | 24.23 | 24.36 | 25.5 | 95.5 | 70 | 130 | * |
| | | | | | | | | |

J:\MS13\ICV_0907.CRT

JA 4/16/08

4/15/08 6:50 AM **242**

INITIAL CALIBRATION VERIFICATION CHECK SHEET

Data File Name: 04140815.D Data File Path: J:\MS13\DATA\2008_04\14\ Operator: WA Date Acquired: 4/14/08 23:45 Acq. Method File: TO15.M Sample Name: 25ng TO-15 ICV Standard Misc Info: S20-04140804/S20-04040804 Instrument Name: GCMS13

| | Name | Ret. | Amt. | Spike % | | Lower | Upper | * OR |
|----------|-----------------------------|-------------|-------------|-----------------|-------------|-------|--------------|------|
| <u>#</u> | Compound | <u>Time</u> | <u>(ng)</u> | <u>Amt.(ng)</u> | <u>Rec.</u> | Limit | <u>Limit</u> | Fail |
| 78) | 4-Ethyltoluene | 24.28 | 25.73 | 26.5 | 97.1 | 70 | 130 | * |
| 79) | 1,3,5-Trimethylbenzene | 24.38 | 24.85 | 26.0 | 95.6 | 70 | 130 | * |
| 80) | alpha-Methylstyrene | 24.56 | 23.55 | 25.5 | 92.4 | 70 | 130 | * |
| 81) | 2-Ethyltoluene | 24.61 | 23.67 | 24.8 | 95.4 | 70 | 130 | * |
| 82) | 1,2,4-Trimethylbenzene | 24.88 | 25.51 | 26.0 | 98.1 | 70 | 130 | * |
| 83) | n-Decane | 24.99 | 25.02 | 26.3 | 95.1 | 70 | 130 | * |
| 84) | Benzyl Chloride | 25.05 | 28.17 | 25.8 | 109.2 | 70 | 130 | * |
| 85) | 1,3-Dichlorobenzene | 25.08 | 23.42 | 25.5 | 91.8 | 70 | 130 | * |
| 86) | 1,4-Dichlorobenzene | 25.16 | 24.12 | 26.3 | 91.7 | 70 | 130 | * |
| 87) | sec-Butylbenzene | 25.21 | 25.53 | 26.8 | 95.3 | 70 | 130 | * |
| 88) | p-Isopropyltoluene | 25.40 | 29.52 | 28.8 | 102.5 | 70 | 130 | * |
| 89) | 1,2,3-Trimethylbenzene | 25.41 | 27.41 | 28.5 | 96.2 | 70 | 130 | * |
| 90) | 1,2-Dichlorobenzene | 25.58 | 23.26 | 25.8 | 90.2 | 70 | 130 | * |
| 91) | d-Limonene | 25.58 | 23.46 | 26.0 | 90.2 | 70 | 130 | * |
| 92) | 1,2-Dibromo-3-Chloropropane | 26.11 | 26.26 | 25.8 | 101.8 | 70 | 130 | * |
| 93) | n-Undecane | 26.50 | 25.02 | 26.5 | 94.4 | 70 | 130 | * |
| 94) | 1,2,4-Trichlorobenzene | 27.63 | 24.47 | 26.0 | 94.1 | 70 | 130 | * |
| 95) | Naphthalene | 27.77 | 25.66 | 26.3 | 97.6 | 70 | 130 | * |
| , 96) | n-Dodecane | 27.74 | 23.67 | 26.5 | 89.3 | 70 | 130 | * |
| 97) | Hexachloro-1,3-butadiene | 28.19 | 24.23 | 26.3 | 92.1 | 70 | 130 | * |
| | | | | | | | | |

Bold = 67 Compound List

A 116/08-

l.65 27.07 9.65 1.27 1.95 %RSD 2.241 0.771 2.738 2.005 2.954 Avg 1.968 1.947 2.190 4.028 3.424 0.774 5.0 = 04140811.D100 3.226 2.195 3.456 0.766 50 2.045 1.980 2.032 3.060 2.226 0.780 2.823 25 TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD - - ISTD---ISTD --ISTD 2.254 0.787 2.744 2.209 1.0 =04140810.D 100 =04140814.D 5.0 2.727 0.763 2.193 2.272 1.0 2.037 0.759 2.277 2.296 2.728 : Mon Apr 28 10:04:59 2008 0.5 Response Via : Initial Calibration 2.025 0.769 2.158 2.766 2.272 0.5 =04140809.D 50 =04140813.D =04140813.D 0.1 Method Path : J:\MS13\METHODS\ (IS3 F) 1,4-Difluorobenzene (1,2-Dichloroethane-Bromochloromethane Bromofluorobenzene : S13041408.M tert-Butylbenzene Chlorobenzene-d5 Toluene-d8 (SS2) n-Butylbenzene Range Calibration Files Compound 0.1 =04140808.D 25 =04140812.D Ú. Ú. Method File •• Last Update Out S IR ДR Title нол 11 (#) $\overrightarrow{\mathsf{N}}$ $\widehat{\mathfrak{S}}$

5 S13041408.M Mon Apr 28 10:05:09 2008

Page:

EN 4/25/05

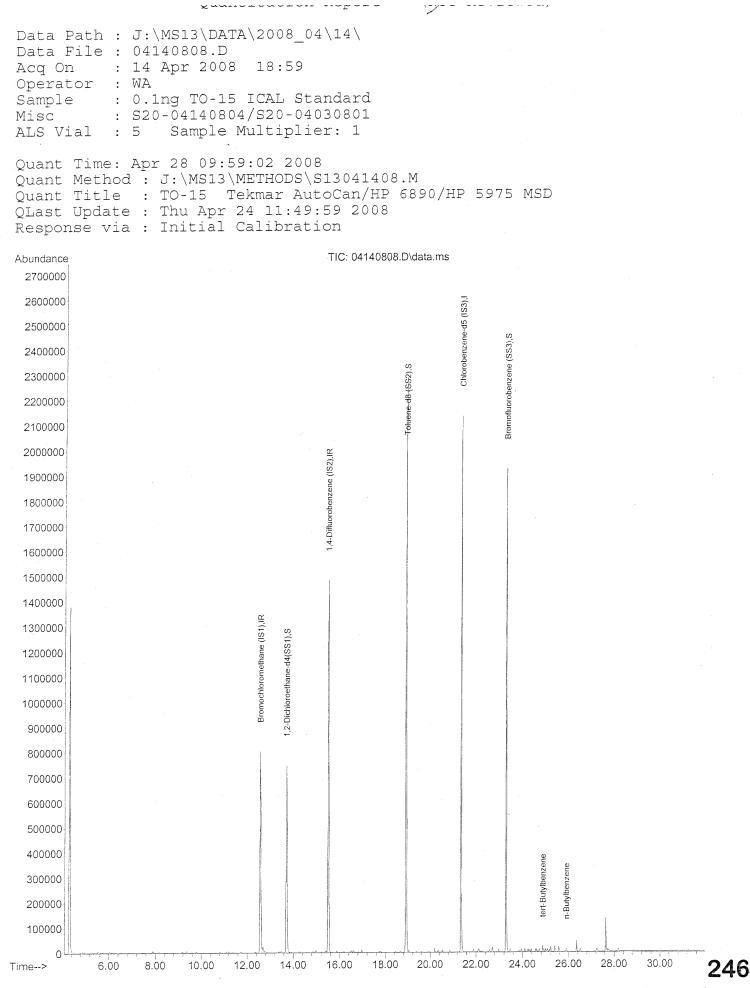
Method Path : J:\MS13\METHODS\ Method File : S13041408.M Title : TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD Last Update : Mon Apr 28 10:06:00 2008 Response Via : Initial Calibration

| #. | ID | Conc | ISTD | Path\File |
|----|-----|------|------|------------------------------------|
| | | | Conc | |
| | | | | |
| 1 | 0.1 | 0 | 25 | J:\MS13\DATA\2008_04\14\04140808.D |
| 2 | 0.5 | 1 | 25 | J:\MS13\DATA\2008_04\14\04140809.D |
| 3 | 1.0 | 1 | 25 | J:\MS13\DATA\2008_04\14\04140810.D |
| 4 | 5.0 | 5 | 25 | J:\MS13\DATA\2008_04\14\04140811.D |
| 5 | 25 | 26 | 25 | J:\MS13\DATA\2008_04\14\04140812.D |
| 6 | 50 | 52 | 25 | J:\MS13\DATA\2008_04\14\04140813.D |
| 7 | 100 | 104 | 25 | J:\MS13\DATA\2008_04\14\04140814.D |

| # | ID | Update Time | Quant Time | Acquisition Time | | | |
|------------------|--------------------------------------|---|---|--|--|--|--|
| 2 3 4 5 | 0.1 0.5 1.0 5.0 25 50 | Apr 28 10:03 2008 Apr 28 10:03 2008 Apr 28 10:03 2008 Apr 28 10:03 2008 Apr 28 10:04 2008 Apr 28 10:04 2008 Apr 28 10:04 2008 | Apr 28 09:59 2008 Apr 28 09:59 2008 Apr 28 10:00 2008 Apr 28 10:00 2008 Apr 28 10:00 2008 Apr 28 10:00 2008 Apr 28 10:02 2008 | 14 Apr 2008 18:59 14 Apr 2008 19:40 14 Apr 2008 20:21 14 Apr 2008 21:01 14 Apr 2008 21:43 14 Apr 2008 22:24 | | | |
| 7 | 100 | Apr 28 10:04 2008 | Apr 28 10:02 2008 | 14 Apr 2008 23:04 | | | |

