(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									
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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		
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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)
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Relinquished by: (Signature)			Date:		Received by: (Signature)	(Signature)			Date:	Time:	Cooler / Blank	ank
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		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
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	Were seals i	intact?								\mathbf{X}
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	Were signat	ture and date included	1?							X
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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}
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	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}
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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	

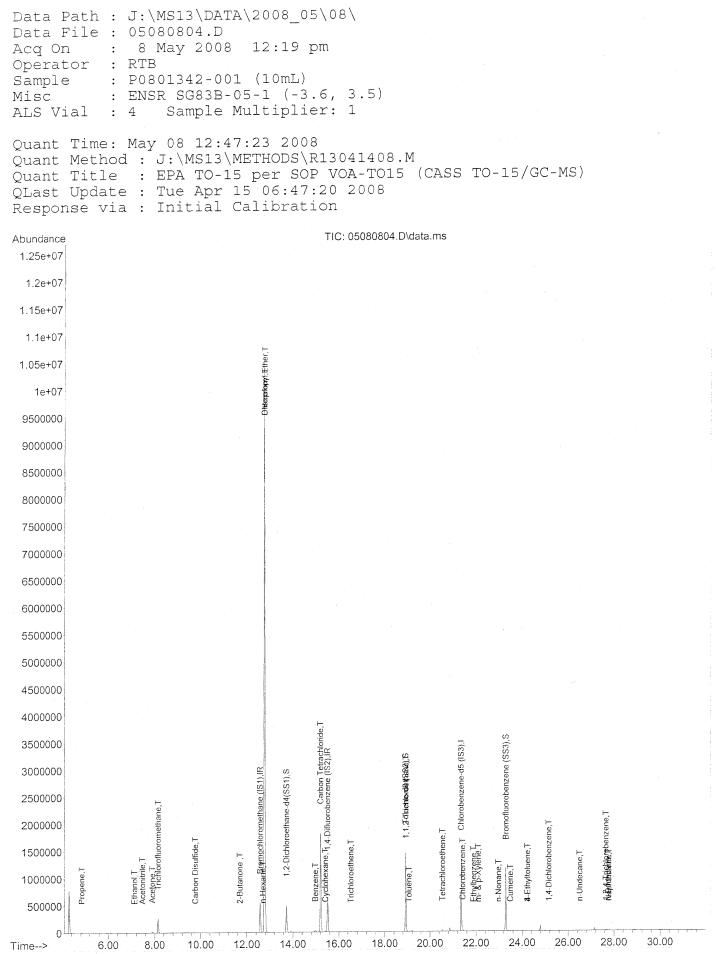
ND = Compound was analyzed for, but not detected above the laboratory detection limit.

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

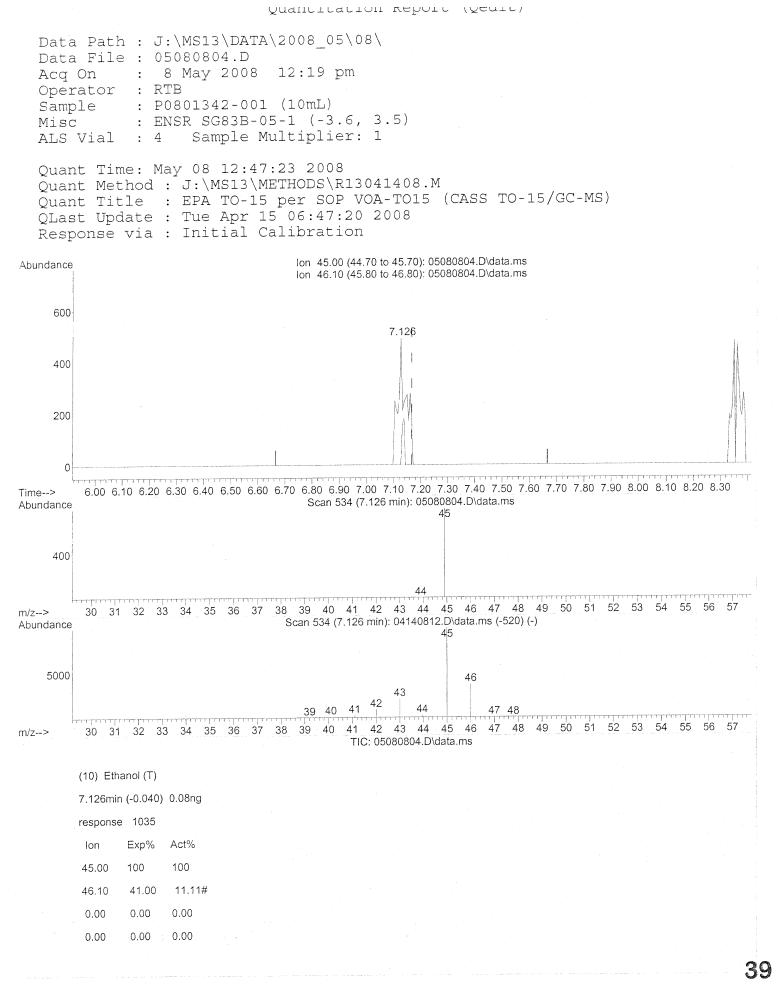
37

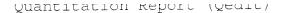
Data Path : J:\MS13\DATA\2008_0 Data File : 05080804.D Acq On : 8 May 2008 12:19 Operator : RTB Sample : P0801342-001 (10mL) Misc : ENSR SG83B-05-1 (-3 ALS Vial : 4 Sample Multipli	pm .6, 3.5)							
Quant Time: May 08 12:47:23 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per 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			
		~~~~~~~						

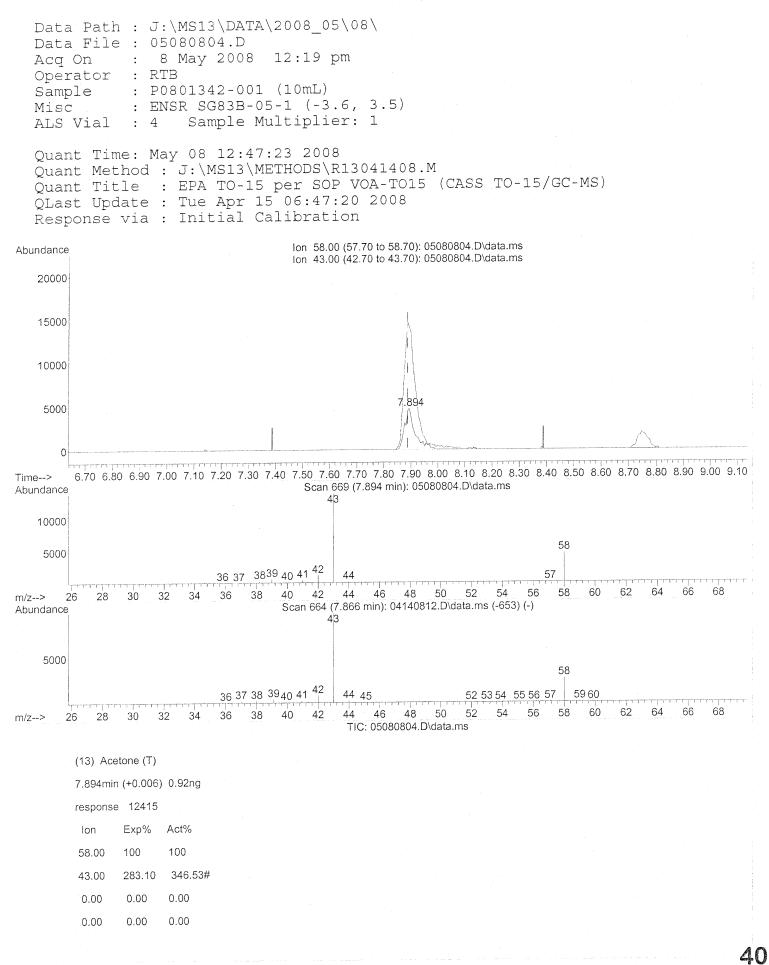
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Fostoglog

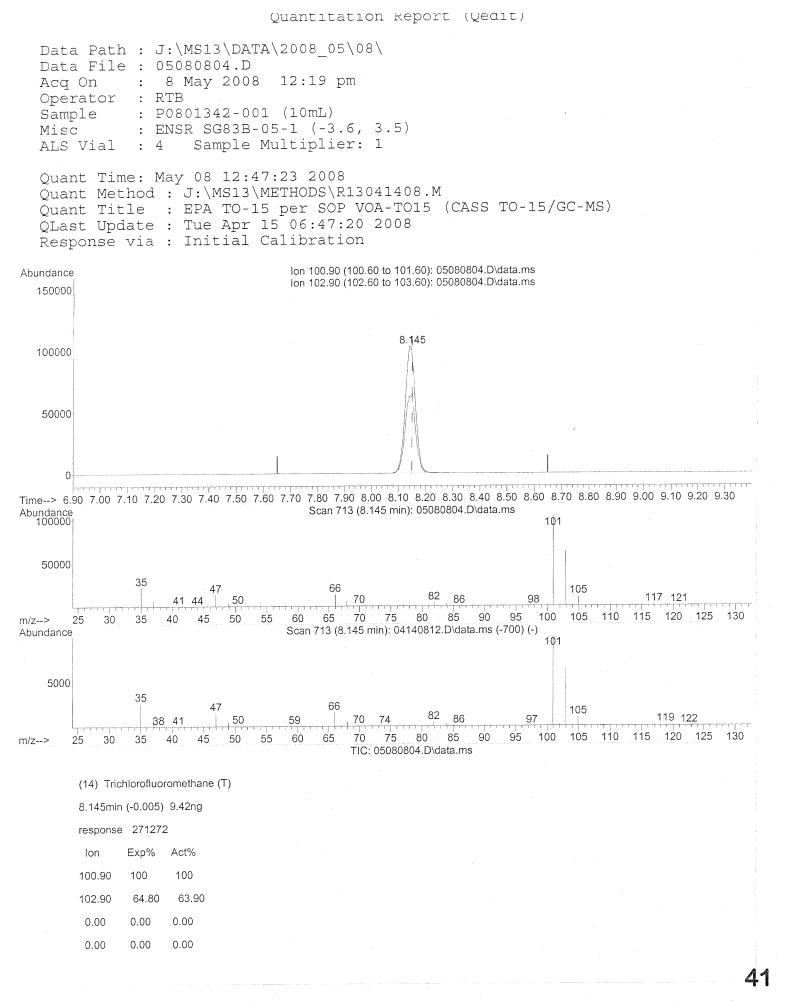
38

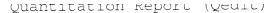


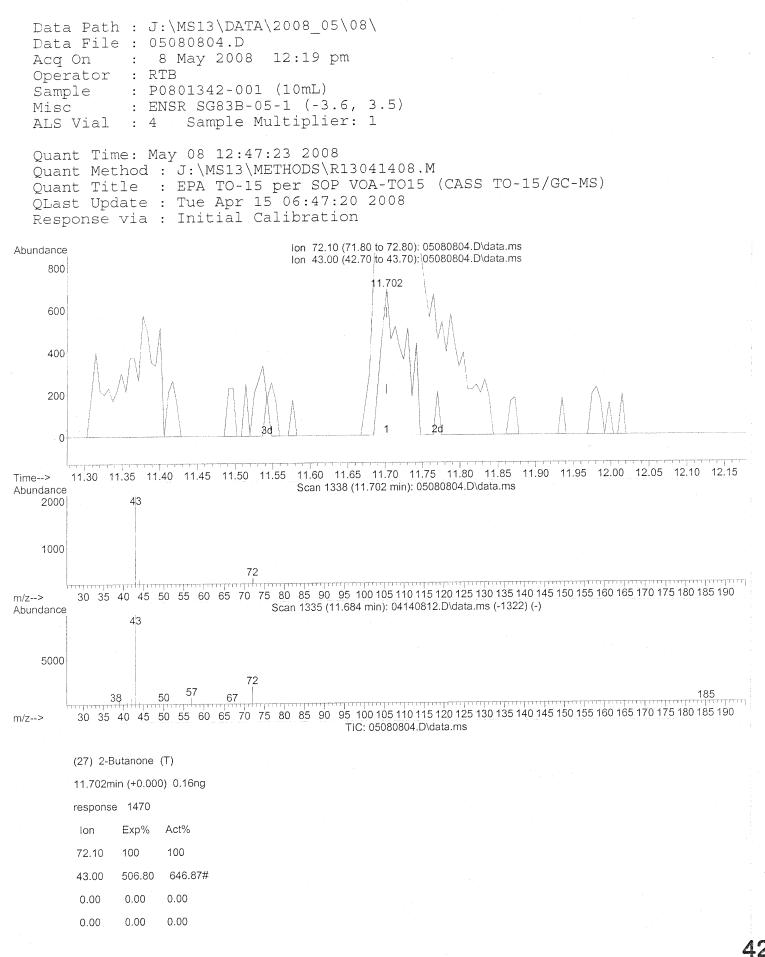




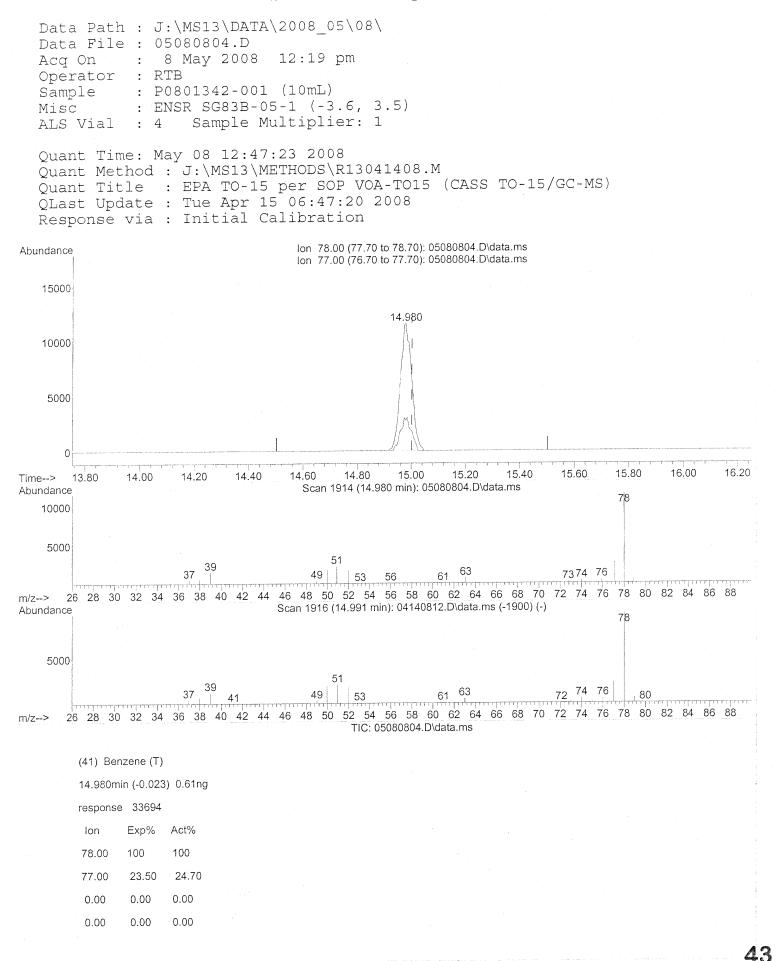
R13041408.M Thu May 08 16:23:52 2008



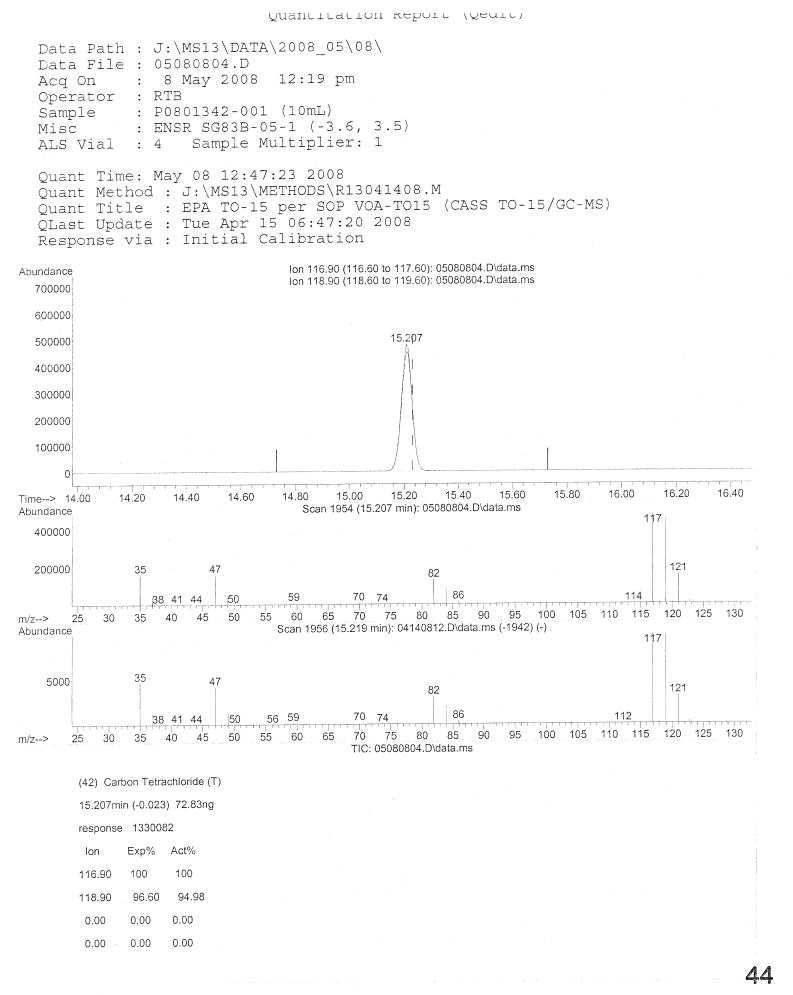


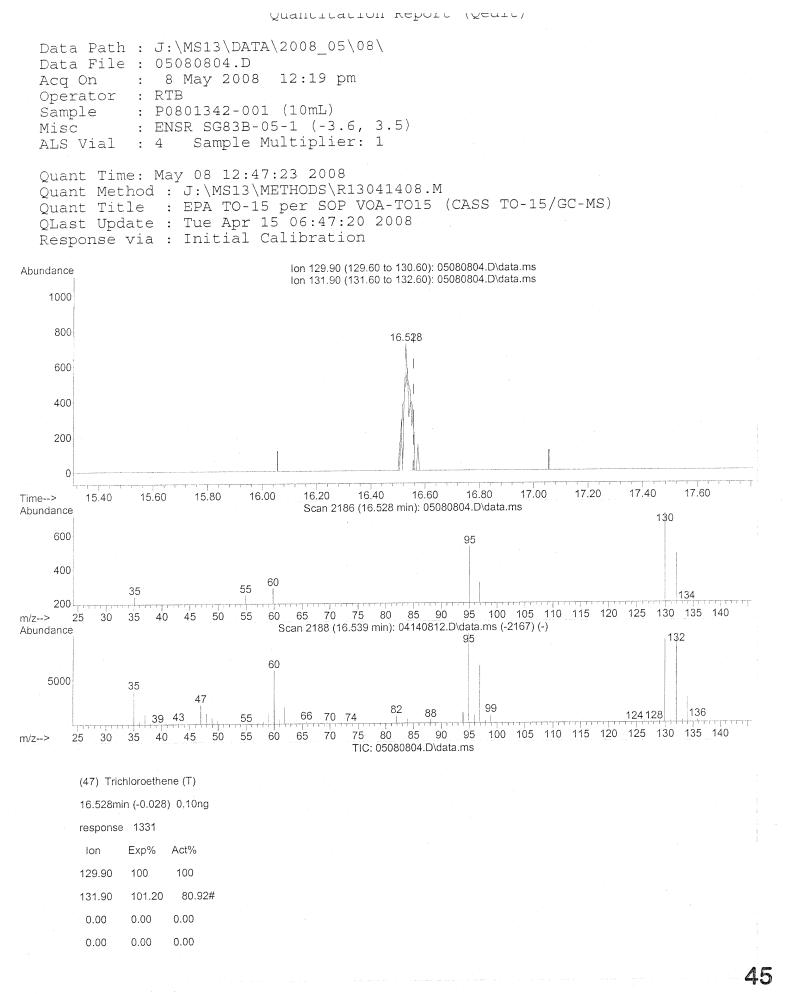


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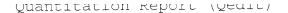


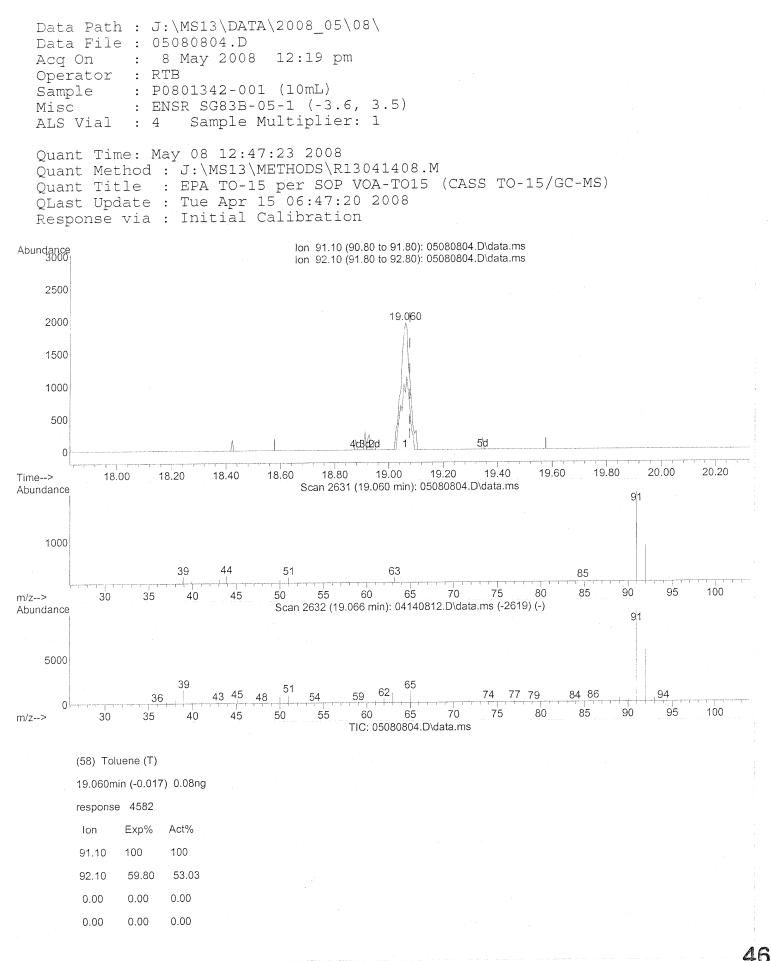
R13041408.M Thu May 08 16:30:01 2008



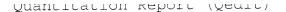


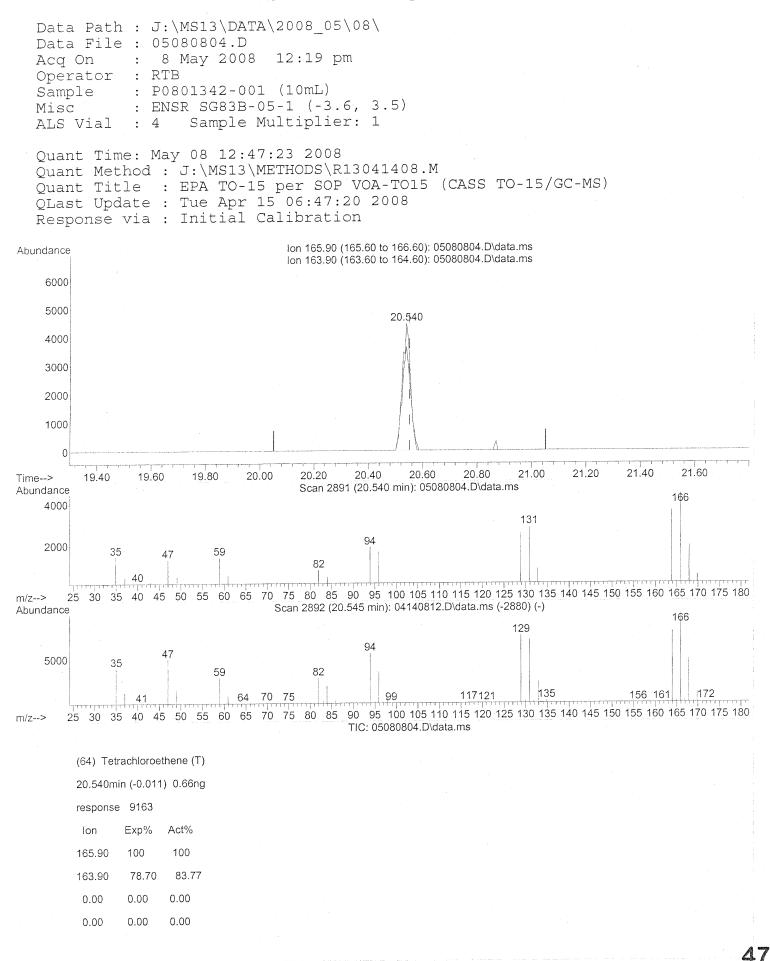
R13041408.M Fri May 09 14:52:13 2008



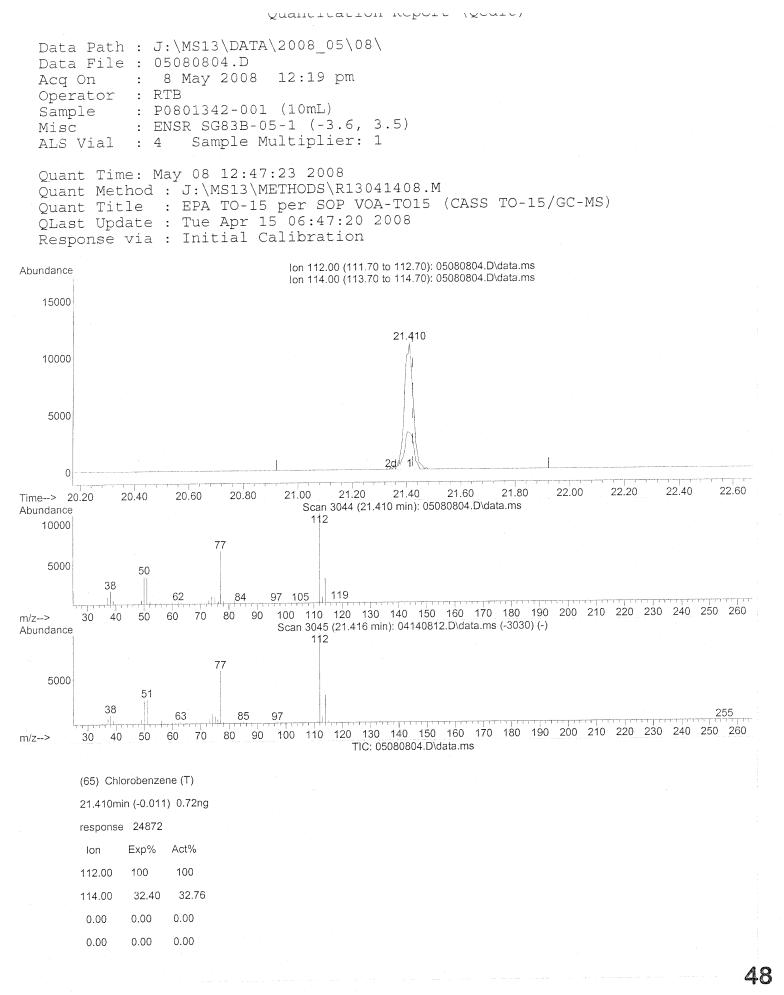


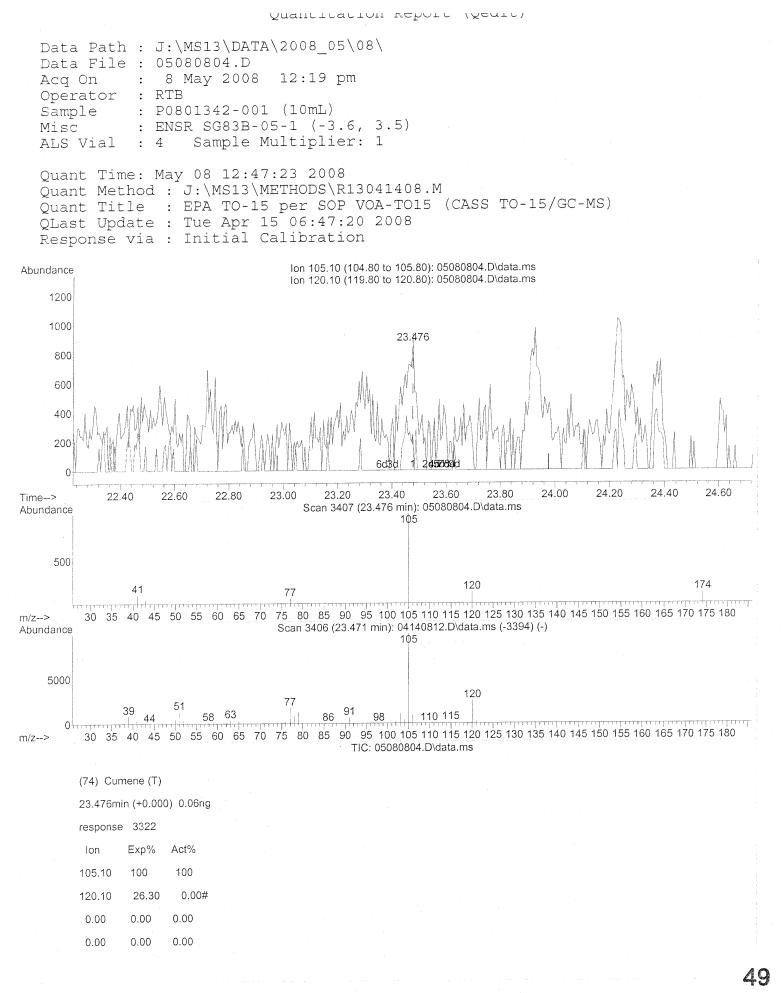
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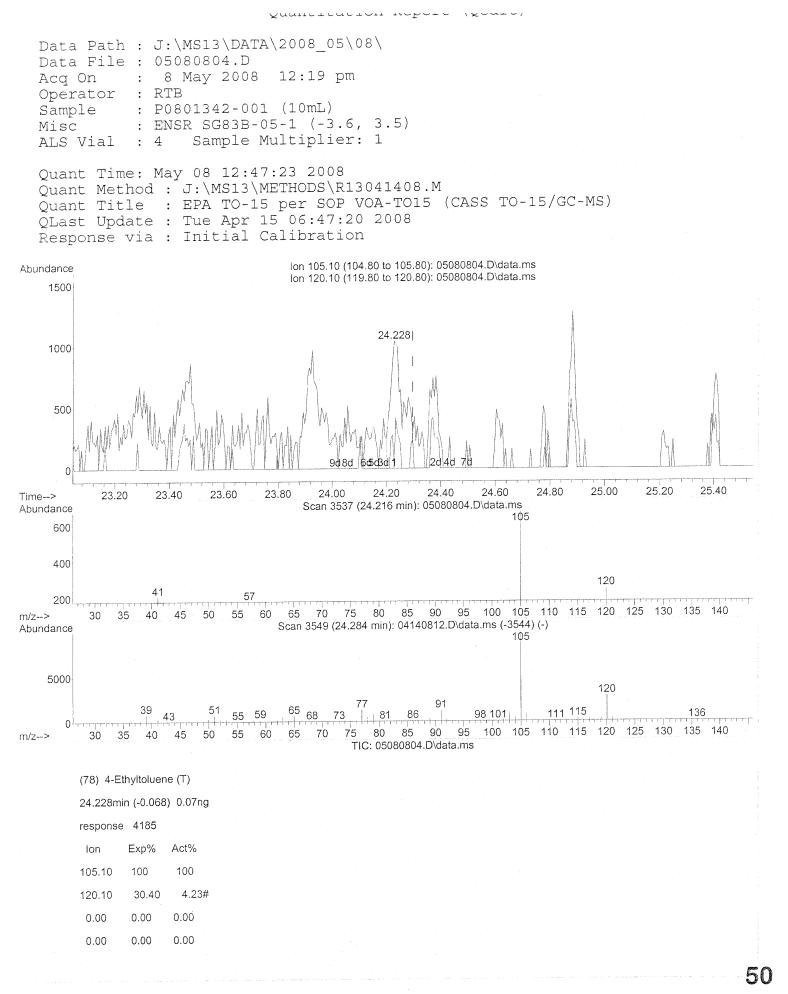


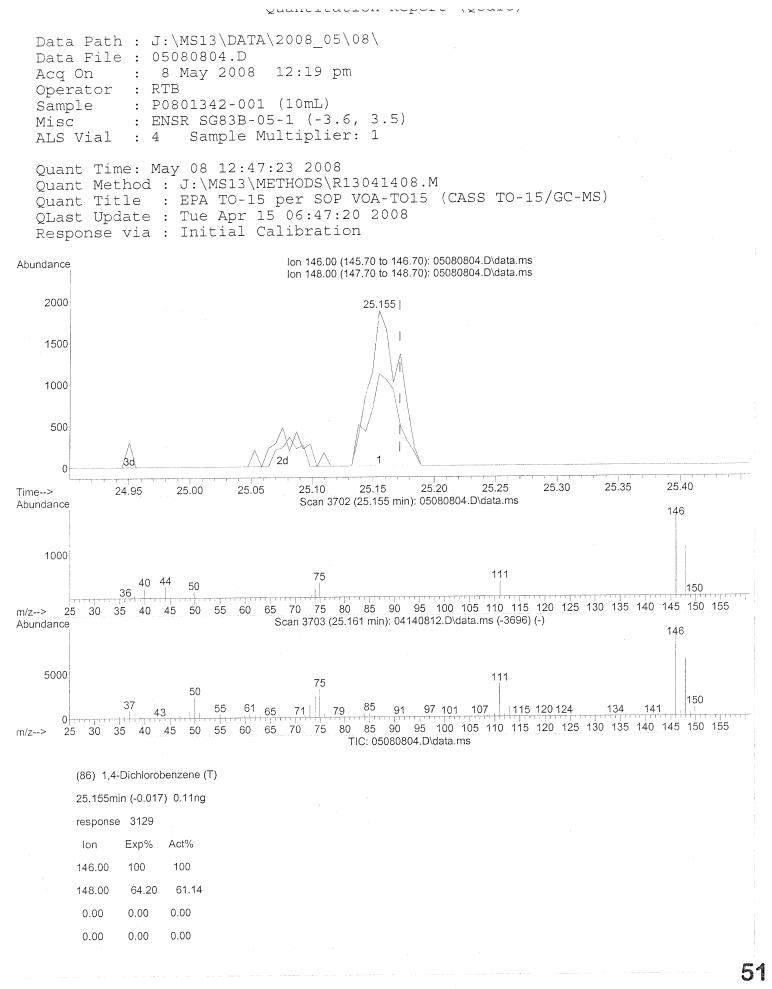


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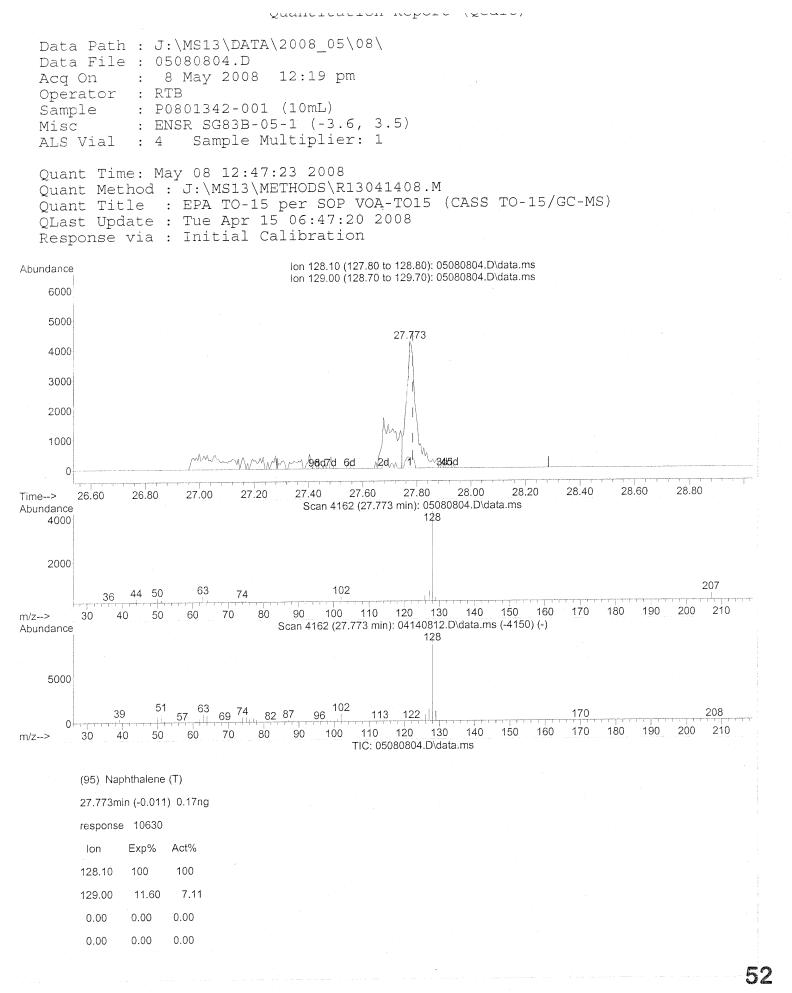


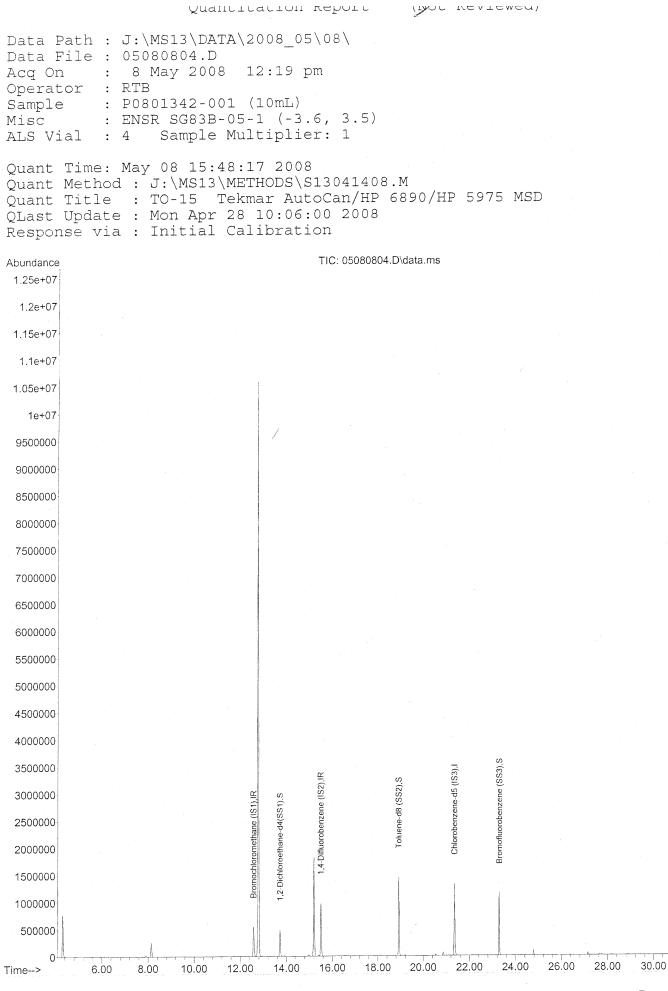






R13041408.M Thu May 08 16:32:02 2008





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

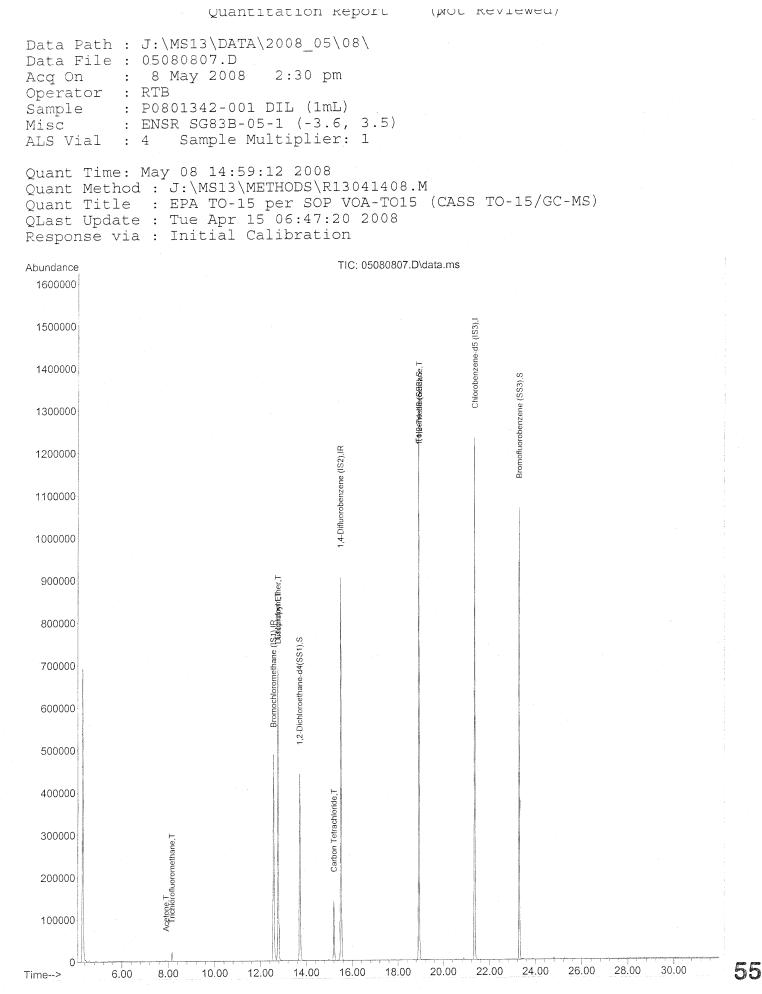
 Spiked Amount
 25.000
 Recovery
 = 101.88%,

 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- &amp; 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.

