

## TECHNICAL MEMORANDUM

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**Date:** November 13, 2008

**Subject:** Technical Memorandum – Screening-Level Indoor Air Health Risk Assessment for the 2008 Tronox Parcels A/B Soil Gas Investigation, BMI Industrial Complex, Clark County, Nevada

### Introduction

The objective of this Technical Memorandum is to present the results of a screening-level indoor air health risk assessment for the Phase 2 soil gas investigation Basic Environmental Company (BEC) and Tronox performed for the Tronox Parcels “A” and “B” (portions of APN Nos. 178-01-401-001, 178-12-101-002, and 178-12-201-006 [Note: Parcel 178-12-601-005, formerly part of Tronox Parcel B, has been sold and is excluded from this analysis]). Parcels A and B will collectively be referred to as the property for the purposes of this Technical Memorandum. The property is located north of Warm Springs Road, 1/4 mile west of the intersection with Boulder Highway, in Henderson, Nevada. Figure 1 shows details of Parcels A and B and the soil gas sampling locations. The Technical Memorandum only presents the methods and results of the screening-level indoor air health risk assessment, and does not present investigation, data summary, data usability, or data adequacy information. This information is provided in the Nevada Division of Environmental Protection (NDEP) approved *Technical Memorandum – Data Review for 2007 Tronox Parcels A/B Investigation* dated February 11, 2008, and the Data Validation Summary Report for the soil gas survey.

### Conceptual Site Model

The conceptual site model (CSM) is used to describe relationships between chemicals and potentially exposed human receptor populations, thereby delineating the relationships between the suspected sources of chemicals identified at the property, the mechanisms by which the chemicals might be released and transported in the environment, and the means by which the

receptors could come in contact with the chemicals. The CSM provides a basis for defining data quality objectives and developing exposure scenarios.

The CSM considers current and potential future land-use conditions. Currently, the property is undeveloped. Current receptors that may use the property include on-site trespassers. Therefore, current exposures to native soils at the property are likely to be minimal. In addition, exposures to future on-site workers will be much greater than current exposures. For example, future receptors include indoor commercial workers who are assumed to be exposed to soil gas emanating from the subsurface for 250 days per year for 25 years which is much greater than any current exposures.

USEPA (1989) guidance states that potential future land use should be considered in addition to current land use when evaluating the potential for human exposure at a site. Therefore, the CSM also considers other future land-uses. For example, the CSM includes the planned use of the property for redevelopment into commercial use.

Given the planned development of the property, potential human receptors include on-site construction workers, on-site indoor commercial workers, on-site outdoor maintenance workers, and on-site visitors. Although several potential human receptors may occur on the property in the future, the screening-level health risk assessment focuses on indoor commercial workers. This receptor is considered to have the highest level of exposure at the property. Other receptors generally have lower exposures, and thus lower risk estimates. Therefore, risk estimates generated for future on-site indoor commercial workers will be protective of other potential receptors at the property.

The previous screening-level health risk assessment evaluated risks from exposure to soil. However, these exposures did not account for potential migration of VOCs from the subsurface into indoor air. In general USEPA does not recommend evaluating the indoor air exposure pathway using soil matrix data (USEPA 2002a). Because groundwater beneath a portion of the property is considered a potential VOC source area, soil gas data were recently collected. These data are further evaluated and are the focus of this screening-level indoor air health risk assessment.

### **Screening-Level Indoor Air Health Risk Assessment**

As discussed above, the previous screening-level health risk assessment did not consider the indoor air pathway. Therefore, soil gas data were collected to specifically evaluate this potential exposure pathway at the property.

Human health risks are represented by estimated theoretical upper-bound cancer risks and non-cancer hazards derived in accordance with standard USEPA methods. The acceptable risk levels defined by USEPA for the protection of human health, and following those discussed previously with NDEP, are:

1. For non-carcinogenic compounds, the acceptable criterion is a cumulative hazard index (HI) of one or less. If the screening HI is determined to be greater than 1.0, target organ-specific HIs will be calculated for primary and secondary organs. The final risk goal will be to achieve target organ-specific non-carcinogenic HIs of less than 1.0; and
2. For known or suspected chemical and radionuclide carcinogens, the acceptable ceiling for a cumulative incremental lifetime cancer risk (ILCR) ranges from  $10^{-6}$  to  $10^{-4}$ . The risk goal established by the NDEP is  $10^{-6}$ .
3. Where background levels exceed risk level goals, metals and radionuclides in property soils are targeted to have risks no greater than those associated with background conditions.