513041408.M Thu May 08 16:19:56 2008

D# 5/8/08



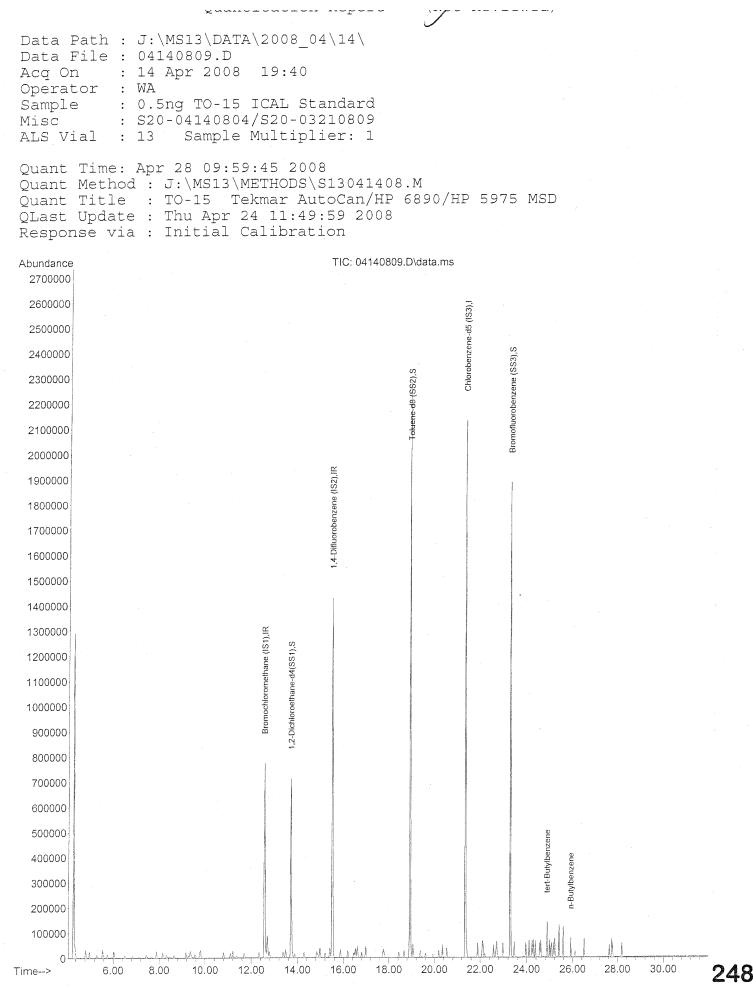
S13041408.M Mon Apr 28 09:59:03 2008

(THE ICVICVICV) Quantituation Report Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140808.D Acq On : 14 Apr 2008 18:59 Operator : WA Sample : 0.1ng TO-15 ICAL Standard Misc : S20-04140804/S20-04030801 ALS Vial : 5 Sample Multiplier: 1 Quant Time: Apr 28 09:59:02 2008 Quant Method : J:\MS13\METHODS\S13041408.M Quant Title : TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD OLast Update : Thu Apr 24 11:49:59 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards 1) Bromochloromethane (IS1)12.5813033067225.000 ng-0.043) 1,4-Difluorobenzene (IS2)15.51114151679925.000 ng-0.034) Chlorobenzene-d5 (IS3)21.358275815225.000 ng-0.01 System Monitoring Compounds 2) 1,2-Dichloroethane-d4(... 13.72 65 669686 31.921 ng -0.03 Spiked Amount25.000Recovery = 127.68%5) Toluene-d8 (SS2)18.9398172218929.398 ng-0.02Spiked Amount25.000Recovery = 117.60%6) Bromofluorobenzene (SS3)23.2917458306920.588 ng-0.01 Recovery = 127.68% Recovery = 82.36% Spiked Amount 25.000 Qvalue Target Compounds 7) tert-Butylbenzene24.8811968050.087ng8) n-Butylbenzene25.919189750.107ng# 99 89

(#) = qualifier out of range (m) = manual integration (+) = signals summed

13041408.M Mon Apr 28 09:59:03 2008

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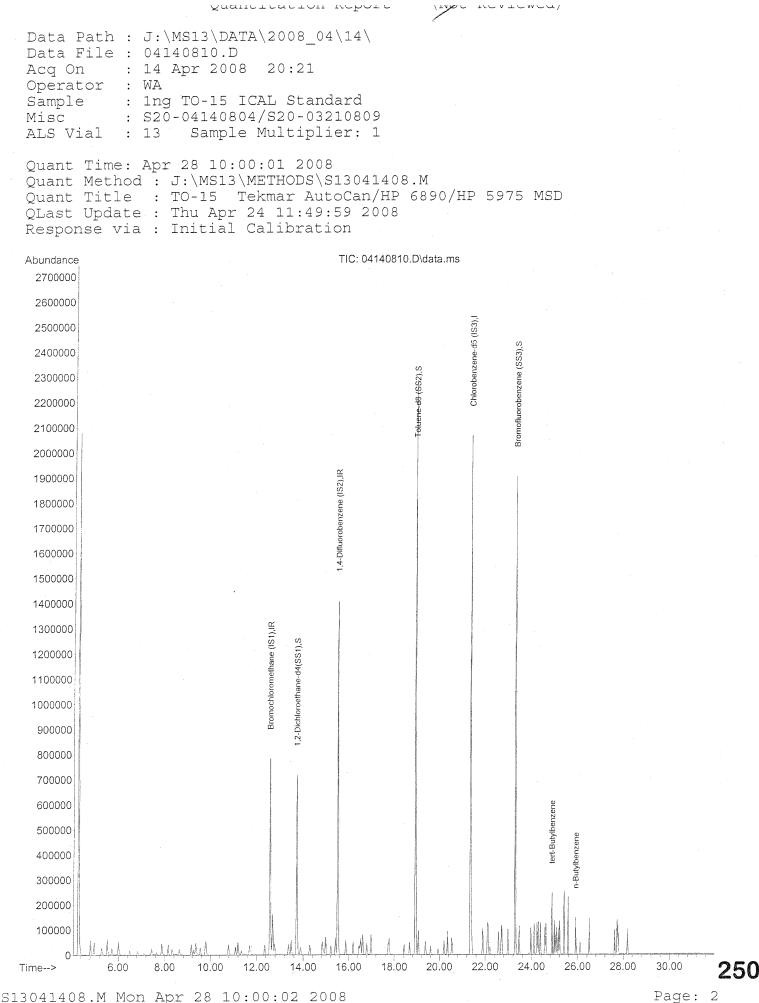


S13041408.M Mon Apr 28 09:59:46 2008

Quanter cacross report Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140809.D Acq On : 14 Apr 2008 19:40 Operator : WA Sample : 0.5ng TO-15 ICAL Standard Misc : S20-04140804/S20-03210809 ALS Vial : 13 Sample Multiplier: 1 Ouant Time: Apr 28 09:59:45 2008 Quant Method : J:\MS13\METHODS\S13041408.M Quant Title : TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD QLast Update : Thu Apr 24 11:49:59 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards 1) Bromochloromethane (IS1) 12.58 130 321232 25.000 ng -0.04 3) 1,4-Difluorobenzene (IS2)15.51114146814225.000 ng-0.024) Chlorobenzene-d5 (IS3)21.358274431125.000 ng-0.01 System Monitoring Compounds 2) 1,2-Dichloroethane-d4(... 13.73 65 654334 32.106 ng -0.03 Recovery = 128.44% Spiked Amount25.000Recovery = 128.44%5) Toluene-d8 (SS2)18.9398169442429.462 ng-0.02Spiked Amount25.000Recovery = 117.84%6) Bromofluorobenzene (SS3)23.2917456526320.330 ng-0.01 Spiked Amount 25.000 Recovery = 81.32% Spiked Amount 25.000 Ovalue Target Compounds 7)tert-Butylbenzene24.88119355500.461ng8)n-Butylbenzene25.9191434520.527ng 98 95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

13041408.M Mon Apr 28 09:59:46 2008



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Xuurrenoucher report Data Path : J:\MS13\DATA\2008_04\14\ Data File : 04140810.D Acq On : 14 Apr 2008 20:21 Operator : WA Sample : 1ng TO-15 ICAL Standard Misc : S20-04140804/S20-03210809 ALS Vial : 13 Sample Multiplier: 1 Ouant Time: Apr 28 10:00:01 2008 Quant Method : J:\MS13\METHODS\S13041408.M Quant Title : TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD QLast Update : Thu Apr 24 11:49:59 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards ______ 1) Bromochloromethane (IS1)12.5813031446125.000 ng-0.043) 1,4-Difluorobenzene (IS2)15.51114145464725.000 ng-0.024) Chlorobenzene-d5 (IS3)21.358273508325.000 ng-0.01 System Monitoring Compounds 2) 1,2-Dichloroethane-d4(... 13.72 65 643137 32.236 ng -0.03 Recovery = 128.96%

 Spiked Amount
 25.000
 Recovery
 = 128.96%

 5) Toluene-d8 (SS2)
 18.93
 98
 1669857
 29.400 ng
 -0.02

 Spiked Amount
 25.000
 Recovery
 = 117.60%

 6) Bromofluorobenzene (SS3)
 23.29
 174
 560818
 20.424 ng
 -0.01

 Recovery

 25.000
 Recovery
 = 81.68%

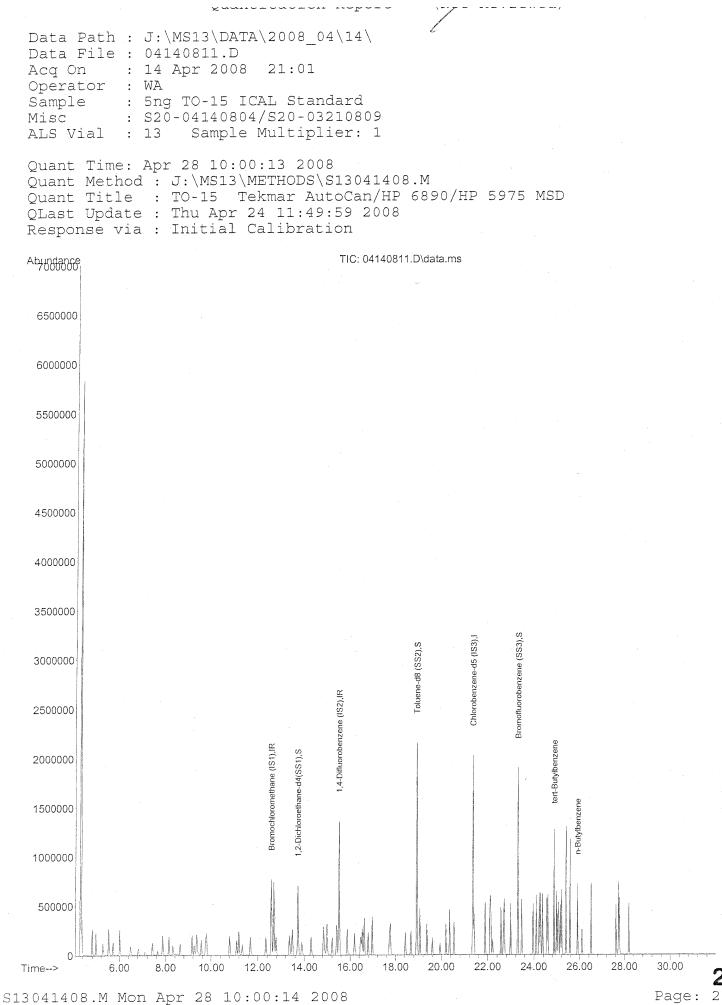
 Spiked Amount 25.000 Recovery = 81.68% Spiked Amount 25.000 Ovalue Target Compounds 7)tert-Butylbenzene24.88119670680.880ng8)n-Butylbenzene25.9191858031.054ng 98 91

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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13041408.M Mon Apr 28 10:00:02 2008