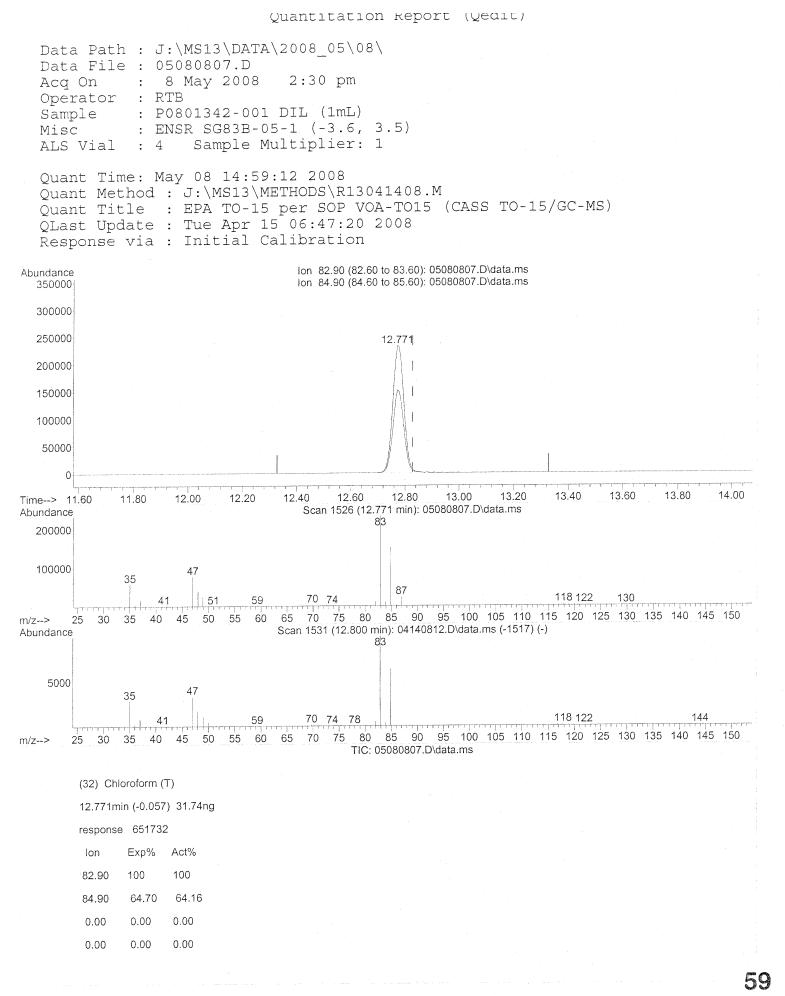
 93) n-Undecane
 26.50
 57
 196
 N.D.

 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
<b>Client Project ID:</b>	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.

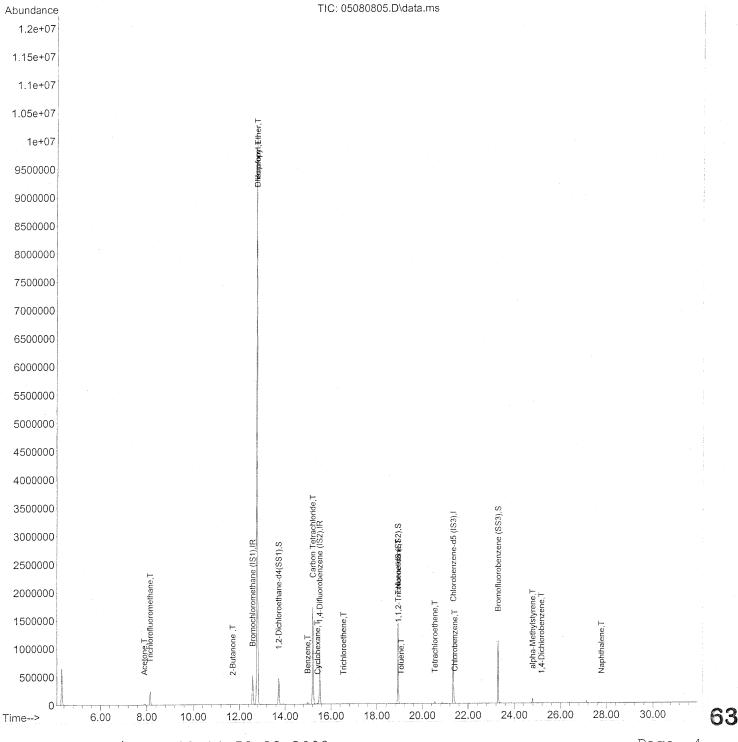
Verified By: <u>Rc-</u>

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 $\pm 5$ cm2 cm2	_1.00			RII-		6
130414	08.M Fri May 09 14:58:29 2	008			205109108	Page	2: 2

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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.

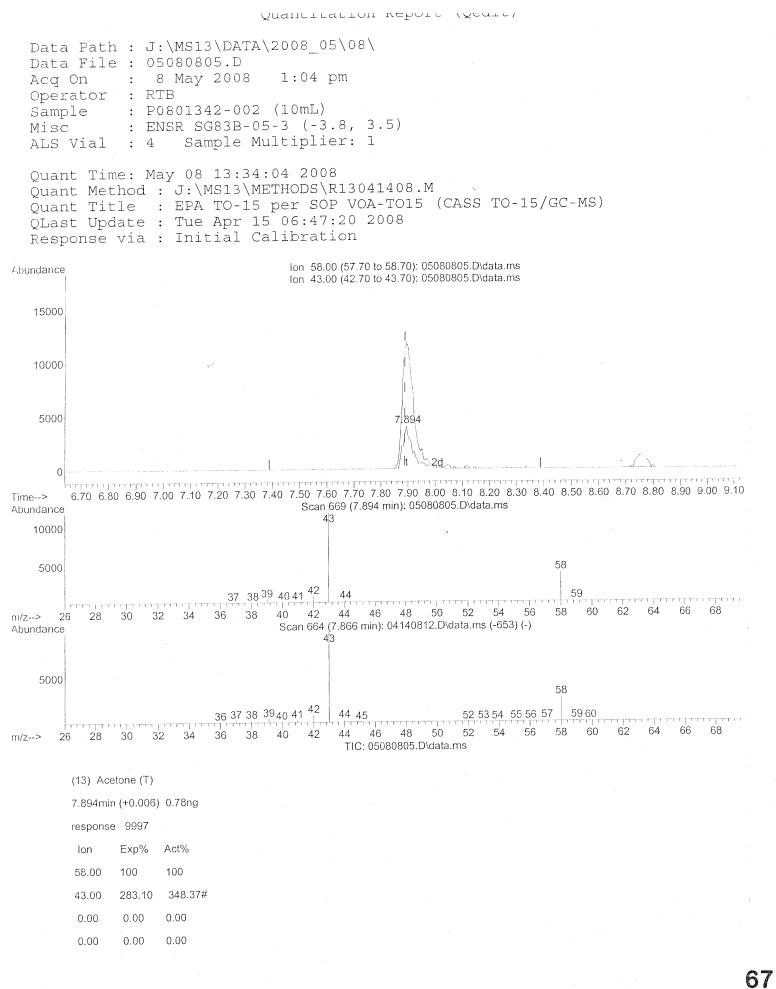
 91) d-Limonene
 26.51
 57
 915
 N.D.
 91
 1,2,4-Trichlorobenzene
 27.64
 180
 229
 N.D.

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

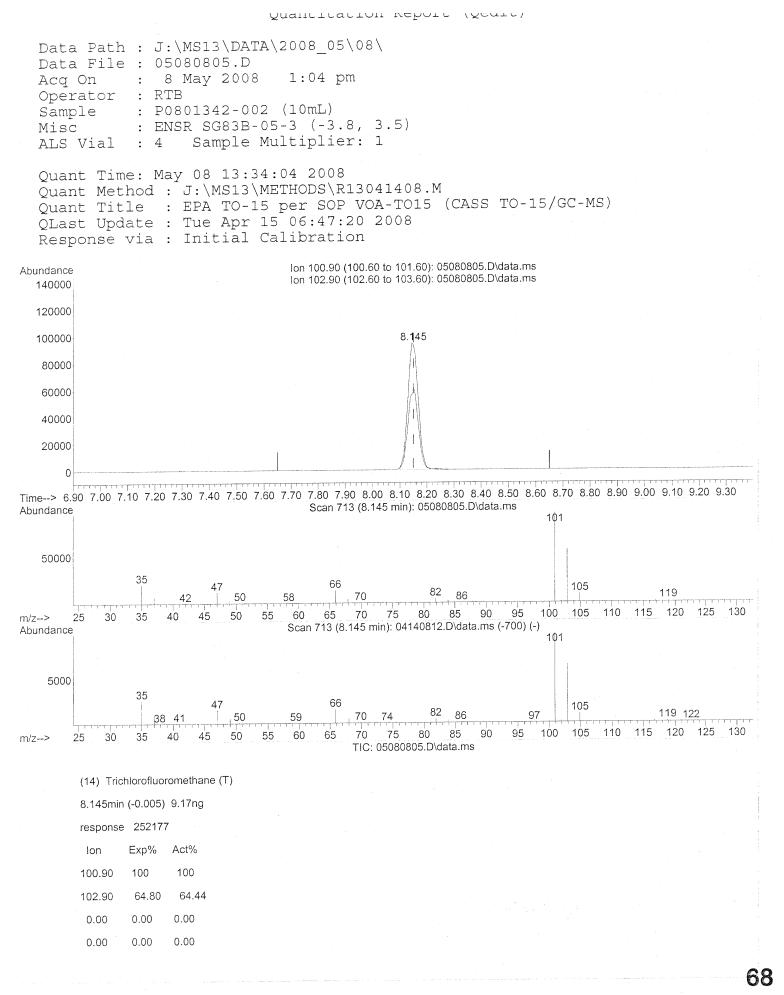
(#) = qualifier out of range (m) = manual integration (+) = signals summed

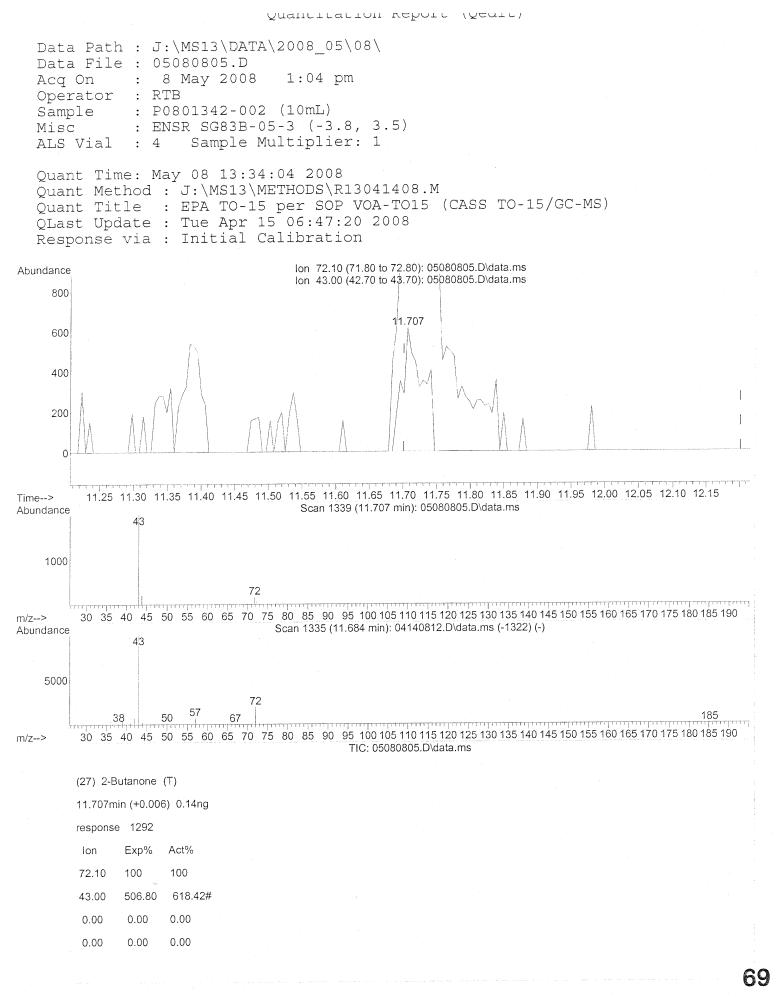
505/09/15

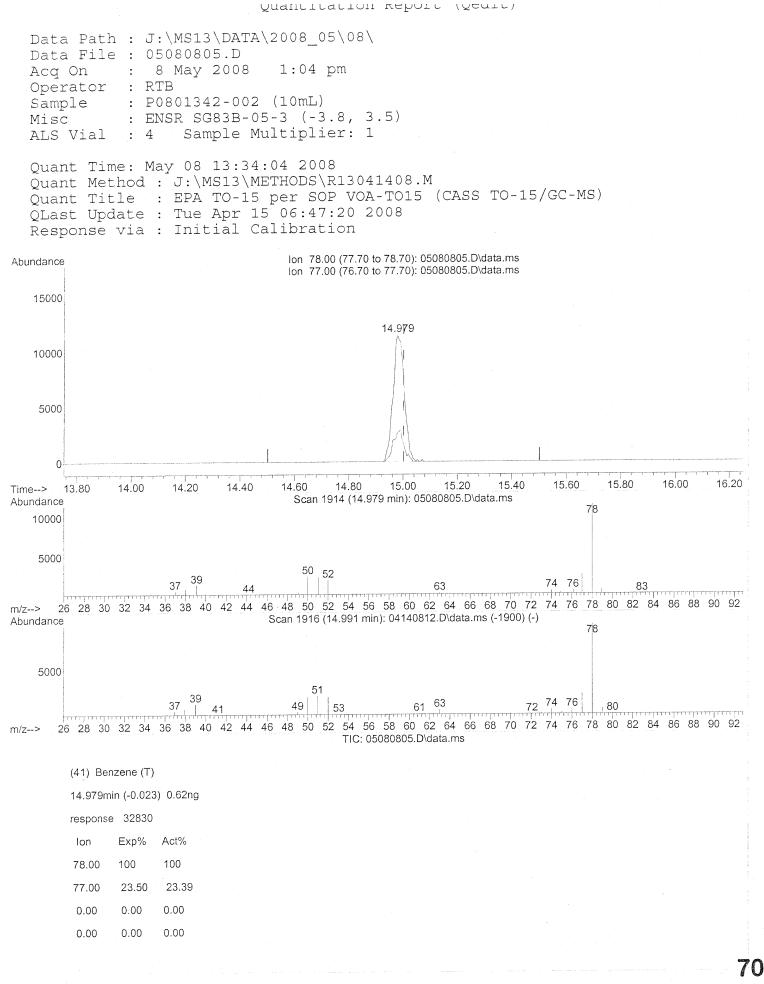
R13041408.M Fri May 09 14:58:29 2008

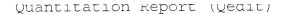


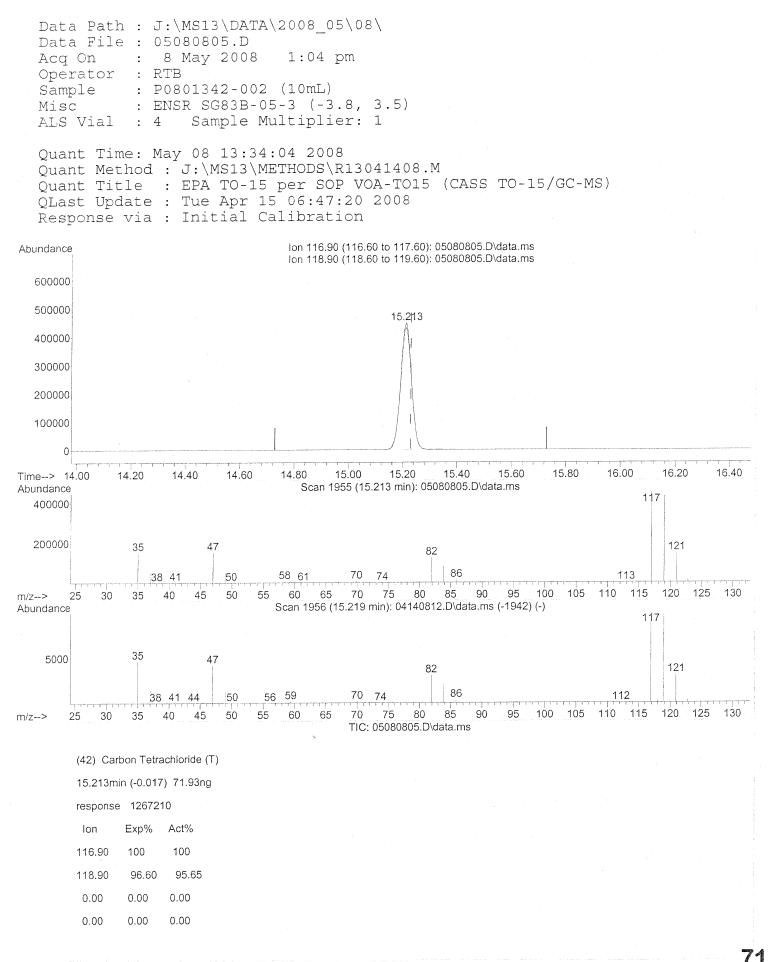
R13041408.M Thu May 08 16:38:23 2008

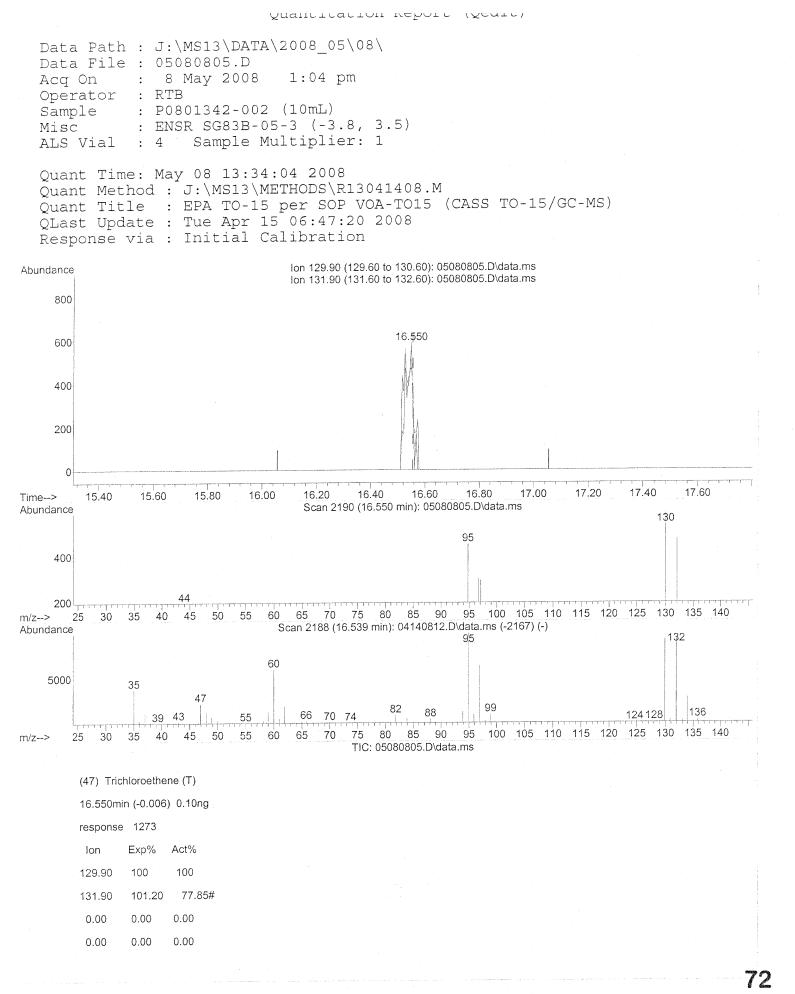




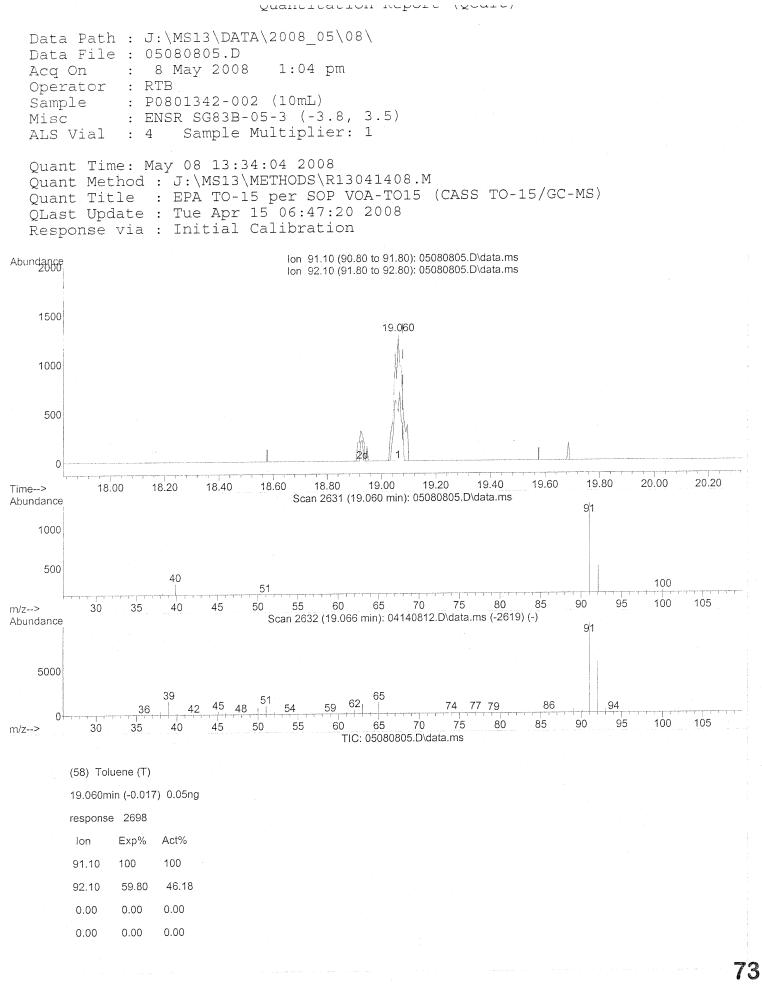


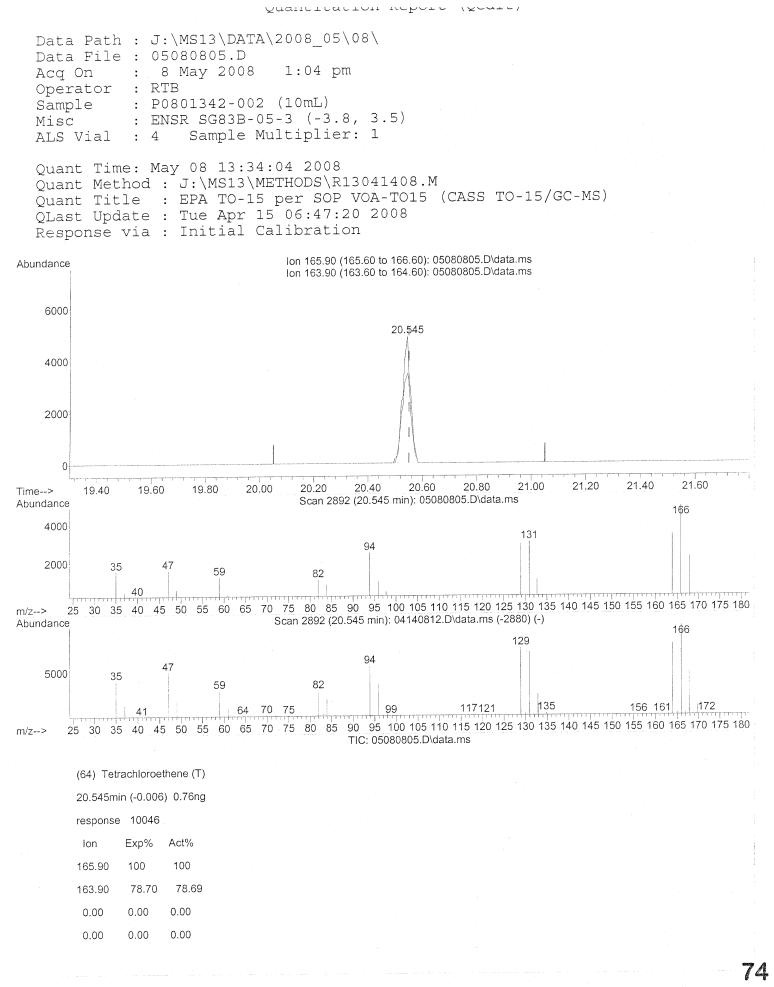


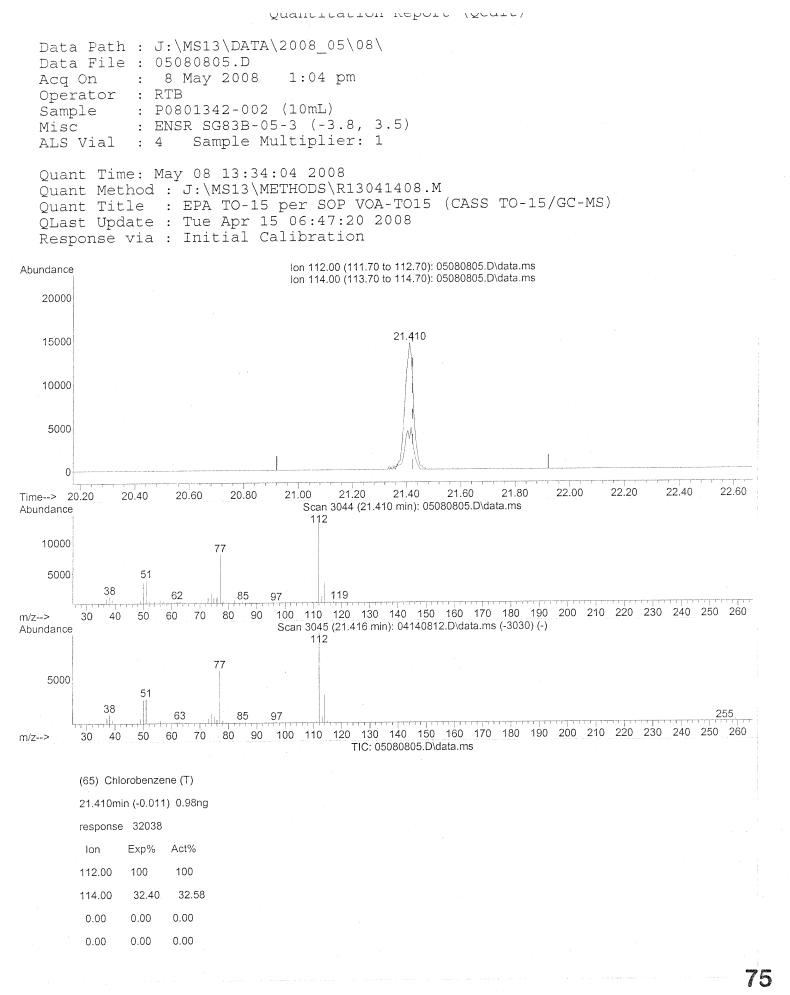


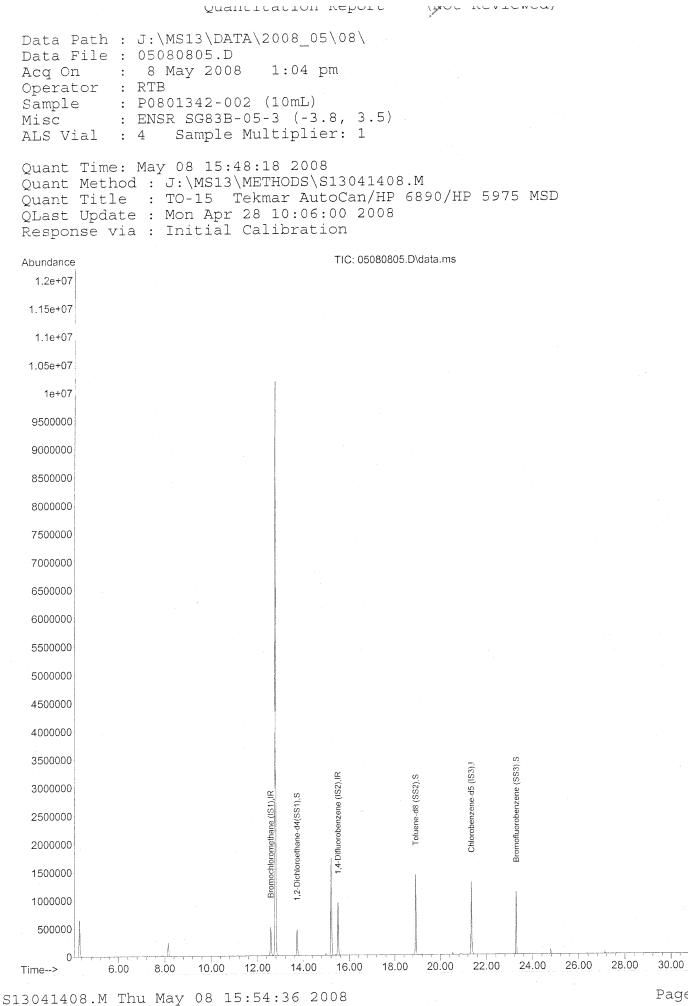


R13041408.M Fri May 09 15:19:21 2008









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

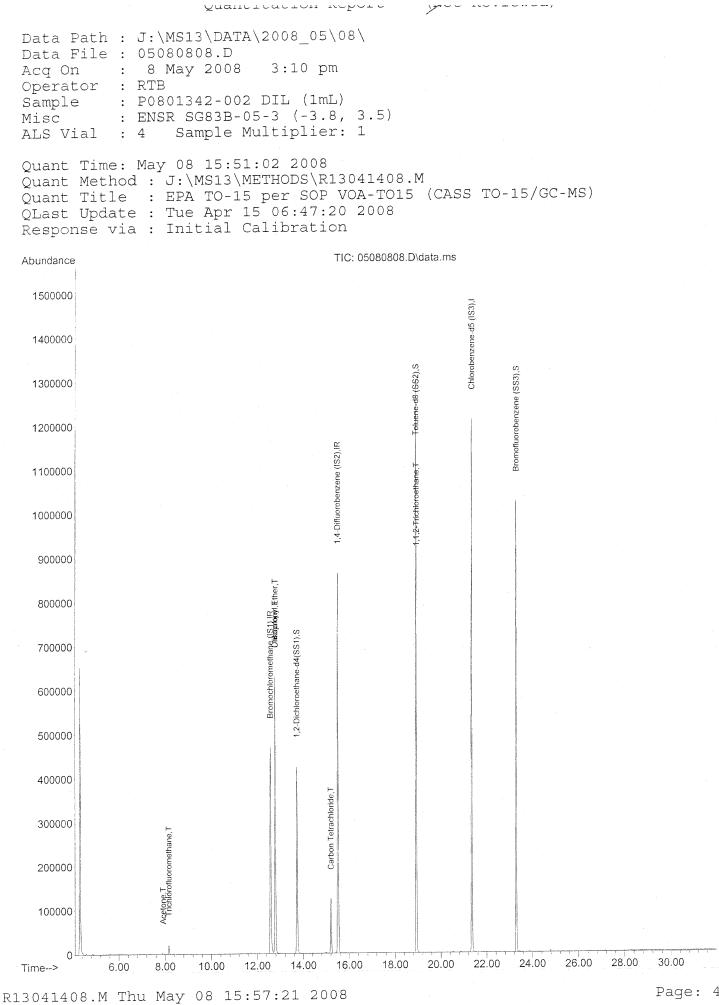
 Spiked Amount
 25.000
 Recovery = 105.96%
 105.96%

 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.	<b>7</b> 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)
<ul> <li>32) Chloroform</li> <li>34) Tetrahydrofuran</li> <li>35) Ethyl tert-Butyl Ether</li> <li>36) 1,2-Dichloroethane</li> <li>38) 1,1,1-Trichloroethane</li> <li>39) Isopropyl Acetate</li> </ul>	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
<ul> <li>40) 1-Butanol</li> <li>41) Benzene</li> <li>42) Carbon Tetrachloride</li> <li>43) Cyclohexane</li> <li>44) tert-Amyl Methyl Ether</li> <li>45) 1,2-Dichloropropane</li> </ul>	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.	
<ul> <li>53) 4-Methyl-2-pentanone</li> <li>54) trans-1,3-Dichloropropene</li> <li>55) 1,1,2-Trichloroethane</li> <li>58) Toluene</li> <li>59) 2-Hexanone</li> <li>60) Dibromochloromethane</li> </ul>		75 97 91 43 129	0 93721 226 0 0	N.D. 7.891 ng N.D. N.D. N.D. N.D.	# 8
<ul> <li>61) 1,2-Dibromoethane</li> <li>62) Butyl Acetate</li> <li>63) n-Octane</li> <li>64) Tetrachloroethene</li> <li>65) Chlorobenzene</li> </ul>	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.	
<ul> <li>66) Ethylbenzene</li> <li>67) m- &amp; p-Xylene</li> <li>68) Bromoform</li> <li>69) Styrene</li> <li>70) o-Xylene</li> </ul>	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.	
<ul> <li>71) n-Nonane</li> <li>72) 1,1,2,2-Tetrachloroethane</li> <li>74) Cumene</li> <li>75) alpha-Pinene</li> <li>76) n-Propylbenzene</li> </ul>	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.	
<ul> <li>77) 3-Ethyltoluene</li> <li>78) 4-Ethyltoluene</li> <li>79) 1,3,5-Trimethylbenzene</li> <li>13041408.M Thu May 08 15:57:20 2</li> </ul>			201 201 201	N.D. N.D. N.D. Fostoslos	<b>8</b> 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

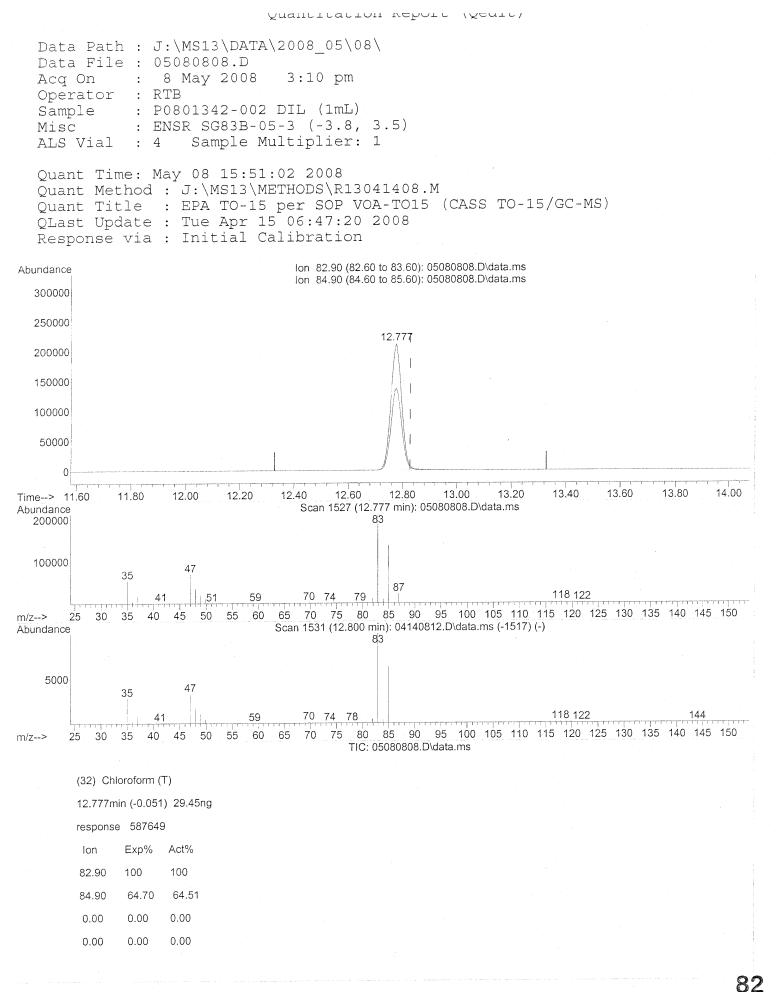
 1.2
 Dichlorobenzene
 25.16
 146

 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
<b>Client Sample ID:</b>	SG83B-05-7
<b>Client Project ID:</b>	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.

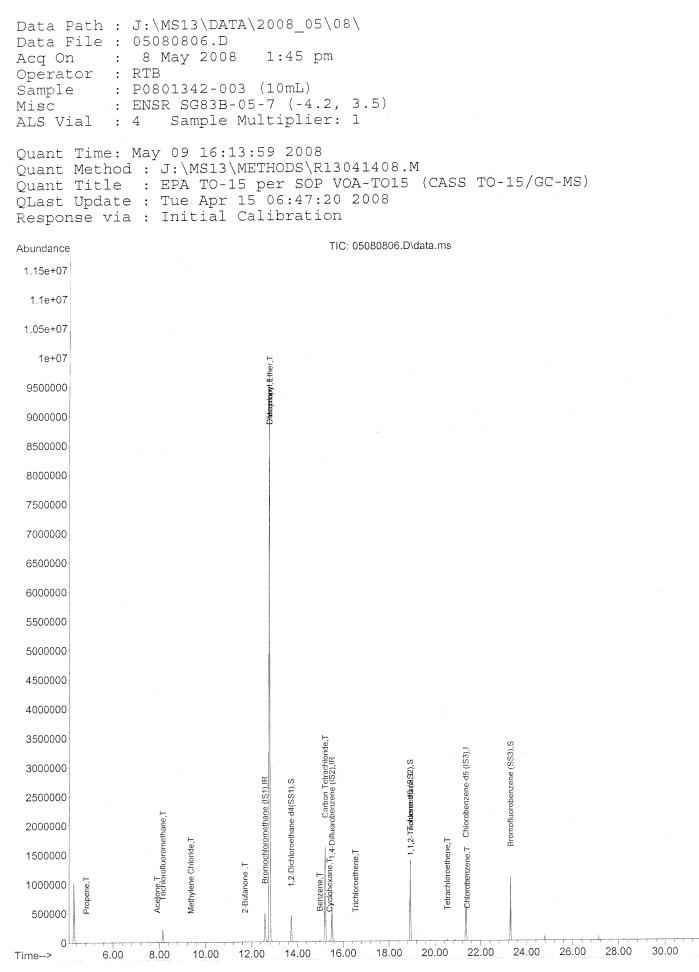
MRL = Method Reporting Limit - The minimum quantity of a target analyte that can be confidently determined by the referenced method.