This screening-level indoor air health risk assessment follows the basic procedures outlined in USEPA *Risk Assessment Guidance for Superfund: Volume I—Human Health Evaluation Manual* (RAGS; USEPA 1989). Other guidance documents were also consulted for the screening-level indoor air health risk assessment.

#### Selection of Chemicals of Potential Concern

The broad suite of analytes sampled for was the initial list of chemicals of potential concern (COPCs) at the property. However, in order to ensure that a risk assessment focuses on those substances that contribute the greatest to the overall risk (USEPA 1989); only one procedure was used to eliminate the chemicals for quantitative evaluation in the screening-level indoor air health risk assessment: identification of chemicals that were not detected in any of the soil gas sample locations within the property. That is, all chemicals that were detected in any soil gas sample within the property was considered a COPC and evaluated in the screening-level indoor air health risk assessment. The identification of those chemicals detected in soil gas samples within the property are presented in Table 1.

#### Determination of Exposure Point Concentrations

A representative exposure concentration is a COPC-specific and media-specific concentration value. In risk assessment, these exposure concentrations are values incorporated into the exposure assessment equations from which potential baseline human exposures are calculated.

Due to the uncertainty associated with determining the true average concentration at a site, where direct measurements of the site average are unavailable, the USEPA recommends using the lower of the maximum detected concentration or the 95 percent upper confidence limit (UCL) as the concentration of a chemical to which an individual could be exposed over time (USEPA 1992). For the 95 percent UCL concentration approach, the 95 percent UCL is typically computed in order to represent the area-wide exposure point concentrations. The 95 percent UCL is defined as the value that, when calculated repeatedly for randomly drawn subsets of site data, equals or exceeds the true mean 95 percent of the time (USEPA 1992). The purpose for using the 95 percent UCL is to take into account the different concentrations a person may be exposed to on any given day. That is, an individual will be exposed to a range of concentrations that exist at an exposure area, from non-detect to the maximum concentration, over an entire exposure period.

The 95 percent UCL statistical calculations were performed using the computer statistical software program GISdT<sup>®</sup> (Neptune and Company 2007). The formulas for calculating the 95 percent UCL COPC concentration (as the representative exposure concentration) are presented in USEPA (1992, 2002b). The representativeness of the 95 percent UCLs for each exposure area, that is, a property-wide mean concentration is valid since concentrations of COPCs are primarily emanating from a sub-surface groundwater source, and localized ‘hot spot’ concentrations within the property are not expected. Therefore each measurement is assumed to be equally representative for that chemical at any point in the property and calculation of the 95 percent UCL is appropriate. The soil gas representative exposure concentrations used in this screening-level indoor air health risk assessment are presented in Table 1.

### *Indoor Air*

The flux of COPCs from the subsurface and dispersion into indoor air were estimated using the USEPA spreadsheet-based Johnson and Ettinger model (USEPA 2004). The model is based on the vapor intrusion model published by Johnson and Ettinger (1991). The Johnson and Ettinger vapor intrusion model is a screening-level model, which incorporates both convective and diffusive mechanisms for estimating the transport of chemical vapors emanating from either subsurface soils or groundwater into indoor spaces located directly above the source of contamination. The model is constructed to calculate steady-state vapor transport (infinite source). Maximum detected VOCs concentrations in soil gas were used as representative exposure concentrations for the indoor air exposure pathway. The default physical properties and building characteristics contained in the USEPA Johnson and Ettinger model were used in

this evaluation. These values are presented in Table 2. Table 3 presents the indoor air concentrations predicted by the Johnson and Ettinger model for each of the COPCs.