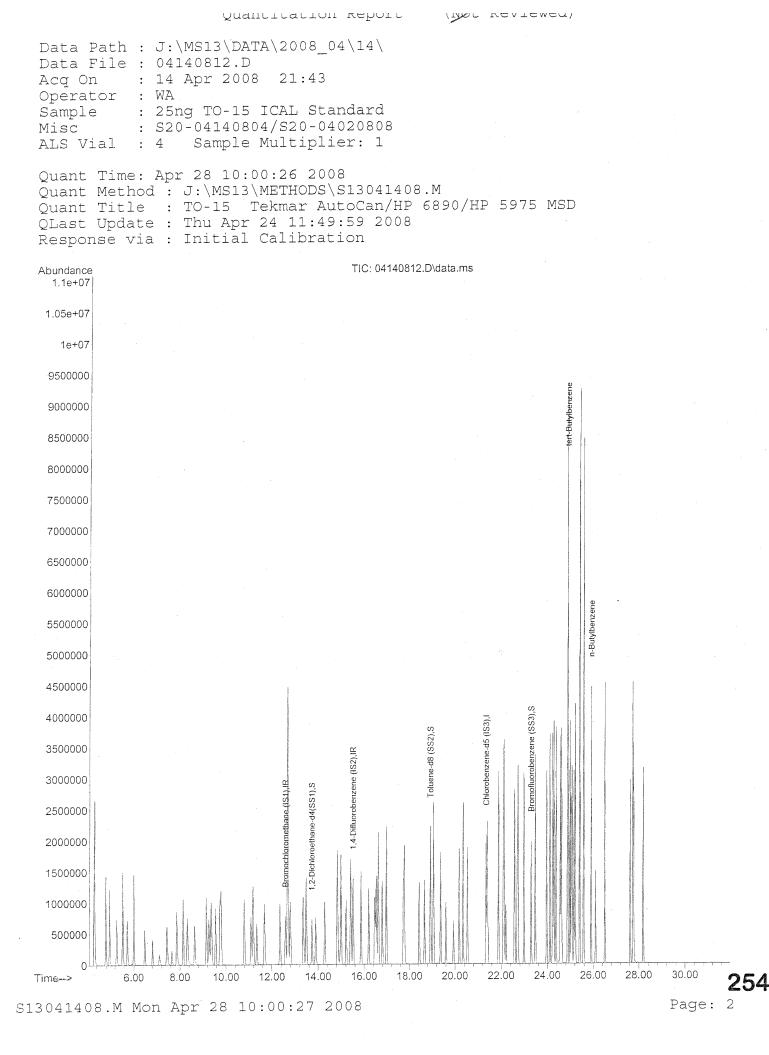
Page: 1



Xaanoroaoron nobee-Data Path : J:\MS13\DATA\2008_04\14\ Data File : 04140811.D Acq On : 14 Apr 2008 21:01 Operator : WA Sample : 5ng TO-15 ICAL Standard Misc : S20-04140804/S20-0321080 : S20-04140804/S20-03210809 ALS Vial : 13 Sample Multiplier: 1 Quant Time: Apr 28 10:00:13 2008 Quant Method : J:\MS13\METHODS\S13041408.M Quant Title : TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD QLast Update : Thu Apr 24 11:49:59 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards 1) Bromochloromethane (IS1)12.5813031358425.000 ng-0.033) 1,4-Difluorobenzene (IS2)15.51114140651525.000 ng-0.024) Chlorobenzene-d5 (IS3)21.358271579925.000 ng-0.01 -0.02 -0.01 System Monitoring Compounds 2) 1,2-Dichloroethane-d4(... 13.73 65 620740 31.200 ng -0.03 Recovery = 124.80% Spiked Amount25.000Recovery= 124.80%5) Toluene-d8 (SS2)18.9398161355629.174 ng-0.02Spiked Amount25.000Recovery= 116.68%6) Bromofluorobenzene (SS3)23.2917456316921.062 ng-0.01Recovery6) Bromofluorobenzene (SS3)23.2917456316921.062 ng-0.01 Spiked Amount 25.000 Recovery = 84.24% Spiked Amount 25.000 Qvalue 99 Target Compounds 7) tert-Butylbenzene24.881193289184.434ng998) n-Butylbenzene25.91914204045.303ng#94 _ __ __ __ __ __ __ __ __ __ __ __

(#) = qualifier out of range (m) = manual integration (+) = signals summed

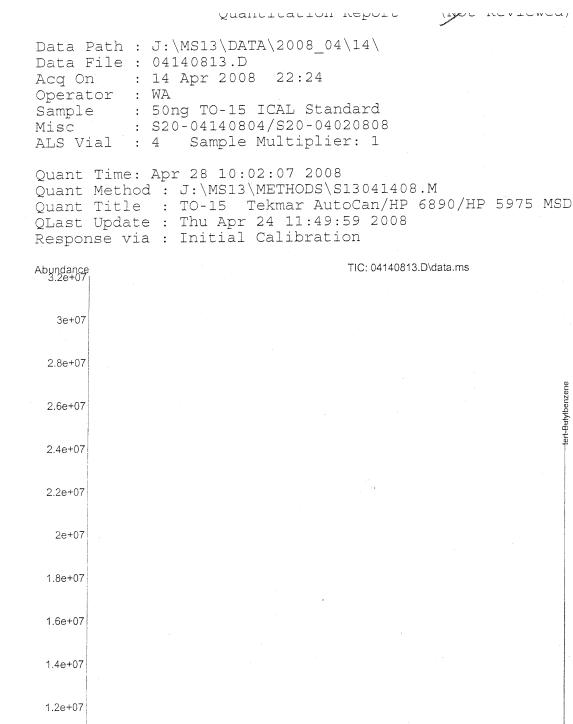
30/26/108

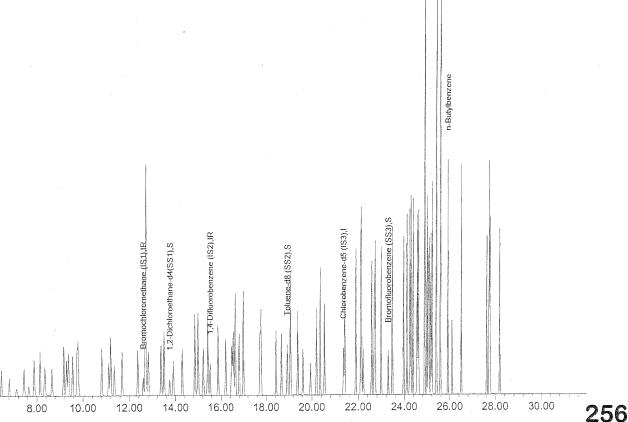


(INCL ICCVICWCC) QUAIILILALIUN NEPULL Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140812.D Acq On : 14 Apr 2008 21:43 Operator : WA Sample : 25ng TO-15 ICAL Standard Misc : S20-04140804/S20-04020808 ALS Vial : 4 Sample Multiplier: 1 Quant Time: Apr 28 10:00:26 2008 Quant Method : J:\MS13\METHODS\S13041408.M Quant Title : TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD QLast Update : Thu Apr 24 11:49:59 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards _____ 1) Bromochloromethane (IS1)12.6013033207025.000 ng-0.023) 1,4-Difluorobenzene (IS2)15.52114146703225.000 ng-0.024) Chlorobenzene-d5 (IS3)21.358276215225.000 ng-0.01 System Monitoring Compounds 2) 1,2-Dichloroethane-d4(... 13.73 65 674919 32.035 ng -0.02 Spiked Amount25.000Recovery = 128.16%5) Toluene-d8 (SS2)18.9398169687528.814 ng-0.02Spiked Amount25.000Recovery = 115.24%6) Bromofluorobenzene (SS3)23.2917459448920.881 ng-0.01 Recovery = 128.16% Recovery = 83.52% Spiked Amount 25.000 Qvalue 99 Target Compounds 7) tert-Butylbenzene24.88119223737728.326ng8) n-Butylbenzene25.9191249999329.618ng# 7) tert-Butylbenzene 93 _____

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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S13041408.M Mon Apr 28 10:02:08 2008

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Qualititation report Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140813.D Acq On : 14 Apr 2008 22:24 Operator : WA Sample : 50ng TO-15 ICAL Standard Misc : S20-04140804/S20-04020808 ALS Vial : 4 Sample Multiplier: 1 Quant Time: Apr 28 10:02:07 2008 Quant Method : J:\MS13\METHODS\S13041408.M Ouant Title : TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD QLast Update : Thu Apr 24 11:49:59 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards 1) Bromochloromethane (IS1)12.6013035913525.000 ng-0.023) 1,4-Difluorobenzene (IS2)15.53114158007725.000 ng-0.014) Chlorobenzene-d5 (IS3)21.368281877225.000 ng0.00 System Monitoring Compounds stem Monitoring Compounds 2) 1,2-Dichloroethane-d4(... 13.74 65 706628 31.013 ng -0.01

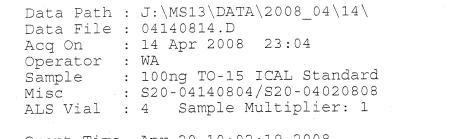
 2) 1,2 Dichlet Cooling and the second sec Recovery = 124.04% Recovery = 81.96% Spiked Amount 25.000 Qvalue Target Compounds 7)tert-Butylbenzene24.89119588564069.362ng8)n-Butylbenzene25.9191565198662.329ng 99 95 ·

(#) = qualifier out of range (m) = manual integration (+) = signals summed

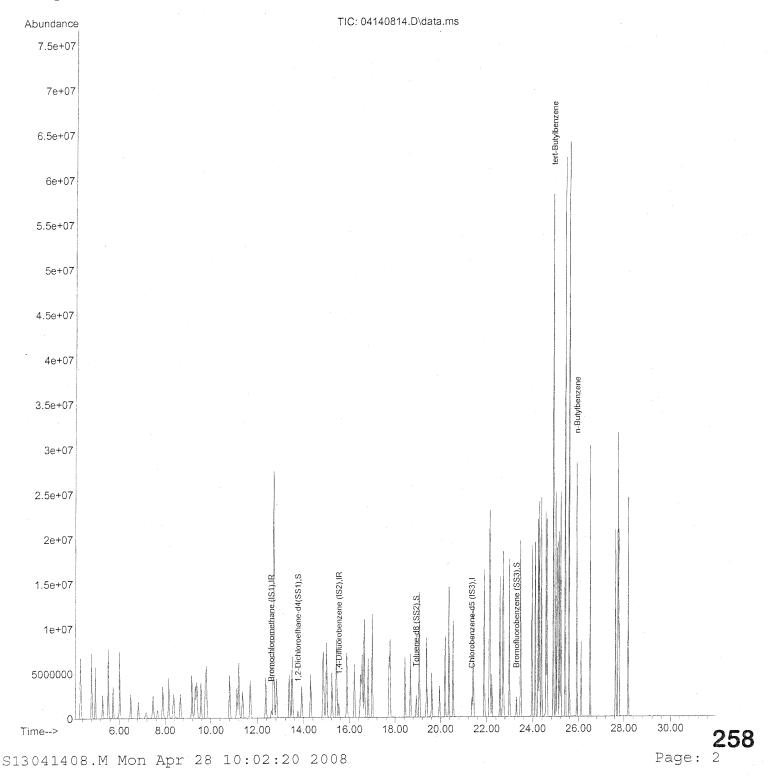
313041408.M Mon Apr 28 10:02:08 2008

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Quant Time: Apr 28 10:02:19 2008 Quant Method : J:\MS13\METHODS\S13041408.M Quant Title : TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD QLast Update : Thu Apr 24 11:49:59 2008 Response via : Initial Calibration