## RESULTS OF ANALYSIS

Page 4 of 4

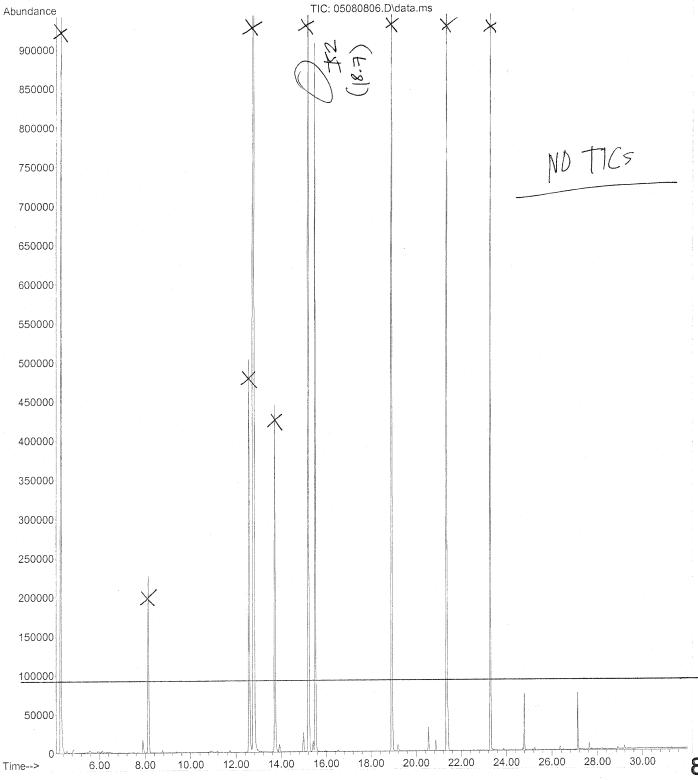
Client:	ENSR		
<b>Client Sample ID:</b>	SG83B-05-7	CAS Project ID: P080	1342
<b>Client Project ID:</b>	Phase B Soil Gas / 04020-023-4311	CAS Sample ID: P080	1342-003
	Tentatively Identified Compo	unds	
Test Code:	EPA TO-15	Date Collected: 5/7/0	8
Instrument ID:	Tekmar AUTOCAN/Agilent 5975Binert/6890N/MS13	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	
<ul> <li>32) Chloroform</li> <li>34) Tetrahydrofuran</li> <li>35) Ethyl tert-Butyl Ether</li> <li>36) 1,2-Dichloroethane</li> <li>38) 1,1,1-Trichloroethane</li> <li>39) Isopropyl Acetate</li> </ul>	13.74 62 0.00 97 0.00 61	0 N. 0 N. 250 N. 0 N. 0 N.	D. D. D. D.
<ul> <li>40) 1-Butanol</li> <li>41) Benzene</li> <li>42) Carbon Tetrachloride</li> <li>43) Cyclohexane</li> <li>44) tert-Amyl Methyl Ether</li> <li>45) 1,2-Dichloropropane</li> <li>46) Bromodichloromethane</li> </ul>	$ \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.	-
<ul> <li>46) Bromodicinoronale finale</li> <li>47) Trichloroethene</li> <li>48) 1,4-Dioxane</li> <li>49) Isooctane</li> <li>50) Methyl Methacrylate</li> <li>51) n-Heptane</li> <li>52) cis-1,3-Dichloropropene</li> <li>53) 4-Methyl-2-pentanone</li> <li>54) trans-1,3-Dichloropropen</li> </ul>	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.
<ul> <li>54) trans-1, 3-Dichiolopiopen</li> <li>55) 1,1,2-Trichloroethane</li> <li>58) Toluene</li> <li>59) 2-Hexanone</li> <li>60) Dibromochloromethane</li> <li>61) 1,2-Dibromoethane</li> <li>62) Butyl Acetate</li> <li>63) n-Octane</li> </ul>	$ \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.
<ul> <li>64) Tetrachloroethene</li> <li>65) Chlorobenzene</li> <li>66) Ethylbenzene</li> <li>67) m- &amp; p-Xylene</li> <li>68) Bromoform</li> <li>69) Styrene</li> <li>70) o-Xylene</li> </ul>	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. <b>90</b>
13041408.M Fri May 09 16:16:03	2008	Rosfiglos	Page: 2

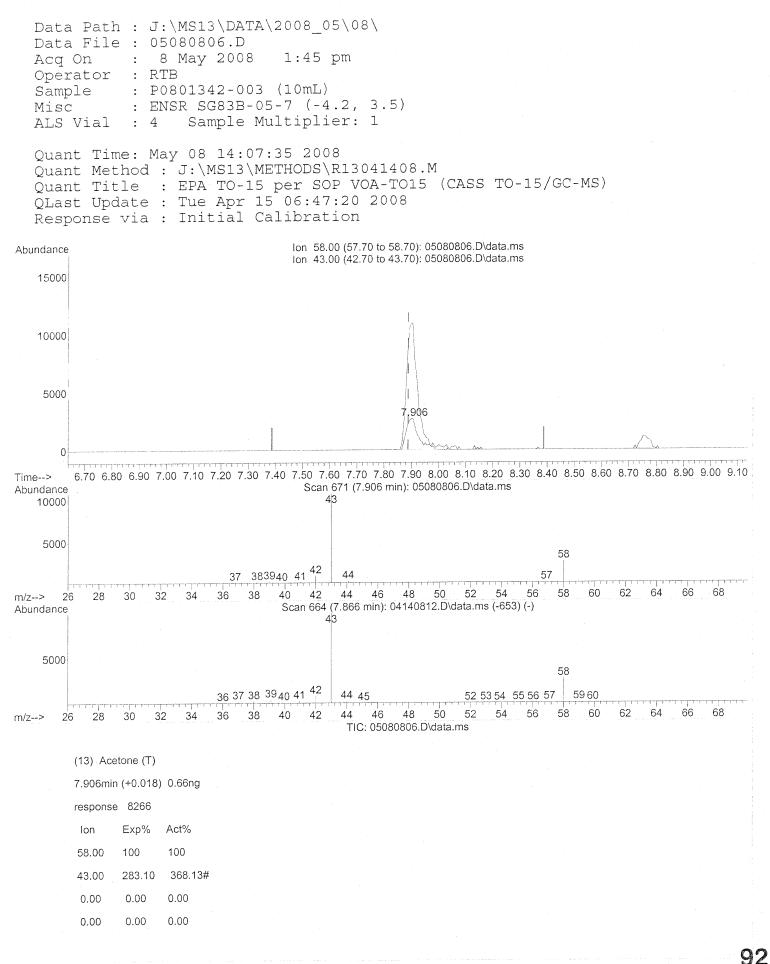
Quantituacion Report (Xr Review,

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. ------

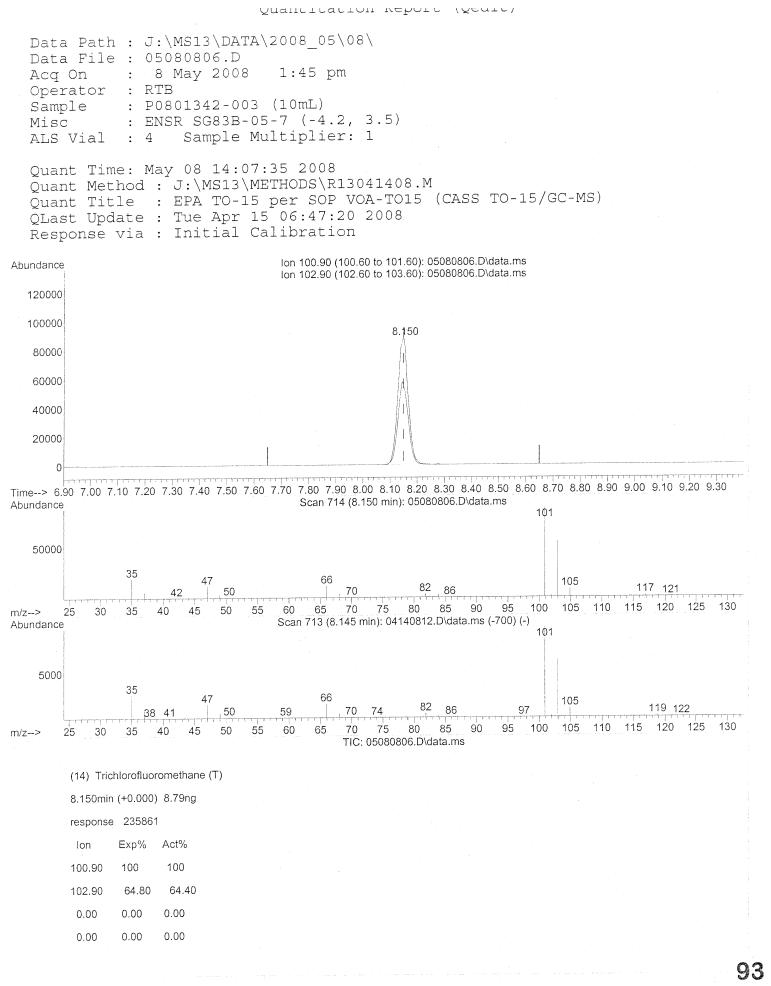
(#) = qualifier out of range (m) = manual integration (+) = signals summed

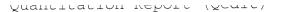
50 stoglor

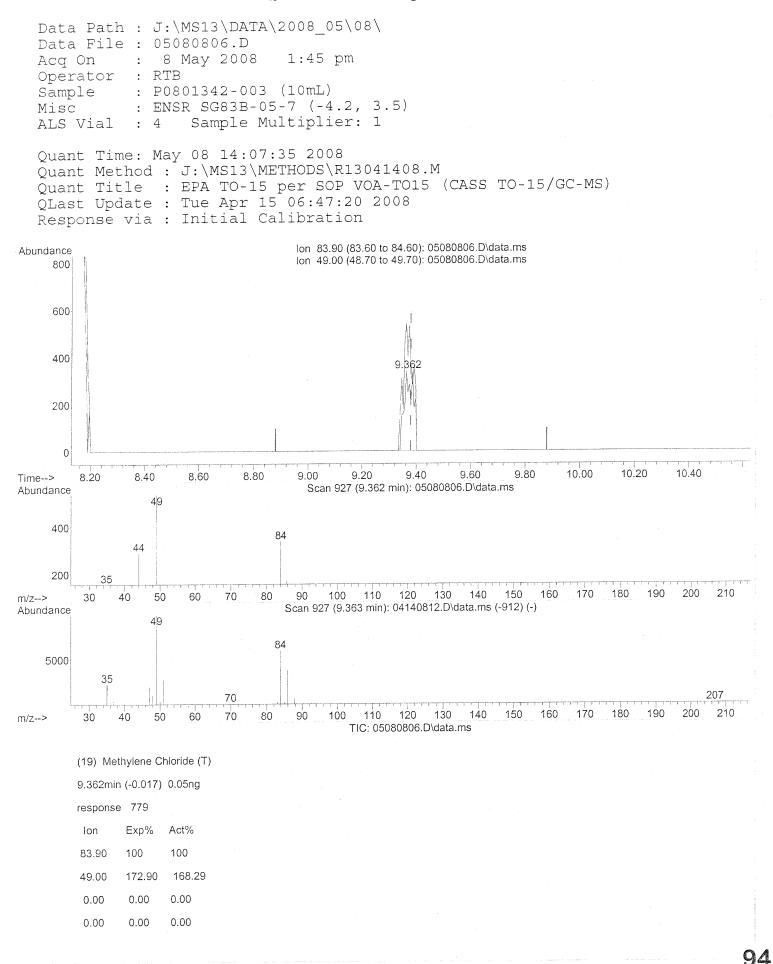




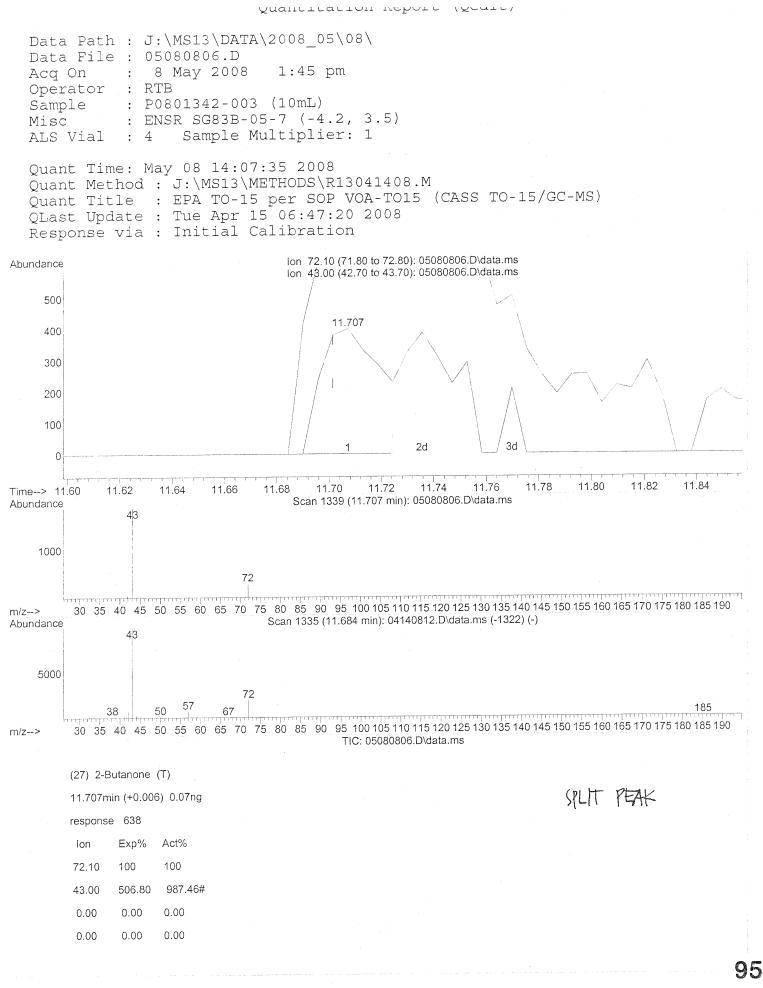
R13041408.M Thu May 08 16:44:44 2008

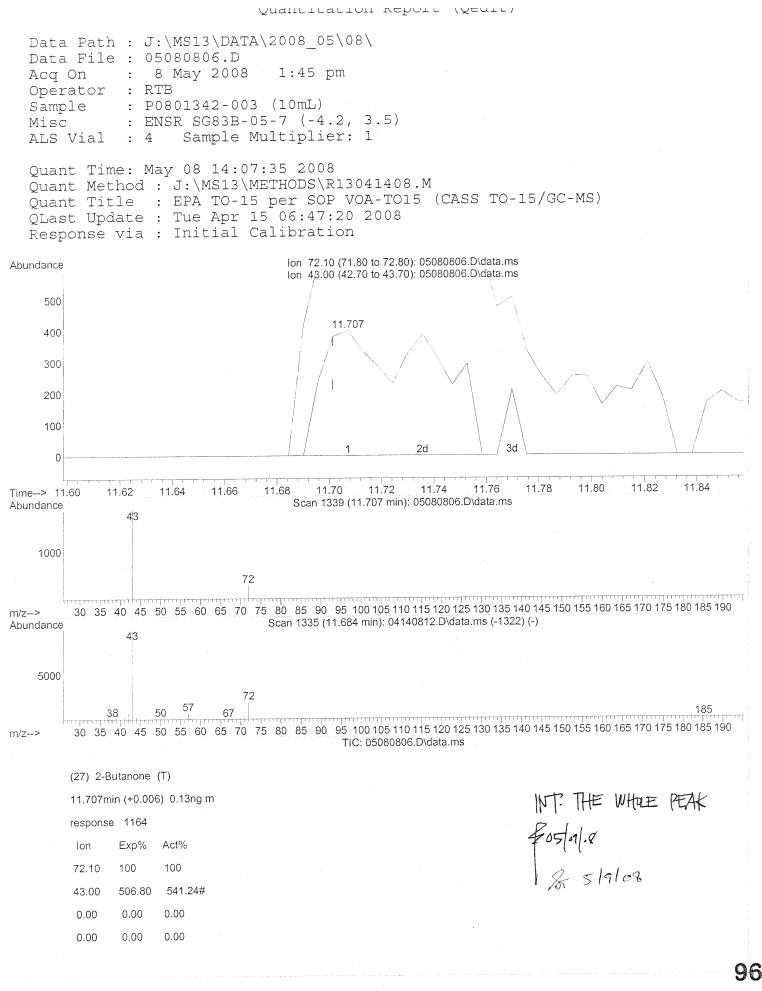


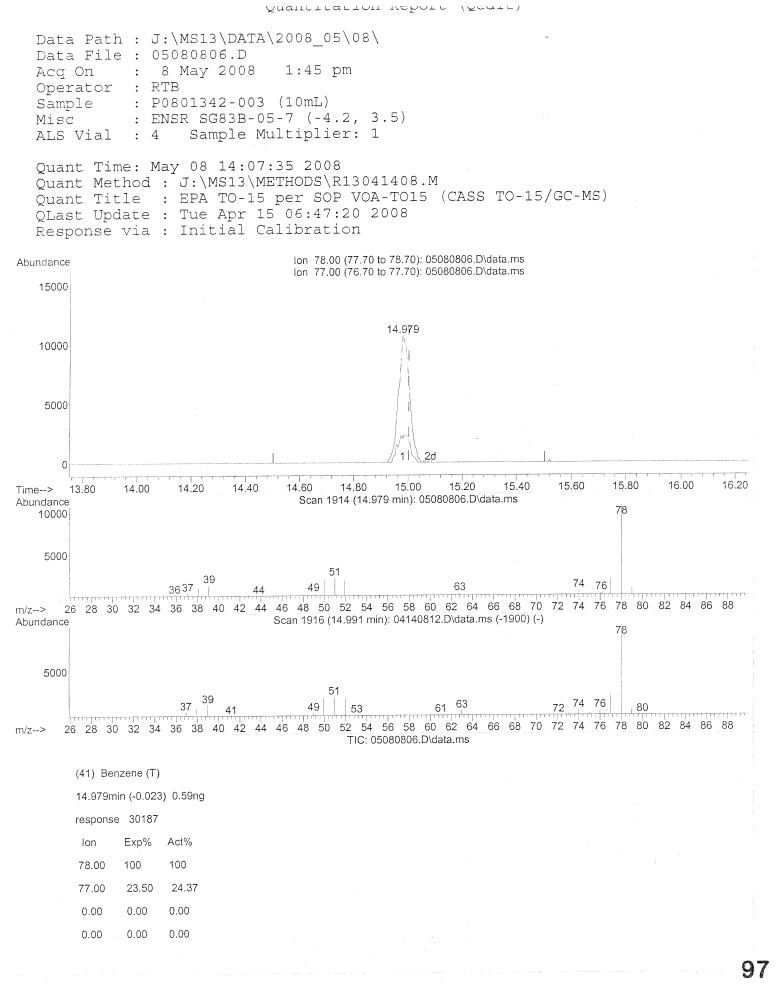




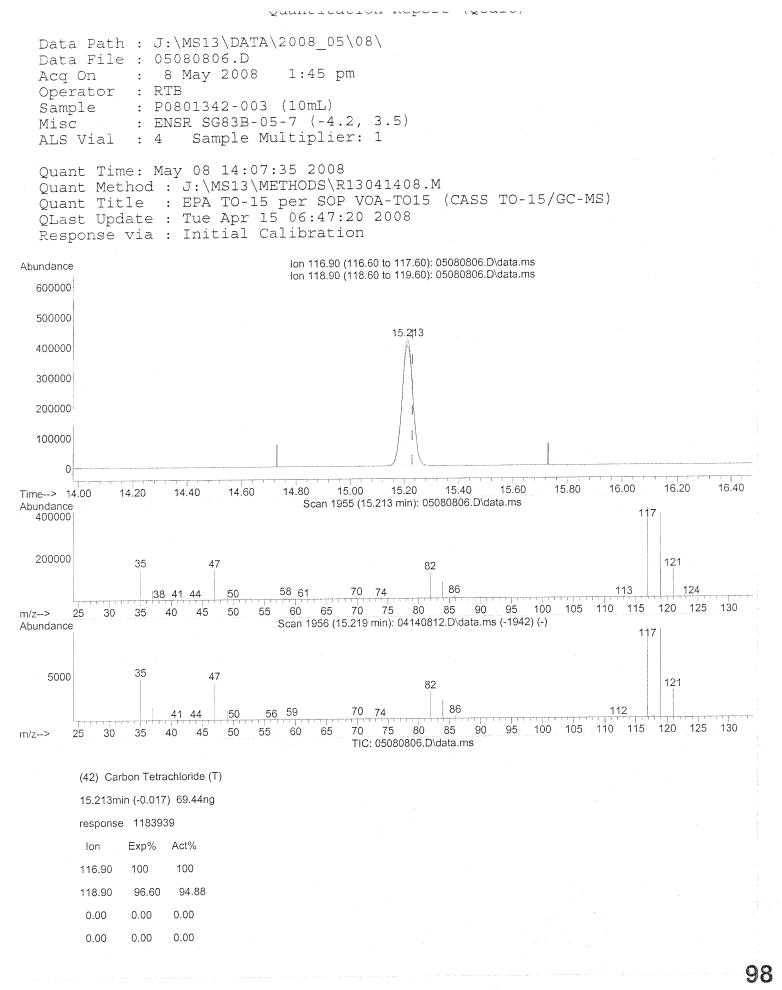
R13041408.M Fri May 09 15:25:06 2008

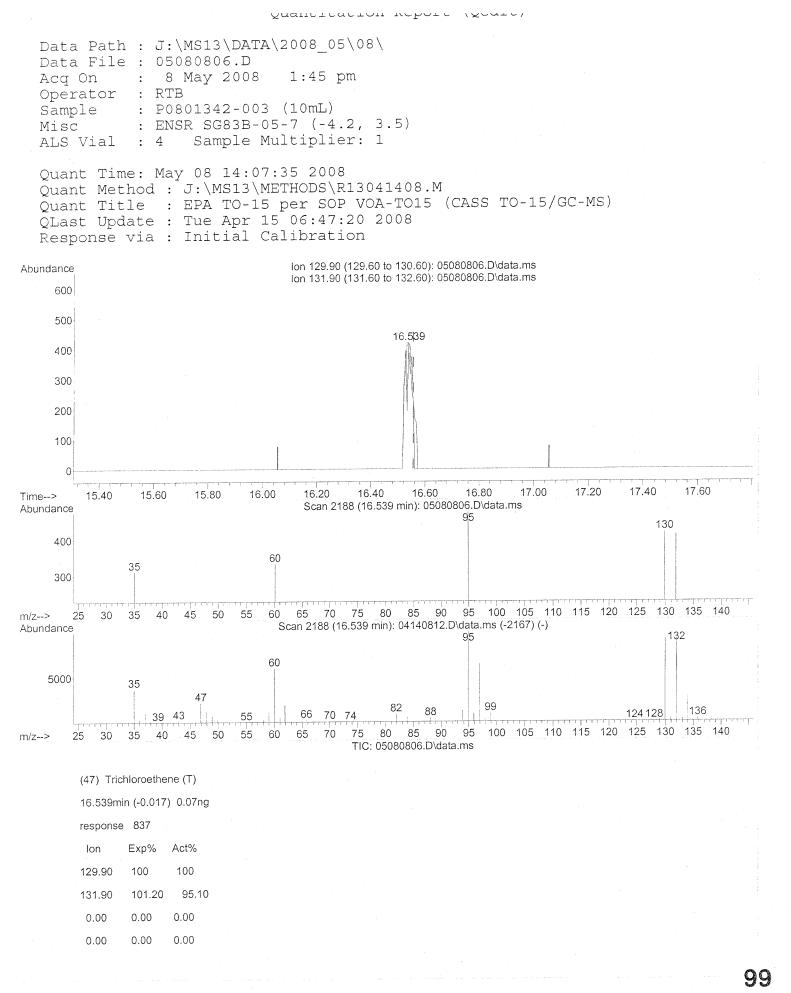




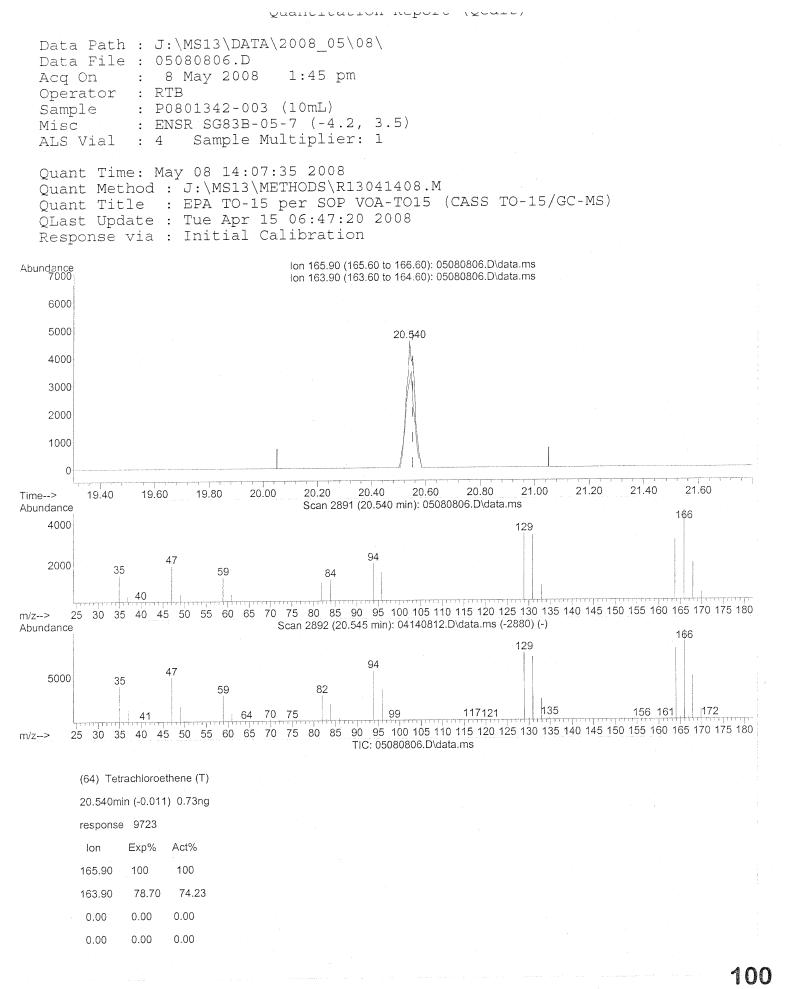


R13041408.M Thu May 08 16:45:54 2008



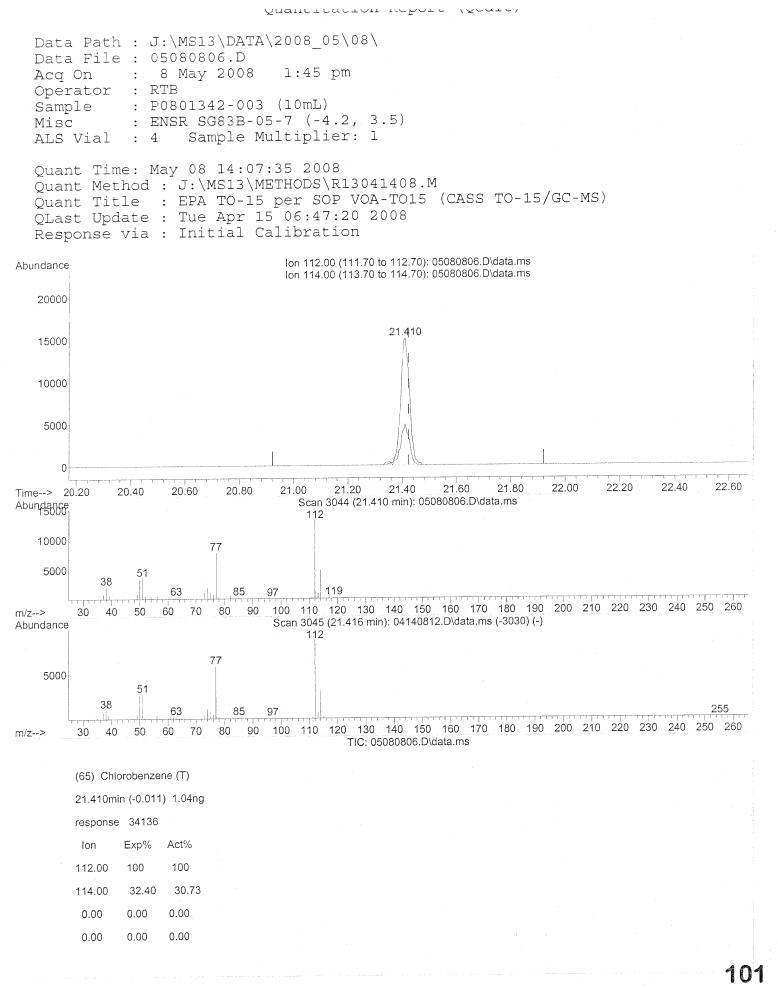


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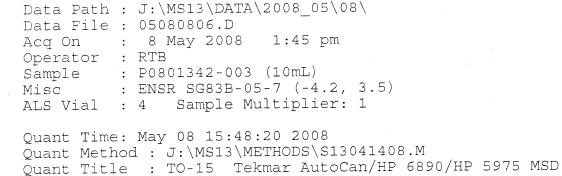


R13041408.M Thu May 08 16:46:40 2008

Page: 1



(INON VENTEMEN)



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

Response via : Initial Calibration

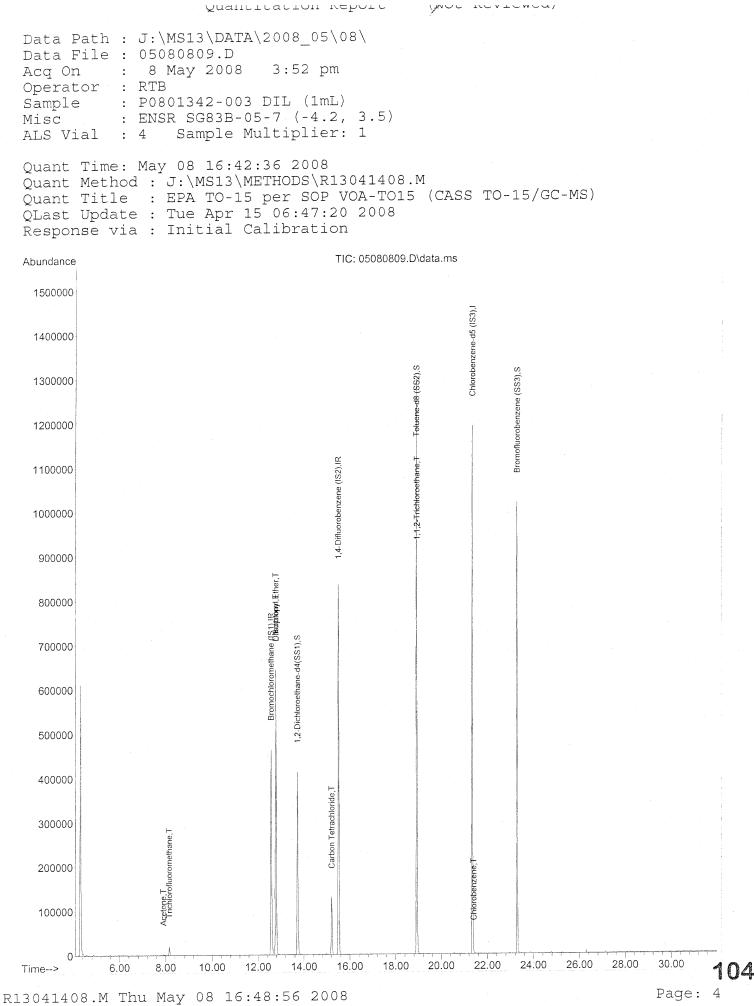
Abundance	TIC: 05080806.D\data.ms	
1.15e+07		
1.1e+07		
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9500000		
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4000000		
3500000	SS SS SS SS SS	
3000000	ne-d5 ((S2)) a (S22)	
2500000	diloromethane (IS1),IR Noroethane-d4(SS1),S Toluene-d8 (SS2),S 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	<b>30.00</b> Page: 2

Yuuncreacton		7	<b>, , , , , , , , , ,</b>				
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
<ol> <li>Acetonitrile</li> <li>Acrolein</li> <li>Acetone</li> <li>Trichlorofluoromethane</li> <li>Isopropanol</li> <li>Acrylonitrile</li> <li>1,1-Dichloroethene</li> <li>tert-Butanol</li> </ol>	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
<ul> <li>19) Methylene Chloride</li> <li>20) Allyl Chloride</li> <li>21) Trichlorotrifluoroethane</li> <li>22) Carbon Disulfide</li> <li>23) trans-1,2-Dichloroethene</li> <li>24) 1,1-Dichloroethane</li> <li>25) Methyl tert-Butyl Ether</li> <li>26) Vinyl Acetate</li> </ul>	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

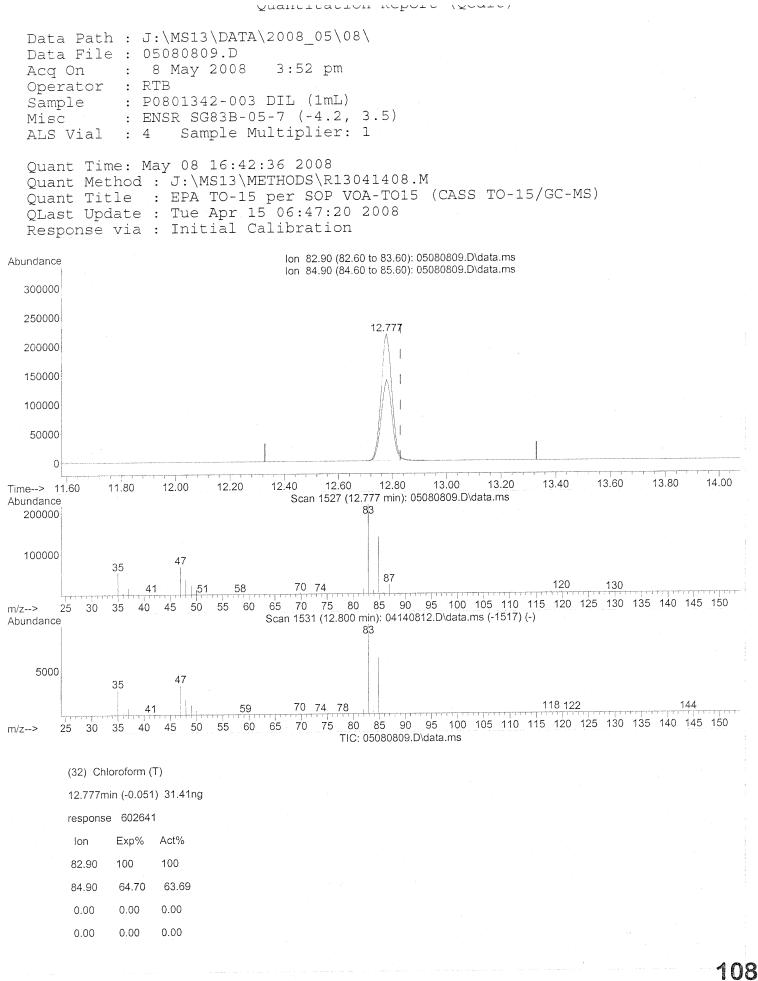
Quantitation	MCDOT C	(1)		- /	
Data Path : J:\MS13\DATA\2008_05 Data File : 05080809.D Acq On : 8 May 2008 3:52 p Operator : RTB Sample : P0801342-003 DIL (1m Misc : ENSR SG83B-05-7 (-4. ALS Vial : 4 Sample Multiplie	om NL) 2, 3.5)				
Quant Time: May 08 16:42:36 2008 Quant Method : J:\MS13\METHODS\R 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 <b>Q</b>	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

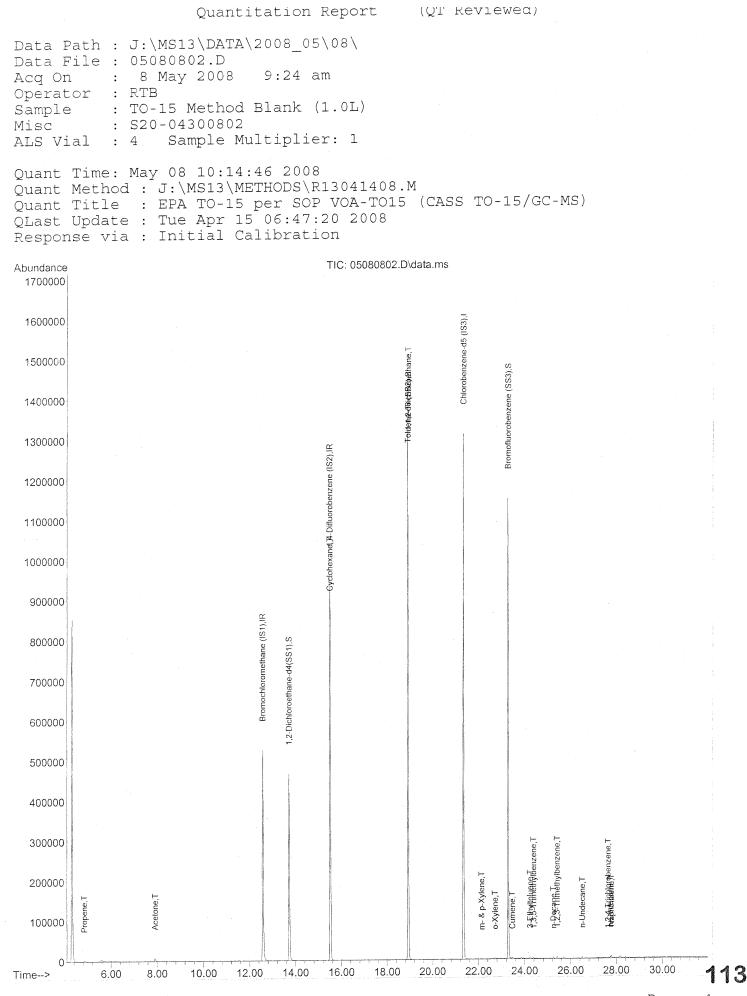
## RESULTS OF ANALYSIS Page 4 of 4

Client:	ENSR		
<b>Client Sample ID:</b>	Method Blank	CAS Project ID: P	0801342
<b>Client Project ID:</b>	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.

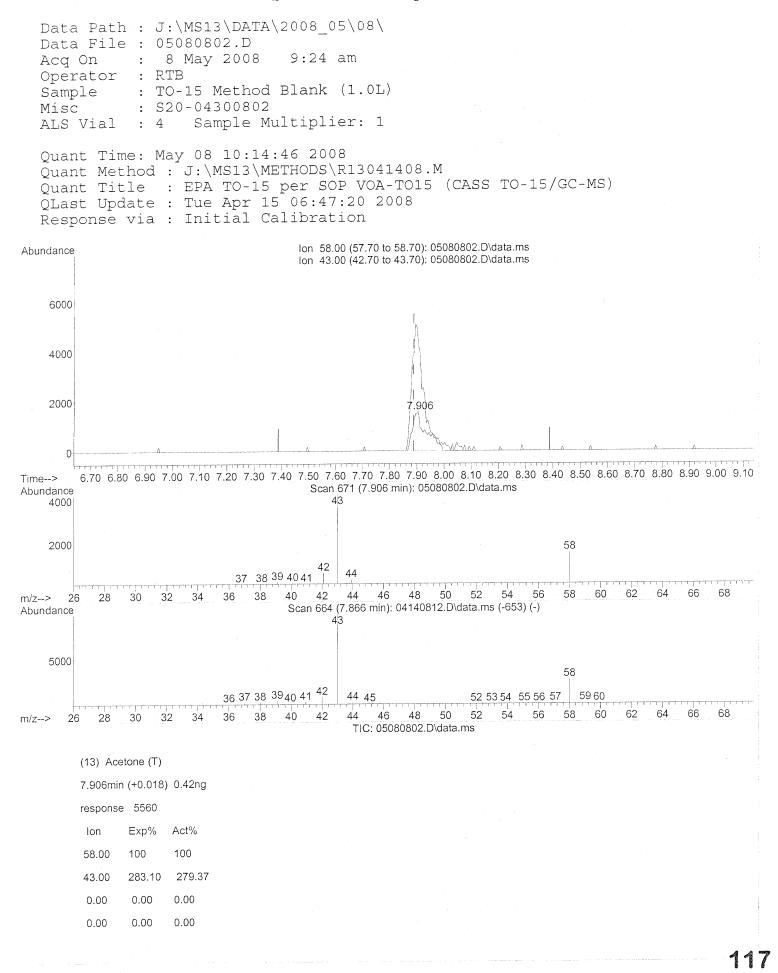
 91)
 d-Limonene
 0.00
 68
 0
 N.D.