### Risk Assessment Methodology

The method used in the screening-level indoor air health risk assessment consists of several steps. The first step is the calculation of exposure point concentrations representative of the particular area (see above). The second step is fate and transport modeling to predict concentrations that may be present when direct measurements are not available. The third step is the exposure assessment for the various receptors present in the particular areas. The next step is to define the toxicity values for each COPC. The final step is risk characterization where theoretical upper-bound ILCRs and non-cancer HIs are calculated. The *BRC Closure Plan* (BRC, ERM, and DBSA 2007) provides a full discussion on the risk assessment methodology for the project, and used in this screening-level indoor air health risk assessment.

Table 2 presents each of the exposure parameters used in the screening-level indoor air health risk assessment. Toxicity values, when available, are published by the USEPA in the on-line Integrated Risk Information System (IRIS; USEPA 2008) and the Health Effects Assessment Summary Tables (HEAST; USEPA 1997). Unit risk factors (URFs) are chemical-specific, experimentally-derived potency values used to calculate the risk of cancer resulting from exposure to carcinogenic chemicals. A higher value implies a more potent carcinogen. Reference concentrations (RfCs) are experimentally derived “no-effect” values used to quantify the extent of adverse non-cancer health effects from exposure to chemicals. Here, a lower RfC implies a more potent toxicant. These criteria are generally developed by USEPA risk assessment work groups and listed in USEPA risk assessment guidance documents and databases. The hierarchy for selecting toxicity criteria presented in the *BRC Closure Plan* (BRC, ERM, and DBSA 2007) was used.

### Uncertainty Analysis

Risk estimates are values that have uncertainties associated with them. These uncertainties, which arise at every step of a risk assessment, are evaluated to provide an indication of the uncertainty associated with a risk estimate. Risk assessments are not intended to estimate the true risk to a receptor associated with exposure to chemicals in the environment. In fact, estimating the true risk is impossible because of the variability in the exposed or potentially exposed populations. Therefore, risk assessment is a means of estimating the probability that an adverse health effect (*e.g.*, cancer, impaired reproduction) will occur in a receptor in order to assist in

decision making regarding the protection of human health. The multitude of conservative assumptions used in risk assessments guard against underestimation of risks.

Risk estimates are calculated by combining site data, assumptions about individual receptor's exposures to impacted media, and toxicity data. The uncertainties in this screening-level indoor air health risk assessment can be grouped into four main categories that correspond to these steps:

- Uncertainties in environmental sampling and analysis
- Uncertainties in fate and transport modeling
- Uncertainties in assumptions concerning exposure scenarios
- Uncertainties in toxicity data and dose-response extrapolations

Additional discussion on the uncertainties associated with the screening-level indoor air health risk assessment is provided below.

The screening-level indoor air health risk assessment for the property was based on the sampling results obtained from an soil gas investigation conducted in 2008. Errors in sampling results can arise from the field sampling, laboratory analyses, and data analyses. Errors in laboratory analysis procedures are possible, although the impacts of these sorts of errors on the risk estimates are likely to be low. The environmental sampling at the property is one source of uncertainty in the evaluation. However, the number of sampling locations and events is large and widespread, and sampling was performed using approved procedures; therefore, the sampling and analysis data is sufficient to characterize the impacts and the associated potential risks.

The amount of COPCs the body absorbs may be different from the amount of a COPC contacted. In this screening-level indoor air health risk assessment, absorption of inhaled COPCs is conservatively assumed to be 100 percent. Actual chemical and site specific values are likely less than this default value.

Toxicity criteria have not been established for many of the chemicals detected at the property. These chemicals were not quantitatively evaluated in the screening-level indoor air health risk assessment. Thus, the risks presented in this assessment could be underestimated as a result.

The selection of exposure pathways is a process, often based on best professional judgement, which attempts to identify the most probable potentially harmful exposure scenarios. In a risk

assessment it is possible that risks are not calculated for all of the exposure pathways that may occur, possibly causing some underestimation of risk. In this assessment, risks were estimated for one receptor; future on-site indoor commercial workers. Risks for the most likely route of exposure to future on-site indoor commercial workers were estimated. Specifically, risks to future on-site indoor commercial workers were estimated for inhalation of indoor air. Although it is possible that other exposure routes could exist, these exposures are expected to be lower than the risks associated with the pathway considered.