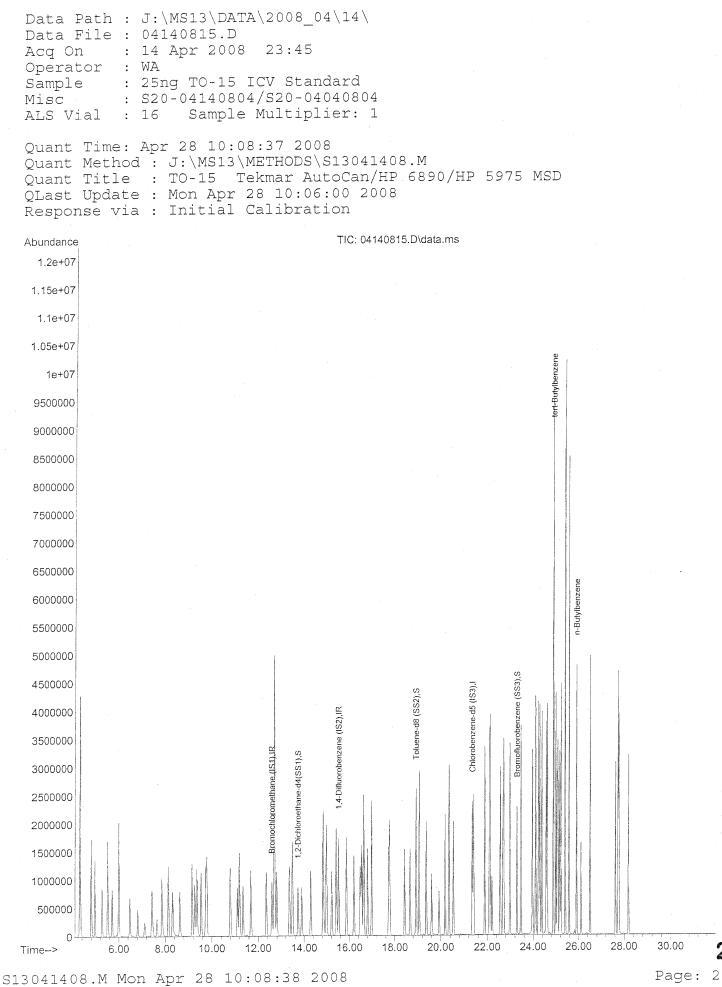


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|--|--------------------------------|-----------|--|---|
| Data Path : J:\MS13\DATA\2008_04 Data File : 04140814.D Acq On : 14 Apr 2008 23:04 Operator : WA Sample : 100ng TO-15 ICAL Sta Misc : S20-04140804/S20-040 ALS Vial : 4 Sample Multiplie | ndard 20808 | | | |
| Quant Time: Apr 28 10:02:19 2008 Quant Method : J:\MS13\METHODS\S Quant Title : TO-15 Tekmar Aut QLast Update : Thu Apr 24 11:49: Response via : Initial Calibratio | 1304140; oCan/HP 59 2008 | 689 | 0/HP 5975 MSD | |
| Internal Standards | R.T. | QIO | n Response Conc Units Dev(Min |) |
| 3) 1.4-Difluorobenzene (IS2) | 15.53 | 114 | 379040 25.000 ng -0.01 1673737 25.000 ng 0.00 871036 25.000 ng 0.00 | |
| 5) Toluene-d8 (SS2) Spiked Amount 25.000 | 18.93 | 98 | 738132 30.694 ng 0.00 Recovery = 122.76% 1908001 28.349 ng -0.01 Recovery = 113.40% 674346 20.725 ng 0.00 Recovery = 82.92% | |
| Target Compounds 7) tert-Butylbenzene 8) n-Butylbenzene | 24.89 25.92 | 119 91 | Qvalue 14594545 161.676 ng # 90 12764675 132.321 ng 93 | _ |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

13041408.M Mon Apr 28 10:02:19 2008

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Qualitication Report

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Quantituation Report STACE THE ATCANCOL Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140815.D Acq On : 14 Apr 2008 23:45 Operator : WA Sample : 25ng TO-15 ICV Standard Misc : S20-04140804/S20-04040804 ALS Vial : 16 Sample Multiplier: 1 Quant Time: Apr 28 10:08:37 2008 Quant Method : J:\MS13\METHODS\S13041408.M Quant Title : TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD QLast Update : Mon Apr 28 10:06:00 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards 1) Bromochloromethane (IS1) 12.59 130 402323 25.000 ng -0.02 3)1,4-Difluorobenzene (IS2)15.52114179919525.000 ng-0.014)Chlorobenzene-d5 (IS3)21.368289926825.000 ng0.00 System Monitoring Compounds stem Monitoring Compounds 2) 1,2-Dichloroethane-d4(... 13.73 65 796514 24.687 ng -0.02 Spiked Amount25.000Recovery =98.76%5) Toluene-d8 (SS2)18.9398200733224.904 ng0.00Spiked Amount25.000Recovery =99.60%6) Bromofluorobenzene (SS3)23.2917467829824.454 ng0.00 Recovery = 98.76% Recovery = 97.80% Spiked Amount 25.000 Qvalue Target Compounds 99

7)tert-Butylbenzene24.88119245327224.914ng998)n-Butylbenzene25.9191272379525.637ng94

(#) = qualifier out of range (m) = manual integration (+) = signals summed

13041408.M Mon Apr 28 10:08:38 2008

INITIAL CALIBRATION VERIFICATION CHECK SHEET

Data File Name: 04140815.D Data File Path: J:\MS13\DATA\2008_04\14\ Operator: WA Date Acquired: 4/14/08 23:45 Acq. Method File: TO15.M Sample Name: 25ng TO-15 ICV Standard Misc Info: S20-04140804/S20-04040804 Instrument Name: GCMS13

| | Name | Ret. | Amt. | Spike | % | Lower | Upper | * OR |
|----|-------------------|-------|-------------|----------|------|--------------|--------------|------|
| # | <u>Compound</u> | Time | <u>(ng)</u> | Amt.(ng) | Rec. | <u>Limit</u> | <u>Limit</u> | Fail |
| 7) | tert-Butylbenzene | 24.88 | 24.91 | 26.3 | 94.7 | 70 | 130 | * |
| 8) | n-Butylbenzene | 25.91 | 25.64 | 26.8 | 95.7 | 70 | 130 | * |

Page 1 of 1

J:\MS13\ICV_S0907_BBENZENES.CRT

5/9/08 4:21 PM 262

CONTINUING CALIBRATION STANDARDS

Data Path : J:\MS13\DATA\2008_05\08\ Data File : 05080801.D Acq On : 8 May 2008 8:41 am Operator : RTB Sample : 25ng TO-15 CCV Standard Misc : S20-04300802/S20-04250805 ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 10:13:43 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration

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Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min Max. RRF Dev : 30% Max. Rel. Area : 200%

| | Compound | AvgRF | CCRF | %Dev Ar | ea% | Dev(min | 1) |
|-------|----------------------------|-------|-------|---------|-----|---------|----|
| 1 IR | Bromochloromethane (IS1) | | 1.000 | 0.0 | 73 | -0.02 | |
| 2 T | Propene | 2.069 | 1.722 | 16.8 | | 0.00 | |
| 3 T | Dichlorodifluoromethane | 3.772 | 3.071 | 18.6 | | 0.00 | |
| 4 T | Chloromethane | 3.153 | 2.549 | 19.2 | 63 | -0.01 | |
| 5 T | Freon 114 | 1.847 | 1.563 | 15.4 | 68 | 0.00 | |
| 6 T | Vinyl Chloride | 2.934 | 2.322 | 20.9 | 63 | 0.00 | |
| 7 T | 1,3-Butadiene | 2.269 | 1.935 | 14.7 | 64 | | |
| 8 T | Bromomethane | 1.409 | 1.246 | 11.6 | 66 | | |
| 9 T | Chloroethane | 1.216 | 1.090 | 10.4 | 68 | | |
| 10 T | Ethanol | 1.411 | | 18.7 | 69 | | |
| 11 T | Acetonitrile | 3.708 | 3.127 | 15.7 | 70 | | |
| 12 T | Acrolein | 0.997 | 0.923 | 7.4 | 69 | | |
| 13 T | Acetone | 1.384 | 1.214 | 12.3 | 69 | -0.03 | |
| 14 T | Trichlorofluoromethane | 2.964 | 2.643 | 10.8 | 68 | 0.00 | |
| 15 T | Isopropanol | 4.683 | 4.232 | 9.6 | 67 | -0.06 | |
| 16 T | Acrylonitrile | 2.147 | 2.080 | 3.1 | 68 | -0.04 | |
| 17 T | 1,1-Dichloroethene | 1.381 | 1.238 | 10.4 | 69 | 0.00 | |
| 18 T | tert-Butanol | 3.884 | 3.749 | 3.5 | 68 | -0.05 | |
| 19 T | Methylene Chloride | 1.590 | 1.338 | 15.8 | 69 | -0.02 | |
| 20 T | Allyl Chloride | 2.124 | | -7.1 | 70 | -0.02 | |
| 21 T | Trichlorotrifluoroethane | 1.267 | 1.136 | 10.3 | 71 | 0.00 | |
| 22 T | Carbon Disulfide | 5.894 | 5.345 | 9.3 | 67 | -0.01 | |
| 23 T | trans-1,2-Dichloroethene | 2.406 | 2.120 | 11.9 | 66 | -0.02 | |
| 24 T | 1,1-Dichloroethane | 2.800 | 2.495 | 10.9 | 66 | -0.02 | |
| 25 T | Methyl tert-Butyl Ether | 4.599 | 4.015 | 12.7 | 68 | -0.02 | |
| 26 T | Vinyl Acetate | 0.275 | 0.298 | -8.4 | 74 | -0.01 | |
| 27 T | 2-Butanone | 0.970 | 0.923 | 4.8 | 68 | -0.02 | |
| 28 T | cis-1,2-Dichloroethene | 2.268 | 1.968 | 13.2 | 65 | -0.02 | |
| 29 T | Diisopropyl Ether | 1.274 | 1.093 | 14.2 | 67 | -0.02 | |
| 30 T | Ethyl Acetate | 0.605 | | 13.1 | 64 | -0.02 | |
| 31 T | n-Hexane | 3.144 | 2.559 | 18.6 | 64 | -0.01 | |
| 32 T | Chloroform | 2.334 | 2.041 | 12.6 | 67 | -0.03 | |
| 33 S | 1,2-Dichloroethane-d4(SS1) | | 1.830 | 8.7 | 66 | -0.02 | |
| 34 T | Tetrahydrofuran | 0.971 | | 7.3 | 67 | -0.02 | |
| 35 T | Ethyl tert-Butyl Ether | | | 9.5 | 70 | -0.02 | |
| 36 T | 1,2-Dichloroethane | 2.312 | 2.003 | 13.4 | 65 | -0.02 | |
| 37 IR | 1,4-Difluorobenzene (IS2) | | 1.000 | 0.0 | 71 | -0.01 | |
| 38 T | 1,1,1-Trichloroethane | 0.521 | 0.486 | 6.7 | 67 | -0.02 | 2 |

213041408.M Thu May 08 10:14:37 2008

\$ 05/08/08

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DVALAACO CONCLINATING CALEDIACION ROPOLO

| Data Path | : | J:\MS13\DATA\2008_05\08\ |
|-----------|---|---------------------------|
| Data File | : | 05080801.D |
| Acq On | : | 8 May 2008 8:41 am |
| Operator | | RTB |
| Sample | | 25ng TO-15 CCV Standard |
| Misc | | S20-04300802/S20-04250805 |
| ALS Vial | : | 4 Sample Multiplier: 1 |
| | | |