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

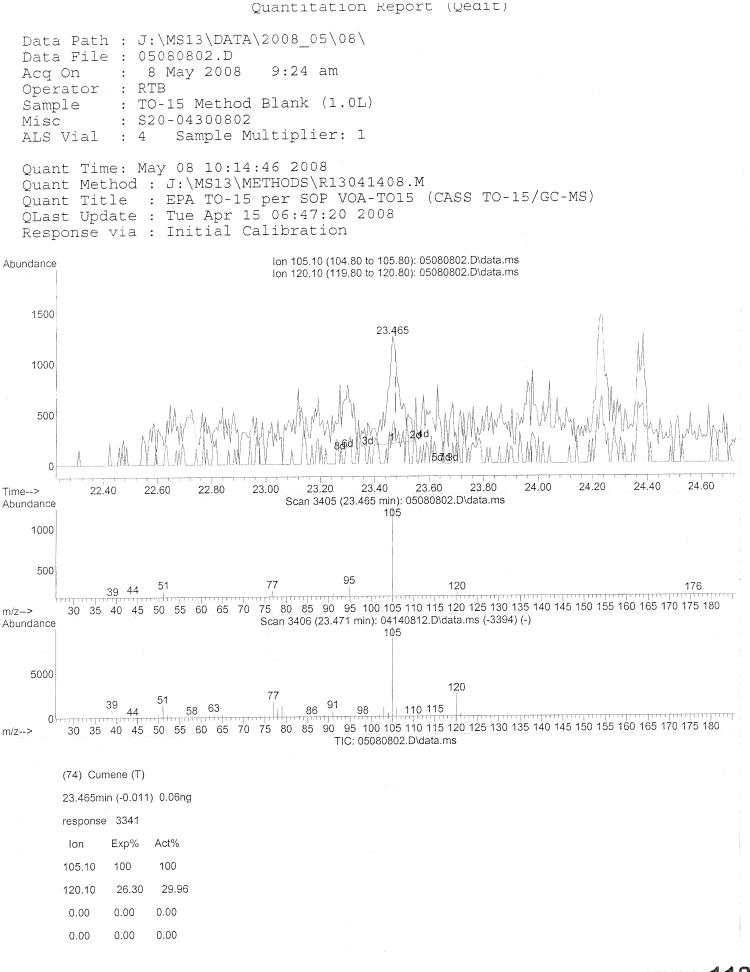
 93)
 n-Undecane
 26.51
 57
 1681
 0.045 ng
 #

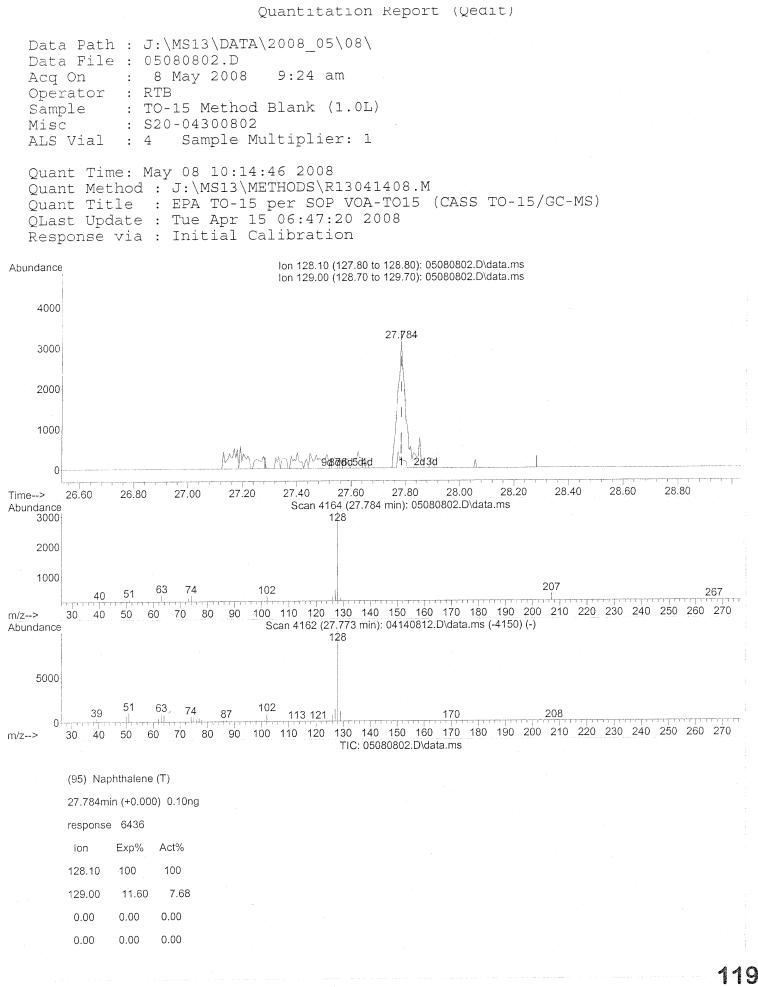
 94)
 1,2,4-Trichlorob 50 90 95 70 90 78 _____

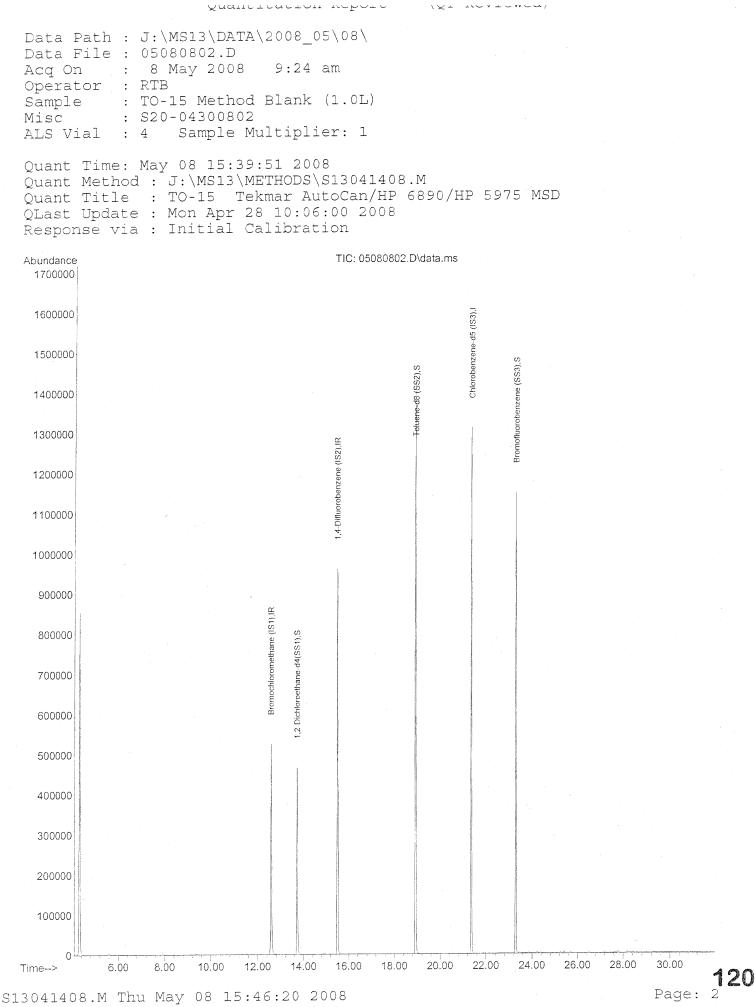
(#) = qualifier out of range (m) = manual integration (+) = signals summed



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:

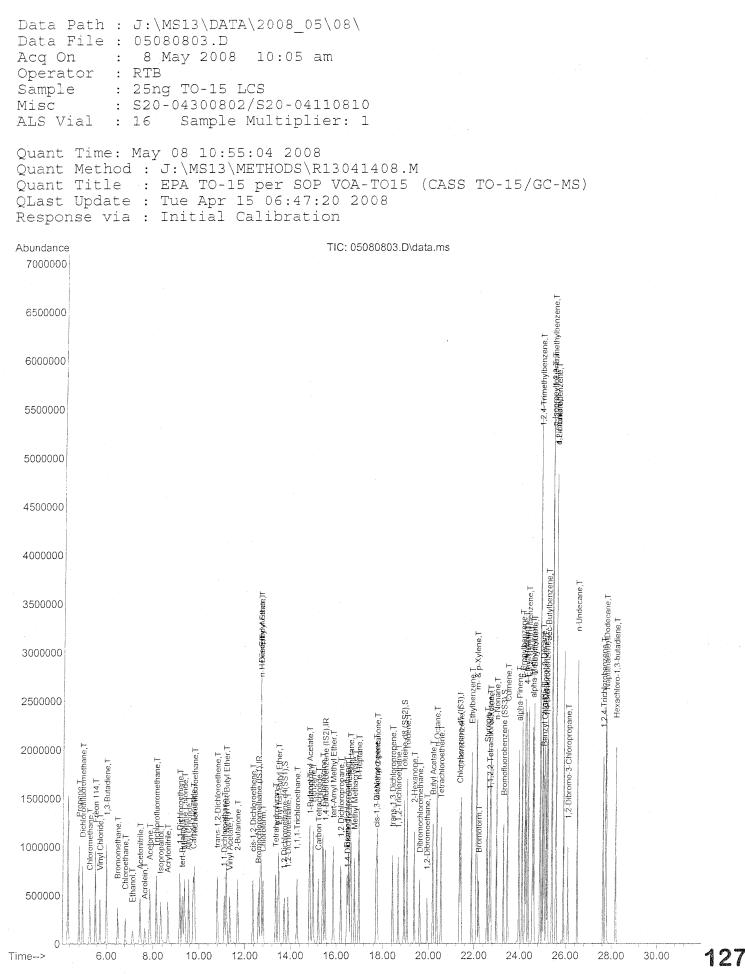
Date Collected: NA Date Received: NA Date Analyzed: 5/08/08 Volume(s) Analyzed: NA Liter(s)

					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 <b>28</b>
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

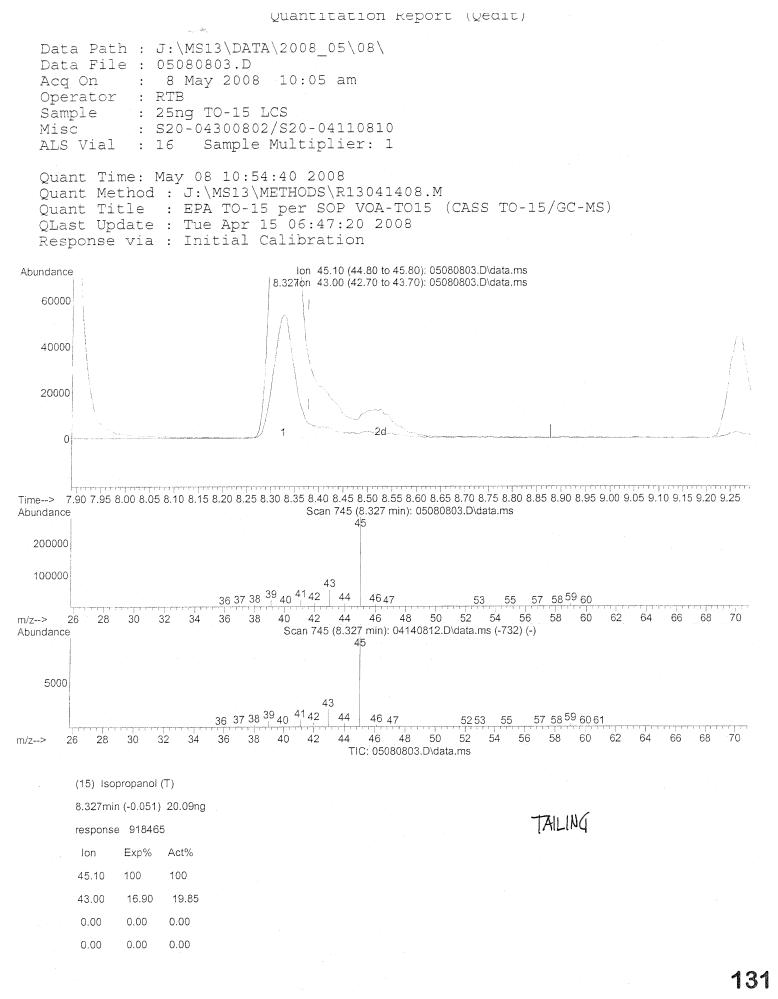
 Internal Standards R.T. QIon Response Conc Units Dev(Min)

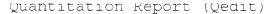
Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080803.D Acq On : 8 May 2008 10:05 am Operator : RTB Sample : 25ng TO-15 LCS Misc : S20-04300802/S20-04110810 ALS Vial : 16 Sample Multiplier: 1 Quant Time: May 08 10:55:04 2008 Ouant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP 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 _____

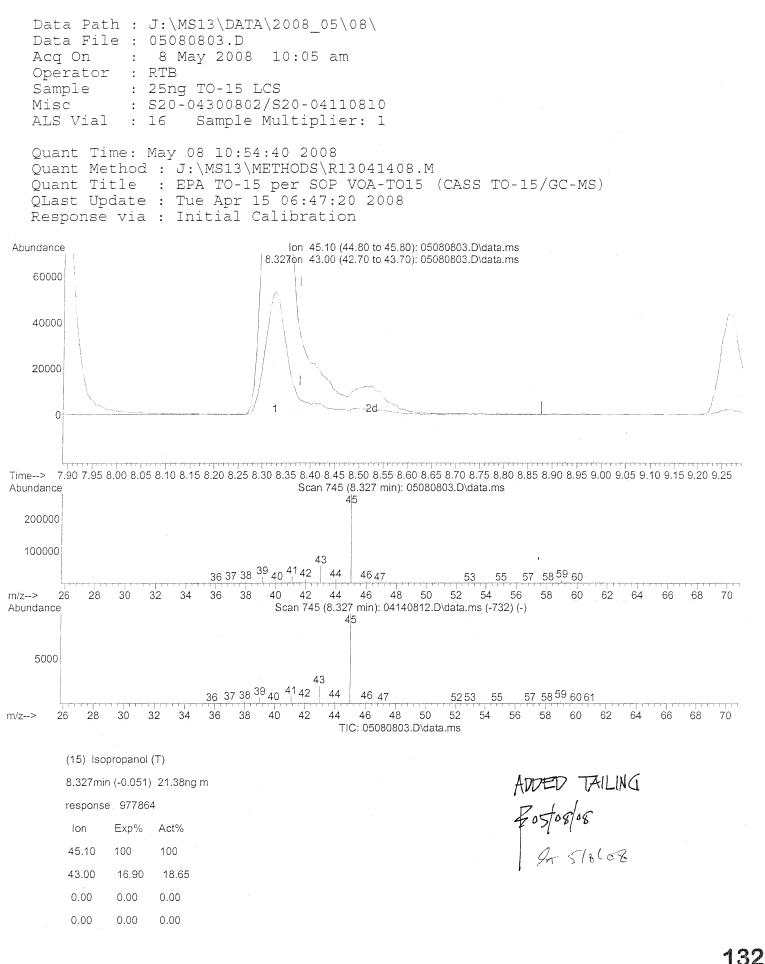
(#) = gualifier out of range (m) = manual integration (+) = signals summed

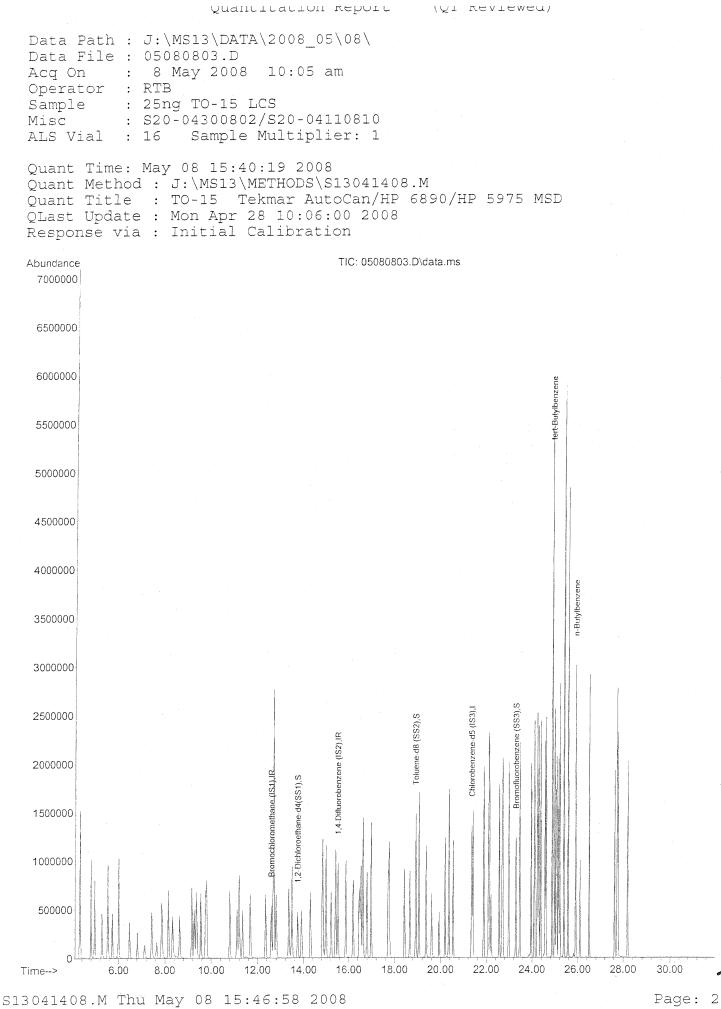
R13041408.M Thu May 08 10:55:26 2008

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	
<b>Client Sample ID:</b>	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

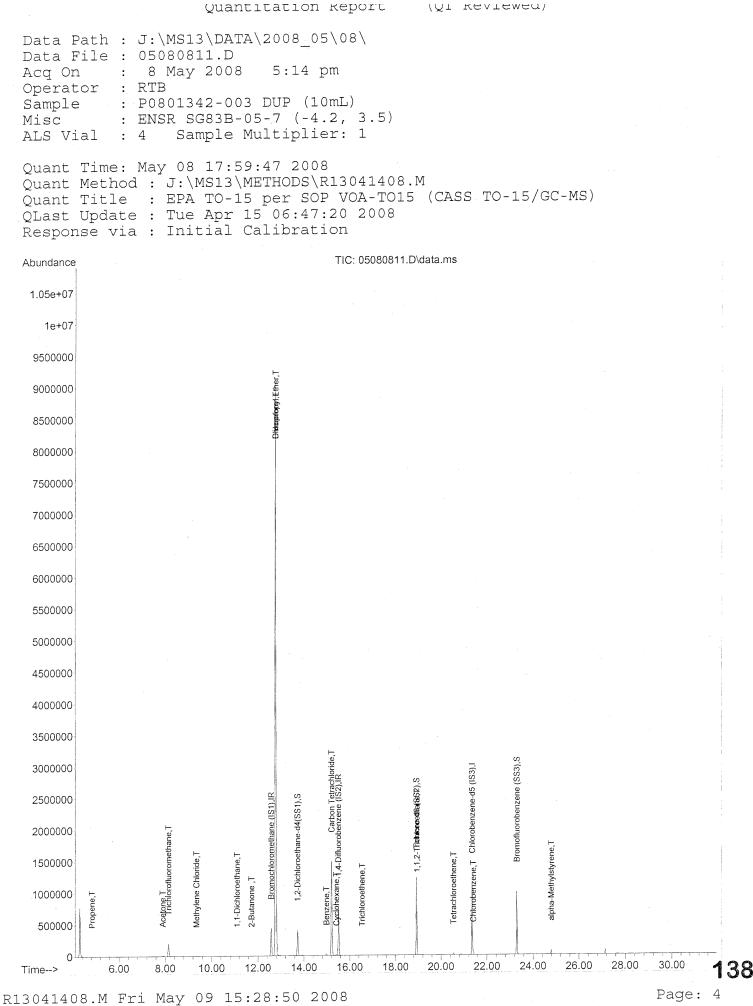
Final Pressure (psig): 3.5

	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	

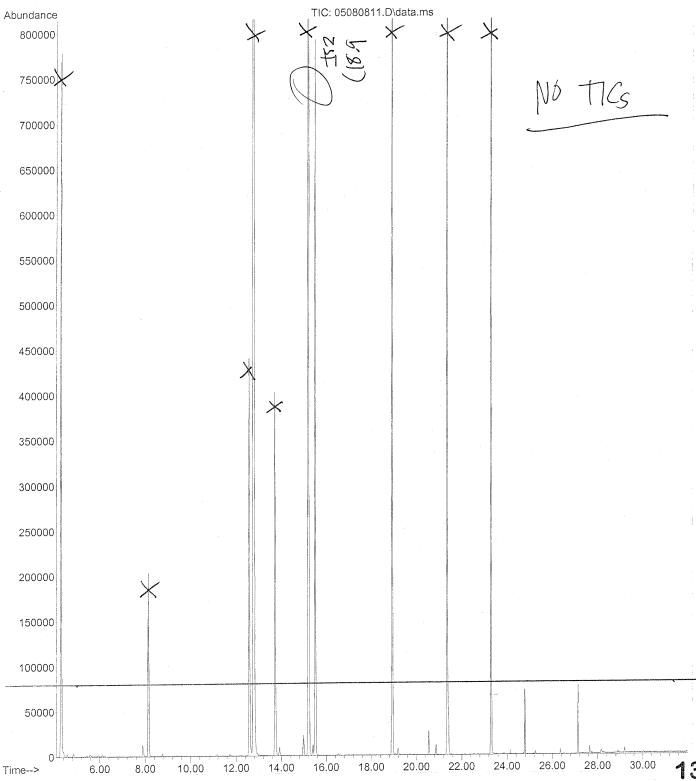
ND = Compound was analyzed for, but not detected above the laboratory detection limit.

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 -<del>0.040</del> 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.	<b>D</b> 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 <b>4</b> # 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.

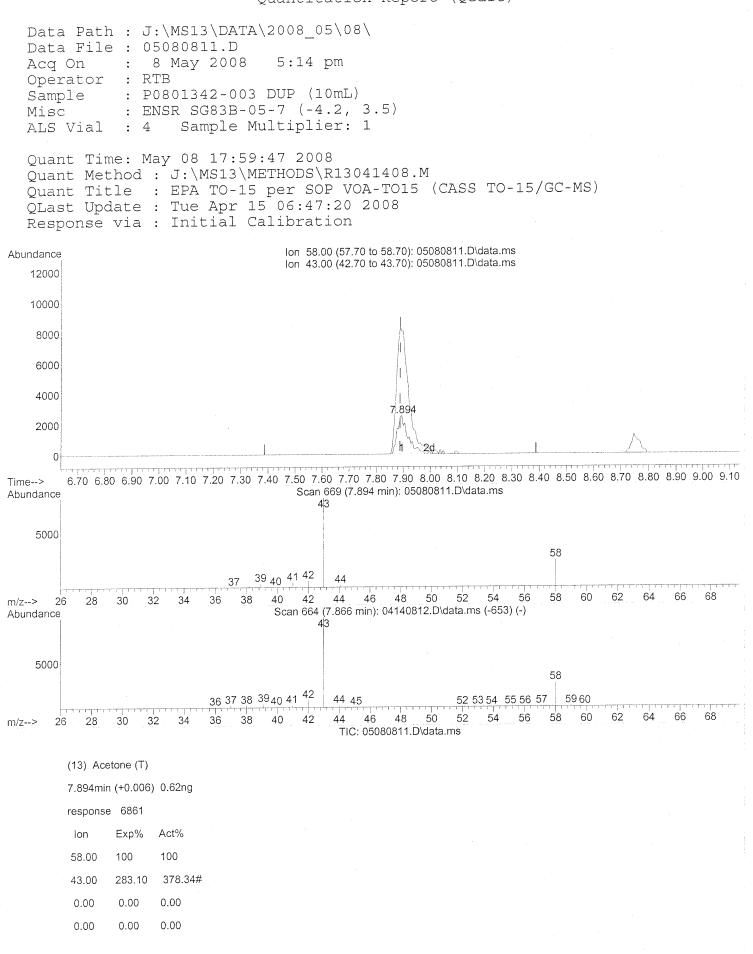
 93) n-Undecane
 26.50
 57
 602
 N.D.

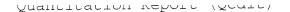
 94) 1,2,4-Trichlorobenzene
 26.50
 57
 602
 N.D.

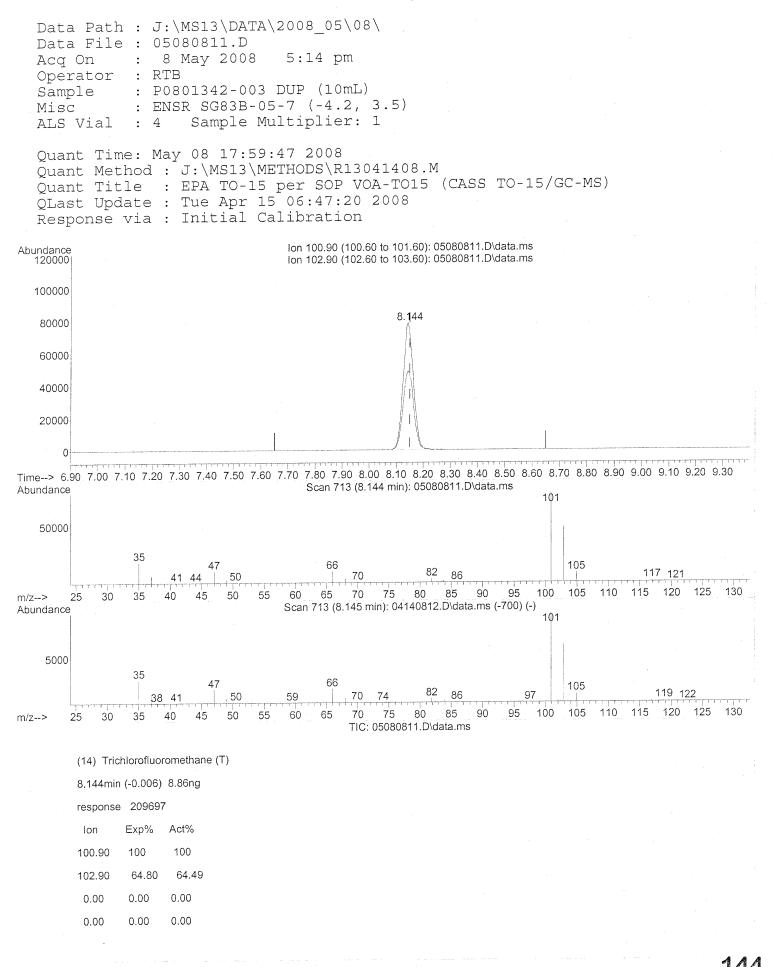
 95) Naphthalene
 27.73
 57
 703
 

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

405/20/108

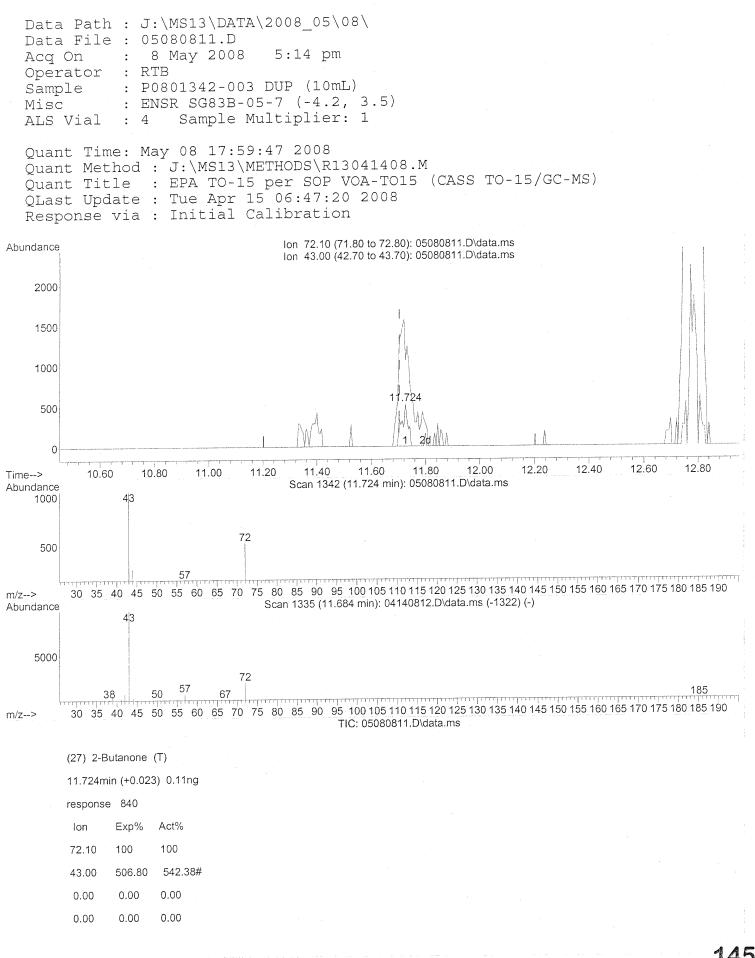


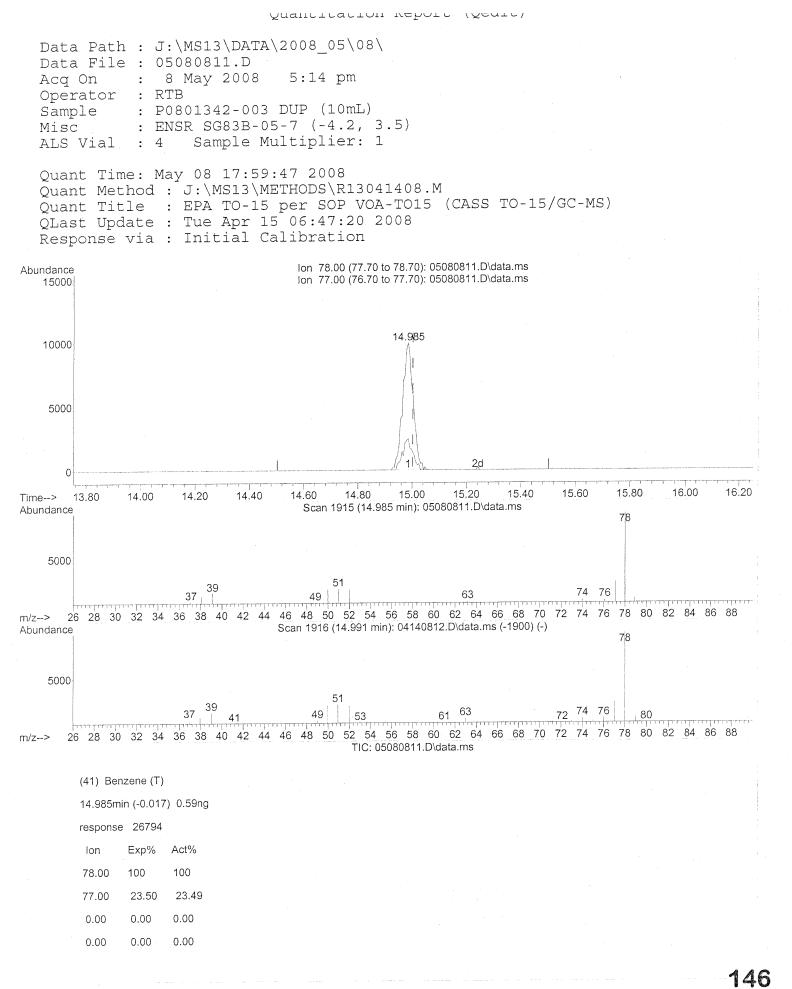




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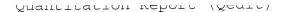
R13041408.M Thu May 08 18:22:43 2008

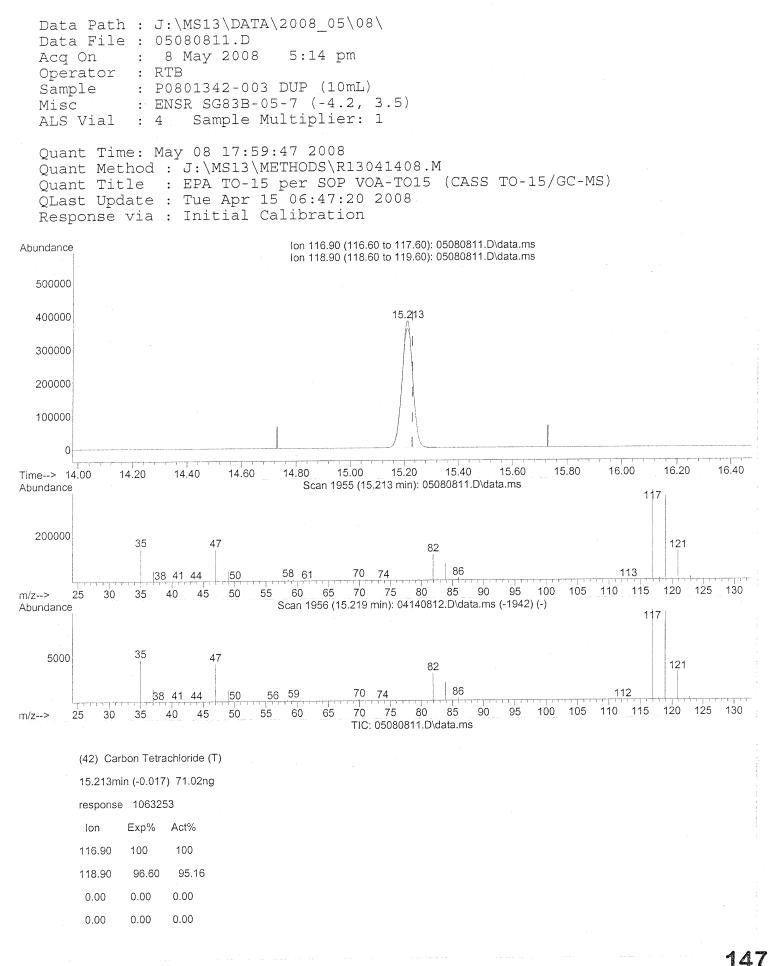




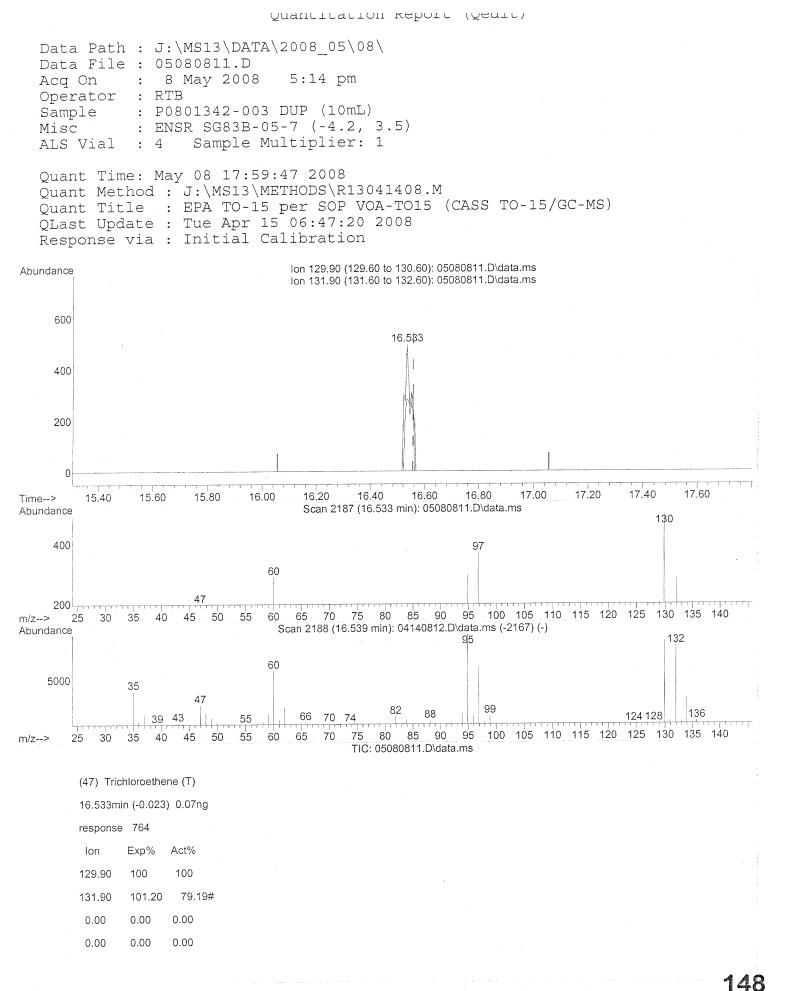
R13041408.M Thu May 08 18:23:39 2008

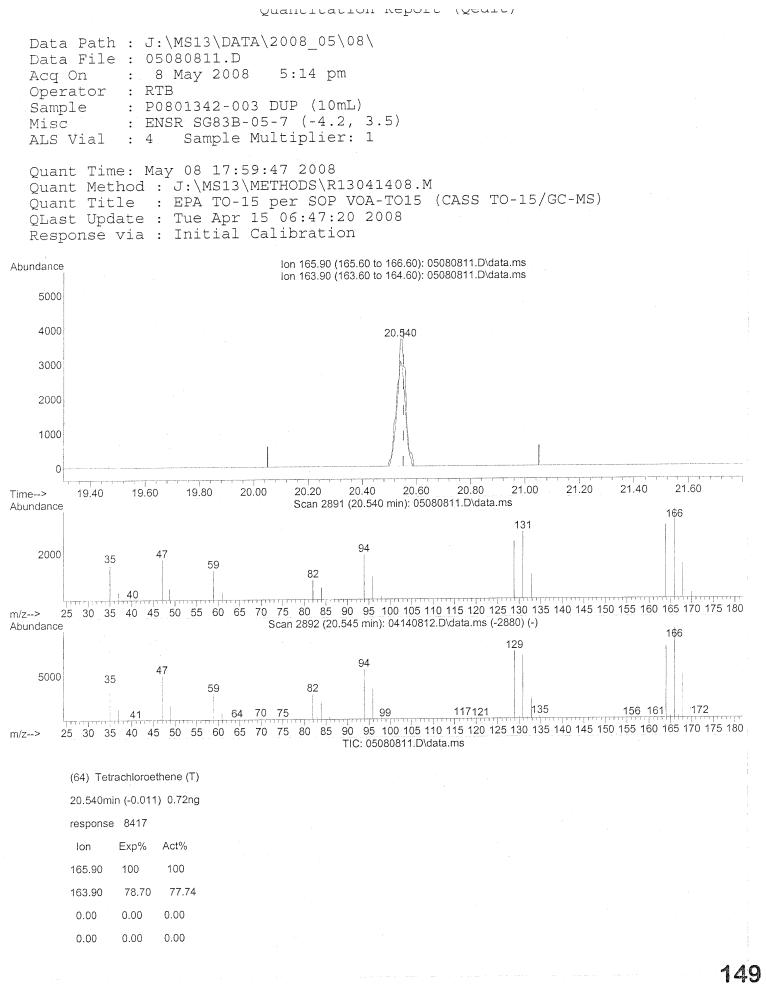
Page: 1

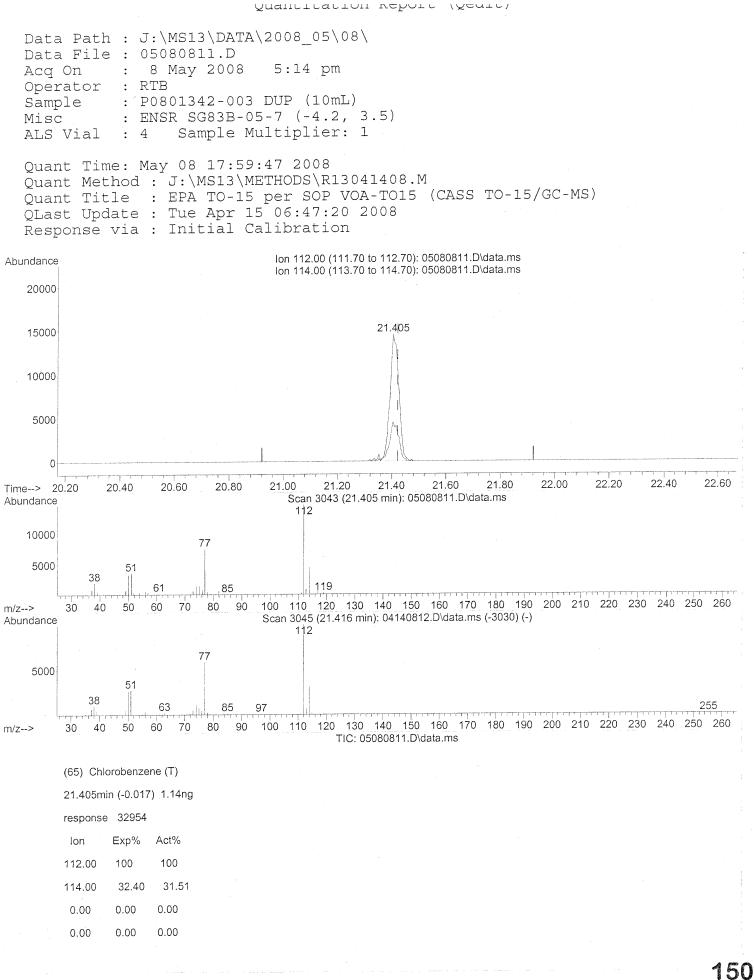




R13041408.M Thu May 08 18:23:44 2008







(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

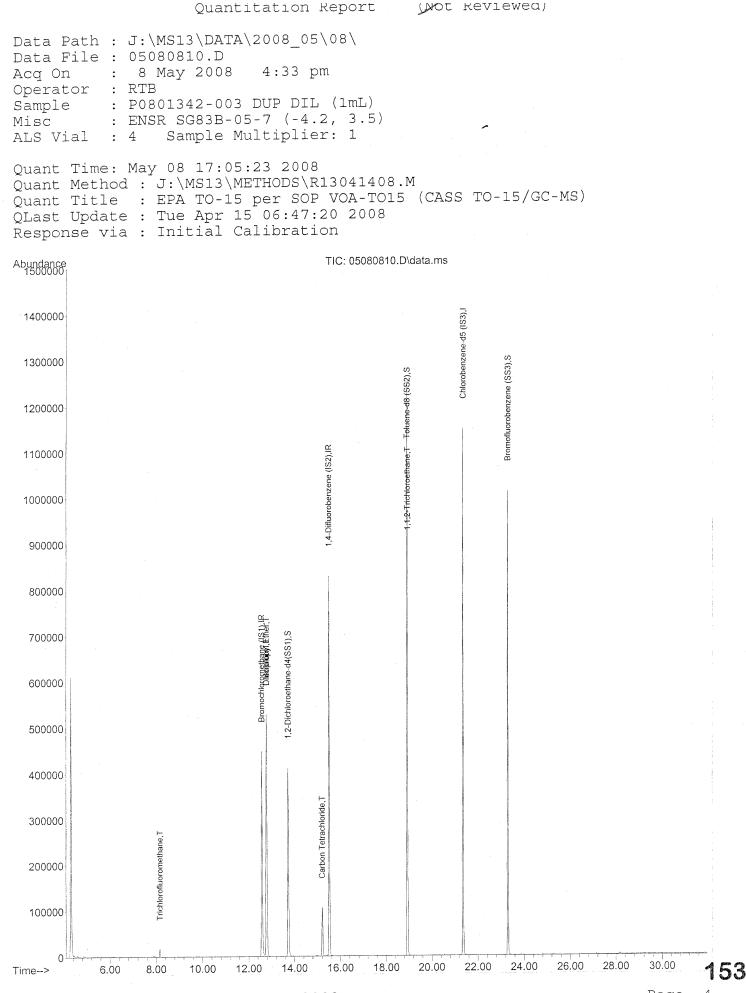
Abundance TIC: 05080811.D\data.ms	
1.05e+07	
1e+07	
950000	
900000	
850000	
800000	
7500000	
700000	
6500000	
600000	
5500000	
500000	
4500000	
400000	
3500000	
3000000	
0000001         0000002         0000002         0000002           1         Toluene-d8 (SS2),S         0000005         0000005           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 <b>151</b> 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\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 0.00\\ 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 <b>15</b> 4
13041408.M Thu May 08 17:11:42 20	08			05/18/18	Page: 1

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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)	
<ul> <li>32) Chloroform</li> <li>34) Tetrahydrofuran</li> <li>35) Ethyl tert-Butyl Ether</li> <li>36) 1,2-Dichloroethane</li> <li>38) 1,1,1-Trichloroethane</li> <li>39) Isopropyl Acetate</li> <li>40) 1-Butanol</li> <li>41) Benzene</li> <li>42) Carbon Tetrachloride</li> </ul>	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- &amp; 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

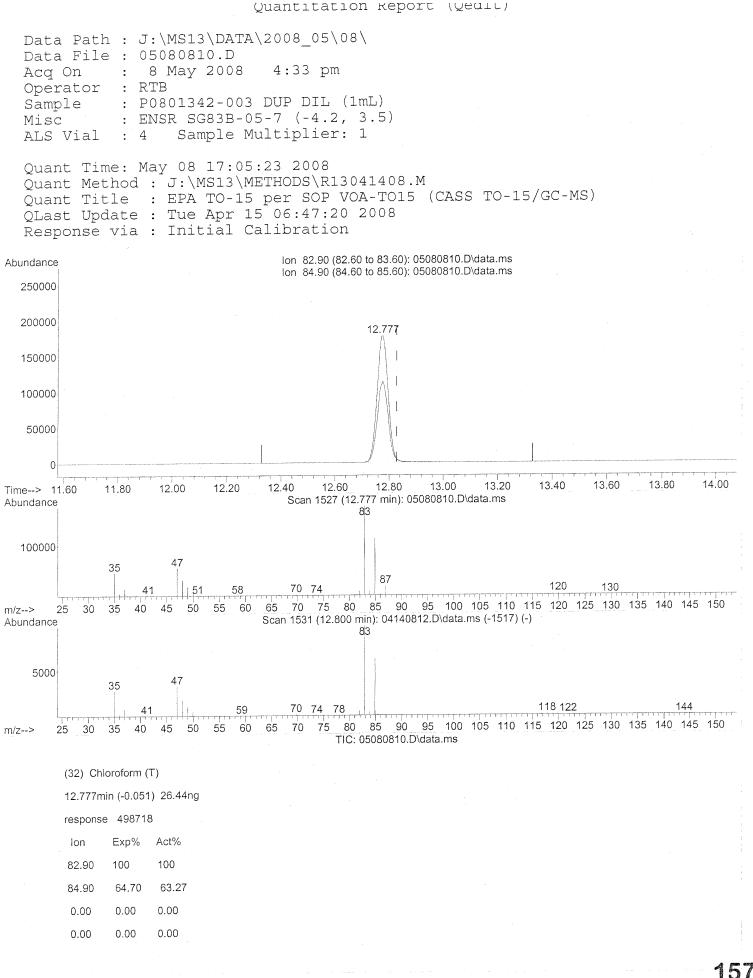
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Prost. 8/08

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 #	<b>RT</b> #	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

1 );;;) ) 5

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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 <b>5</b> <b>0</b> <b>5</b> <b>0</b>	.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

4

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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)	<del>[</del> 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	<b>l</b> 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)	<del>[</del> 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)	<del>[</del> 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