Uncertainties from different sources are compounded in the screening-level indoor air health risk assessment. For example, if a person's daily intake rate for a chemical is compared to an RfC to determine potential health risks, the uncertainties in the concentration measurements, exposure assumptions, and toxicities will all be expressed in the result. Because the exposure assumptions and toxicity criteria are considered conservative, the risk estimates calculated in this screening-level indoor air health risk assessment are likely to overestimate rather than underestimate potential risks.

#### Screening-Level Indoor Air Health Risk Assessment Results

This screening-level indoor air health risk assessment has evaluated potential risks to human health associated with chemicals detected in soil gas at the Tronox Parcels A and B property. The calculation of chemical theoretical upper-bound ILCRs and non-cancer health effects are presented in Table 4. All calculation spreadsheets for this screening-level indoor air health risk assessment are included in Attachment A.

The total cumulative non-cancer HI for future on-site indoor commercial workers at the property is 0.01, which is below the target HI of 1.0. Therefore, because the total cumulative HI is below 1.0, the potential for adverse health effects is considered unlikely.

The theoretical upper-bound ILCR for future on-site indoor commercial workers at the property is  $4 \times 10^{-6}$ . The risks are primarily driven by chloroform, which contributes 95 percent of the theoretical upper-bound ILCR. Although the ILCR is above the risk goal of  $1 \times 10^{-6}$ , it is within the acceptable risk range from  $10^{-6}$  to  $10^{-4}$ . Therefore, these results indicate that future receptor exposures at the property should not result in unacceptable carcinogenic risks.

#### **Summary**

Based on the results of the 2008 soil gas investigation, this data review, and the screening-level indoor air health risk assessment, concentration levels of chemicals in soil gas at the Tronox

Parcels A and B property are not at levels of concern for human health risk for an indoor commercial scenario. In summary, BEC concludes that an NFAD for the property is further supported by these results.

## REFERENCES

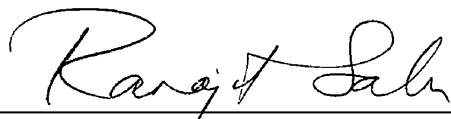
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Attachments: Figure 1 – Tronox Parcels A/B Phase B Soil Gas Sample Locations  
Table 1 – Chemicals of Potential Concern and Representative Exposure  
Concentrations in Soil Gas Table 2 – Johnson and Ettinger Model Input  
Parameters  
Table 3 – Model Estimated Indoor Air Concentrations  
Table 4 – Screening-Level Indoor Air Health Risk Assessment Results  
Attachment A – Screening-Level Indoor Air Health Risk Assessment Calculation  
Spreadsheets (on CD)

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I hereby certify that I am responsible for the services described in this document and for the preparation of this document. The services described in this document have been provided in a manner consistent with the current standards of the profession and to the best of my knowledge comply with all applicable federal, state and local statutes, regulations and ordinances. I hereby certify that all laboratory analytical data was generated by a laboratory certified by the NDEP for each constituent and media presented herein.



November 13, 2008

Dr. Ranajit Sahu, C.E.M. (No. EM-1699, Exp. 10/07/2009)      Date  
BRC Project Manager