Quant Time: May 08 10:13:43 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min Max. RRF Dev : 30% Max. Rel. Area : 200%

| | Compound | AvgRF | CCRF | %Dev Ar | ea% | Dev(min) |
|------|---------------------------|-------|-------|---------|-----|----------|
| 39 T | Isopropyl Acetate | 0.227 | 0.217 | 4.4 | 67 | -0.02 |
| 40 T | 1-Butanol | 0.341 | 0.324 | 5.0 | 69 | -0.03 |
| 41 T | Benzene | 1.326 | 1.188 | 10.4 | 66 | |
| 42 T | Carbon Tetrachloride | 0.439 | 0.445 | -1.4 | 70 | |
| 43 T | Cyclohexane | 0.491 | 0.457 | 6.9 | 69 | |
| 44 T | tert-Amyl Methyl Ether | 0.944 | 0.883 | 6.5 | 67 | |
| 45 T | 1,2-Dichloropropane | 0.380 | 0.339 | 10.8 | 67 | |
| 46 T | Bromodichloromethane | 0.451 | 0.420 | 6.9 | 67 | |
| 47 T | Trichloroethene | 0.326 | 0.307 | 5.8 | 70 | |
| 48 T | 1,4-Dioxane | 0.235 | 0.231 | 1.7 | 70 | -0.02 |
| 49 T | Isooctane | 1.570 | 1.441 | 8.2 | 67 | |
| 50 T | Methyl Methacrylate | 0.120 | 0.121 | -0.8 | 71 | -0.01 |
| 51 T | n-Heptane | 0.367 | 0.328 | 10.6 | 66 | |
| 52 T | cis-1,3-Dichloropropene | 0.517 | 0.514 | 0.6 | 68 | |
| 53 T | 4-Methyl-2-pentanone | 0.363 | 0.347 | 4.4 | 66 | |
| 54 T | trans-1,3-Dichloropropene | 0.446 | 0.459 | -2.9 | 68 | |
| 55 T | 1,1,2-Trichloroethane | 0.319 | 0.293 | 8.2 | 70 | -0.01 |
| 56 I | Chlorobenzene-d5 (IS3) | 1.000 | 1.000 | 0.0 | 66 | 0.00 |
| 57 S | Toluene-d8 (SS2) | 2.241 | 2.288 | -2.1 | 68 | 0.00 |
| 58 T | Toluene | 2.815 | 2.724 | 3.2 | 67 | |
| 59 T | 2-Hexanone | 2.097 | 2.169 | -3.4 | 64 | -0.02 |
| 60 T | Dibromochloromethane | 0.679 | 0.722 | -6.3 | 69 | |
| 61 T | 1,2-Dibromoethane | 0.662 | | -7.9 | 70 | -0.01 |
| 62 T | Butyl Acetate | 2.097 | 2.254 | -7.5 | 66 | -0.01 |
| 63 T | n-Octane | 0.660 | 0.649 | 1.7 | 66 | -0.01 |
| 64 T | Tetrachloroethene | 0.705 | 0.708 | -0.4 | 71 | 0.00 |
| 65 T | Chlorobenzene | 1.744 | 1.732 | 0.7 | 69 | |
| 66 T | Ethylbenzene | 3.144 | 3.178 | -1.1 | 66 | -0.01 |
| 67 T | m- & p-Xylene | 2.101 | 2.126 | -1.2 | 66 | -0.02 |
| 68 T | Bromoform | 0.464 | 0.526 | -13.4 | 69 | -0.01 |
| 69 T | Styrene | 1.814 | 1.859 | -2.5 | 67 | -0.01 |
| 70 T | o-Xylene | 2.261 | 2.238 | 1.0 | 66 | -0.01 |
| 71 T | n-Nonane | 1.811 | 1.734 | 4.3 | 63 | -0.01 |
| 72 T | 1,1,2,2-Tetrachloroethane | 1.079 | | 0.4 | 65 | -0.02 |
| 73 S | Bromofluorobenzene (SS3) | 0.771 | 0.795 | -3.1 | 67 | 0.00 |
| 74 T | Cumene | 2.869 | 2.986 | -4.1 | 67 | |
| 75 T | alpha-Pinene | 1.525 | 1.549 | -1.6 | 66 | -0.01 |
| 76 T | n-Propylbenzene | 3.839 | 4.034 | -5.1 | 67 | -0.01 2 |

13041408.M Thu May 08 10:14:37 2008

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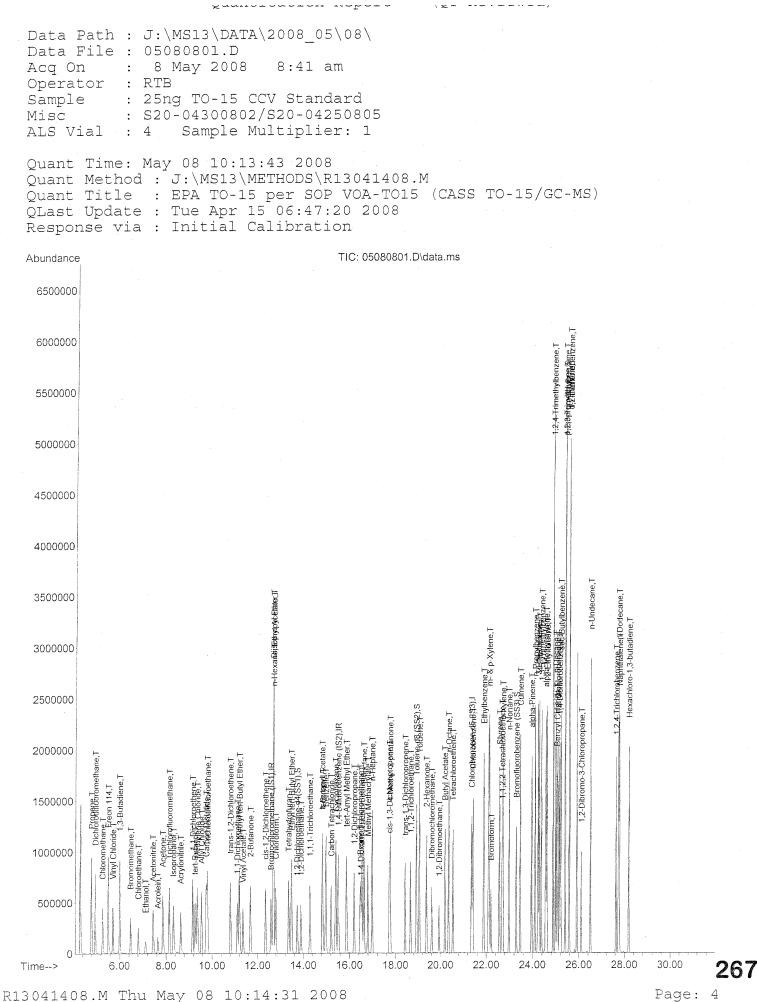
Page: 2

265

waraace concentanty caresesered tope-Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080801.D Acq On : 8 May 2008 8:41 am Operator : RTB Sample: 25ng TO-15 CCV StandardMisc: S20-04300802/S20-04250805 ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 10:13:43 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min Max. RRF Dev : 30% Max. Rel. Area : 200% AvgRF CCRF %Dev Area% Dev(min) Compound 77 T3-Ethyltoluene3.1283.222-3.067-0.0178 T4-Ethyltoluene2.8802.995-4.067-0.0179 T1,3,5-Trimethylbenzene2.5582.618-2.3670.0080 Talpha-Methylstyrene1.3551.421-4.967-0.0181 T2-Ethyltoluene3.1453.254-3.567-0.0182 T1,2,4-Trimethylbenzene2.8902.895-0.264-0.0283 Tn-Decane1.5981.610-0.866-0.0284 TBenzyl Chloride1.9702.266-15.067-0.0285 T1,3-Dichlorobenzene1.4971.545-3.269-0.0286 T1,4-Dichlorobenzene1.4971.545-3.269-0.0287 Tsec-Butylbenzene3.3943.547-4.567-0.0188 Tp-Isopropyltoluene2.9593.109-5.164-0.0189 T1,2.3-Trimethylbenzene1.6041.5771.765-0.0190 T1,2-Dichlorobenzene1.6041.5771.765-0.0191 Td-Limonene1.3111.2335.961-0.0192 T1,2.4-Trichlorobenzene0.4250.511-20.2720.0093 Tn-Undecane1.6781.689-0.7650.0094 T1,2.4-Trichlorobenzene0.9941.036-4.2690.00</ ______

(#) = Out of Range SPCC's out = 0 CCC's out = 0

F05/08/08



Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080801.D Acq On : 8 May 2008 8:41 am Operator : RTB Sample : 25ng TO-15 CCV Standard Misc : S20-04300802/S20-04250805 ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 10:13:43 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration Internal Standards R.T. QIon Response Conc Units Dev(Min) _____ 1) Bromochloromethane (IS1)12.5913024247725.000 ng-0.0237) 1,4-Difluorobenzene (IS2)15.52114104841125.000 ng-0.0156) Chlorobenzene-d5 (IS3)21.358250073025.000 ng0.00

 System Monitoring Compounds

 33) 1,2-Dichloroethane-d4(...
 13.73
 65
 443692
 22.817 ng
 -0.02

 Spiked Amount
 25.000
 Recovery
 =
 91.28%

 57) Toluene-d8 (SS2)
 18.93
 98
 1145691
 25.527 ng
 0.00

 Spiked Amount
 25.000
 Recovery
 =
 102.12%

 73) Bromofluorobenzene (SS3)
 23.29
 174
 397980
 25.767 ng
 0.00

 Spiked Amount
 25.000
 Recovery
 =
 103.08%

 System Monitoring Compounds

 Spiked Amount
 25.000
 Recovery
 Lint

 Target Compounds
 Qvalue

 2) Propene
 4.79
 42
 451017
 22.473 ng
 91

 3) Dichlorodifluoromethane
 4.95
 85
 774448
 21.67 ng
 100

 4) Chloromethane
 5.27
 50
 630543
 20.621 ng
 97

 5) Freon 114
 5.52
 135
 406287
 22.679 ng
 99

 6) Vinyl Chloride
 5.72
 62
 581139
 20.421 ng
 96

 7) 1, 3-Butadiene
 6.48
 94
 317718
 23.255 ng
 97

 9) Chloroethane
 6.48
 64
 317718
 23.255 ng
 97

 10 Ethanol
 7.12
 45
 253592
 18.529 ng
 97

 11 Acetonitrile
 7.43
 41
 742494
 20.655 ng
 97

 12 Acrolein
 7.64
 56
 32779
 24.383 ng
 65

 14
 Trichlorofluoromethane
 8.14
 101
 665952
 23.314 ng
 94

 16 Acrylonitrile
 8.63
 53
 510483
 24.511 ng
 98< Recovery = 103.08% 🗸 Spiked Amount 25.000 ⁹¹268

13041408.M Thu May 08 10:14:30 2008

Quantituation report (AT VENTEMEN) Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080801.D Acq On : 8 May 2008 8:41 am Operator : RTB Sample: 25ng TO-15 CCV StandardMisc: S20-04300802/S20-04250805 ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 10:13:43 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Tue Apr 15 06:47:20 2008 Response via : Initial Calibration

 Internal Standards
 R.T. Qion Response Conc Units Dev(Min)