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(#) = Out of Range

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

4

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	$\chi$	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)

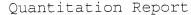
Method Path : J:\MS13\METHODS\ Method File : R13041408.M Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) Last Update : Tue Apr 15 06:47:20 2008 Response Via : Initial Calibration

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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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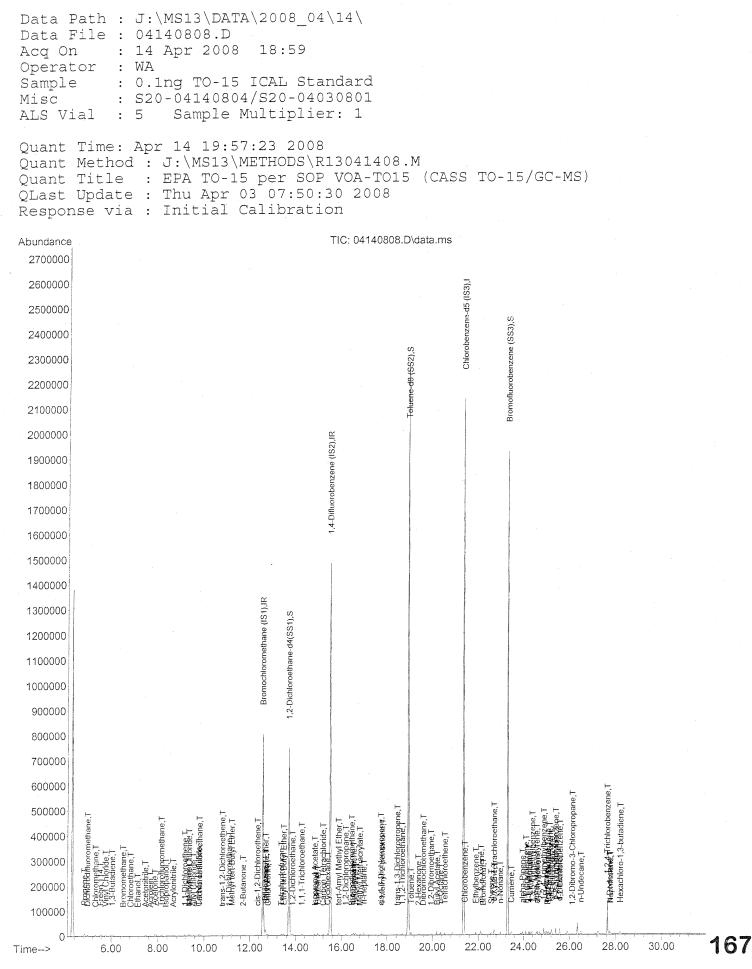
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2	0.5	Apr 15 06:33 2008	Apr 14 20:24 2008	14 Apr 2008 19:40			
3	1.0	Apr 15 06:33 2008	Apr 15 06:20 2008	14 Apr 2008 20:21			
4	5.0	Apr 15 06:34 2008	Apr 15 06:22 2008	14 Apr 2008 21:01			
5	25	Apr 15 06:34 2008	Apr 15 06:24 2008	14 Apr 2008 21:43			
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
<ul> <li>6) Vinyl Chloride</li> <li>7) 1,3-Butadiene</li> <li>8) Bromomethane</li> <li>9) Chloroethane</li> <li>10) Ethanol</li> <li>11) Acetonitrile</li> <li>12) Acrolein</li> <li>13) Acetone</li> <li>14) Trichlorofluoromethane</li> <li>15) Isopropanol</li> <li>16) Acrylonitrile</li> <li>17) 1,1-Dichloroethene</li> <li>18) tert-Butanol</li> <li>19) Methylene Chloride</li> <li>20) Allyl Chloride</li> <li>21) Trichlorotrifluoroethane</li> </ul>	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) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 42) 43) 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) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 52) 53) 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 <b>169</b>

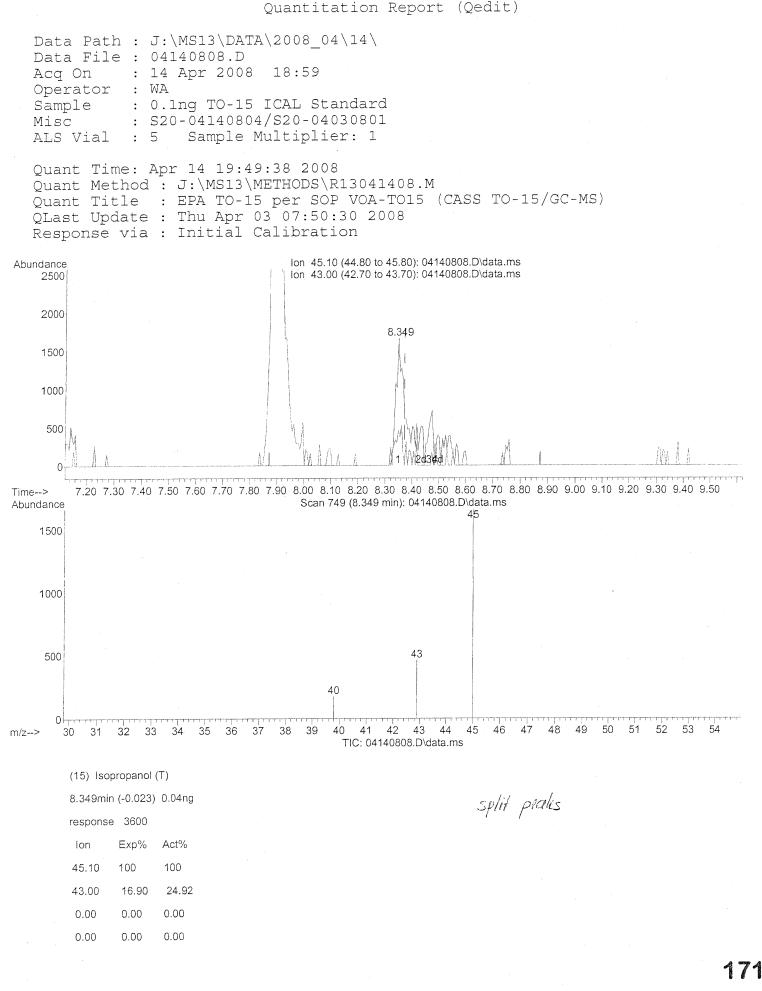
213041408.M Mon Apr 14 20:02:23 2008

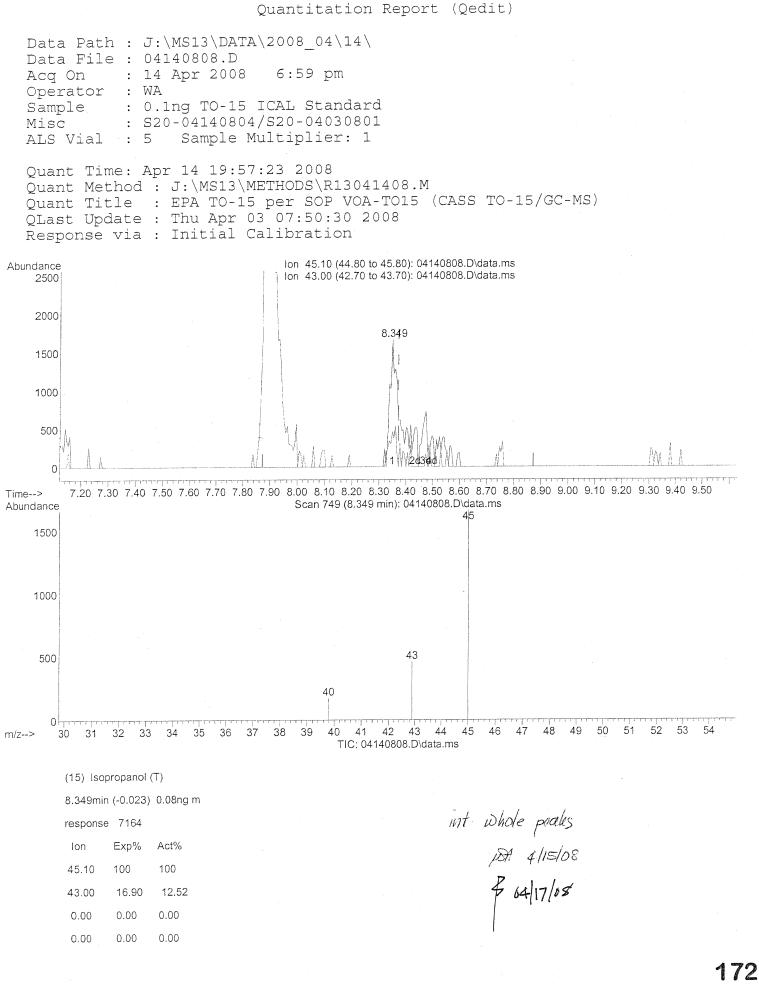
jor 4/15/08

Page: 2

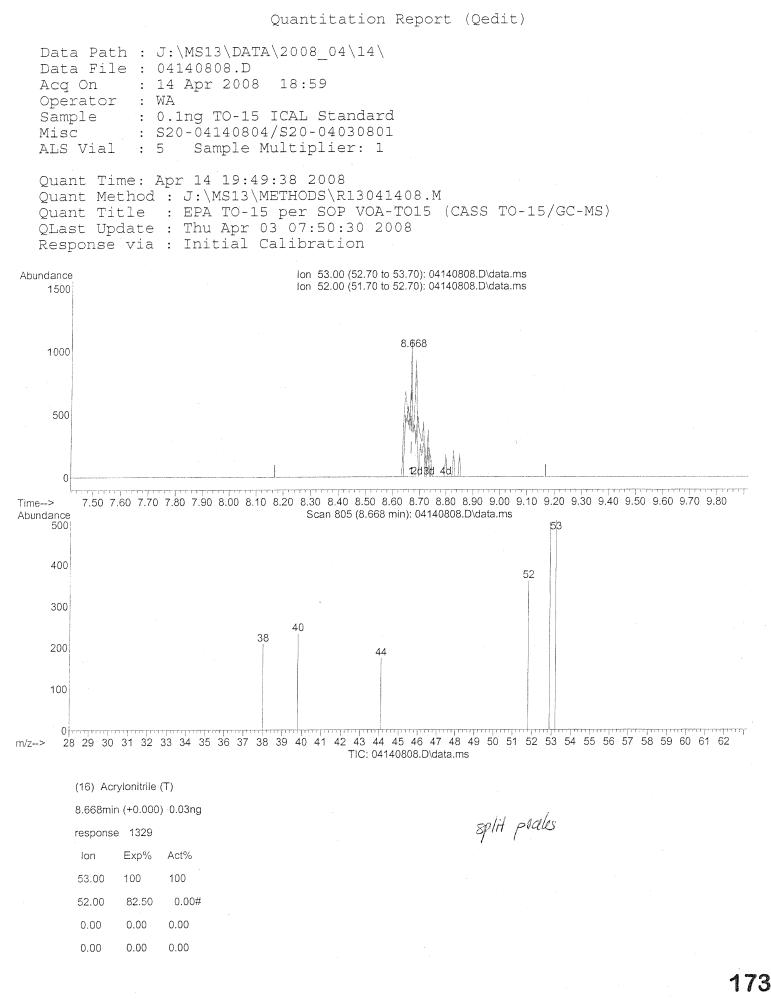
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Quant Time: Apr 14 19:57:23 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Thu Apr 03 07:50:30 2008 Response via : Initial Calibration							
Internal Standards	R.T.	QIon	Response	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	

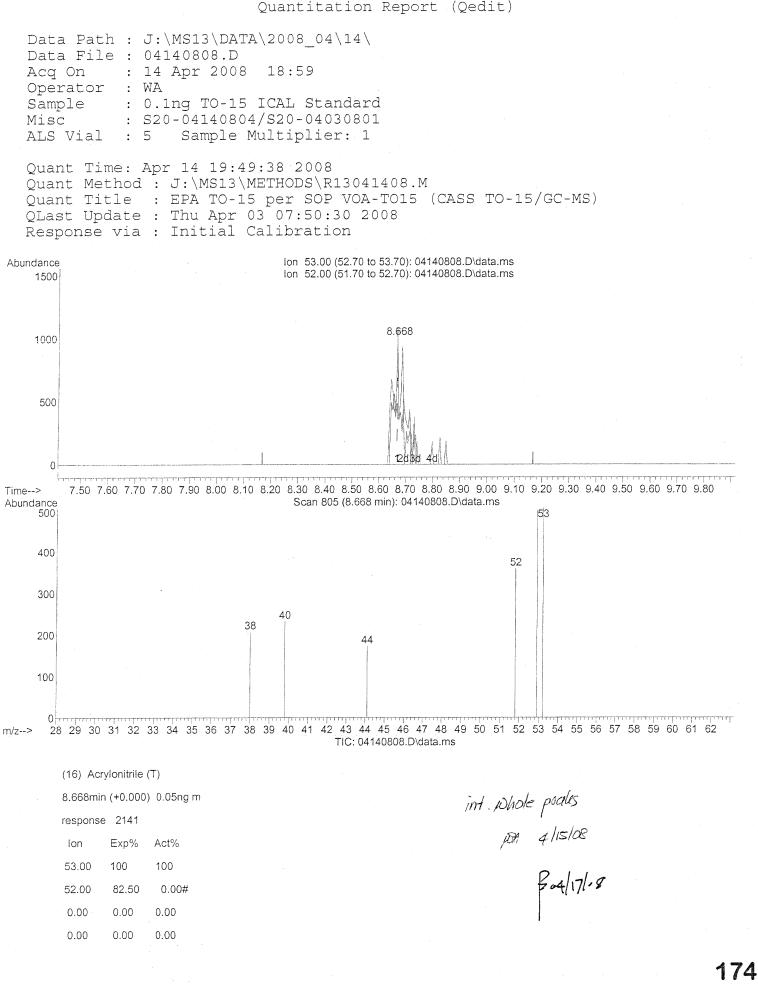
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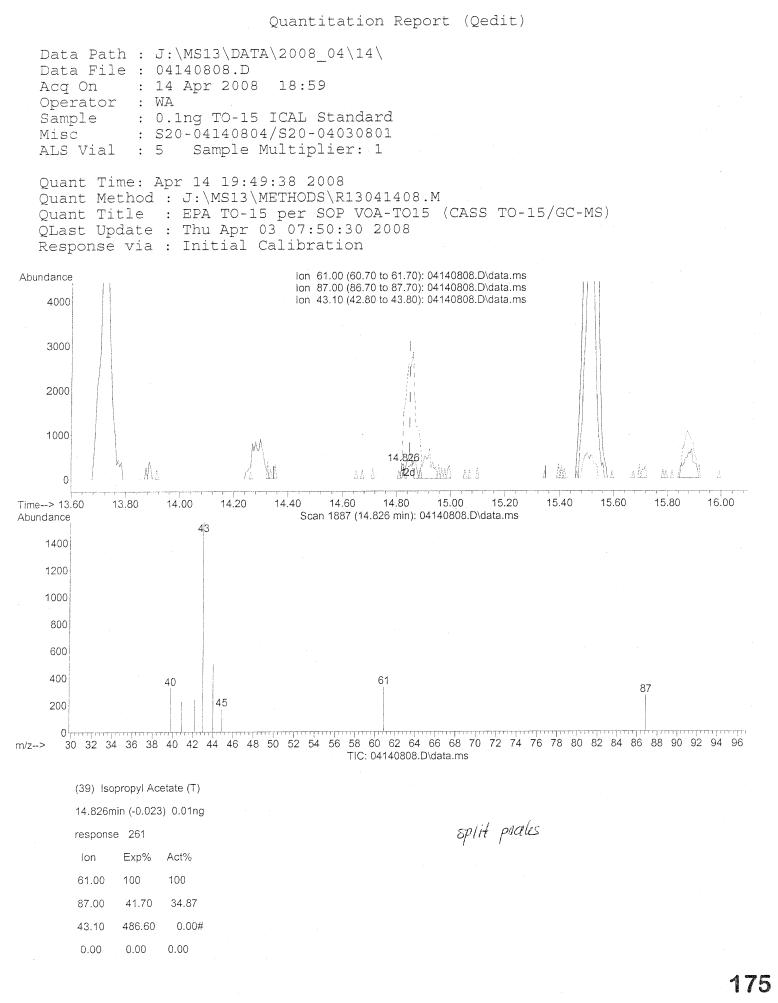


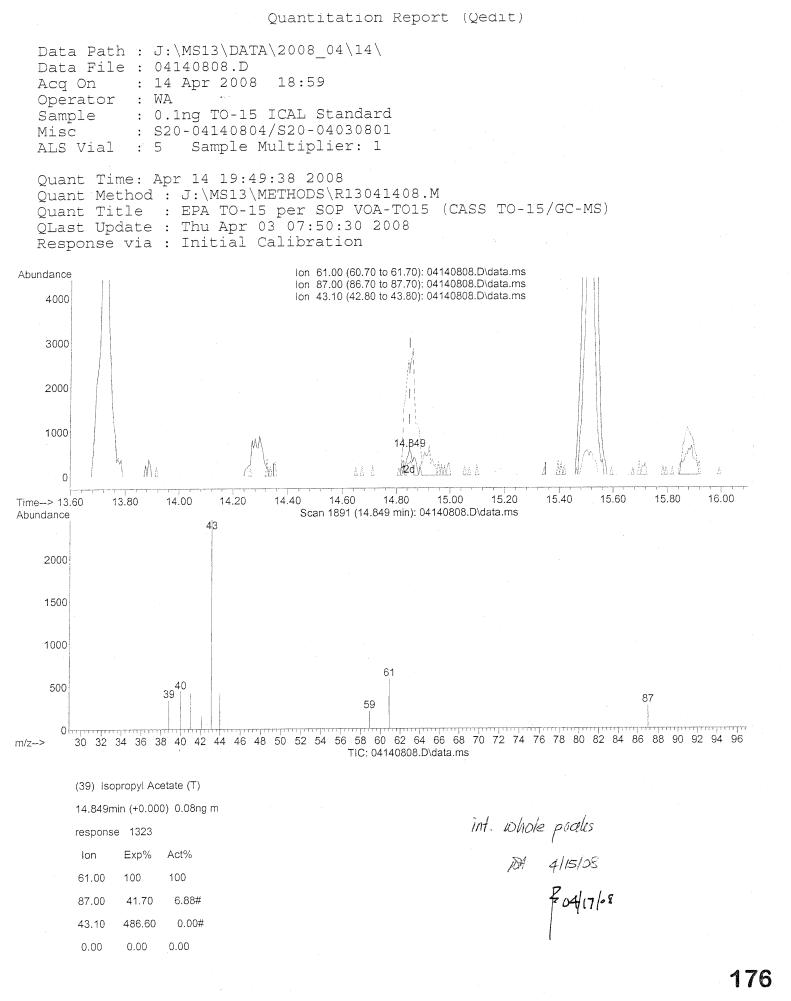
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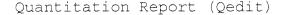


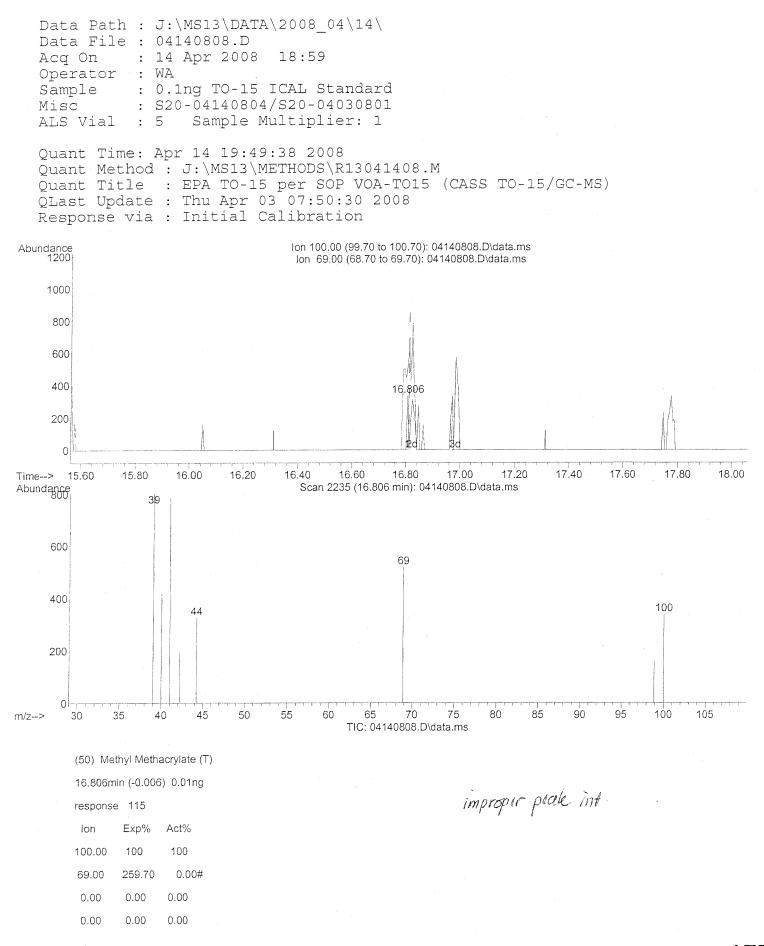


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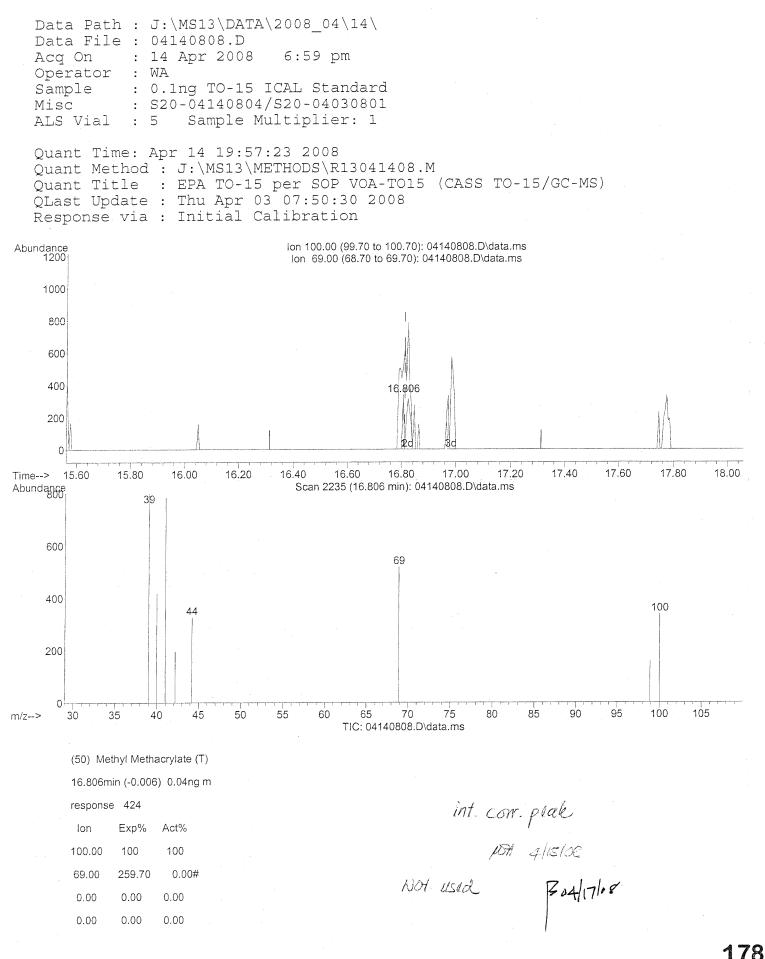




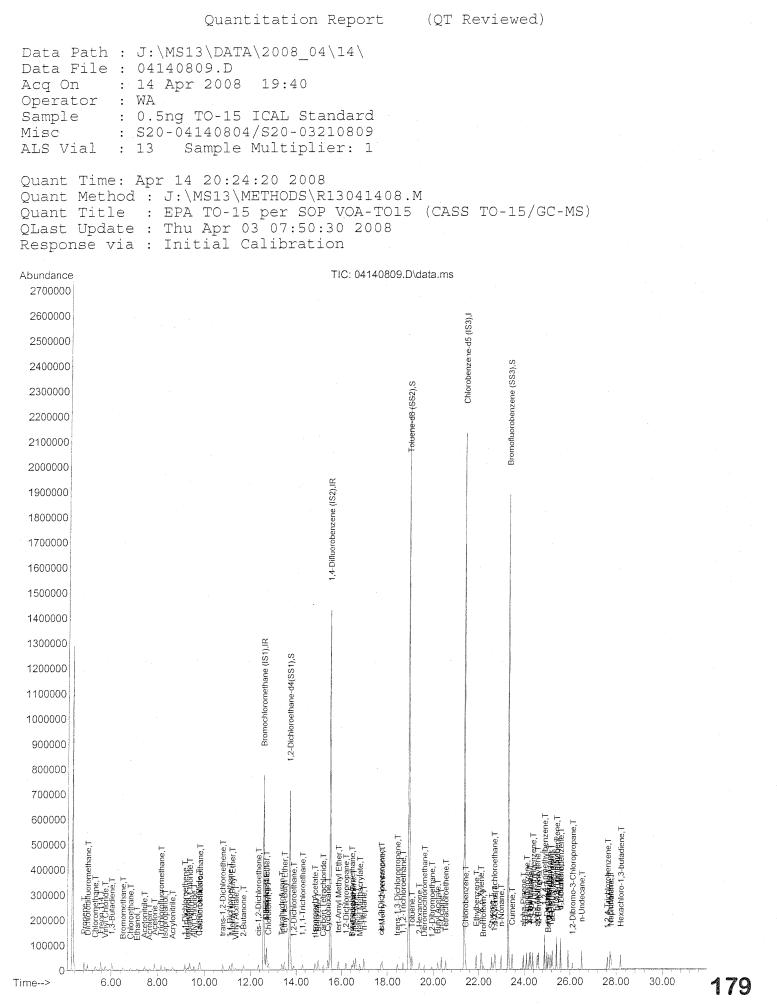


Page: 1





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



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

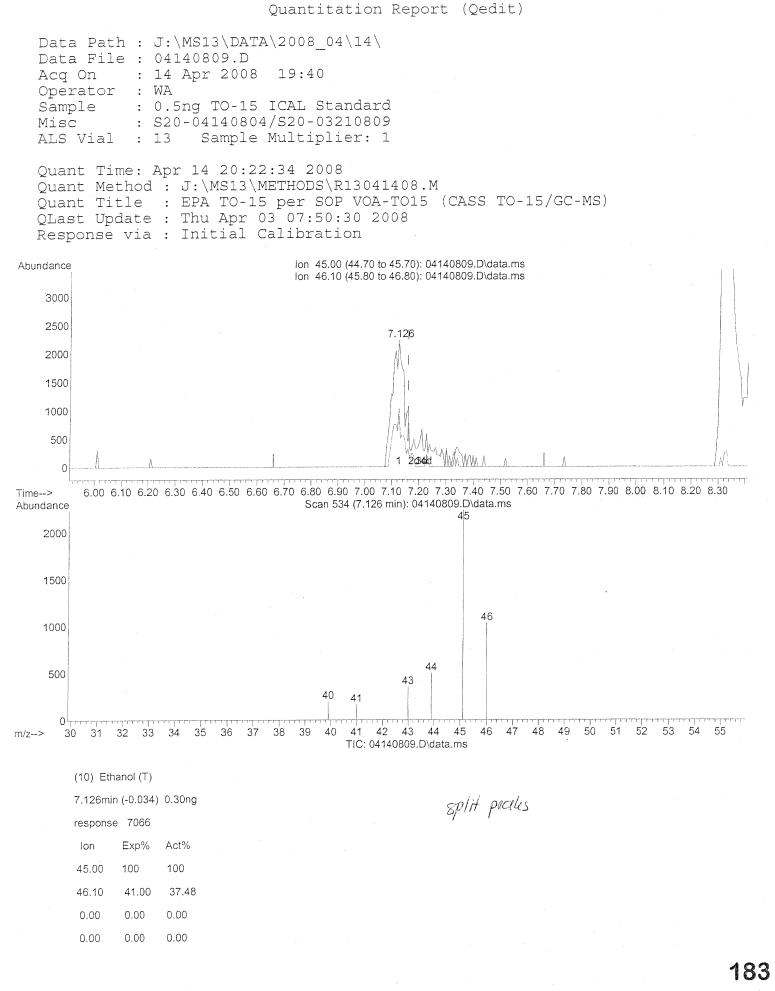
Data Path : J:\MS13\DATA\2008_04 Data File : 04140809.D Acq On : 14 Apr 2008 19:40 Operator : WA Sample : 0.5ng TO-15 ICAL Sta Misc : S20-04140804/S20-032 ALS Vial : 13 Sample Multipli Quant Time: Apr 14 20:24:20 2008 Quant Method : J:\MS13\METHODS\R Quant Title : EPA TO-15 per SOP QLast Update : Thu Apr 03 07:50: Response via : Initial Calibrati	ndard 10809 er: 1 1304140 VOA-TO 30 2008	8.M	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	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 0.387 0.385 0.387 0.385 0.387 0.385 0.387 0.385 0.387 0.385 0.387 0.385 0.387 0.385 0.387 0.385 0.387 0.385 0.387 0.385 0.387 0.385 0.387 0.385 0.387 0.385 0.387 0.385 0.387 0.385 0.387 0.385 0.387 0.385 0.387 0.385 0.387 0.385 0.387 0.385 0.387 0.385 0.387 0.385 0.385 0.385 0.387 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.385 0.396	ng ng ng ng ng ng ng ng ng ng ng ng ng ng ng ng ng ng ng ng ng n	99 97 96 98 # 71 98

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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- &amp; 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 83 99 83 99 83 99 83 99 83 99 83 99 83 99 83 99 83 99 83 99 83 99 83 99 83 99 83 99 83 99 83 99 83 99 83 99 83 99 83 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

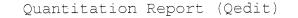
Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140809.D Acq On : 14 Apr 2008 19:40 Operator : WA Sample : 0.5ng TO-15 ICAL Standard Misc : S20-04140804/S20-03210809 ALS Vial : 13 Sample Multiplier: 1 Quant Time: Apr 14 20:24:20 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 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 