## FIGURES





Tronox Parcels A/B Boundary

**Soil Gas Sample Locations**

- Location within Parcels A/B
- Other Sample Location

BEC / Tronox Parcels A/B Data Review  
 BMI Common Areas, Henderson, Nevada

**FIGURE 1**

**TRONOX PARCELS A/B  
 PHASE B SOIL GAS  
 SAMPLE LOCATIONS**



Prepared by: MKJ  
 Date: 11/03/08

JOB No. 0069073  
 FILE: GIS/BEC/TRONOX\_AB/FIGURE\_1.MXD



## TABLES

**TABLE 1**  
**CHEMICALS OF POTENTIAL CONCERN AND REPRESENTATIVE EXPOSURE CONCENTRATIONS IN SOIL GAS**  
**TRONOX PARCELS A/B SOIL GAS INVESTIGATION**  
**CLARK COUNTY, NEVADA**  
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| Chemical                       | Number of Samples | Number of Detections | Frequency of Detects | Minimum DL | Maximum DL | Minimum Detection | Median <sup>a</sup> | Mean <sup>a</sup> | Maximum Detection | 95% UCL | EPC   |
|--------------------------------|-------------------|----------------------|----------------------|------------|------------|-------------------|---------------------|-------------------|-------------------|---------|-------|
| 1,1,1-Trichloroethane          | 9                 | 0                    | 0%                   | 0.15       | 0.17       | --                | 0.08                | 0.079             | --                | --      | --    |
| 1,1,2,2-Tetrachloroethane      | 9                 | 0                    | 0%                   | 0.15       | 0.17       | --                | 0.08                | 0.079             | --                | --      | --    |
| 1,1,2-Trichloroethane          | 9                 | 1                    | 11%                  | 0.15       | 0.17       | 0.12              | 0.08                | 0.083             | 0.12              | 0.093   | 0.093 |
| 1,1,2-Trichlorotrifluoroethane | 9                 | 9                    | 100%                 | --         | --         | 0.45              | 0.49                | 0.5               | 0.63              | 0.55    | 0.55  |
| 1,1-Dichloroethane             | 9                 | 7                    | 78%                  | 0.15       | 0.16       | 0.11              | 0.41                | 8                 | 27                | 15.6    | 15.6  |
| 1,1-Dichloroethene             | 9                 | 2                    | 22%                  | 0.15       | 0.17       | 0.1               | 0.08                | 0.086             | 0.12              | 0.097   | 0.097 |
| 1,2,4-Trichlorobenzene         | 9                 | 3                    | 33%                  | 0.15       | 0.17       | 0.2               | 0.08                | 0.21              | 0.75              | 0.37    | 0.37  |
| 1,2,4-Trimethylbenzene         | 9                 | 9                    | 100%                 | --         | --         | 0.12              | 0.37                | 0.87              | 3.5               | 1.8     | 1.8   |
| 1,2-Dibromo-3-chloropropane    | 9                 | 0                    | 0%                   | 0.74       | 0.85       | --                | 0.39                | 0.4               | --                | --      | --    |
| 1,2-Dichlorobenzene            | 9                 | 0                    | 0%                   | 0.15       | 0.17       | --                | 0.08                | 0.079             | --                | --      | --    |
| 1,2-Dichloroethane             | 9                 | 3                    | 33%                  | 0.15       | 0.16       | 0.32              | 0.08                | 0.27              | 1.1               | 0.56    | 0.56  |
| 1,2-Dichloropropane            | 9                 | 4                    | 44%                  | 0.15       | 0.17       | 0.085             | 0.085               | 0.13              | 0.47              | 0.25    | 0.25  |
| 1,2-Dichlorotetrafluoroethane  | 9                 | 5                    | 56%                  | 0.77       | 0.85       | 0.085             | 0.1                 | 0.23              | 0.1               | 0.33    | 0.1   |
| 1,3,5-Trimethylbenzene         | 9                 | 5                    | 56%                  | 0.77       | 0.85       | 0.09              | 0.385               | 0.49              | 1.9               | 0.99    | 0.99  |