 32) Chloroform
 12.60
 83
 639410
 28.244 ng
 ng
 95

 34) Tetrahydrofuran
 13.35
 72
 242667
 25.775 ng
 95

 35) Ethyl tert-Butyl Ether
 13.46
 87
 387048
 23.824 ng
 # 00

 35) Icopropyl Acctate
 14.83
 61
 230639
 24.192 ng
 # 35

 40) I-Bucanol
 14.84
 56
 308652
 21.634 ng
 95

 41) Benzene
 15.41
 84
 53345
 25.92 ng
 # 79

 43) Cyclohexane
 15.41
 84
 53345
 25.92 ng
 # 79

 43) Cyclohexane
 16.46
 83
 507625
 28.297 ng
 92

 44) Isooctane
 16.64
 93
 507625
 28.297 ng
 92

 49) Isooctane
 16.62
 57
 1570604
 23.854 ng
 75

 50) Methyl Methacrylate
 16.60
 100
 134316
 26.593 ng
 # 75

 51) r.-Heptane
 17.77
 58
 382318
 21.131 ng
 86

 52) clar.1.3.-Dic Internal Standards R.T. QIon Response Conc Units Dev(Min) ⁹²69

.13041408.M Thu May 08 10:14:30 2008

| × • • • • • • • • • • • • • • • • • • • | <u>P</u> | \ × | | ~ / | | |
|--|--|---|---|--|---|--|
| Data Path : J:\MS13\DATA\2008_05 Data File : 05080801.D Acq On : 8 May 2008 8:41 a Operator : RTB Sample : 25ng TO-15 CCV Stand Misc : S20-04300802/S20-042 ALS Vial : 4 Sample Multiplie | m .ard 50805 | | | | | |
| Quant Time: May 08 10:13:43 2008 Quant Method : J:\MS13\METHODS\R Quant Title : EPA TO-15 per SOP QLast Update : Tue Apr 15 06:47: Response via : Initial Calibrati Internal Standards | 1304140 VOA-TO 20 2008 on | 015 (C | | | its De | ev(Min) |
| <pre>80) alpha-Methylstyrene 81) 2-Ethyltoluene 82) 1,2,4-Trimethylbenzene 83) n-Decane 84) Benzyl Chloride 85) 1,3-Dichlorobenzene 86) 1,4-Dichlorobenzene 87) sec-Butylbenzene 88) p-Isopropyltoluene 89) 1,2,3-Trimethylbenzene 90) 1,2-Dichlorobenzene 91) d-Limonene 92) 1,2-Dibromo-3-Chloropr 93) n-Undecane 94) 1,2,4-Trichlorobenzene 95) Naphthalene</pre> | 24.56 24.61 24.88 24.98 25.05 25.15 25.21 25.40 25.41 25.58 25.58 26.11 26.50 27.63 27.77 27.74 | 118 105 57 91 146 146 105 119 105 146 68 157 57 180 128 57 | 725626 1616291 1594582 838565 1216575 841352 850779 1903738 1836968 1576581 853071 654184 265979 889951 580788 1842526 873340 | 26.745 m 25.662 m 27.543 m 26.192 m 30.828 m 26.844 m 28.366 m 28.008 m 30.990 m 27.692 m 26.558 m 24.921 m 31.256 m 26.477 m 29.175 m 28.787 m 25.609 m | a a a a a a a a a a a a a a a a a a a | 96 97 97 94 99 98 96 92 97 100 97 100 94 70 87 96 99 85 |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

205/08/0¢

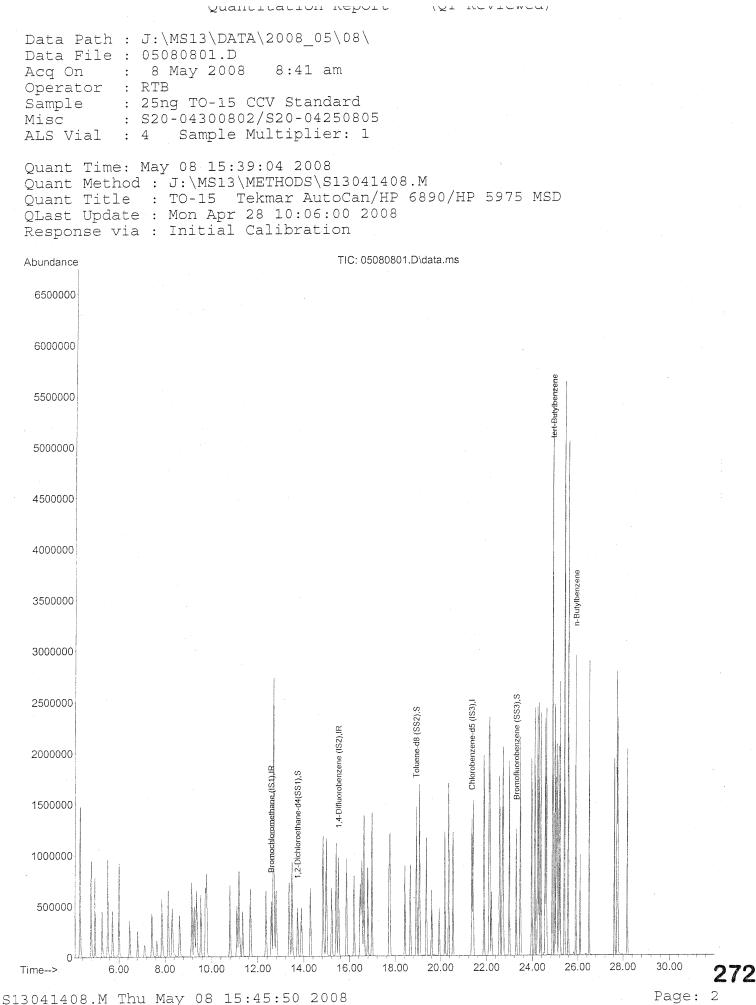
waraaco oonornanig oarraractor neper-Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080801.D Acq On : 8 May 2008 8:41 am Operator : RTB Sample : 25ng TO-15 CCV Standard Misc : S20-04300802/S20-04250805 ALS Vial : 4 Sample Multiplier: 1 Ouant Time: May 08 15:39:04 2008 Quant Method : J:\MS13\METHODS\S13041408.M Quant Title : TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD QLast Update : Mon Apr 28 10:06:00 2008 Response via : Initial Calibration Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.33min Max. RRF Dev : 30% Max. Rel. Area : 200% AvgRF CCRF %Dev Area% Dev(min) Compound -1 IRBromochloromethane (IS1)1.0001.0000.073-0.022 S1,2-Dichloroethane-d4(SS1)2.0051.8308.766-0.02 3 IR 1,4-Difluorobenzene (IS2) 1.000 1.000 0.0 71 -0.01 4 IChlorobenzene-d5 (IS3)1.0001.0000.0660.005 SToluene-d8 (SS2)2.2412.288-2.1680.006 SBromofluorobenzene (SS3)0.7710.795-3.1670.007tert-Butylbenzene2.7382.753-0.564-0.018n-Butylbenzene2.9543.108-5.267-0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

313041408.M Thu May 08 15:49:05 2008

F15708/08



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Zaamereacton metore Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080801.D Acq On : 8 May 2008 8:41 am Operator : RTB Sample : 25ng TO-15 CCV Standard Misc : S20-04300802/S20-04250805 ALS Vial : 4 Sample Multiplier: 1 Quant Time: May 08 15:39:04 2008 Quant Method : J:\MS13\METHODS\S13041408.M Quant Title : TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD QLast Update : Mon Apr 28 10:06:00 2008 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Internal Standards _______ 1) Bromochloromethane (IS1)12.5913024247725.000 ng-0.023) 1,4-Difluorobenzene (IS2)15.52114104841125.000 ng-0.014) Chlorobenzene-d5 (IS3)21.358250073025.000 ng0.00 System Monitoring Compounds 2) 1,2-Dichloroethane-d4(... 13.73 65 443692 22.817 ng -0.02

 Spiked Amount
 25.000
 Recovery = 91.28% /

 5) Toluene-d8 (SS2)
 18.93
 98
 1145691
 25.527 ng
 0.00

 Spiked Amount
 25.000
 Recovery = 102.12% /

 6) Bromofluorobenzene (SS3)
 23.29
 174
 397980
 25.767 ng
 0.00

 Recovery = 103.08% 🗸 Spiked Amount 25.000 Ovalue Target Compounds 7) tert-Butylbenzene24.88119143345726.143ng8) n-Butylbenzene25.9191166851628.204ng 99 94 _____

(x= ======,

(#) = qualifier out of range (m) = manual integration (+) = signals summed

205/08/08

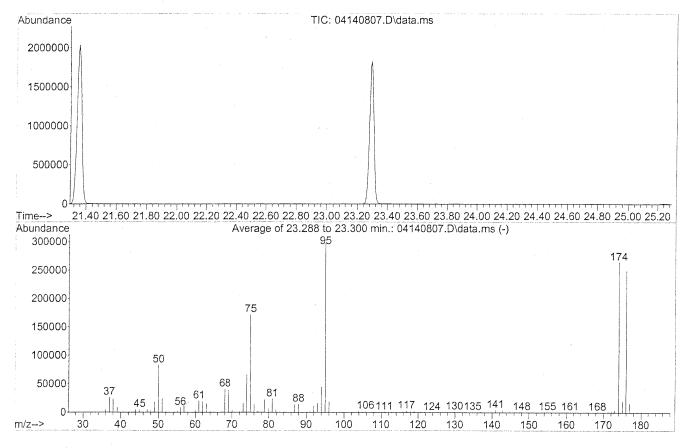
273

BFB TUNING & MASS CALIBRATIONS

Data Path : J:\MS13\DATA\2008_04\14\ Data File : 04140807.D Acq On : 14 Apr 2008 18:18 Operator : WA Sample : BFB Tune Standard (200ml) Misc : S20-04140804 ALS Vial : 4 Sample Multiplier: 1

Integration File: RTEINT.P

Method : J:\MS13\METHODS\R13041408.M Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) Last Update : Tue Apr 15 06:47:20 2008



AutoFind: Scans 3374, 3375, 3376; Background Corrected with Scan 3363

| Target Mass | Rel. to Mass | Lower Limit% | Upper Limit% | Rel. Abn% | Raw Abn | Result Pass/Fail |
|----------------|-----------------|-----------------|-----------------|--------------|------------|---------------------|
| 50 | 95 | 8 | 40 | 27.9 | 83021 | PASS |
| 75 | 95 | 30 | 66 | 57.6 | 171669 | PASS |
| 95 | 95 | 100 | 100 | 100.0 | 297792 | PASS |
| 96 | 95 | 5 | 9 | 6.3 | 18818 | PASS |
| 173 | 174 | 0.00 | 2 | 1.5 | 4055 | PASS |
| 174 | 95 | 50 | 120 | 88.7 | 264171 | PASS |
| 175 | 174 | 4 | 9 | 7.6 | 20123 | PASS |
| 176 | 174 | 93 | 101 | 94.5 | 249664 | PASS |
| 177 | 176 | 5 | 9 | 6.3 | 15749 | PASS |

R13041408.M Thu Apr 17 09:03:26 2008

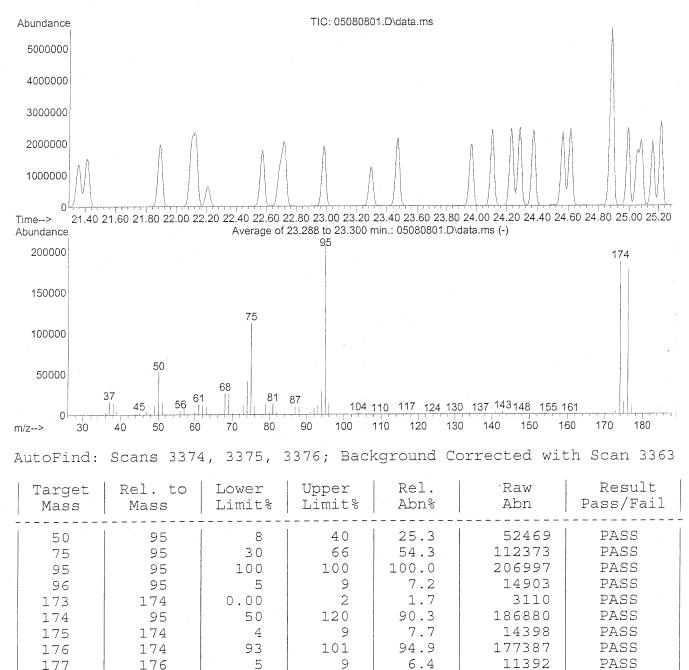
DH 4/17/08

275

Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080801.D 8 May 2008 8:41 am Acq On : Operator : RTB : 25ng TO-15 CCV Standard Sample : S20-04300802/S20-04250805 Misc Sample Multiplier: 1 : 4 ALS Vial

Integration File: RTEINT.P

```
Method
          : J:\MS13\METHODS\R13041408.M
          : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS)
Title
Last Update : Tue Apr 15 06:47:20 2008
```