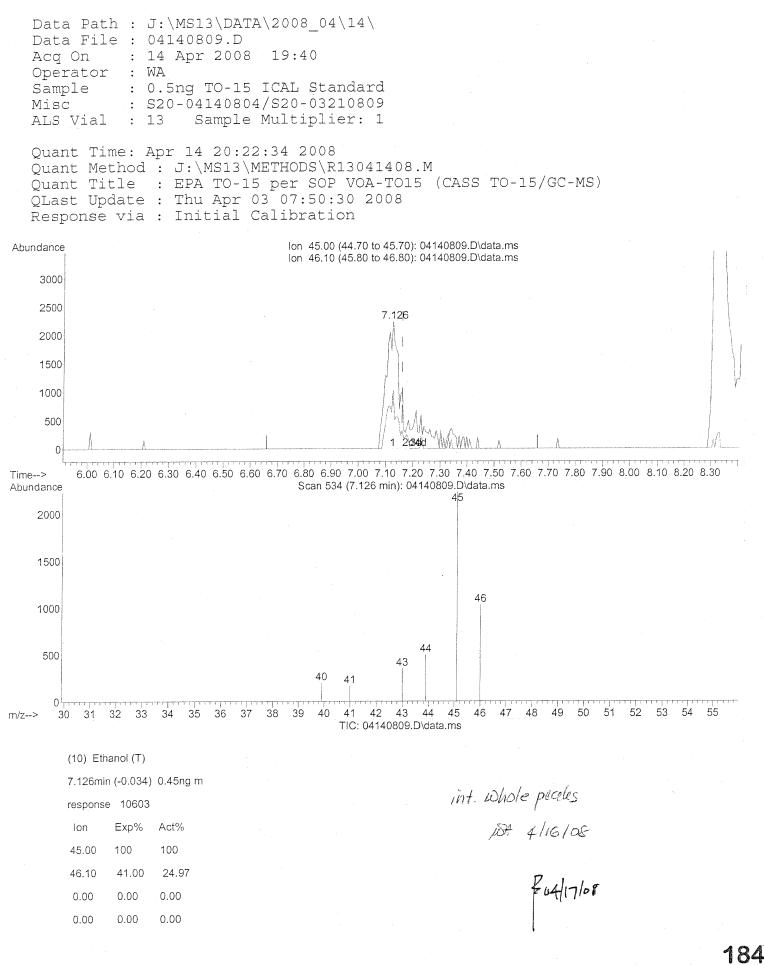
(#) = qualifier out of range (m) = manual integration (+) = signals summed

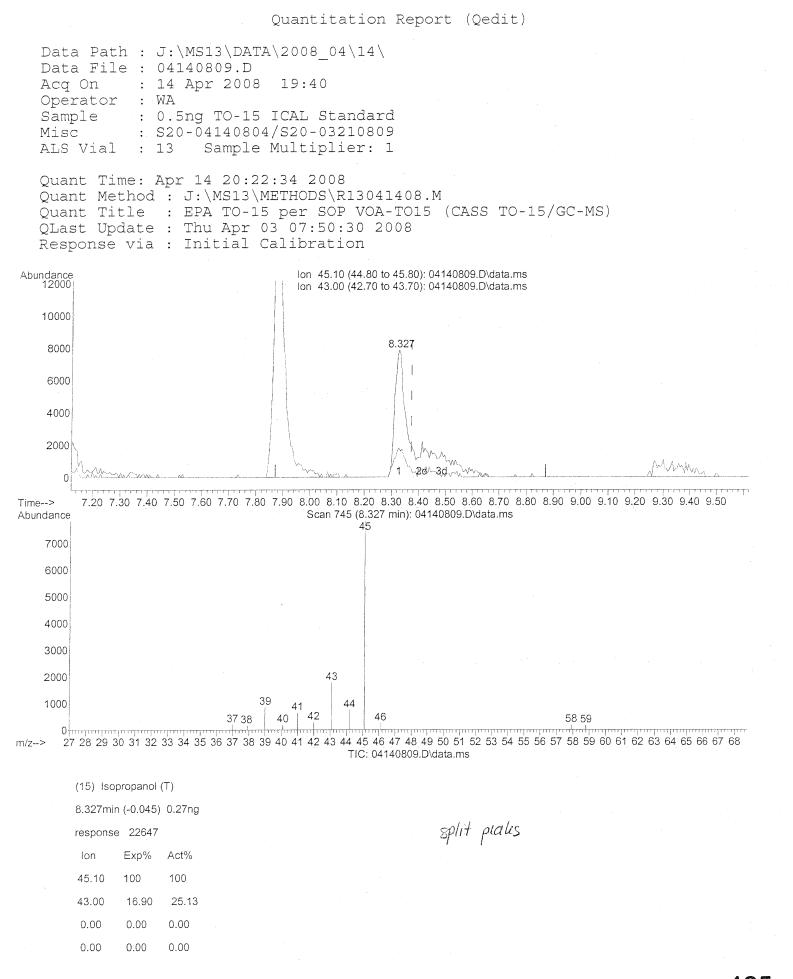


R13041408.M Mon Apr 14 20:23:25 2008

Page: 1

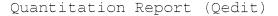


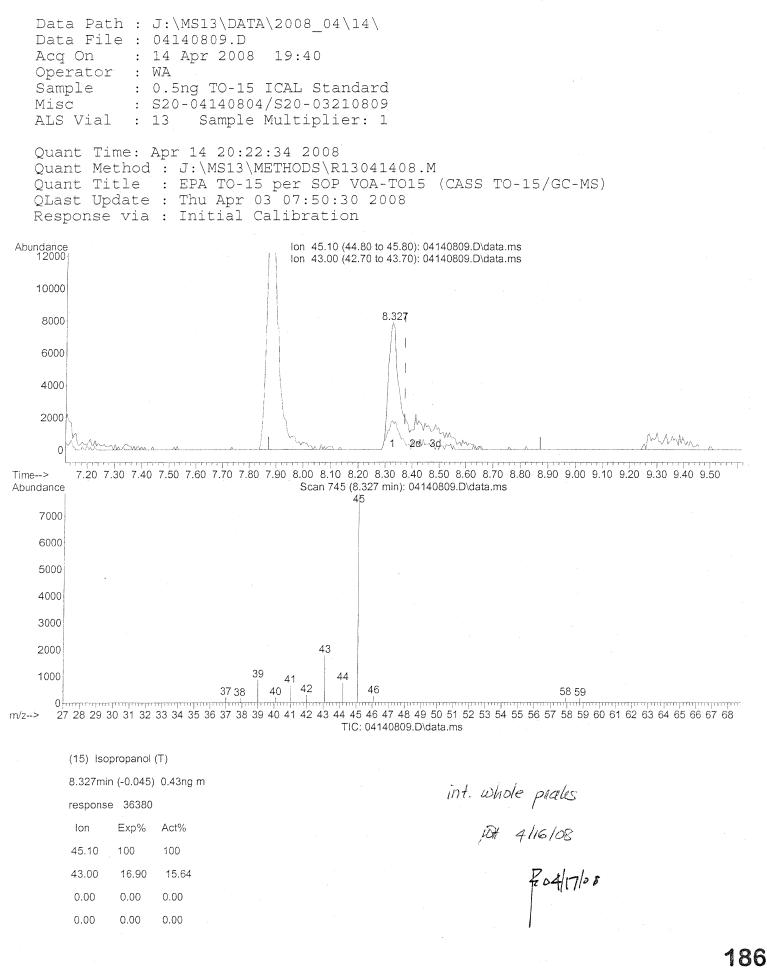




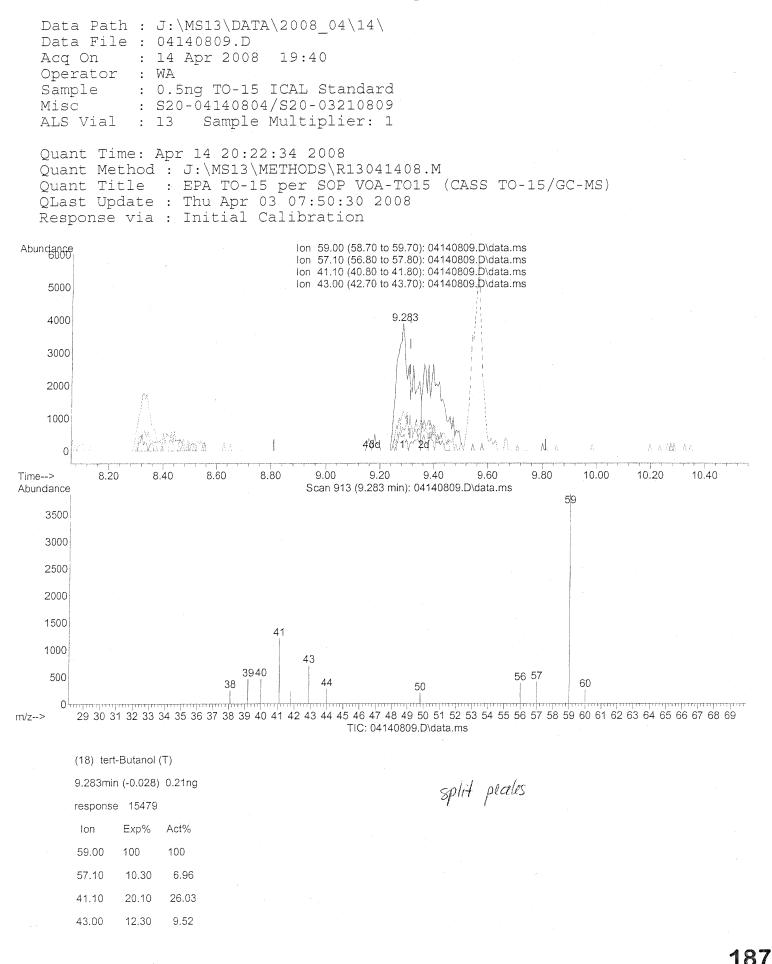
R13041408.M Mon Apr 14 20:24:03 2008

Page: 1

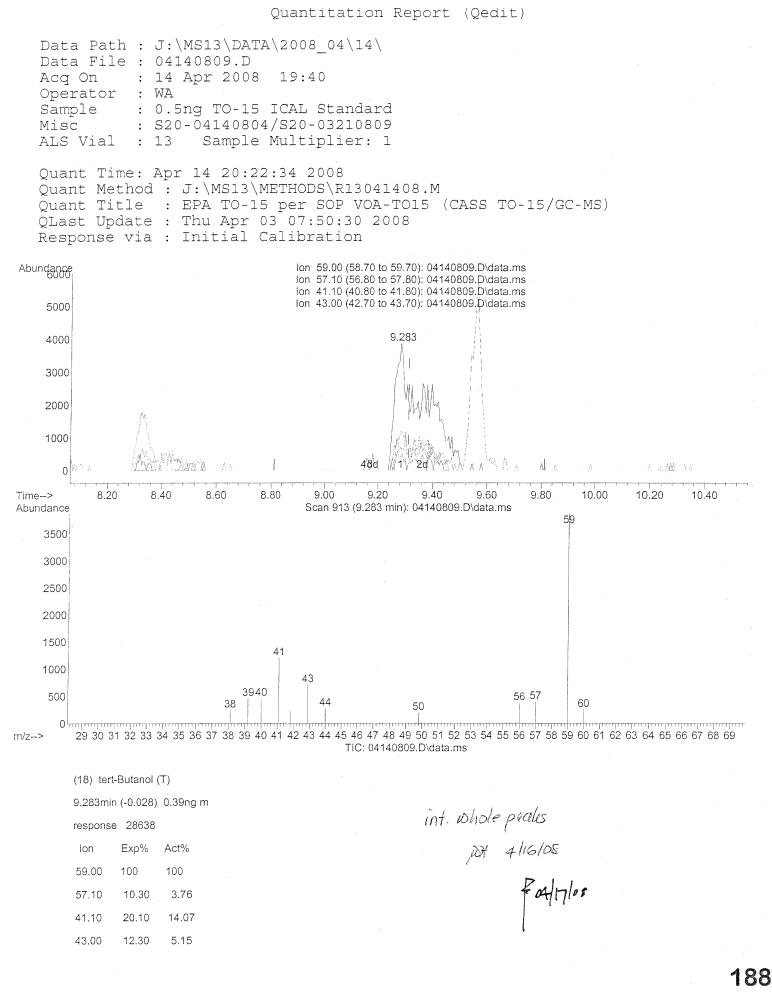


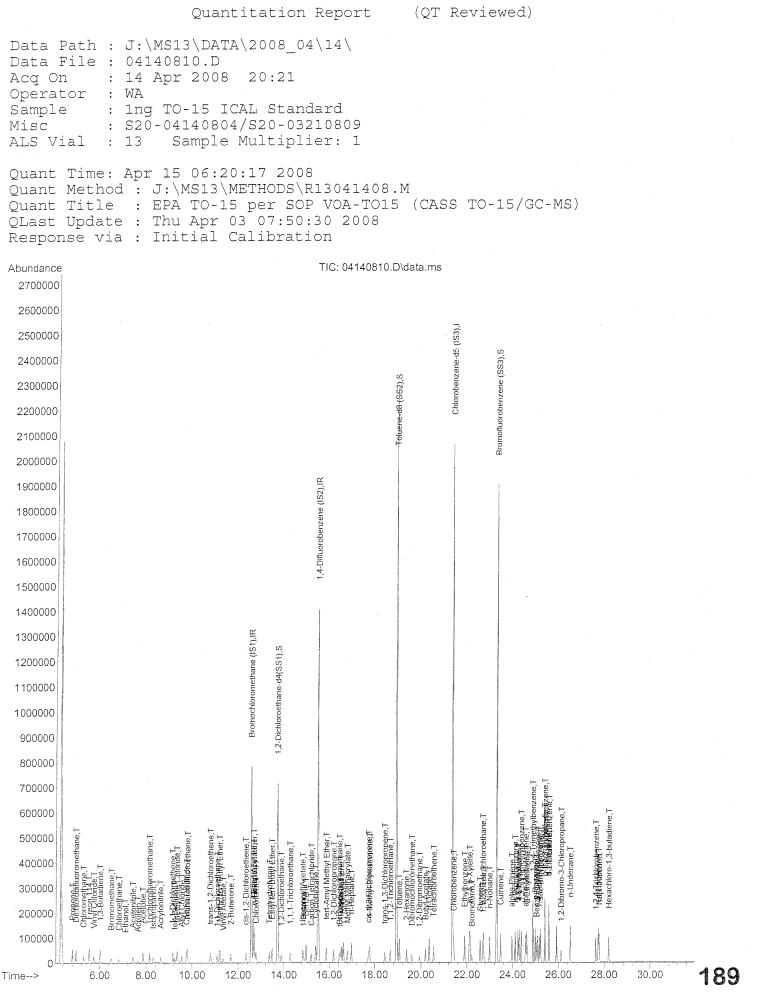


R13041408.M Mon Apr 14 20:24:11 2008



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- &amp; 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 <b>191</b>

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

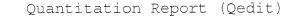
101 4/16/08

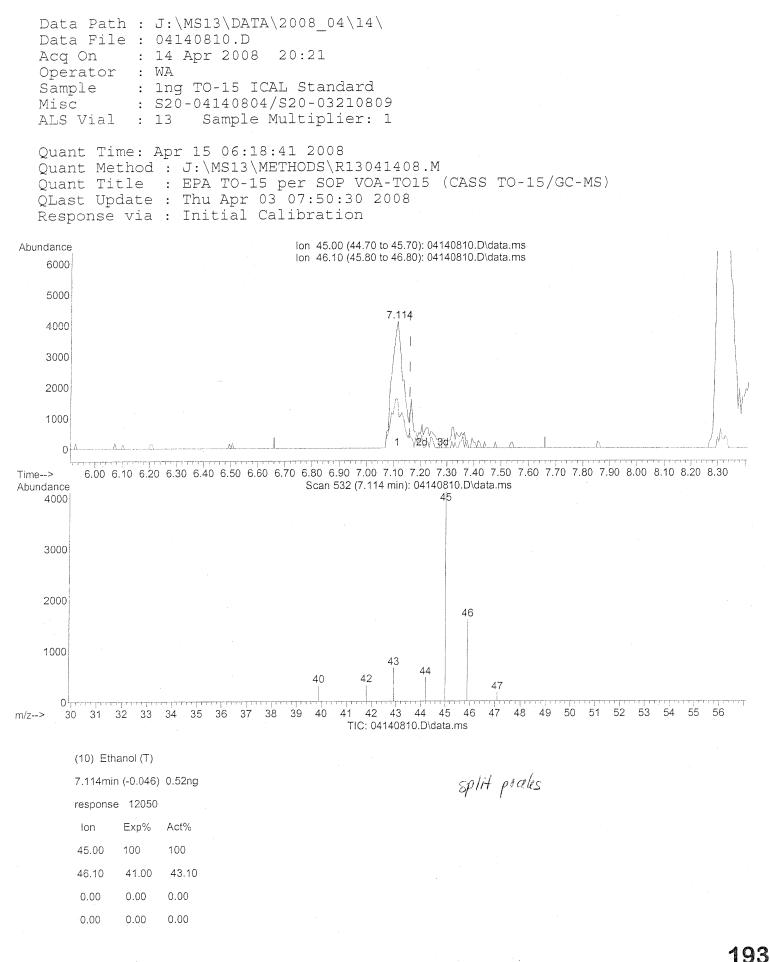
Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140810.D Acq On : 14 Apr 2008 20:21 Operator : WA Sample : 1ng TO-15 ICAL Standard Misc : S20-04140804/S20-03210800 : S20-04140804/S20-03210809 ALS Vial : 13 Sample Multiplier: 1 Quant Time: Apr 15 06:20:17 2008 Quant Method : J:\MS13\METHODS\R13041408.M 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

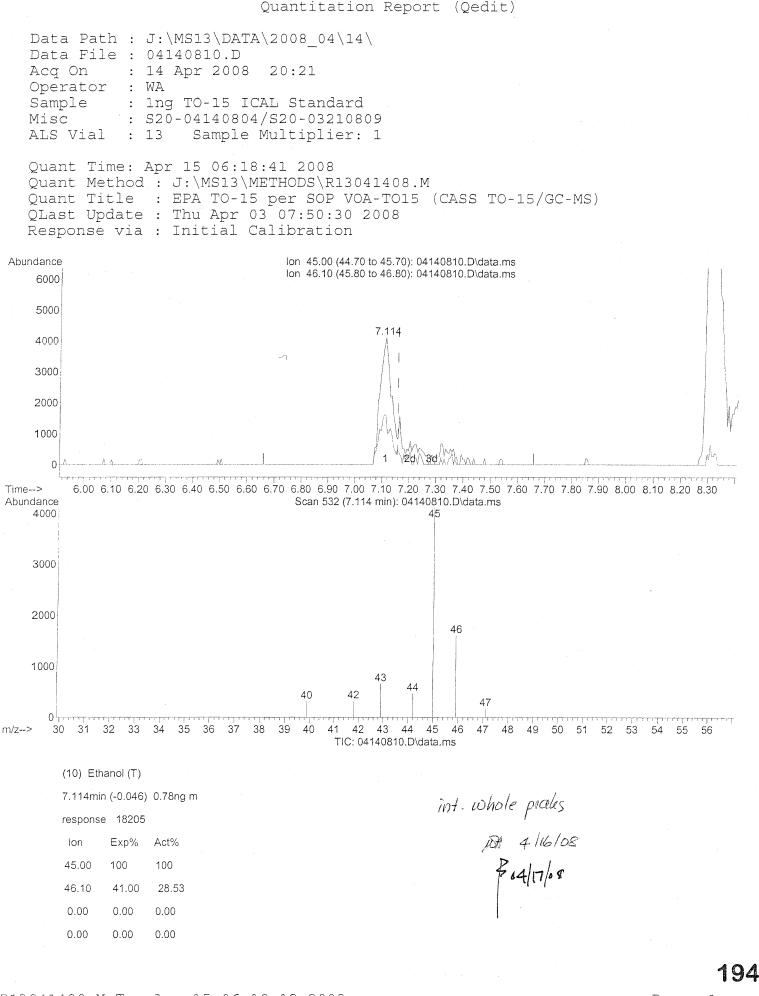
181 4/16/08



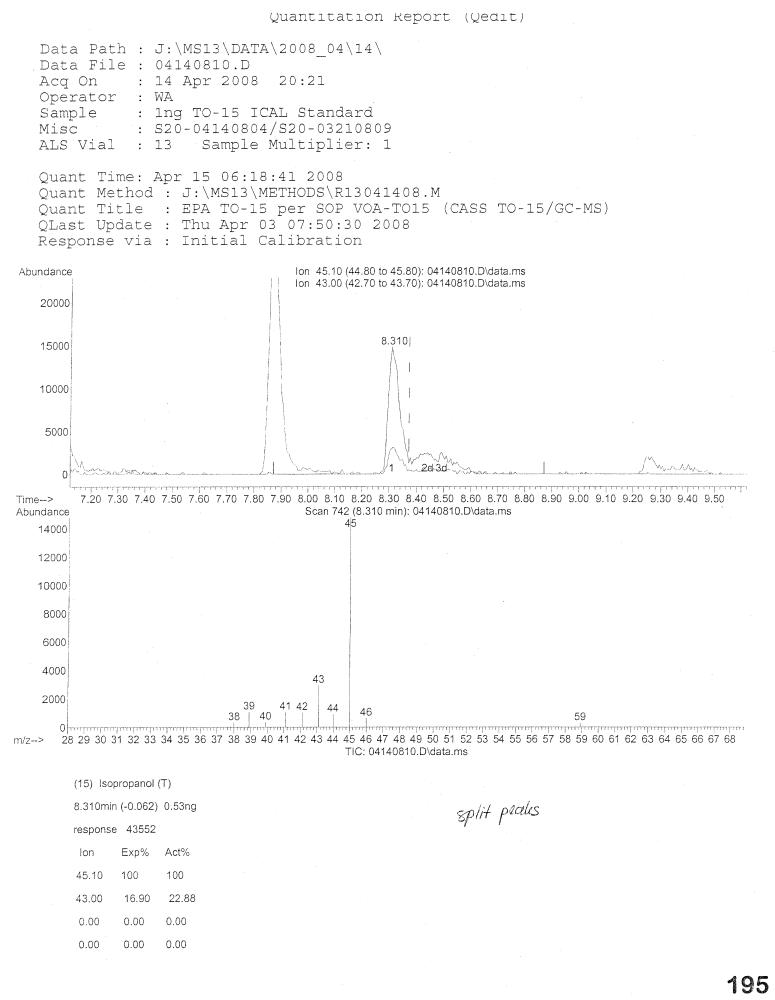


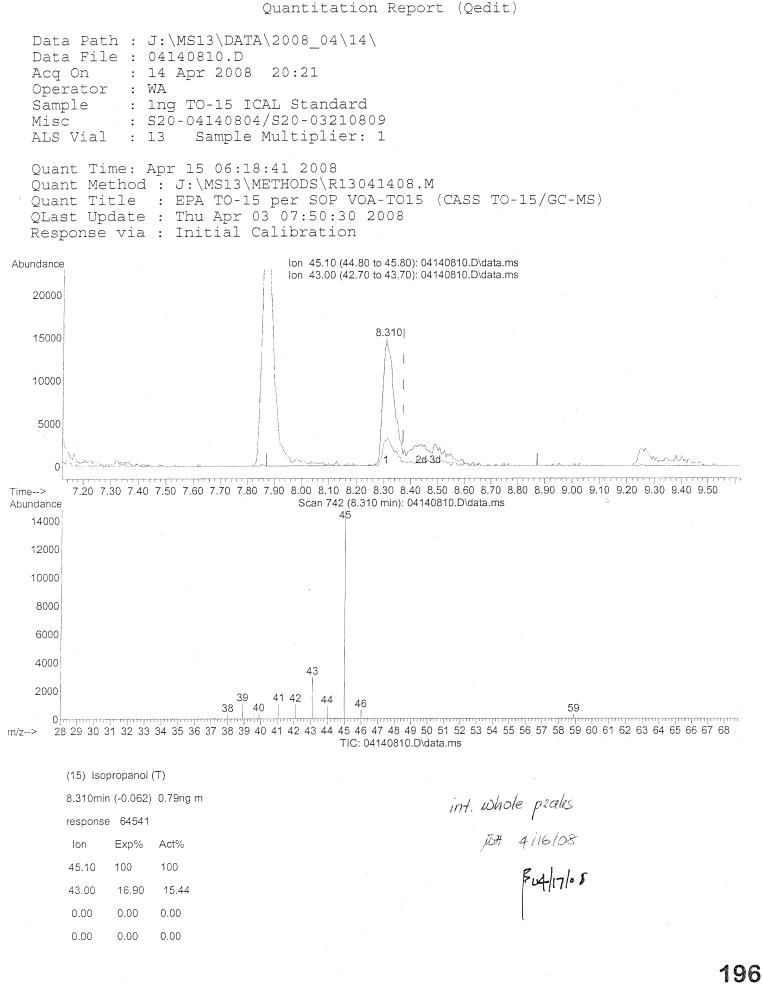
R13041408.M Tue Apr 15 06:19:06 2008

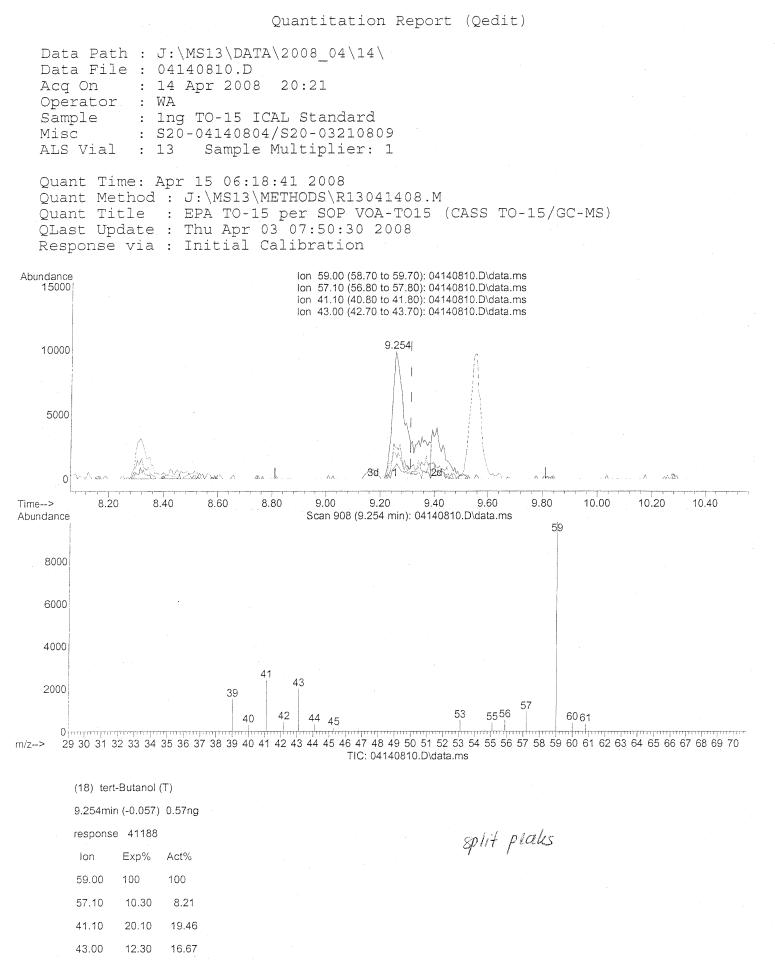
Page: 1



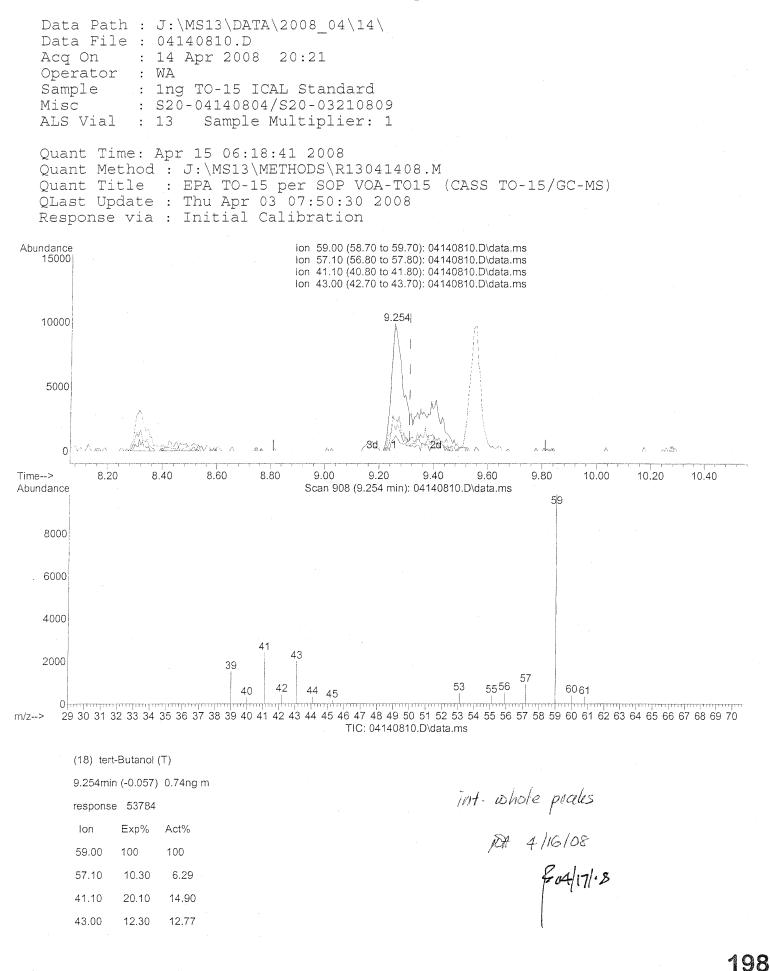
R13041408.M Tue Apr 15 06:19:12 2008





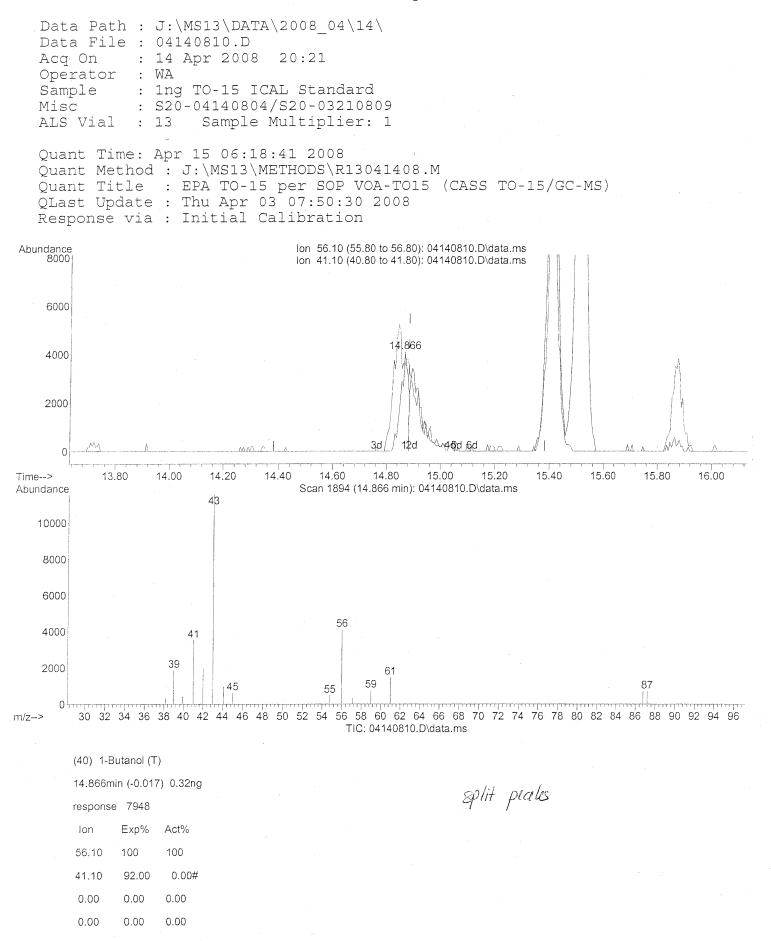


R13041408.M Tue Apr 15 06:19:36 2008

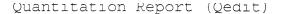


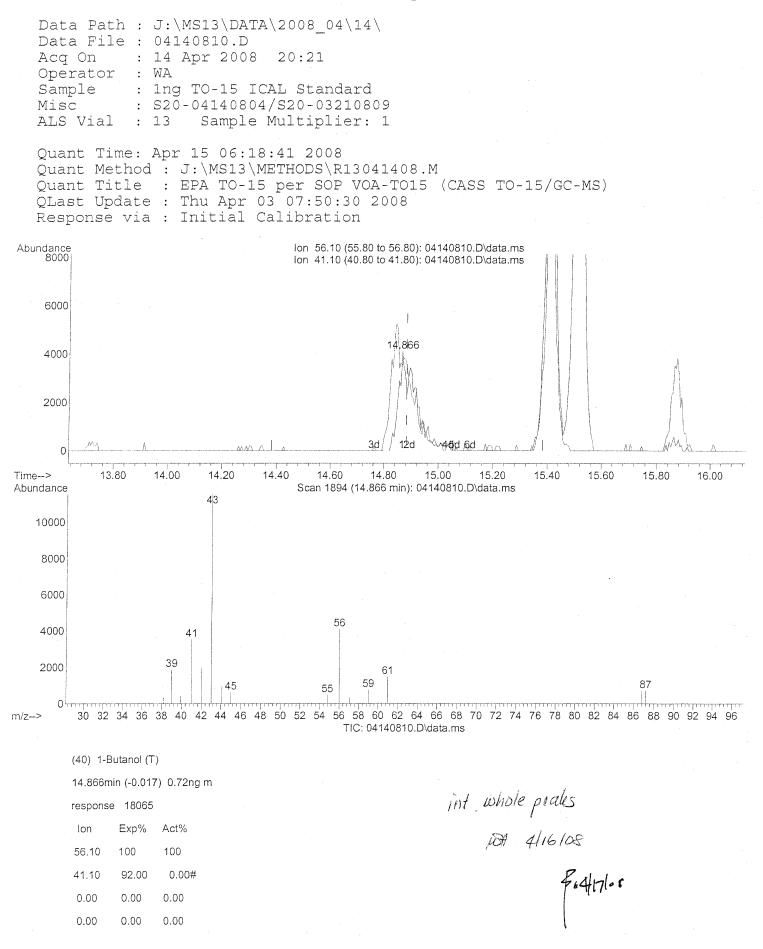
Page: 1



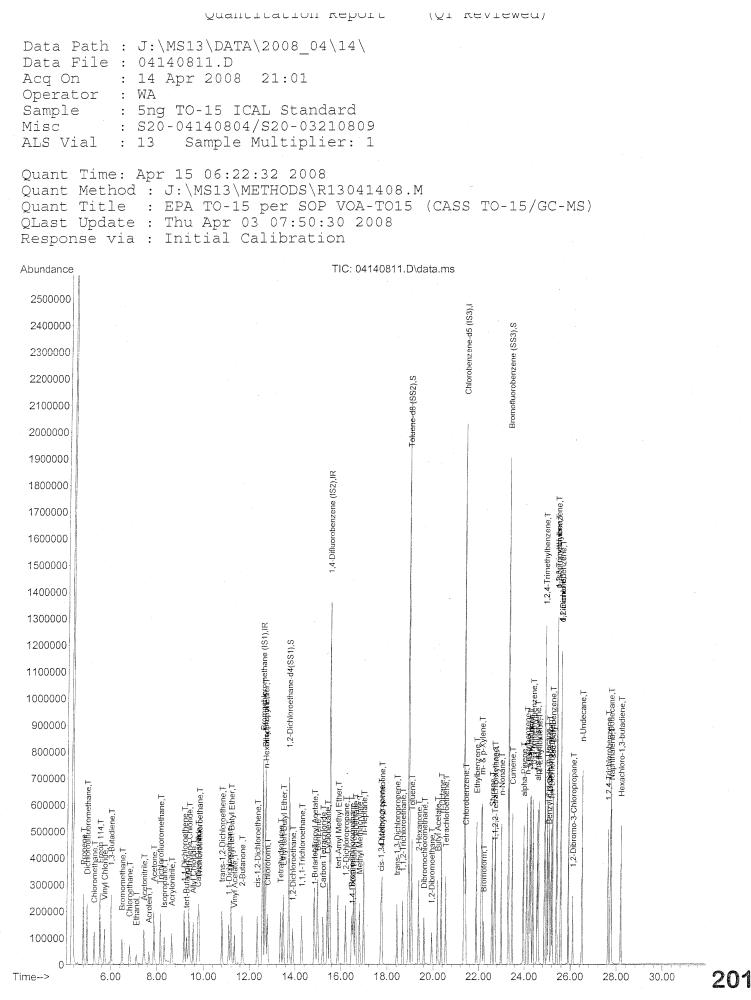


R13041408.M Tue Apr 15 06:20:15 2008





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

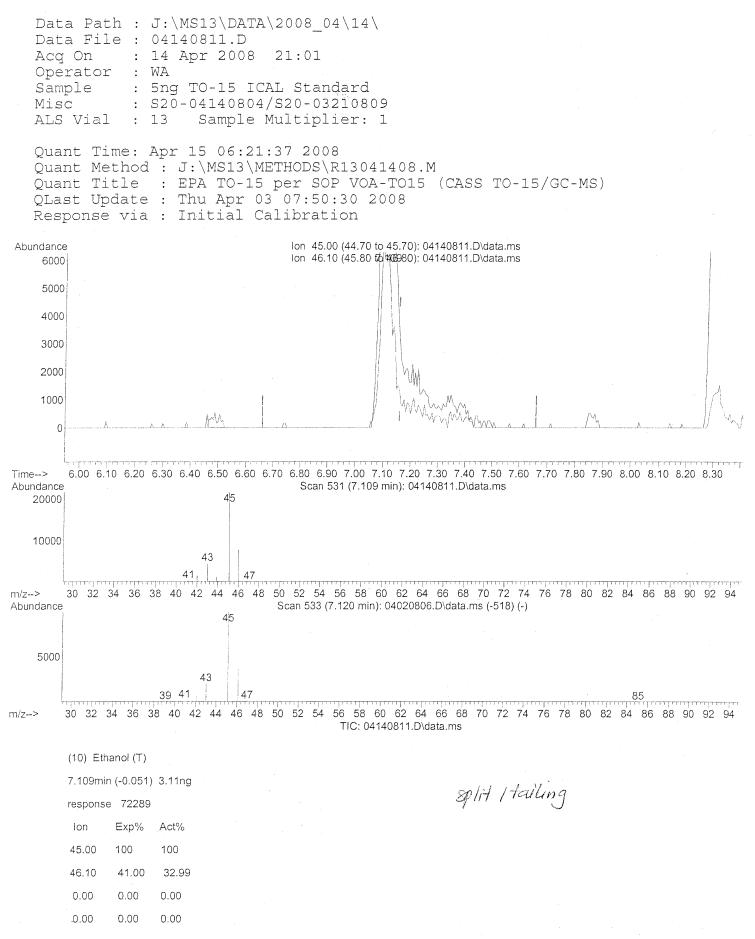
 51) n-Heptane
 Internal Standards R.T. QIon Response Conc Units Dev(Min)

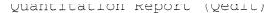
13041408.M Tue Apr 15 06:24:08 2008

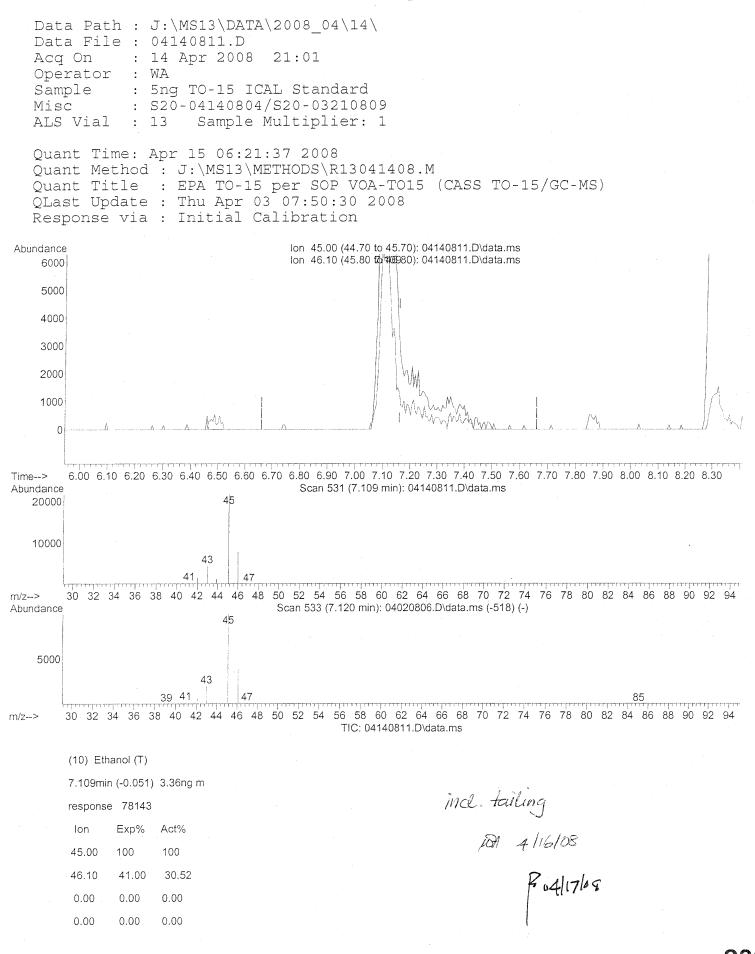
Data Path : J:\MS13\DATA\2008_04 Data File : 04140811.D Acq On : 14 Apr 2008 21:01 Operator : WA Sample : 5ng TO-15 ICAL Standa Misc : S20-04140804/S20-0323 ALS Vial : 13 Sample Multiplie	ard 10809				
Quant Time: Apr 15 06:22:32 2008 Quant Method : J:\MS13\METHODS\R Quant Title : EPA TO-15 per SOP QLast 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



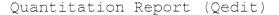


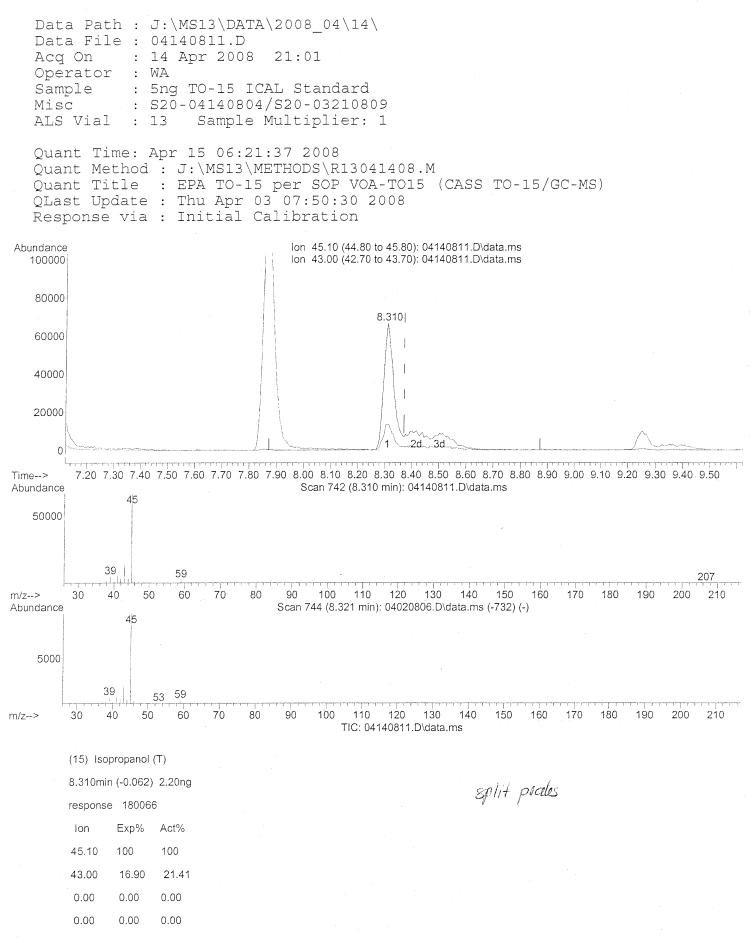




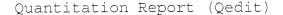
R13041408.M Tue Apr 15 06:22:07 2008

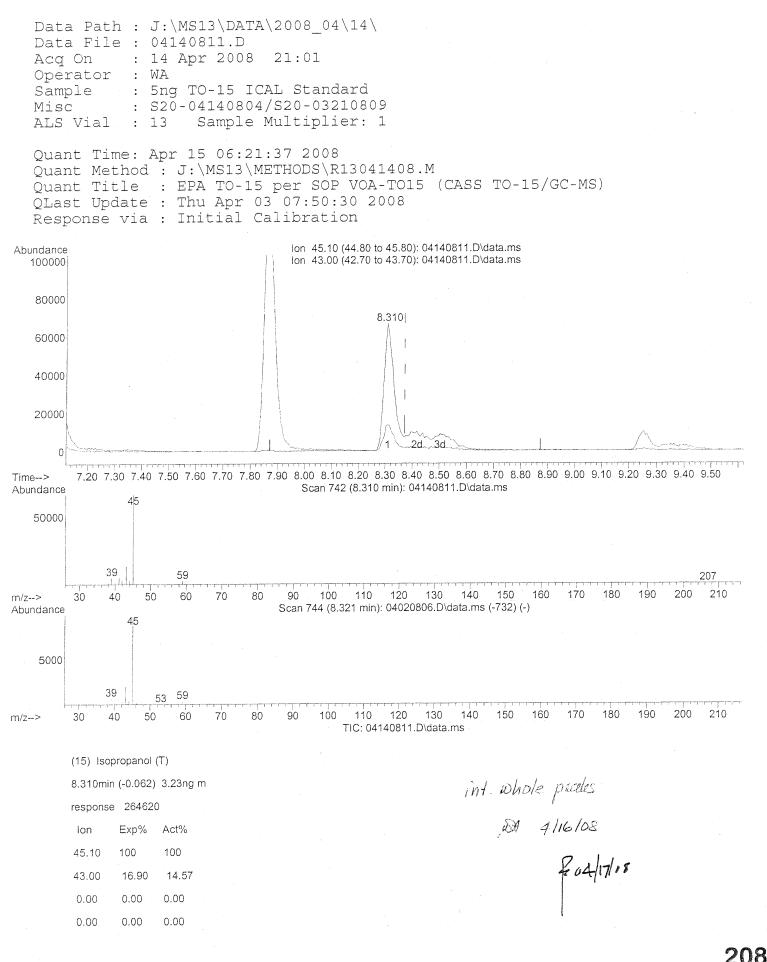
Page: 1

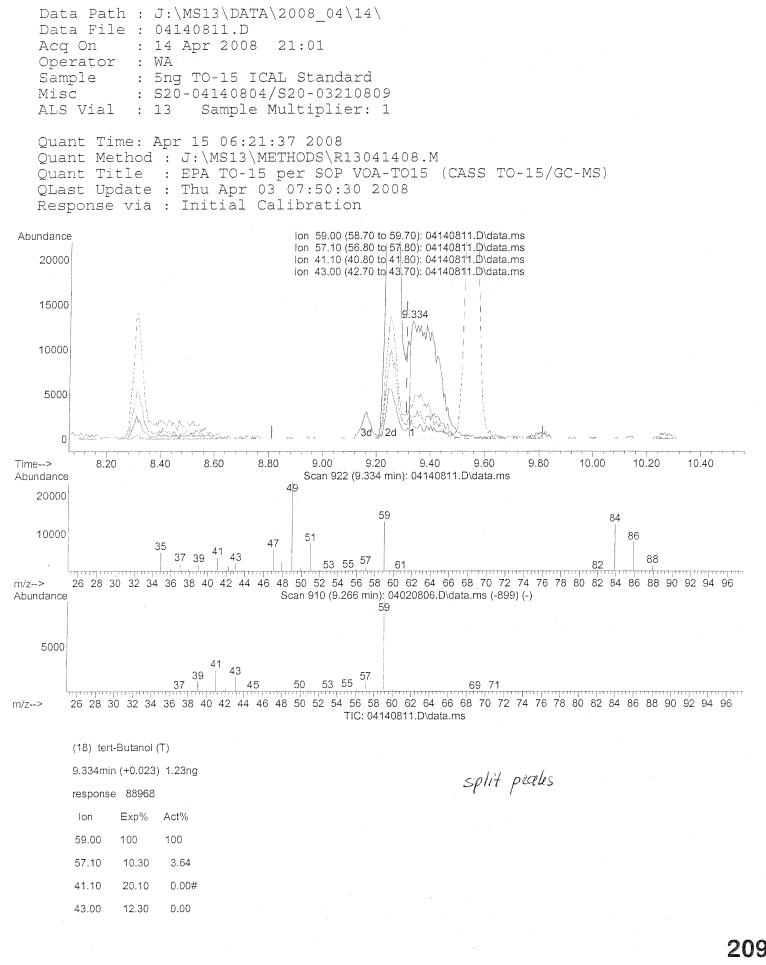


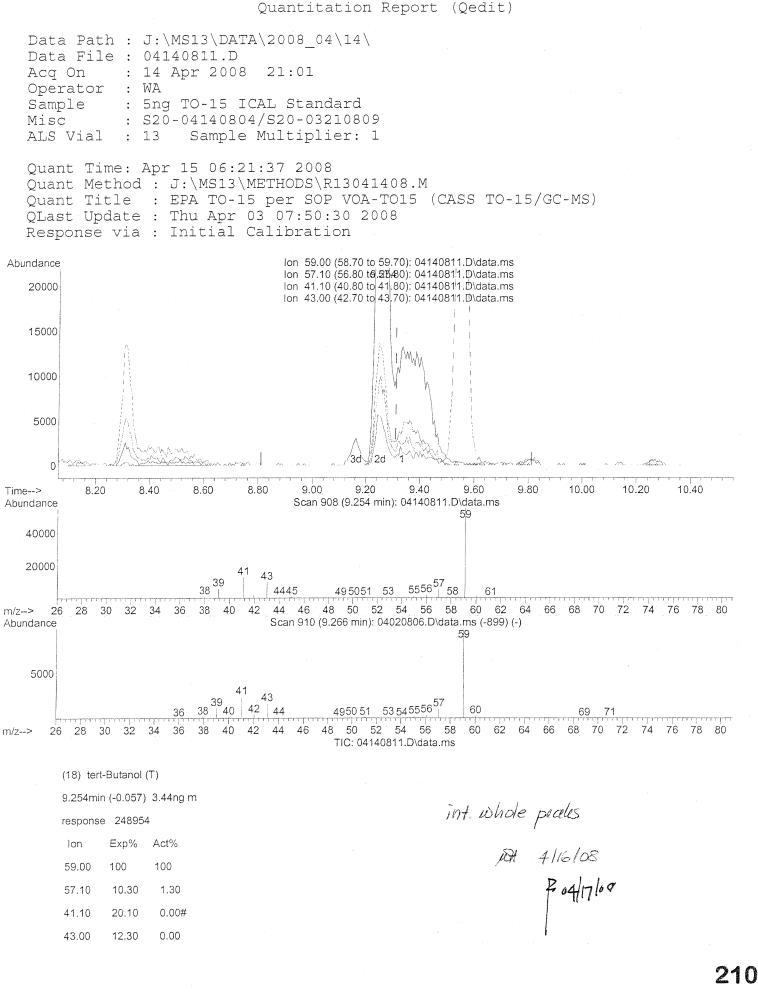


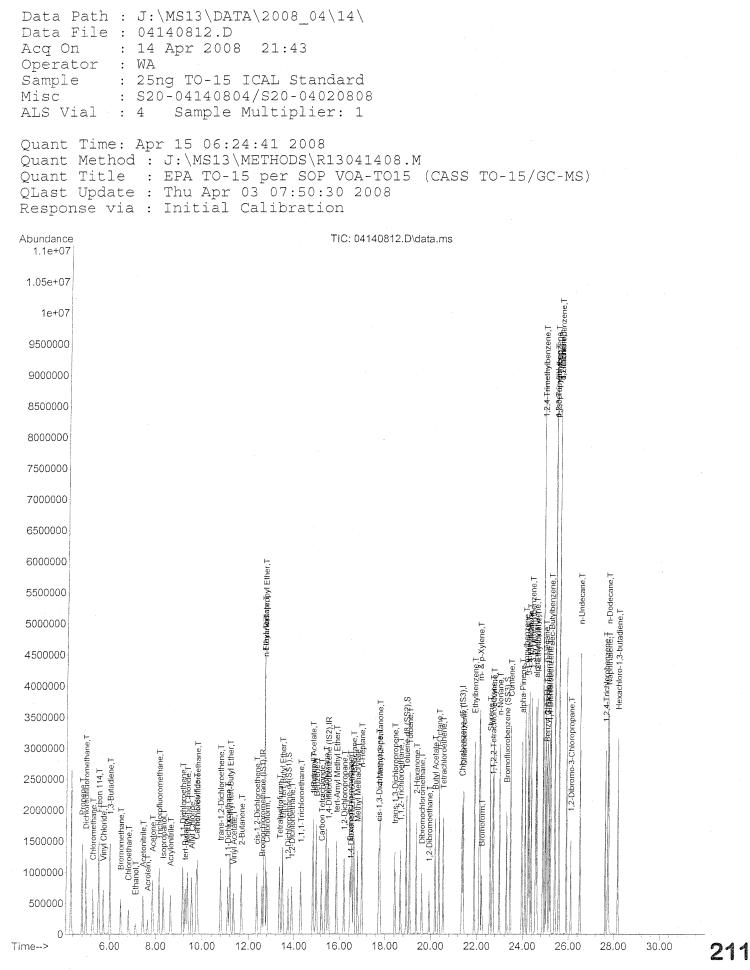
R13041408.M Tue Apr 15 06:22:16 2008











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- &amp; 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

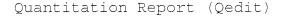
Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140812.D Acg On : 14 Apr 2008 21:43 Operator : WA 
 Sample
 : 25ng TO-15 ICAL Standard

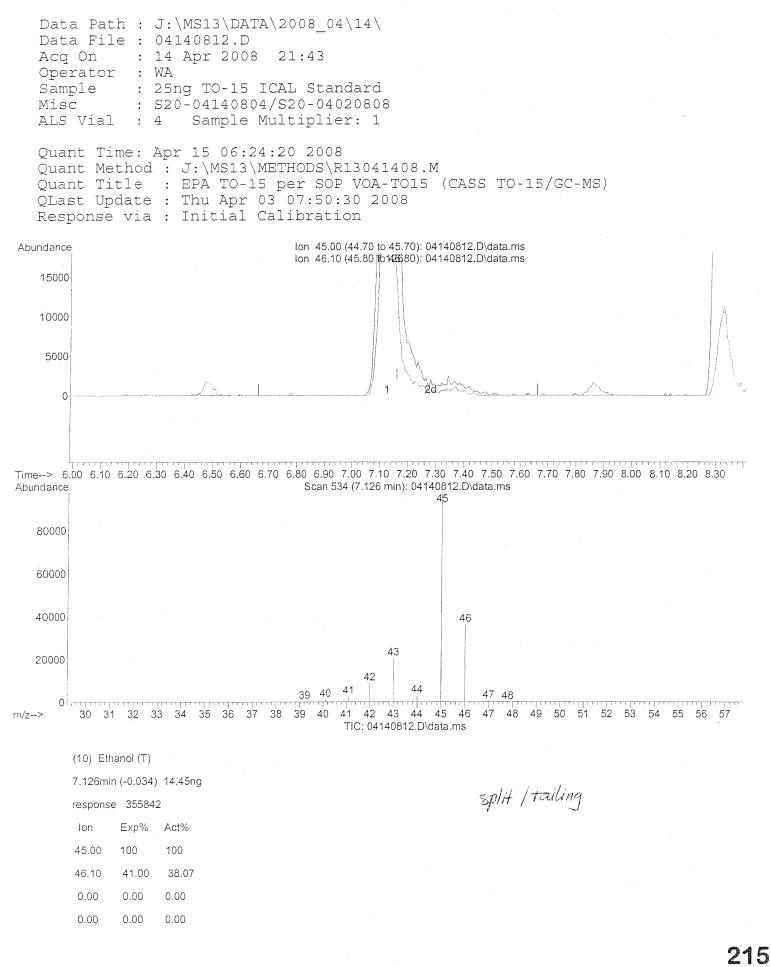
 Misc
 : S20-04140804/S20-04020808
 ALS Vial : 4 Sample Multiplier: 1 Quant Time: Apr 15 06:24:41 2008 Quant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 per SOP VOA-TO15 (CASS TO-15/GC-MS) QLast Update : Thu Apr 03 07:50:30 2008 Response via : Initial Calibration R.T. QIon Response Conc 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

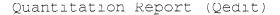
13041408.M Tue Apr 15 06:25:46 2008

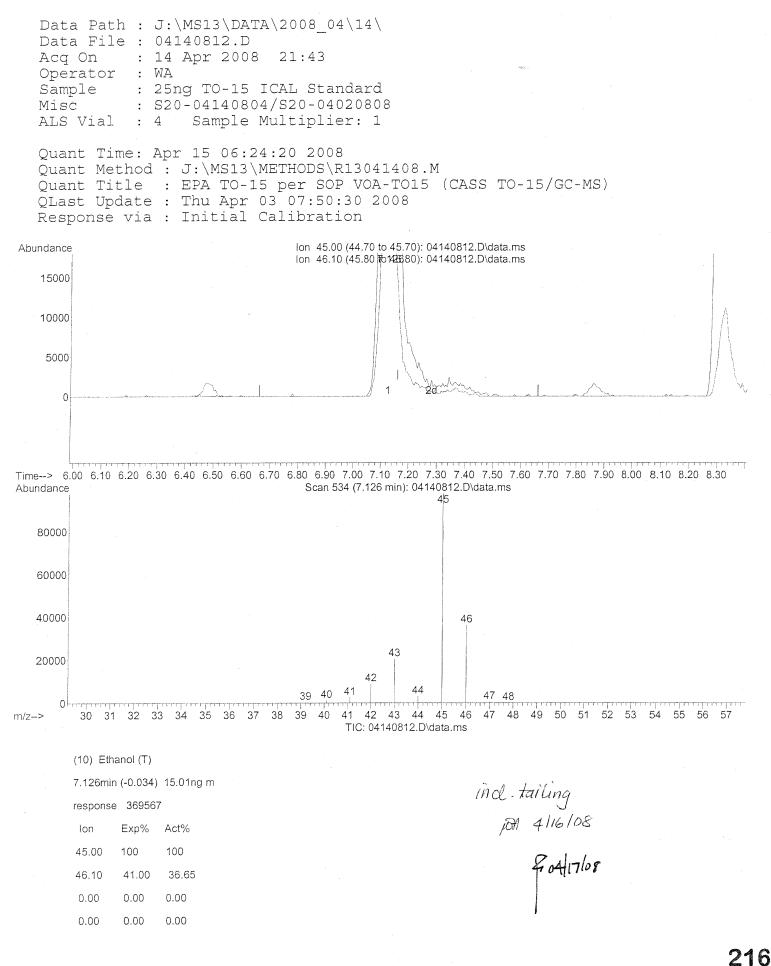
PA 4/16/08

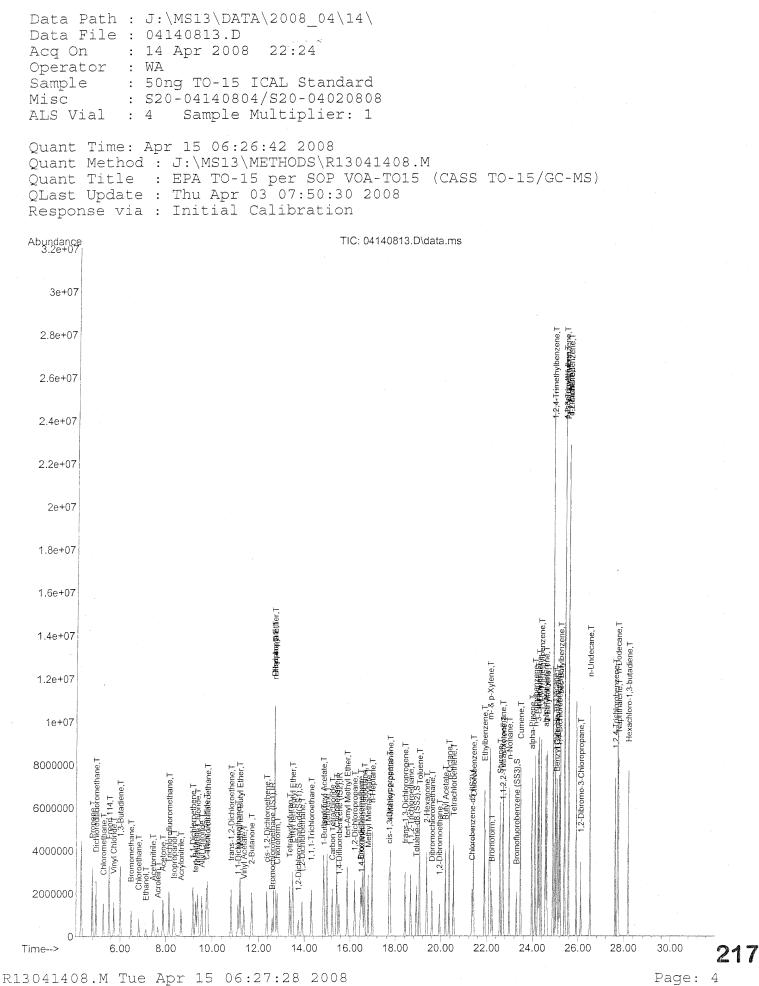




R13041408.M Tue Apr 15 06:24:38 2008







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			
<ul> <li>13) Acetone</li> <li>14) Trichlorofluoromethane</li> <li>15) Isopropanol</li> <li>16) Acrylonitrile</li> <li>17) 1,1-Dichloroethene</li> <li>18) tert-Butanol</li> <li>19) Methylene Chloride</li> </ul>	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

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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

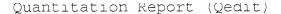
 92) 1,2-Dibromo-3-Chloropr...
 26.11
 157
 833376
 63.938
 ng
 #
 62

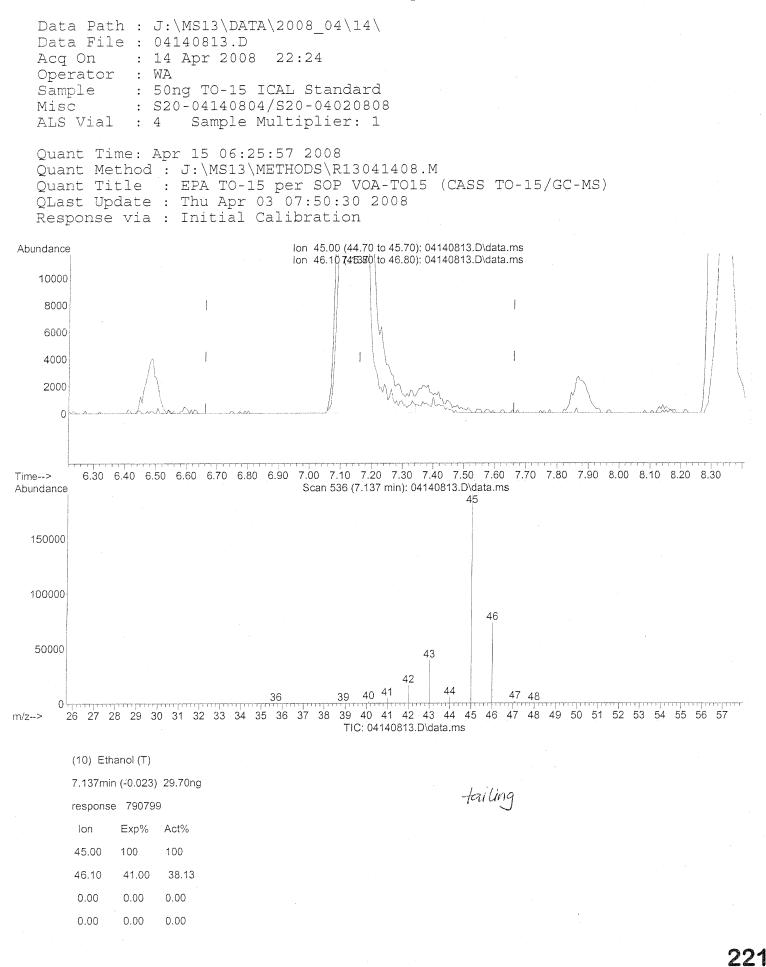
 93) n-Undecane26.5057314646550.655 ng94) 1,2,4-Trichlorobenzene27.63180201784663.304 ng95) Naphthalene27.78128617126064.890 ng96) n-Dodecane27.7457317722552.528 ng97) Hexachloro-1,3-butadiene28.19225131524169.543 ng 87 96 99 86 100 _____

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

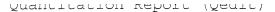
13041408.M Tue Apr 15 06:27:27 2008