| 1,3-Dichlorobenzene            | 9                 | 3                    | 33%                  | 0.15       | 0.17       | 0.098             | 0.085               | 0.12              | 0.32              | 0.19    | 0.19  |
| 1,4-Dichlorobenzene            | 9                 | 9                    | 100%                 | --         | --         | 0.31              | 0.84                | 8                 | 43                | 21.1    | 21.1  |
| 1,4-Dioxane                    | 9                 | 5                    | 56%                  | 0.77       | 0.85       | 0.14              | 0.385               | 0.29              | 0.39              | 0.37    | 0.37  |
| 2-Butanone                     | 9                 | 9                    | 100%                 | --         | --         | 4.6               | 7                   | 7.3               | 13                | 9.1     | 9.1   |
| 2-Hexanone                     | 9                 | 9                    | 100%                 | --         | --         | 0.26              | 0.43                | 0.42              | 0.52              | 0.46    | 0.46  |
| 2-Methoxy-2-methyl-butane      | 9                 | 0                    | 0%                   | 0.74       | 0.85       | --                | 0.39                | 0.4               | --                | --      | --    |
| 4-Ethyltoluene                 | 9                 | 6                    | 67%                  | 0.77       | 0.85       | 0.11              | 0.385               | 0.41              | 1.5               | 0.77    | 0.77  |
| 4-Isopropyltoluene             | 9                 | 7                    | 78%                  | 0.77       | 0.85       | 0.13              | 0.385               | 0.8               | 4.4               | 1.8     | 1.8   |
| 4-Methyl-2-pentanone           | 9                 | 9                    | 100%                 | --         | --         | 0.14              | 0.29                | 1.3               | 9.2               | 4.2     | 4.2   |
| Acetone                        | 9                 | 7                    | 78%                  | 15         | 24         | 12                | 18                  | 21                | 50                | 30.9    | 30.9  |
| Acrylonitrile                  | 9                 | 3                    | 33%                  | 0.77       | 0.85       | 0.11              | 0.385               | 0.31              | 0.12              | 0.40    | 0.12  |
| Allyl chloride                 | 9                 | 1                    | 11%                  | 0.15       | 0.17       | 0.17              | 0.08                | 0.089             | 0.17              | 0.11    | 0.11  |
| alpha-Methylstyrene            | 9                 | 4                    | 44%                  | 0.74       | 0.85       | 0.13              | 0.39                | 1.1               | 7.7               | 3.6     | 3.6   |
| Benzene                        | 9                 | 9                    | 100%                 | --         | --         | 1.2               | 1.9                 | 1.9               | 2.7               | 2.2     | 2.2   |
| Benzyl Chloride                | 9                 | 0                    | 0%                   | 0.15       | 0.17       | --                | 0.08                | 0.079             | --                | --      | --    |
| Bromodichloromethane           | 9                 | 6                    | 67%                  | 0.16       | 0.17       | 0.098             | 0.18                | 0.22              | 0.67              | 0.38    | 0.38  |
| Bromoform                      | 9                 | 1                    | 11%                  | 0.74       | 0.85       | 0.27              | 0.39                | 0.39              | 0.27              | 0.41    | 0.27  |
| Bromomethane                   | 9                 | 1                    | 11%                  | 0.15       | 0.17       | 0.11              | 0.08                | 0.082             | 0.11              | 0.091   | 0.091 |
| Carbon disulfide               | 9                 | 7                    | 78%                  | 1.1        | 1.4        | 1.5               | 2                   | 4.9               | 14                | 8.2     | 8.2   |