R13041408.M Thu May 08 10:13:34 2008

177

F05/08/05

276

RUN LOGS

| 1.97 | 45 | 1.00 |
|------|-----|------|
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| ··· U | Comment | leiV | Operator | Misc Info | al alqms2 | emsN elia | amiT\ats0 | |
|--|--|----------|----------|----------------------------|----------------------------------|-------------|------------------|-------|
| | | 4 | AW | 220-04140804 | BFB Tune Standard (200ml) | G.70804140 | | |
| | | 9 | AW | 250-04140804/250-04030801 | | | 81:81 80/41/40 | |
| /70 67 | राज्याम्लः स सिर्ह्याप्त वर | 13- | AW | 250-04140804/250-03510806 | 0.5nd TO-15 ICAL Standard | G.80804140 | 65:81 80/41/40 | |
| | | | AW | | 0.5ng TO-15 ICAL Standard | G.00804140 | | |
| | HOLD AL AL ALE ALCHING TO ALL | 13 El | | 90801250-022/40804140-022 | TADI St-OT put | 0.01804140 | 12:02 80/41/40 | |
| | alt so to the state | 13 | AW AW | S20-04140804/S20-03210809 | 5ng TO-15 ICAL Standard | G.11804140 | 10:12 80/11/0 | |
| | | * | AW | S20-04140804/S20-04020808 | bisbrief JAOI 21-OT gn2S | G.21804140 | 54:12 80/41/40 | |
| | 1947 | * | AW | | 50ng TO-15 ICAL Standard | 0.21804140 | 04/14/08 22:24 | |
| | | 91 | AW | S20-04140804/S20-04020808 | Toppets VOI at OT product | C.41804140 | | |
| Commentation and a commentation of the second se | frezer all (milage | 4 | AW | S20-04140804/S20-04040804 | 25ng TO-15 ICV Standard | 0.31804140 | 04/14/08 23:45 | |
| | | 3 | AW | 1891 | (Im002) /0802/04022 | D.91804140 | 92:0 80/31/40 | |
| | · []:::::::::::::::::::::::::::::::::::: | 3 | AW | 1651 | (1m25) 10807120-041 | 0.71804140 | 70:1 80/31/40 | |
| | 4 Marine 1997 - | 3 | AW | 1201 | 250-04120804 (220MI) | G.81804140 | 84:1 80/31/10 | |
| | 294. 87. 98 [°] | <u>۱</u> | AW | fest Test | | G.07804140 | 62:2 80/31/40 | |
| | | ۴ | AW |] 189⊥ | 250-04120805 (220ml) | 0.12804140 | | |
| | | | | 100 | (10007) 7000 / 10-070 | D.12804140 | 75:9 80/21/40 | 12 |
| 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - | | L | AW | S20-04140804/S20-04140802 | 25ng MAPH CCV Standard | Q.10802140 | 85:7 80/31/40 | 1 |
| | | L | AW | S20-04140804 | (Im0001) Ansi8 botteM H9AM/81-OT | 0.20802140 | 04/15/08 9:33 | z |
| and the second | | 5 | AW | S20-04140804/S20-04140803 | 25ng MAAM bras | G.E0803140 | 41:01 80/S1/40 | 3 (|
| | | 8 | AW | Test | (1m001) 500-79600809 | 0.40802140 | 98:01 80/91/40 | * (|
| | | 9 | AW | (8,5,8,0-) | (Im0001) 100-78600809 | 0.20802140 | 75:11 80/21/40 | 2 (|
| | | L | AW | Test | (Im0001) 200-78600809 | D.30802140 | 81:21 80/21/40 | 9 |
| | | 6 | AW | (S.E., e. h- | (IM0001) 400-73600809 | D.70802140 | 10:51 80/31/40 |) [] |
| | | 01 | AW | (j. 1. 5, 3. 9) | (Im0001) 200-73600809 | Q.80802140 | 04/12/08 13:42 | 0 8 |
| | | 9 | AW | (8.5, 8.0-) | (1m0001) AUG 100-78600809 | Q.60803140 | 04/12/08 14:23 | 5 6 |
| | | 11 | AW | (7.6.,4.6.) | (Im0001) 900-79600809 | 0.01802140 | \$0:51 80/51/\$0 | 10 C |
| | | 15 | AW | (5.5.,5.5.) | (Im0001) 700-79600809 | 0.11802140 | 24:31 80/31/40 |) LL |
| | | 8 | AW | (9.2, 7.1) | (Im0001) 200-79600809 | 0.2150812.D | 82:91 80/91/70 | 15 C |
| | ે જિલ્લ | + * | втя | S20-04140804/S20-04020808 | 25ng TO-15 CCV Standard | 04150813.D | 04/16/08 17:29 | 13 (|
| | passa) | + * | ВТЯ | S20-04140804 | TO-15 Method Blank (1.0L) | 0.41508140 | 06:81 80/31/40 | 14 C |
| and a second | 1658#2W -17 P25502 - | + + L | втя | 867009VA\7620007\288000A | (J0.f) eguse/OF/nsO OD SAO | CI.21802140 | £1:61 80/31/40 |) St. |
| | | 2 | втя | 767003VA/47r000374/880000A | (J0.1) egusƏ/ƏAnsƏ OQ SAƏ | 0.9150816.D | 95:01 80/51/40 | 0 91 |
| | | E | RTB | 967009VA\82900071\462000A | (10.1) agusð/04/nsð 00 SAO | 0.71802140 | 04/15/08 20:39 | 0 21 |
| | | 9 | ВТЯ | S27009VA\76800031\F66000A | (10.1) eguse/C7/nsC CC 2AC | 04150818.D | 04/16/08 21:22 | 0 81 |
| and a second | | 9 | втя | AC00582/FC00540/AVG00588 | CAS QC Can/FC/Gauge (1.0L) | Q.61802140 | 90:22 80/91/70 | 0 61 |
| and a second | | 2 | втя | AC01136/FC00269/A/G00808 | CAS QC Can/FC/Gauge (1.0L) | 04150820.D | 84:22 80/21/40 | 0 0Z |
| ····· | | 8 | 878 | AC00292/FC00536/AVG00809 | (10.1) aguso/Can/FC/Gauge (1.0L) | C1.12803140 | 14/15/08 23:31 | 54 0 |
| the second s | | 6 | RTB | AC01273/FC00565/A\G00804 | (J0.1) aguso/C3n/FC/Cauge (J0.1) | 04150822.D | \$1:0 80/91/70 | o zz |
| and the second | | 01 | втя | AC00803/FC00676/AVG00800 | (10.1) 9gusÐ\JA\nsJ OQ SAD | 04150823.D | 29:0 80/91/70 | 53 C |
| and the second | | 11 | ВТЯ | AC01320/FC00548/AVG00777 | (10.1) 9gusð/07/ns0 00 2A0 | 0.4150824.D | 04:1 80/91/40 | 5¢ C |
| | | 15 | RTB | 85000VA/64200031637000A | (J0.1) agusa/CAnso OD SAO | 04160825.D | 04/16/08 2:23 | 52 0 |
| | | 13 | RTB | AC01157/FC00583/AVG00776 | CAS QC Can/FC/Gauge (1.0L) | | | 56 C |