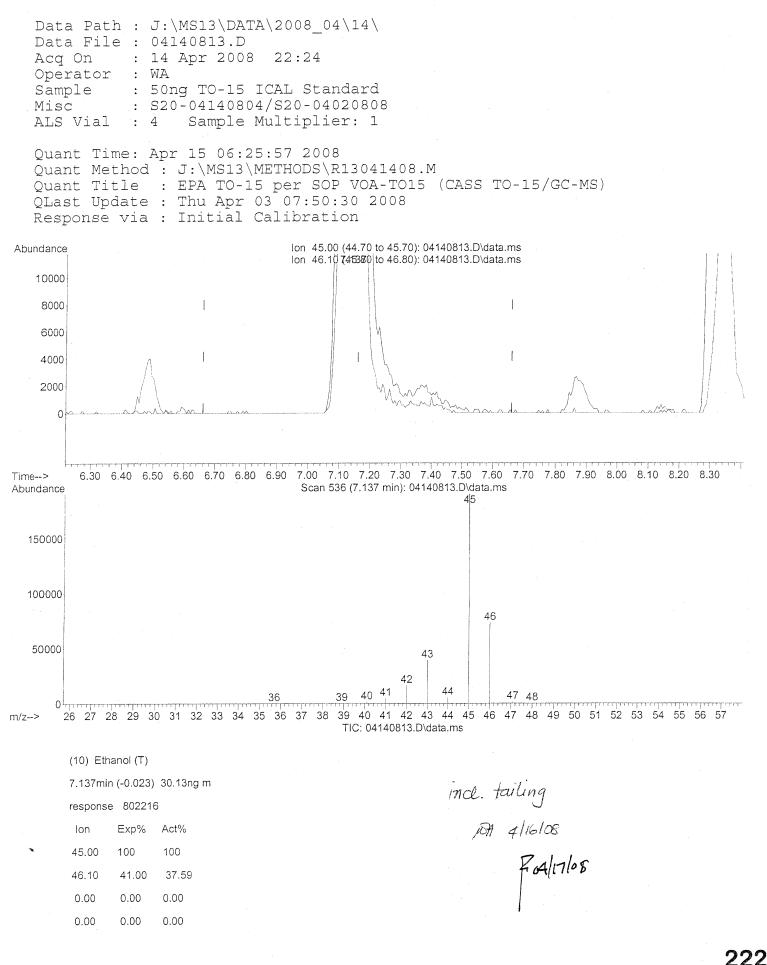
184 4/16/08

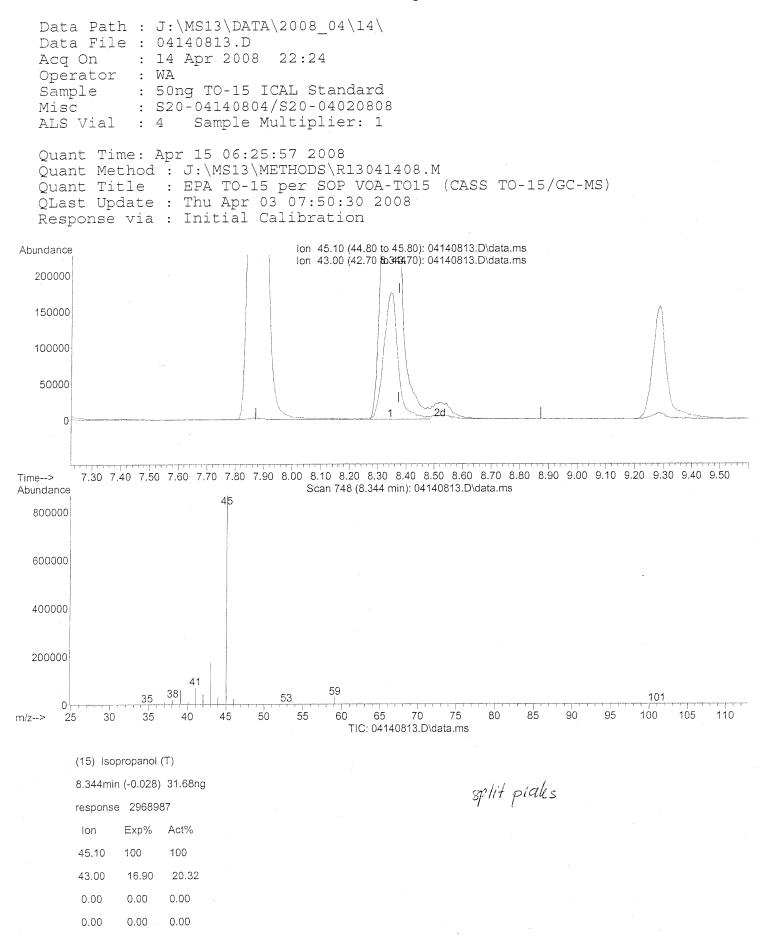


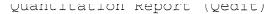


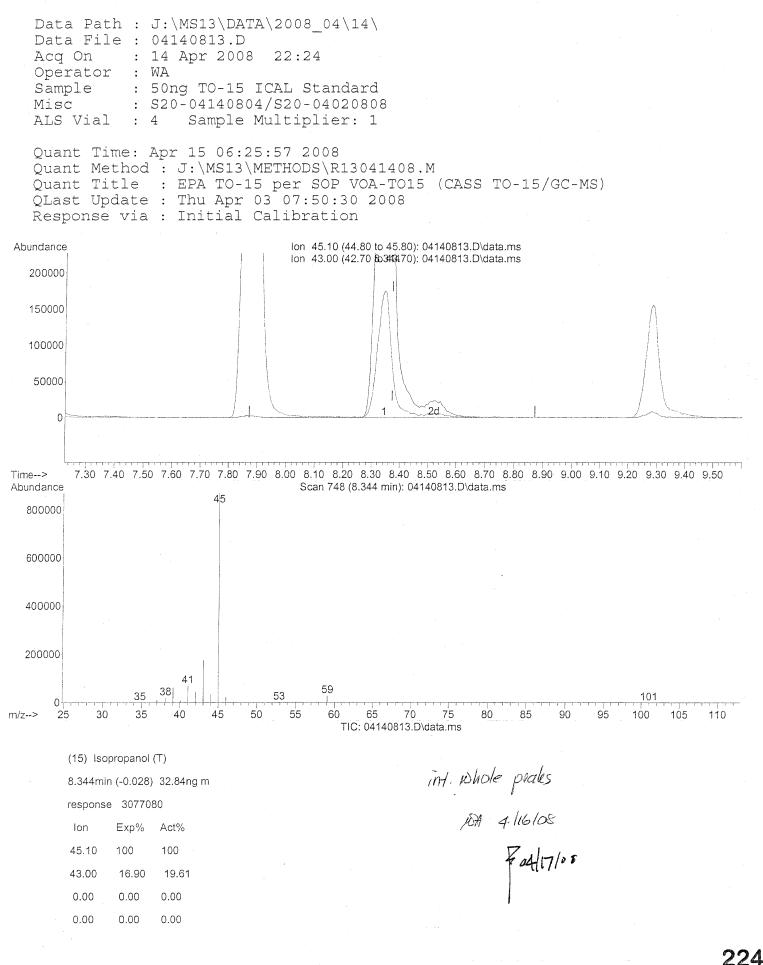
R13041408.M Tue Apr 15 06:26:20 2008

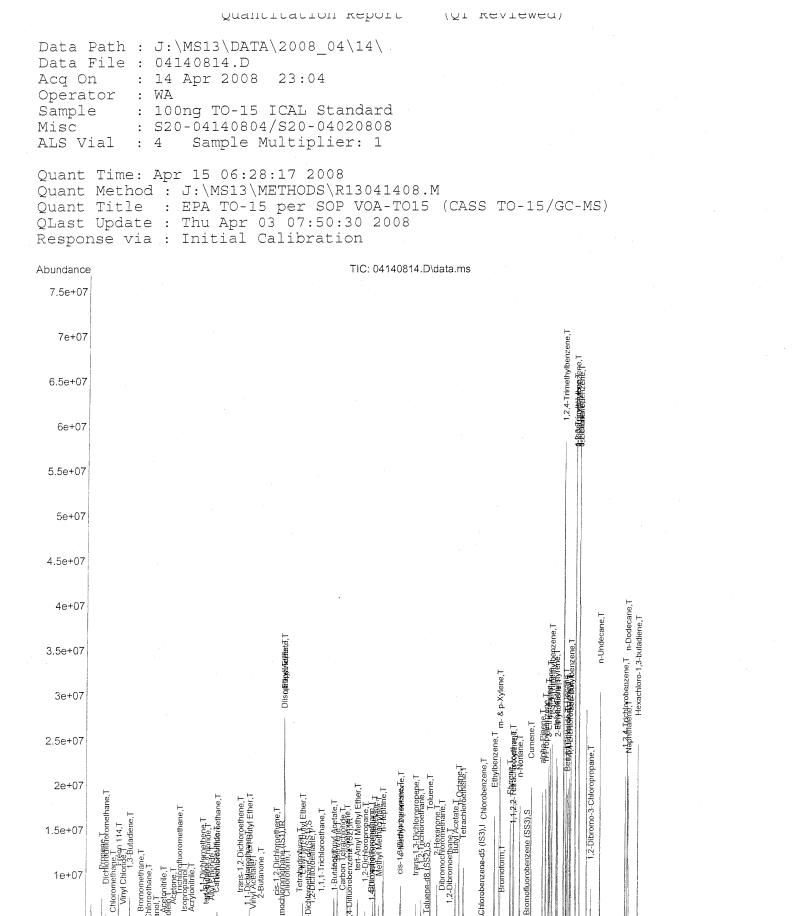


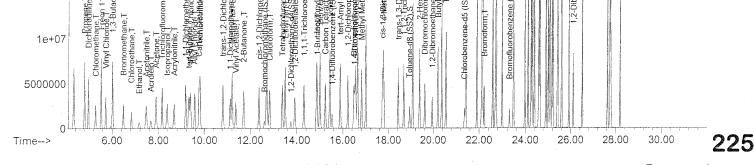












pon

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

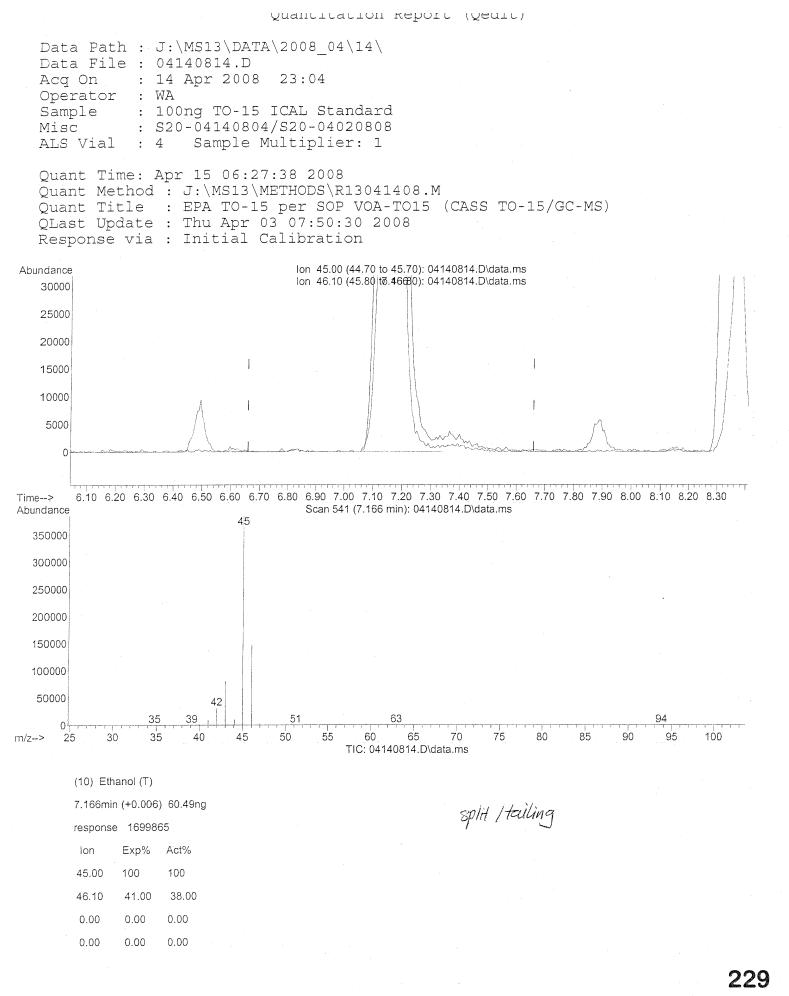
 44: tert-Amyl Methyl Ether
 15.88
 73<700525</td>
 82.635 ng
 91

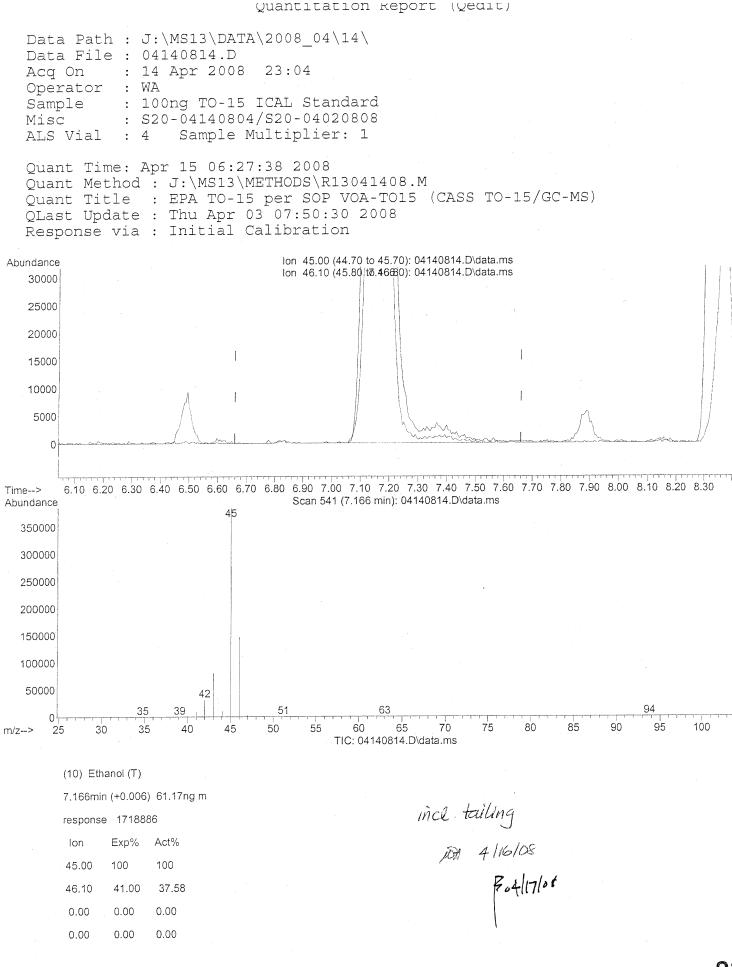
 45: Isooctane
 16.51
 82
 011 ng
 100

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

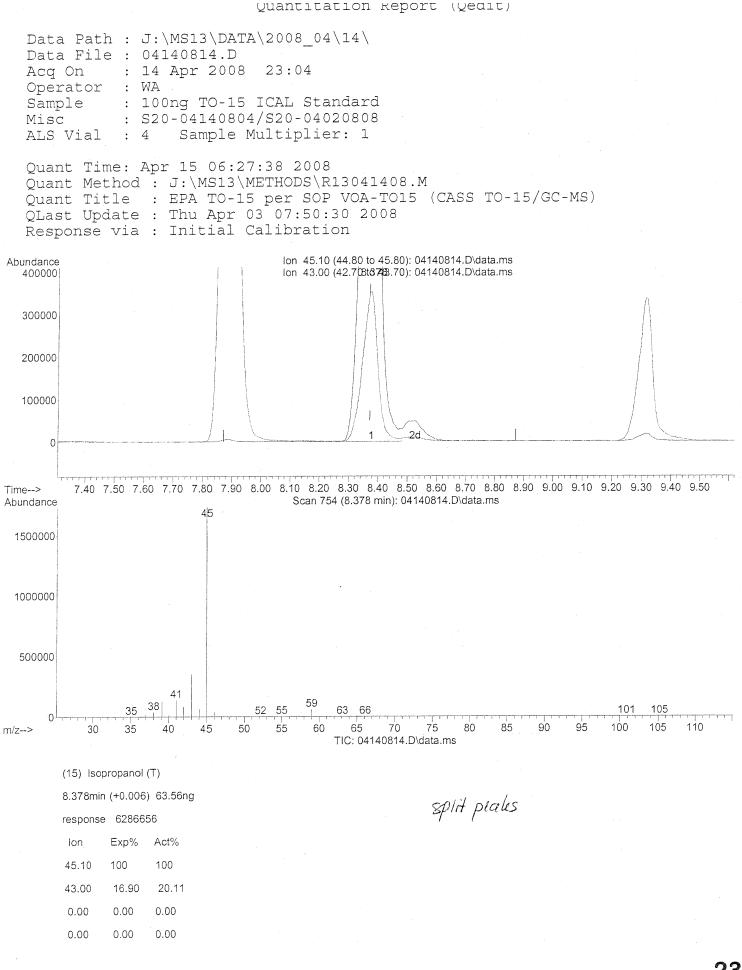
Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140814.D Acq On : 14 Apr 2008 23:04 Operator : WA Sample : 100ng TO-15 ICAL Standard Misc : S20-04140804/S20-04020808 ALS Vial : 4 Sample Multiplier: 1 Quant Time: Apr 15 06:28:17 2008 Ouant Method : J:\MS13\METHODS\R13041408.M Quant Title : EPA TO-15 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 _____

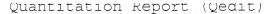
(#) = qualifier out of range (m) = manual integration (+) = signals summed

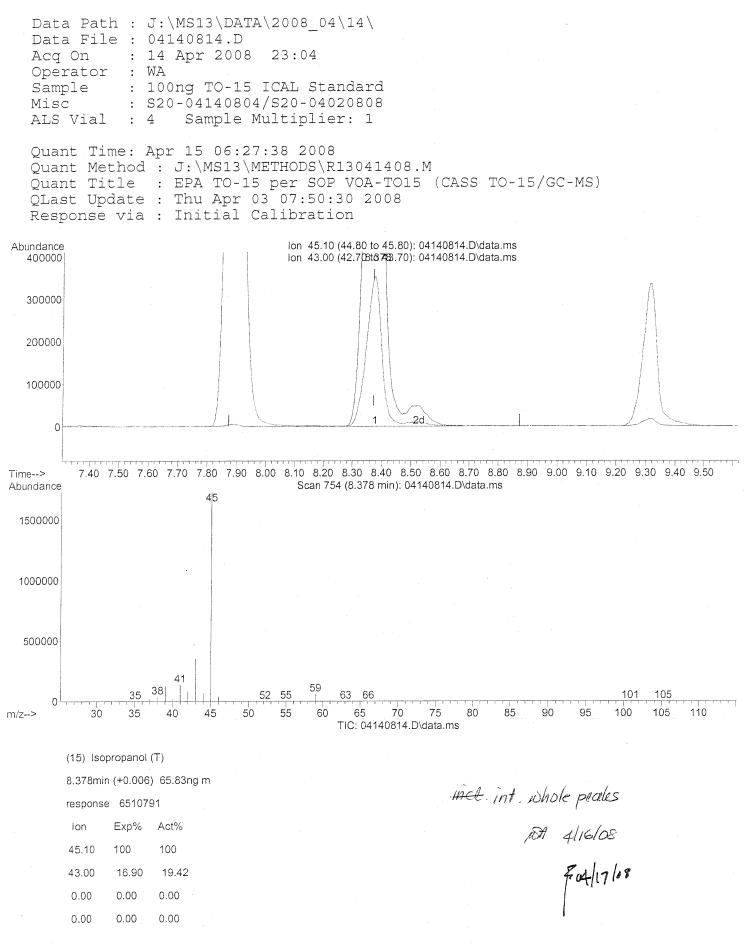


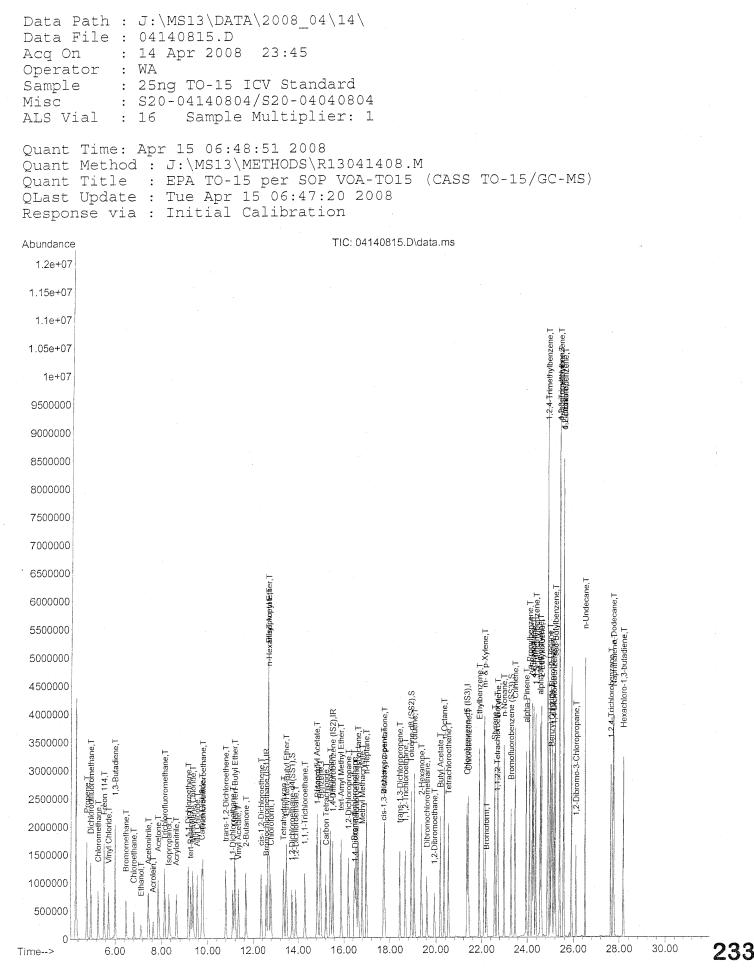


R13041408.M Tue Apr 15 06:28:05 2008









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

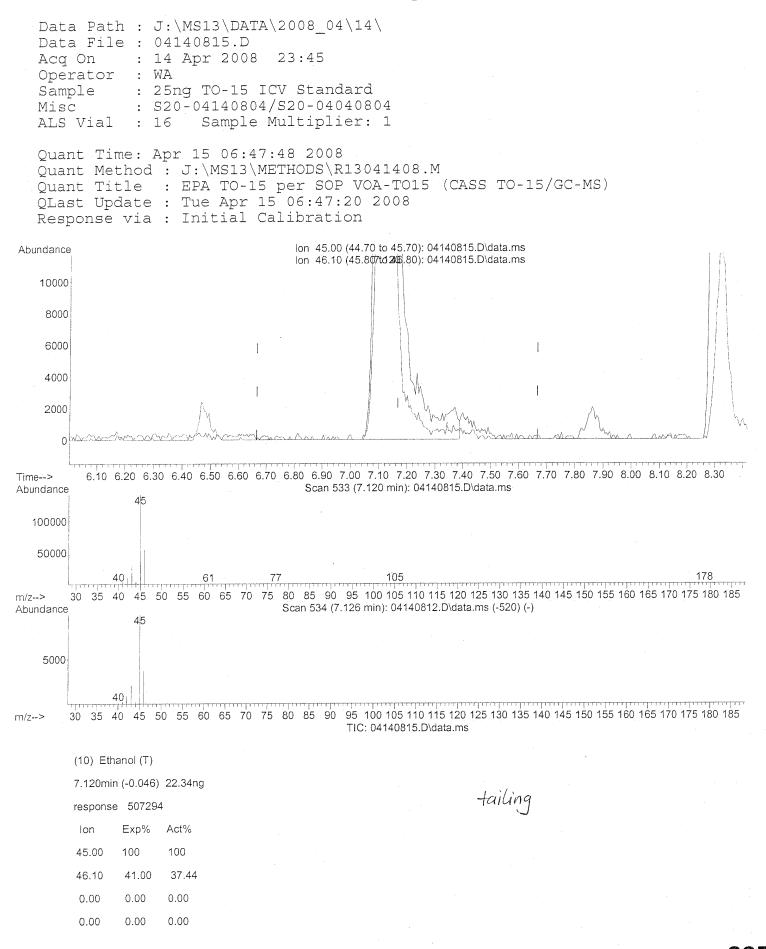
 53) Metyl Methacrylate Internal Standards R.T. QIon Response Conc Units Dev(Min)

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 #

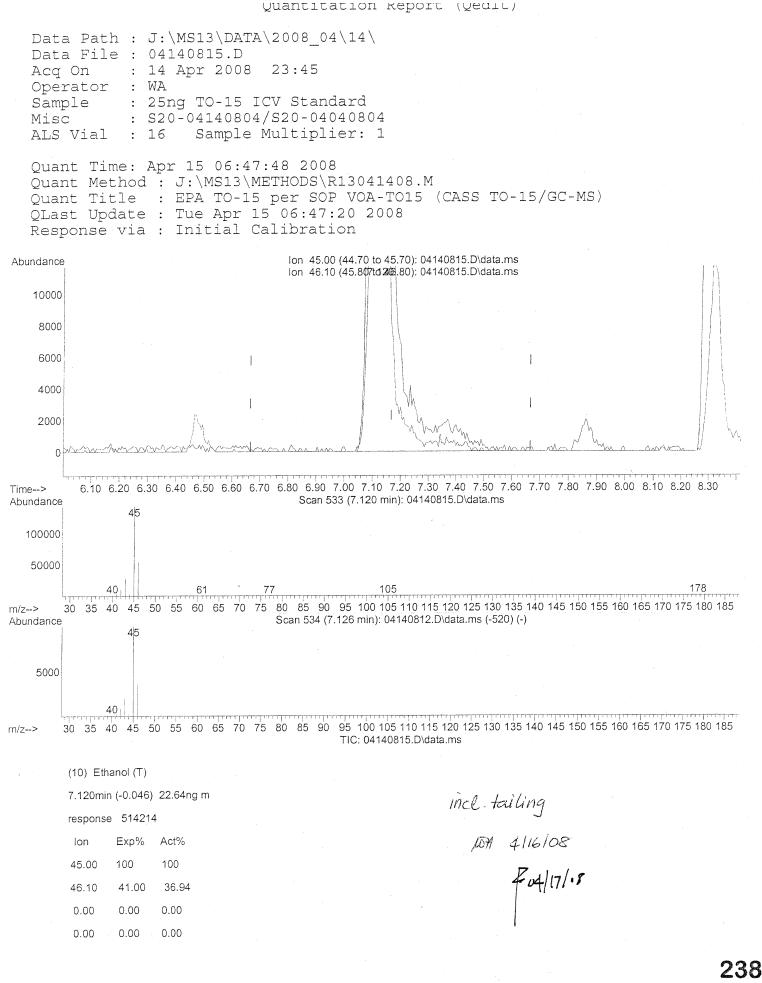
(#) = qualifier out of range (m) = manual integration (+) = signals summed

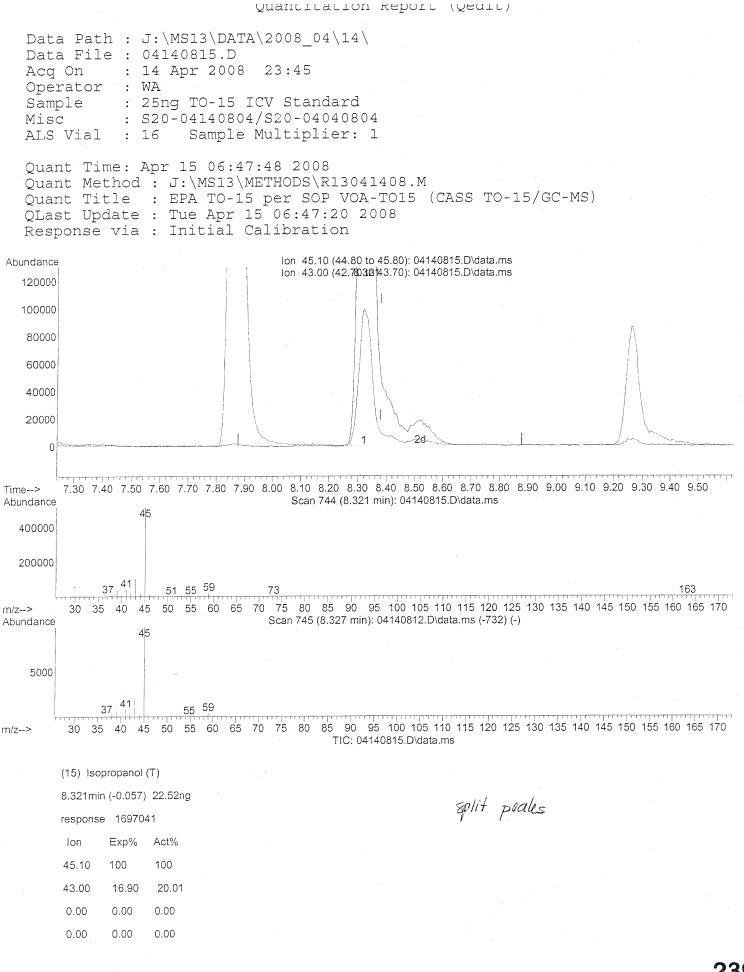
13041408.M Tue Apr 15 06:49:39 2008

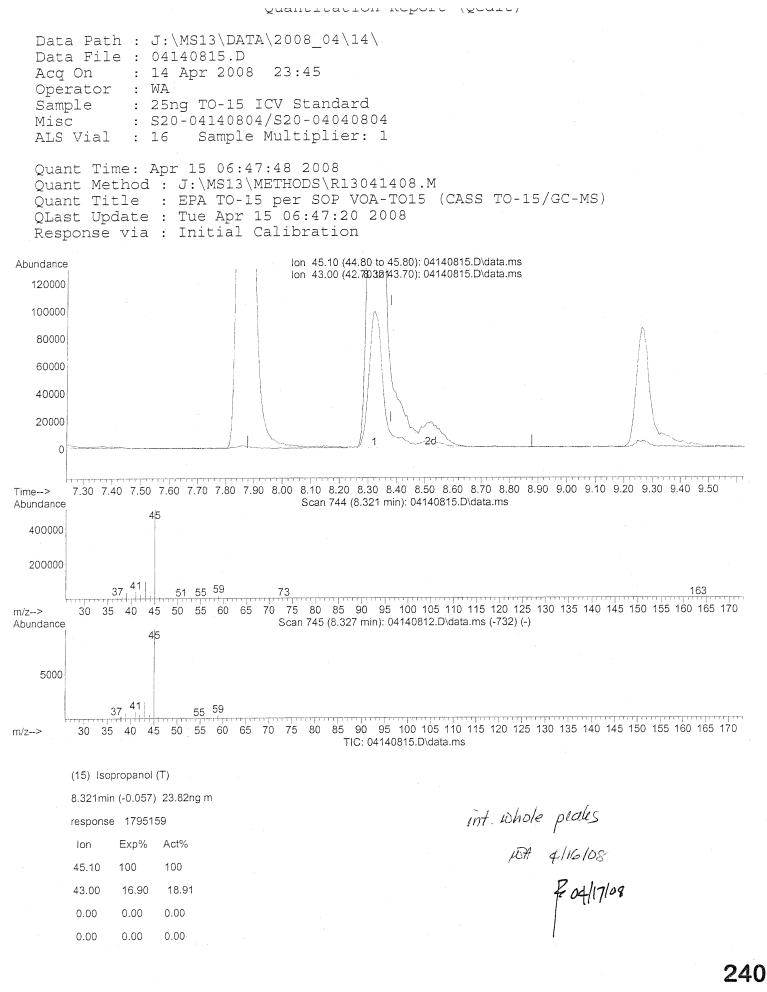
A116108



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	<del>9</del> 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

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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

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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

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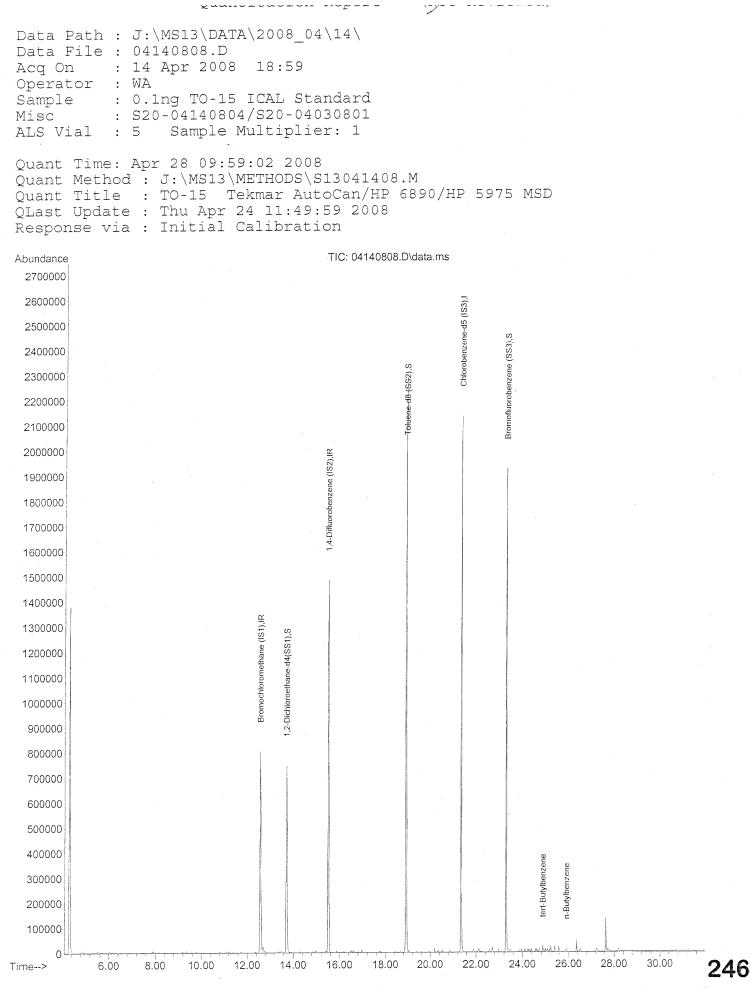
Method Path : J:\MS13\METHODS\ Method File : S13041408.M Title : TO-15 Tekmar AutoCan/HP 6890/HP 5975 MSD Last Update : Mon Apr 28 10:06:00 2008 Response Via : Initial Calibration

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

D# 5/8/08



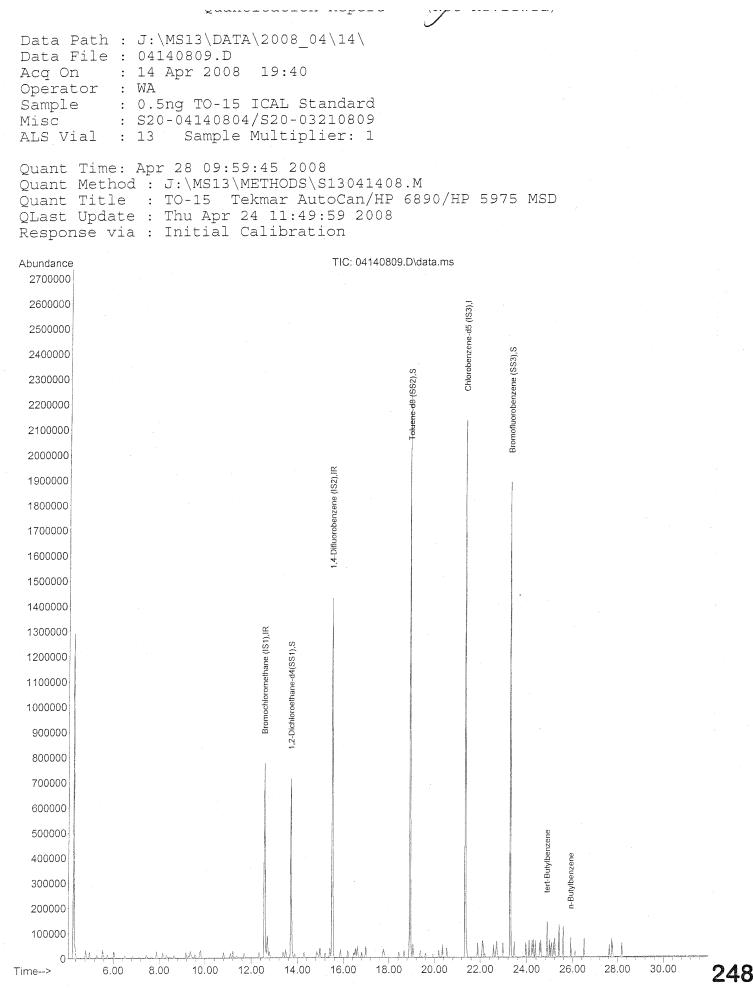
S13041408.M Mon Apr 28 09:59:03 2008

(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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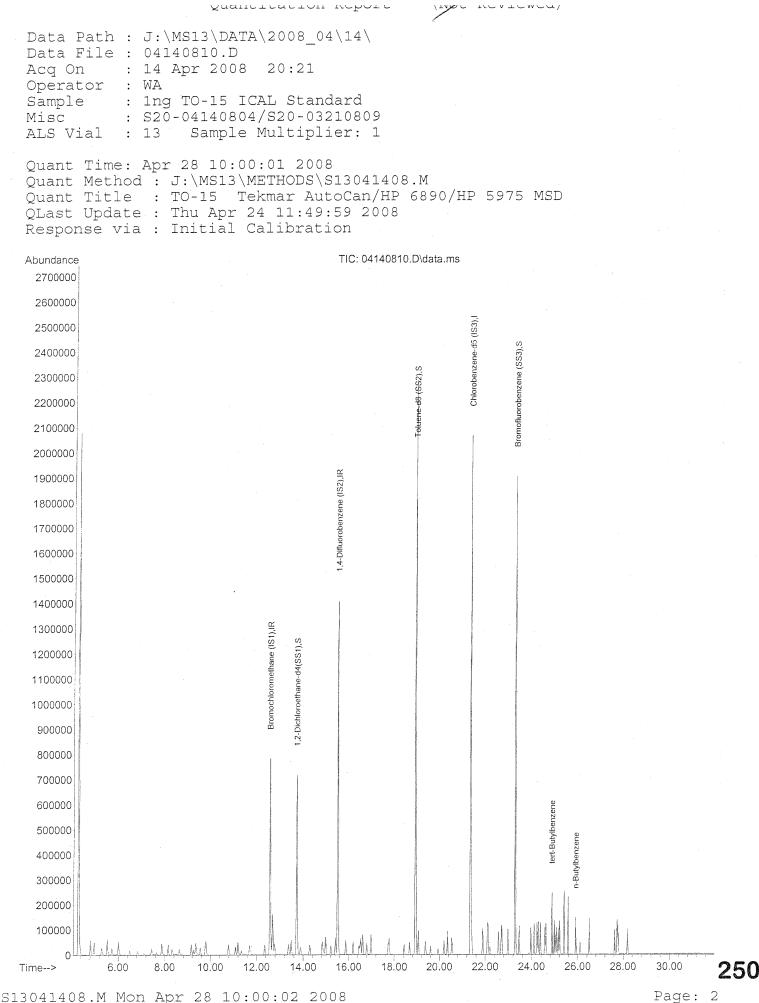


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Quanter cacross report Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140809.D Acq On : 14 Apr 2008 19:40 Operator : WA Sample : 0.5ng TO-15 ICAL Standard Misc : S20-04140804/S20-03210809 ALS Vial : 13 Sample Multiplier: 1 Ouant Time: Apr 28 09:59:45 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%

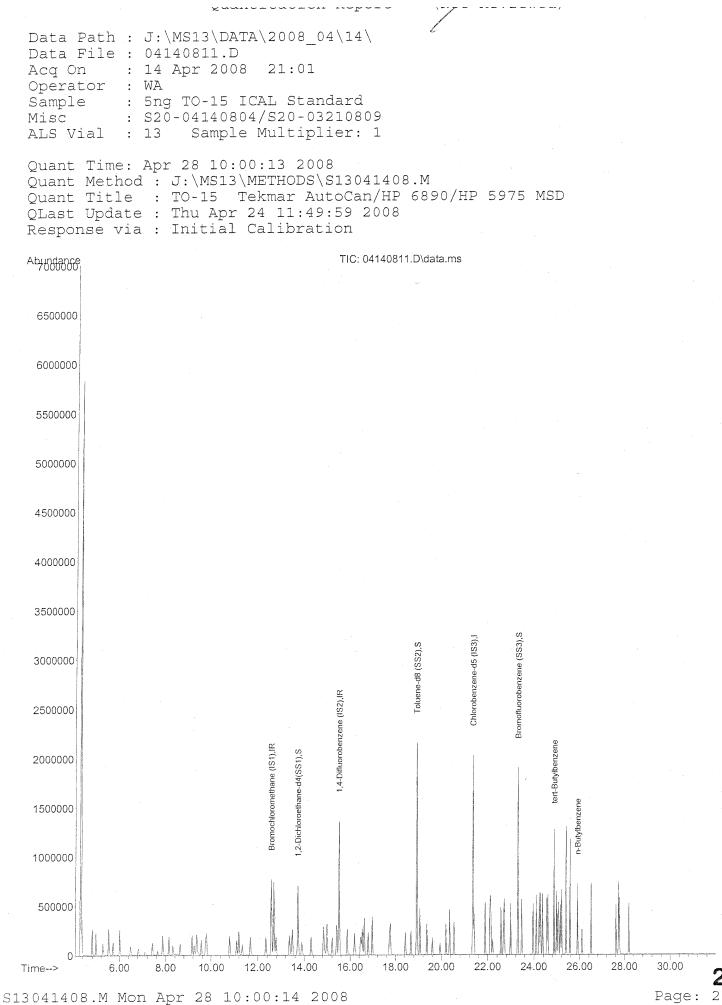
 Spiked Amount 25.000 Recovery = 81.68% Spiked Amount 25.000 Ovalue Target Compounds 7)tert-Butylbenzene24.88119670680.880ng8)n-Butylbenzene25.9191858031.054ng 98 91 

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

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

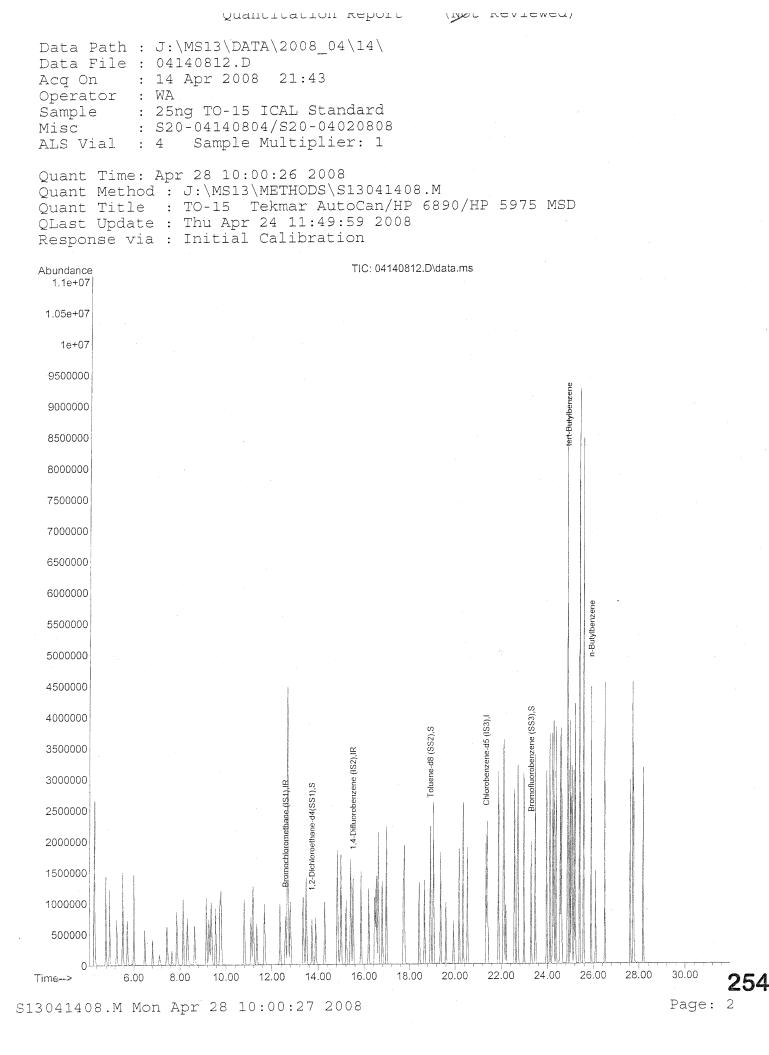
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Xaanoroaoron nobee-Data Path : J:\MS13\DATA\2008_04\14\ Data File : 04140811.D Acq On : 14 Apr 2008 21:01 Operator : WA Sample : 5ng TO-15 ICAL Standard Misc : S20-04140804/S20-0321080 : S20-04140804/S20-03210809 ALS Vial : 13 Sample Multiplier: 1 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

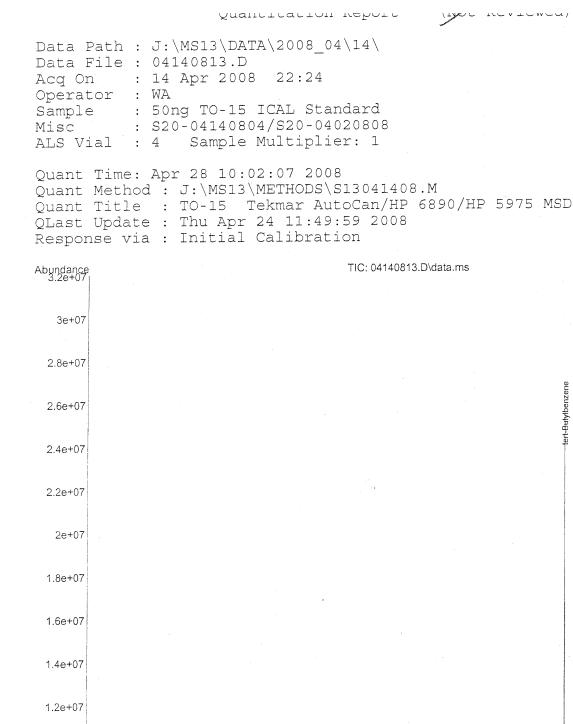
30/26/108

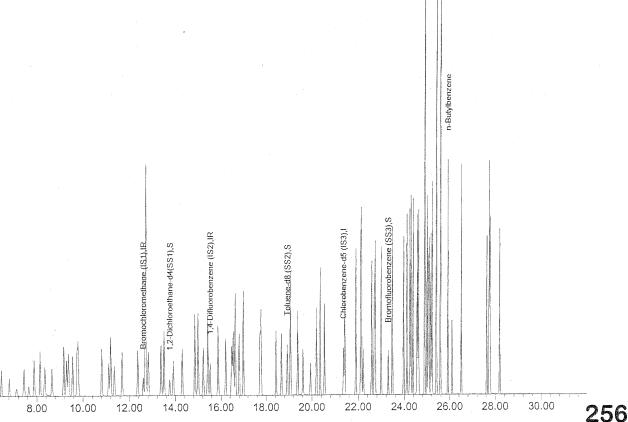


(INCL ICCVICWCC) QUAIILILALIUN NEPULL Data Path : J:\MS13\DATA\2008 04\14\ Data File : 04140812.D Acq On : 14 Apr 2008 21:43 Operator : WA Sample : 25ng TO-15 ICAL Standard Misc : S20-04140804/S20-04020808 ALS Vial : 4 Sample Multiplier: 1 Quant Time: Apr 28 10:00:26 2008 Quant Method : J:\MS13\METHODS\S13041408.M 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

255





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

1e+07

8000000

6000000

4000000

2000000

Time-->

0

6.00

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 

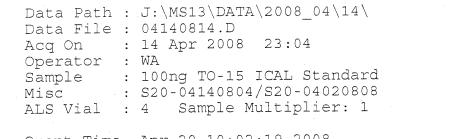
 2) 1,2 Dichlet Cooling and the second sec Recovery = 124.04% Recovery = 81.96% Spiked Amount 25.000 Qvalue Target Compounds 7)tert-Butylbenzene24.89119588564069.362ng8)n-Butylbenzene25.9191565198662.329ng 99 95 ·