**TABLE 1**  
**CHEMICALS OF POTENTIAL CONCERN AND REPRESENTATIVE EXPOSURE CONCENTRATIONS IN SOIL GAS**  
**TRONOX PARCELS A/B SOIL GAS INVESTIGATION**  
**CLARK COUNTY, NEVADA**  
 (Page 2 of 3)

| <b>Chemical</b>            | <b>Number of Samples</b> | <b>Number of Detections</b> | <b>Frequency of Detects</b> | <b>Minimum DL</b> | <b>Maximum DL</b> | <b>Minimum Detection</b> | <b>Median<sup>a</sup></b> | <b>Mean<sup>a</sup></b> | <b>Maximum Detection</b> | <b>95% UCL</b> | <b>EPC</b> |
|----------------------------|--------------------------|-----------------------------|-----------------------------|-------------------|-------------------|--------------------------|---------------------------|-------------------------|--------------------------|----------------|------------|
| Carbon tetrachloride       | 9                        | 9                           | 100%                        | --                | --                | 0.25                     | 0.39                      | 3                       | 11                       | 5.8            | 5.8        |
| Chlorobenzene              | 9                        | 3                           | 33%                         | 0.15              | 0.17              | 0.16                     | 0.08                      | 0.12                    | 0.31                     | 0.18           | 0.18       |
| Chloroethane               | 9                        | 7                           | 78%                         | 0.15              | 0.16              | 0.14                     | 0.87                      | 3.1                     | 11                       | 5.9            | 5.9        |
| Chloroform                 | 9                        | 9                           | 100%                        | --                | --                | 8.6                      | 34                        | 140                     | 440                      | 259            | 259        |
| Chloromethane              | 9                        | 1                           | 11%                         | 0.15              | 0.17              | 0.076                    | 0.08                      | 0.079                   | 0.076                    | 0.082          | 0.076      |
| cis-1,2-Dichloroethene     | 9                        | 2                           | 22%                         | 0.15              | 0.17              | 0.15                     | 0.08                      | 1.5                     | 13                       | 5.8            | 5.8        |
| cis-1,3-Dichloropropene    | 9                        | 0                           | 0%                          | 0.74              | 0.85              | --                       | 0.39                      | 0.4                     | --                       | --             | --         |
| Dibromochloromethane       | 9                        | 1                           | 11%                         | 0.15              | 0.17              | 0.12                     | 0.08                      | 0.084                   | 0.12                     | 0.094          | 0.094      |
| Dichlorodifluoromethane    | 9                        | 9                           | 100%                        | --                | --                | 1.8                      | 2                         | 2                       | 2.1                      | 2.1            | 2.1        |
| Ethanol                    | 9                        | 9                           | 100%                        | --                | --                | 2.3                      | 11                        | 14                      | 32                       | 20.5           | 20.5       |
| Ethyl t-butyl ether        | 9                        | 0                           | 0%                          | 0.74              | 0.85              | --                       | 0.39                      | 0.4                     | --                       | --             | --         |
| Ethylbenzene               | 9                        | 7                           | 78%                         | 0.77              | 0.85              | 0.1                      | 0.385                     | 0.44                    | 1.2                      | 0.70           | 0.70       |
| Ethylene dibromide         | 9                        | 0                           | 0%                          | 0.15              | 0.17              | --                       | 0.08                      | 0.079                   | --                       | --             | --         |
| Hexachlorobutadiene        | 9                        | 5                           | 56%                         | 0.15              | 0.17              | 0.49                     | 0.49                      | 0.66                    | 2.4                      | 1.2            | 1.2        |
| isopropyl ether            | 9                        | 0                           | 0%                          | 0.74              | 0.85              | --                       | 0.39                      | 0.4                     | --                       | --             | --         |
| Isopropylbenzene           | 9                        | 3                           | 33%                         | 0.74              | 0.85              | 0.088                    | 0.385                     | 0.31                    | 0.19                     | 0.40           | 0.19       |
| m,p-Xylene                 | 9                        | 8                           | 89%                         | 0.77              | 0.77              | 0.22                     | 0.8                       | 1.4                     | 5.9                      | 2.8            | 2.8        |
| Methyl methacrylate        | 9                        | 1                           | 11%                         | 0.74              | 0.85              | 0.42                     | 0.39                      | 0.4                     | 0.42                     | 0.41           | 0.41       |
| Methyl tert butyl ether    | 9                        | 6                           | 67%                         | 0.15              | 0.16              | 0.1                      | 0.33                      | 1.4                     | 7.8                      | 3.7            | 3.7        |
| Methylene chloride         | 9                        | 8                           | 89%                         | 0.77              | 0.77              | 0.23                     | 0.63                      | 1.2                     | 3.7                      | 2.0            | 2.0        |
| Naphthalene                | 9                        | 9                           | 100%                        | --                | --                | 0.42                     | 0.83                      | 1.2                     | 4.2                      | 2.1            | 2.1        |
| N-Butylbenzene             | 9                        | 9                           | 100%                        | --                | --                | 0.12                     | 0.26                      | 0.31                    | 0.68                     | 0.44           | 0.44       |
| n-Heptane                  | 9                        | 6                           | 67%                         | 0.77              | 0.85              | 0.24                     | 0.425                     | 0.42                    | 0.72                     | 0.52           | 0.52       |
| n-Octane                   | 9                        | 4                           | 44%                         | 0.77              | 0.85              | 0.23                     | 0.385                     | 0.49                    | 1.5                      | 0.86           | 0.86       |
| N-Propylbenzene            | 9                        | 5                           | 56%                         | 0.77              | 0.85              | 0.084                    | 0.385                     | 0.31                    | 0.52                     | 0.41           | 0.41       |
| o-Xylene                   | 9                        | 7                           | 78%                         | 0.77              | 0.85              | 0.12                     | 0.42                      | 0.61                    | 2.1                      | 1.1            | 1.1        |
| sec-Butylbenzene           | 9                        | 1                           | 11%                         | 0.74              | 0.85              | 0.097                    | 0.385                     | 0.36                    | 0.097                    | 0.43           | 0.10       |
| Styrene                    | 9                        | 5                           | 56%                         | 0.77              | 0.85              | 0.16                     | 0.385                     | 0.38                    | 0.6                      | 0.45           | 0.45       |
| t-Butyl alcohol            | 9                        | 9                           | 100%                        | --                | --                | 0.2                      | 0.45                      | 0.44                    | 0.67                     | 0.53           | 0.53       |
| tert-Butylbenzene          | 9                        | 1                           | 11%                         | 0.29              | 0.34              | 0.14                     | 0.155                     | 0.16                    | 0.14                     | 0.16           | 0.14       |
| Tetrachloroethene          | 9                        | 9                           | 100%                        | --                | --                | 1.1                      | 5.3                       | 7.4                     | 30                       | 13.8           | 13.8       |
| Toluene                    | 9                        | 9                           | 100%                        | --                | --                | 1.2                      | 2                         | 4.4                     | 19                       | 9.8            | 9.8        |
| trans-1,2-Dichloroethylene | 9                        | 0                           | 0%                          | 0.15              | 0.17              | --                       | 0.08                      | 0.079                   | --                       | --             | --         |