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| | | · · · · · · · · · · · · · · · · · · · | 88 | 51 | 1991 | S20-03100803 (50mL) | 0.a1808020 | 0:8 80/60/50 | 0 91 |
| | | | | 51 | 159_ | Blank (100mL) S20-03100803 (50mL) | | | |
| | | | 88 88 88 | | | | Q.21808020 | £2:7 80/60/20 | 15 1 |
| | | | ая | 4 | JzeT | Blank (100mL) | 0.51808080 0.51808080 | 02\09\08 \:23 | 12 10 |
| | | | 89 89 | * | JseT TesT | Biank (100mL) Biank (100mL) | 0.51808020 0.41808020 0.21808020 | 05/08/08/20 5:7 80/80/20 5:7 80/80/20 | 42 (47 (43 (|
| | | ≠235 0 J | 88 88 88 | 7 7 7 | r Test Test | Blank (100mL) Blank (100mL) DIL (10mL) | 0.51808080 0.51808080 0.41808080 0.41808080 0.51808080 0.51808080 | 22:3 80/80/20 92:08/08 2:54 5:7 80/80/20 5:08/08 2:5 | 12 (C |
| | | م م ا ا ا ا ا ا ا ا ا ا ا ا ا ا ا ا ا ا | 88 88 88 | * * * | (2.5.2.) (2.5.2.) (2.5.35.) JseT JzeT | P0801305-001 DIL (10mL) P0801305-001 DIL (10mL) P0801305-001 DIL (10mL) P0801305-001 DIL (10.50mL) | G.11808020 G.21808020 G.21808020 G.21808020 G.21808020 | CZ:7 80\60\20 42:7 80\80\20 25:8 80\80\20 25:8 80\80\20 25:8 80\80\20 | 12 (13 (15 (15 (15 (11 (|
| | | \$232 nJ - | 92 92 92 92 92 | * * * * * | Test Test Test Test | P0801342-003 DUP (10mL) P0801305-001 DIL (10mL) Blank (100mL) Blank (100mL) | 0.01808070 0.51808020 0.51808020 0.51808020 0.51808020 0.51808020 0.51808020 | 22:7 80/80/20 42:7 80/80/20 25:8 80/80/20 41:2 80/80/20 55:4 80/80/20 55:4 80/80/20 | 12 (0 13 (0 13 (0 11 (0 11 (0 |
| | | | 82 87 87 87 87 87 87 87 87 87 87 87 87 87 | | ENSR SG838-05-7 (-4.2.3.5) ENSR SG838-05-7 (-4.2.3.5) Test Test | P0801342-003 DUP DIL (10mL) P0801305-001 DIL (0.50mL) P0801305-001 DIL (0.50mL) P0801305-001 DIL (0.50mL) P0801305-001 DIL (0.50mL) | G.01808020 G.01808020 G.11808020 G.51808020 G.51808020 G.51808020 G.21808020 | 25:2 80/80/50 25:2 80/80/50 20/80/50 20/80/50 20/80/50 20/80/50 | 12 (C 13 (C 12 (C)))))))))))))))))))))))))))))))))))) |
| | | | 88 88 88 88 88 88 88 88 88 88 88 88 88 | t t t t t t t t t t t t | Test ENSR SG83B-05-7 (-4, 2, 3, 5) ENSR SG83B-05-7 (-4, 2, 3, 5) ENSR SG83B-05-7 (-4, 2, 3, 5) Test Test | Blank (100mL) Blank (100mL) P0801342-003 DUP (10mL) P0801342-003 DUP (10mL) P0801342-003 DUP (10mL) P0801342-003 DUP (10mL) P0801342-003 DUP (1mL) | 0.8080808 0.5080809 0.5080804 0.51808080 0.51808080 0.51808080 0.51808080 0.51808080 0.51808080 0.51808080 0.51808080 0.508080 0.50808080 0.508080000000000 | C:: 2 80/80/50 C:: 2 80/80/50 | 12 (0 13 (0 15 (0 15) (0 15 (0 15) (0 15 (0 15) (0)) (0)) (0)) (0)) (0)) (0)) (0)) (0 |
| | | | 88 88 88 88 88 88 88 88 88 88 88 88 88 | v | Test ENSR SG83B-05-7 (-4,2,3,5) ENSR SG83B-05-7 (-4,2,3,5) ENSR SG83B-05-7 (-4,2,3,5) ENSR SG83B-05-7 (-4,2,3,5) ENSR SG83B-05-7 (-4,2,3,5) Test | אפטנואל (100mL) 10801342-003 DIL (10mL) 10801342-003 DIL (10mL) 10801342-003 DIL (10mL) 10801342-003 DIL (10mL) 10801342-003 DIL (10mL) 10801342-003 DIL (1mL) 10801342-003 DIL (1mL) | G. 50808020 G. 50808020 G. 51808020 G. 51808020 G. 51808020 G. 51808020 G. 51808020 G. 51808020 G. 51808020 | 05/08/08 2:30 05/08/08 2:22 02/08/08 2:22 02/08/08 2:22 02/08/08 2:22 02/08/08 2:23 02/08/08 2:23 02/08/08 2:32 02/08/08 2:32 | 12 0 12 0 12 0 12 0 12 0 12 0 12 0 12 0 |
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| | | | 82 82 82 82 82 82 82 82 82 82 82 82 82 8 | t t t t t t t t t t t t t t t t | ENGR SG33B-05-7 (-4.2.3.5) ENGR SG33B-05-1 (-3.6.3.5) ENGR SG33B-05-7 (-4.2.3.5) ENGR SG33B-05-7 (-4.2.3.5) ENGR SG33B-05-7 (-4.2.3.5) ENGR SG33B-05-7 (-4.2.3.5) Test Test | פוסטוניקב-2003 (10mL) Blank (100mL) הספטוניקב-2003 בוור (ניניתר) הספטוניקב-2003 בוור (ניניתר) הספטוניקב-2003 בוור (ניתר) הספטוניקב-2003 בוור (ניתר) הספטוניקב-2003 בוור (ניתר) בוור (ניתר) הייניקב-2003 בוור (ניתר) הייניקב-2003 בוור (ניתר) הייניקב-2003 נוור (ניתר) הייניק | G \$1808050 G \$1808050 G \$1808050 G \$1808050 G \$1808050 G \$1808050 G \$0808050 G \$0808050 G \$0808050 G \$0808050 G \$0808050 | 22:2 80/80/92 22:2 80/80/92 20:08/08 22 20:08/08 22 20:08/08 22 20:08/08 22 20:08/08 22 20:08/08 20 20:2 80/80/50 20:2 80/80/50 20:2 80/80/50 20:2 80/80/50 | 12 0 12 0 14 0 15 0 15 0 15 0 15 0 15 0 15 0 15 0 15 |
| | | - Fussed | 82 82 82 82 82 82 82 82 82 82 82 82 82 8 | t | ENGR SG838-05-3 (-3.8, 3.5) ENGR SG838-05-7 (-4.2, 3.5) Test Test | Blank (100mL) P0801342-002 (10mL) P0801342-003 DLP (10mL) P0801342-003 DLP (10mL) P0801342-003 DLP (10mL) P0801342-003 DLP (1mL) P0801342-003 DLP (1mL) P0801342-003 DLP (1mL) P0801342-003 DLP (1mL) P0801342-003 (10mL) P0801342-002 (10mL) P0801342-002 (10mL) P0801342-002 (10mL) | G \$1 808050 G \$1 808050 G \$1 808050 G \$1 808050 G \$1 808050 G \$1 808050 G \$0808050 G \$0808050 G \$0808050 G \$0808050 G \$0808050 G \$0808050 G \$0808050 | C2:2 80/80/50 C2:2 80/ | 12 (0 14 (0 15 (0 14 (0 15 (0 10 (0 14 (0)))))))))))))))))))))))))))))))))))) |
| | | разли] - | 82 82 82 82 82 82 82 82 82 82 82 82 82 8 | t | ENSR SG838-05-1 (-3.6, 3.5) ENSR SG838-05-7 (-4.2, 3.5) <td>P0801342-001 (10mL) P0801342-002 (10mL) P0801342-002 (10mL) P0801342-003 DUP DIL (1mL) P0801342-003 D</td> <td>G \$1 808050 G \$0808050 G \$0808050 G \$0808050 G \$0808050 G \$0808050 G \$0808050 G \$0808050</td> <td>20:01 80/80/50 20:01 80/80/50</td> <td>12 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 0 11 0 0 11 0 0 0 11 0 0 0 11 0</td> | P0801342-001 (10mL) P0801342-002 (10mL) P0801342-002 (10mL) P0801342-003 DUP DIL (1mL) P0801342-003 D | G \$1 808050 G \$0808050 G \$0808050 G \$0808050 G \$0808050 G \$0808050 G \$0808050 G \$0808050 | 20:01 80/80/50 20:01 80/80/50 | 12 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 11 0 0 0 11 0 0 11 0 0 0 11 0 0 0 11 0 |
| | | | 92 92 92 92 92 92 92 92 92 92 92 92 92 9 | t | Test Test Test Test Test Test Test Test Test Test Test | Sang TO-16 LOS Poso1342-001 (10mL) Poso1342-001 (10mL) Poso1342-003 DIL (10mL) Poso1342-003 DIL (10mL) Poso1342-003 DIL (1mL) Poso1342-003 DIL (10mL) Poso1342-003 DIL (10mL) Poso1342-00 | G 20806020 G 20806020 G 2000000 G 20000000 G 200000000 G 200000000 G 200000000 G 200000000 G 200000000 G 200000000 G 200000000 G 200000000 G 200000000 G 2000000000 G 20000000000 | x2:6 80/80/20 x2:7 80/80/20 x2:7 80/80/20 x2:8 80/80/20 x2:9 80/80/20 x2:1 80/80/20 x2:2 80/80/20 x2:2 80/80/20 x2:5 80/80/20 x2:6 80/80/20 x2:7 80/80/20 | 2 0 0 2 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 |
| | | Þ∂5502 - Þ∂5502 - Þ∂5502 - | 92 92 92 92 92 92 92 92 92 92 92 92 92 9 | t | Test Test Test Test Test Test Test Test Test Test Test Test | TO-15 MadiMa Damia (1,0L) Sang TO-15 LCS Posorst-2001 (10mL) Posorst-2003 (10mL) Posorst-2003 (10mL) Posorst-2003 DIL (1mL) Posorst-2003 | G 20806020 G 20806020 G 2000000 G 20000000 G 200000000 G 200000000 G 200000000 G 200000000 G 200000000 G 200000000 G 200000000 G 200000000 G 200000000 G 2000000000 G 20000000000 | \$2.6 80/80/20 \$2.6 80/80/20 \$2.5 80/ | 2 0 0 2 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 |
| | | | 92 92 92 92 92 92 92 92 92 92 92 92 92 9 | t | Test Test Test Test Test Test Test Test Test Test Test Test | ۲۵-۱5 CCV Standard الا الحالة الحالة الحالة (10.1) الا الحالة الحالة الحالة (10.1) الا الحالة الحالة الحالة (10mL) الحالة (10mL) الحال | G.10808020 G.10808020 G.2080800 G.20808000 G.20808000 G.20808000 G.208080000 G.208080000 G.20808000000000000000000000000000000000 | L+:6 80/80/20 L+:6 80/ | 117 (117 (11 |
| | | | 83 83 83 93 93 93 93 93 93 93 93 93 93 93 93 93 | t | Test Test Test Test Test Test Test Test Test Test Test Test | S20-03100803 (50mL) S50-03100803 (50mL) 25ng TO-15 CCV Standard P0801342-001 (10mL) P0801342-003 DLL (1mL) P0801342-003 DLL (1mL) P0801342-003 DLL (1mL) P0801342-003 DLL (1mL) P0801342-003 DLP (1mL) P0801342-003 DLL (1mL) P0801342-003 DLP (1mL) P0801342-003 DLP (1mL) P0801342-003 DLL (1mL) P0801342-003 DLP (10mL) P0801342-003 DLP (10mL) P08 | G. 1 5 5 0 7 0 2 0 2 0 2 0 2 0 2 0 2 0 2 0 2 0 2 | +0:8 80/80/20 +2:2 80/80/20 +2:2 80/80/20 +2:2 80/80/20 +1:2 80/80/20 +1:2 80/80/20 +1:2 80/80/20 +0:1 80/20/20 50:01 80/20/20 +0:1 80/20/20 +0:1 80/20/20 +0:1 80/80/20 +0:2 | 112 0 0 113 0 0 114 0 115 0 11 |
| | | | 83 83 83 93 93 93 93 93 93 93 93 93 93 93 93 93 | t t | S20-04300802/S20-04250805 S20-04300802/S20-04250805 ENSR SG83B-05-7 (-4.2.3.5) ENSR SG83B-05-7 (-4.2.3.5) Tast Tast Tast | ۲۵-۱5 CCV Standard الا الحالة الحالة الحالة (10.1) الا الحالة الحالة الحالة (10.1) الا الحالة الحالة الحالة (10mL) الحالة (10mL) الحال | G. 1 5 5 0 7 0 2 0 2 0 2 0 2 0 2 0 2 0 2 0 2 0 2 | 22:7 80/80/20 22:7 80/80/20 22:7 80/80/20 22:2 80/80/20 20:05/05 20:05/ | 12 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 |
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| | | | 83 94 95 95 95 95 95 95 95 95 95 95 95 95 | t t t <td>Теят Теят Теят Емая Scasae-05-7 (-4.2, 3, 3, 5) Емая Scasae-05-7 (-4.2, 5) Емая Scasae-05-7 (-4.2, 5) Емая Sca</td> <td>المالة (100mL) المالة (100mL) Sac.03100803 (50mL) Sac.03100803 (50mL) Sac.03100803 (50mL) Sang TO-15 CCV Standard Poso1342-003 (10mL) Poso1342-003 DIL (1mL) Poso1342-003 DIL (1mL) Poso1342</td> <td>G 1+2202000 G 2+220200 G 2+220200 G 2+220200 G 2+220200 G 2000000 G 2000000 G 2000000 G 2000000 G 2000000 G 20000000 G 2000000 G 2000000 G 2000000 G 2000000 G 2000000 G 2000000 G 2000000 G 2000000 G 2000000 G 20000000 G 2000000 G 20000000 G 200000000000 G 200000000 G 2000000000000000000000000000000000000</td> <td>EX: 8 80/20/50 EX: 8 80/20/50 EX: 8 80/20/50 EX: 8 80/80/50 EX: 8 80/20/50 EX: 8 80/20/50</td> <td>112 112 112 113 114 114 114 115 114</td> | Теят Теят Теят Емая Scasae-05-7 (-4.2, 3, 3, 5) Емая Scasae-05-7 (-4.2, 5) Емая Scasae-05-7 (-4.2, 5) Емая Sca | المالة (100mL) المالة (100mL) Sac.03100803 (50mL) Sac.03100803 (50mL) Sac.03100803 (50mL) Sang TO-15 CCV Standard Poso1342-003 (10mL) Poso1342-003 DIL (1mL) Poso1342-003 DIL (1mL) Poso1342 | G 1+2202000 G 2+220200 G 2+220200 G 2+220200 G 2+220200 G 2000000 G 2000000 G 2000000 G 2000000 G 2000000 G 20000000 G 2000000 G 2000000 G 2000000 G 2000000 G 2000000 G 2000000 G 2000000 G 2000000 G 2000000 G 20000000 G 2000000 G 20000000 G 200000000000 G 200000000 G 2000000000000000000000000000000000000 | EX: 8 80/20/50 EX: 8 80/20/50 EX: 8 80/20/50 EX: 8 80/80/50 EX: 8 80/20/50 EX: 8 80/20/50 | 112 112 112 113 114 114 114 115 114 |