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

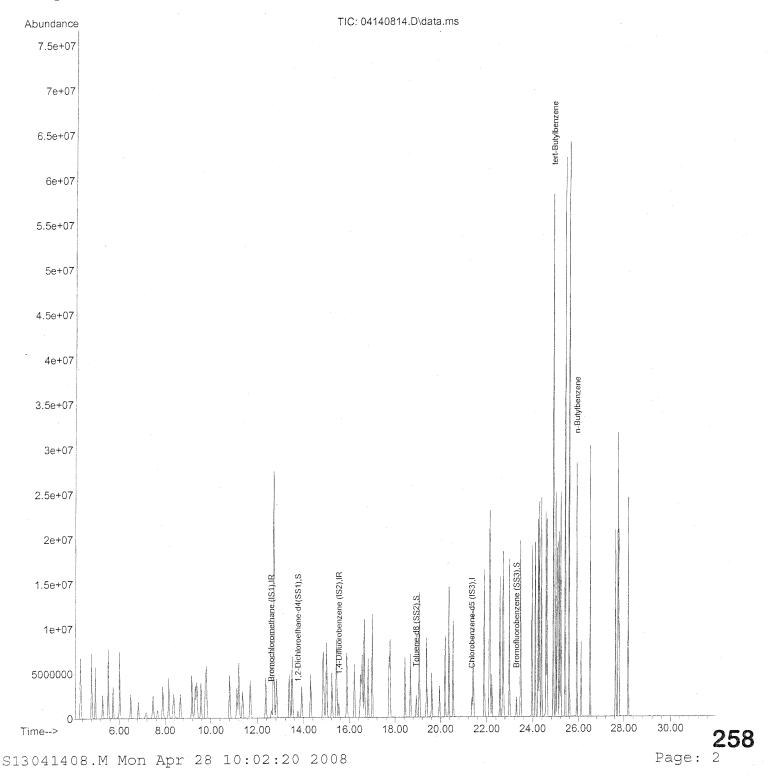
313041408.M Mon Apr 28 10:02:08 2008

\$# 4/28/08

The Treatest



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

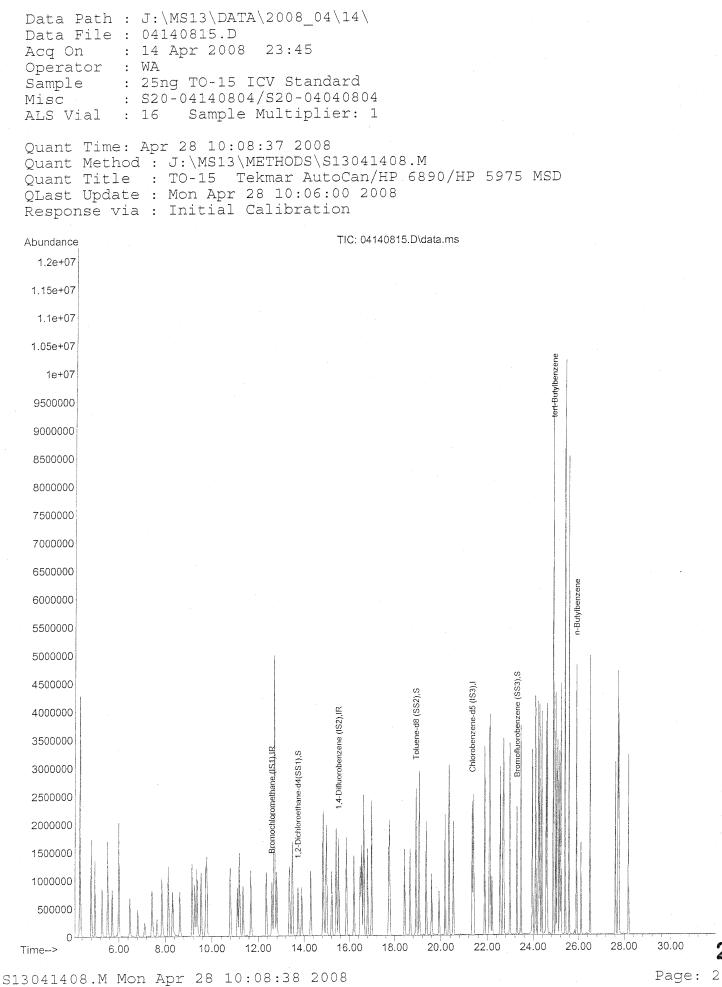


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Data Path : J:\MS13\DATA\2008_04 Data File : 04140814.D Acq On : 14 Apr 2008 23:04 Operator : WA Sample : 100ng TO-15 ICAL Sta Misc : S20-04140804/S20-040 ALS Vial : 4 Sample Multiplie	ndard 20808			
Quant Time: Apr 28 10:02:19 2008 Quant Method : 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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260

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

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264

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

F05/05/05/08

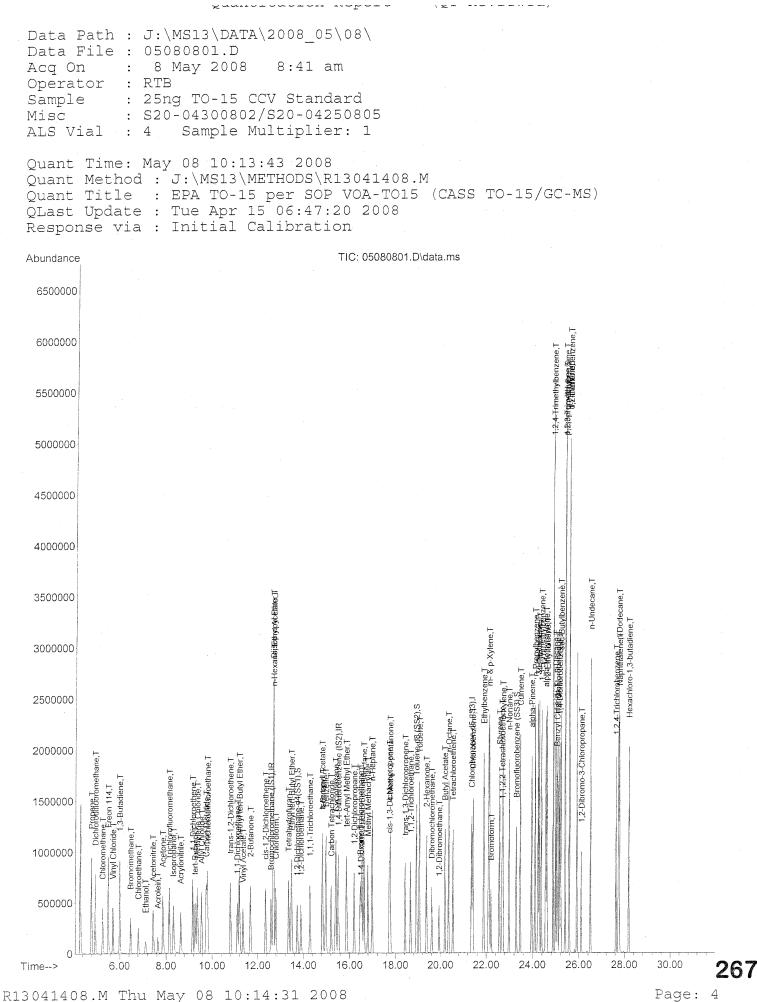
Page: 2

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¢

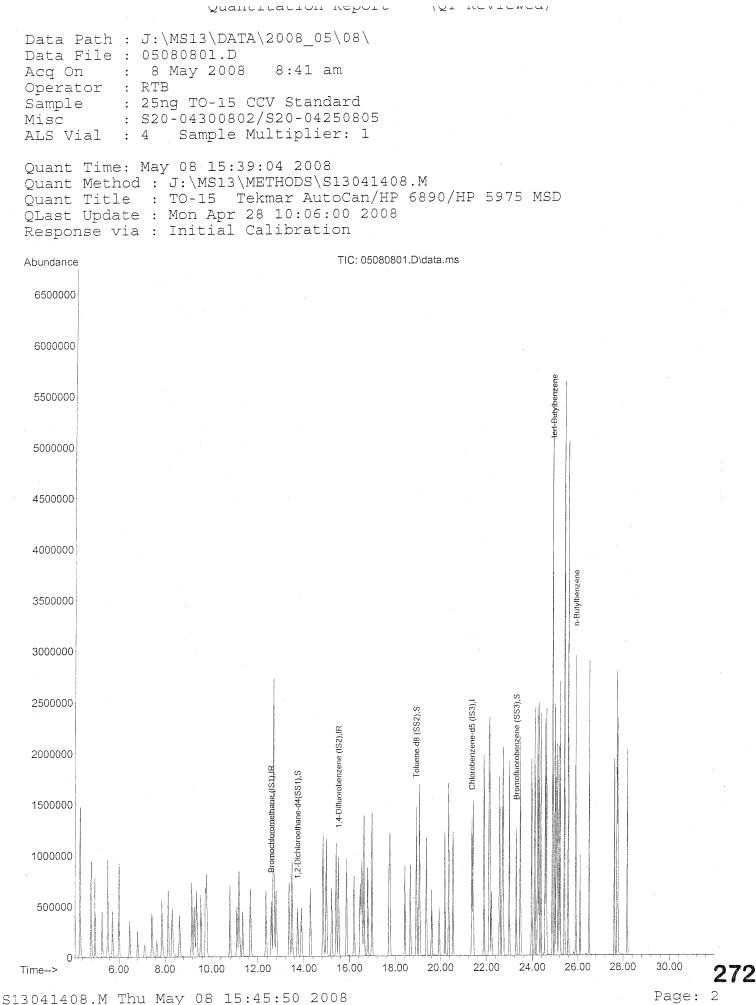
waraaco oonornanig oarraractor neper-Data Path : J:\MS13\DATA\2008 05\08\ Data File : 05080801.D Acq On : 8 May 2008 8:41 am Operator : RTB Sample : 25ng TO-15 CCV Standard Misc : S20-04300802/S20-04250805 ALS Vial : 4 Sample Multiplier: 1 Ouant Time: May 08 15:39:04 2008 Quant Method : J:\MS13\METHODS\S13041408.M 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



1×+

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

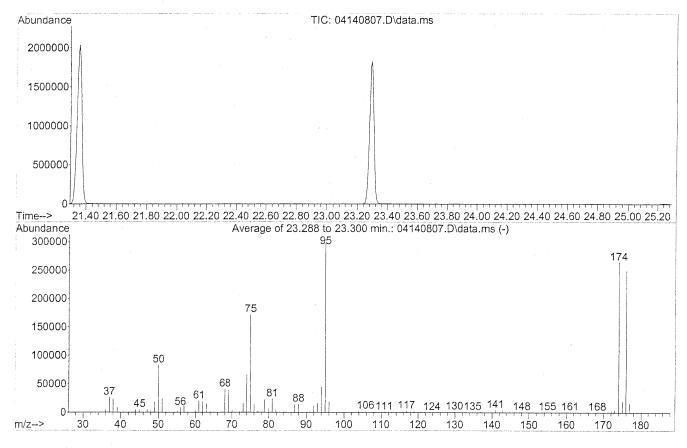
273

## 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

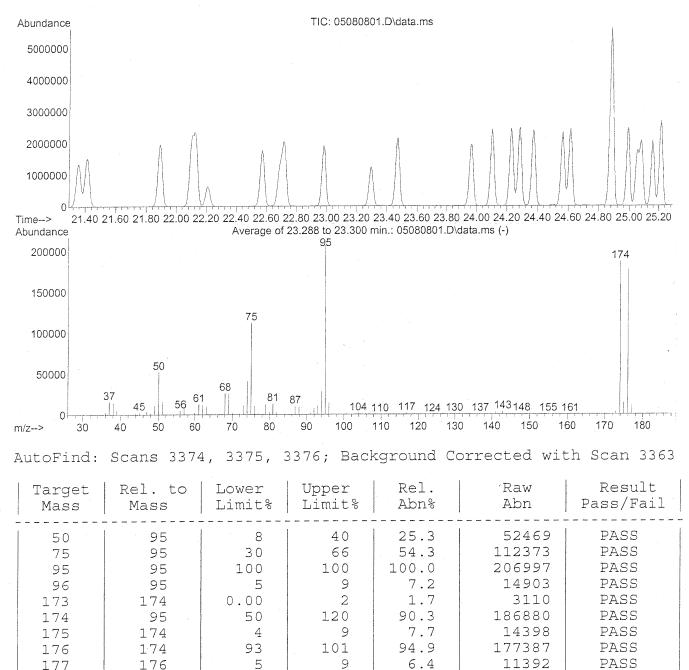
DH 4/17/08

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
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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 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 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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				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)	<ul> <li>G.11808020</li> <li>G.21808020</li> <li>G.21808020</li> <li>G.21808020</li> <li>G.21808020</li> </ul>	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 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.50808080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.50808080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 0.508080 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       v       v       v       v       v       v       v       v       v       v       v       v       v       v       v       v       v       v       v       v       v       v       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
			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       t       t       t	Test ENGR SG83B-05-1 (-3.6, 3.5) ENGR SG83B-05-7 (-4.2, 3.5) ENGR SG83B-05-7 (-4.2, 3.5) ENGR SG83B-05-7 (-4.2, 3.5) ENGR SG83B-05-7 (-4.2, 3.5) Test Test	P0801342-001 DIL (1mL) P0801342-002 DIL (10mL) P0801342-003 DUP (10mL) P0801342-003 DUP (10mL) P0801342-003 DUP (10mL) P0801342-003 DUP (10mL) P0801342-003 DUP (10mL) P0801342-003 DUP (10mL) P0801342-001 DIL (1mL) P0801342-001 DIL (1mL)	G 21808050 G 21808050 G 21808050 G 21808050 G 21808050 G 20808050 G 20808050 G 20808050 G 30808050 G 30808050 G 30808050	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 0 12 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0
			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       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       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	<ul> <li>C2:2 80/80/50</li> <li>C2:2 80/</li></ul>	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       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       t       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 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 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 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0
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			92 92 92 92 92 92 92 92 92 92 92 92 92 9	t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           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.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.20808020 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.2080800 G.20808000 G.20808000 G.20808000 G.208080000 G.208080000 G.20808000000000000000000000000000000000	<ul> <li>L+:6 80/80/20</li> <li>L+:6 80/</li></ul>	117 ( 117 ( 11
			83 83 83 93 93 93 93 93 93 93 93 93 93 93 93 93	t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           t           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 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 115 0 11
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