**TABLE 1**  
**CHEMICALS OF POTENTIAL CONCERN AND REPRESENTATIVE EXPOSURE CONCENTRATIONS IN SOIL GAS**  
**TRONOX PARCELS A/B SOIL GAS INVESTIGATION**  
**CLARK COUNTY, NEVADA**  
**(Page 3 of 3)**

| <b>Chemical</b>           | <b>Number of Samples</b> | <b>Number of Detections</b> | <b>Frequency of Detects</b> | <b>Minimum DL</b> | <b>Maximum DL</b> | <b>Minimum Detection</b> | <b>Median<sup>a</sup></b> | <b>Mean<sup>a</sup></b> | <b>Maximum Detection</b> | <b>95% UCL</b> | <b>EPC</b> |
|---------------------------|--------------------------|-----------------------------|-----------------------------|-------------------|-------------------|--------------------------|---------------------------|-------------------------|--------------------------|----------------|------------|
| trans-1,3-Dichloropropene | 9                        | 0                           | 0%                          | 0.74              | 0.85              | --                       | 0.39                      | 0.4                     | --                       | --             | --         |
| Trichloroethene           | 9                        | 9                           | 100%                        | --                | --                | 0.96                     | 1.3                       | 6.5                     | 42                       | 19.4           | 19.4       |
| Trichlorofluoromethane    | 9                        | 9                           | 100%                        | --                | --                | 0.95                     | 1.1                       | 1.1                     | 1.4                      | 1.2            | 1.2        |
| Vinyl acetate             | 9                        | 7                           | 78%                         | 7.7               | 7.8               | 0.99                     | 3.5                       | 3.4                     | 5                        | 4.2            | 4.2        |
| Vinyl chloride            | 9                        | 2                           | 22%                         | 0.15              | 0.16              | 0.12                     | 0.08                      | 0.087                   | 0.12                     | 0.099          | 0.099      |

Note: All units in  $\mu\text{g}/\text{m}^3$ .

a - Includes both detect values and non-detect values, with one-half the DL used for non-detect values.

DL = detection limit

UCL = upper confidence limit

EPC = exposure point concentration

-- = Not applicable or statistic not evaluated because all results were non-detect..

**TABLE 2**  
**JOHNSON AND ETTINGER MODEL INPUT PARAMETERS**  
**TRONOX PARCELS A/B SOIL GAS INVESTIGATION**  
**CLARK COUNTY, NEVADA**  
**(Page 1 of 1)**

| Parameter  | Value | Reference/Rationale           |
|--|-------|-------------------------------|
| Depth below grade to bottom of enclosed floor space (cm)       | 15    | Model default (slab on grade) |
| Average soil temperature (°C)                                  | 10    | Model default                 |
| Soil gas sampling depth (cm)                                   | 200   | Site-specific (five feet bgs) |
| Thickness of soil stratum (cm)                                 | 200   | Site-specific (five feet bgs) |
| Soil stratum used to calculate soil vapor permeability         | S     | Sand                          |
| Vadose zone dry bulk density (g/cm <sup>3</sup> )              | 1.66  | Model default                 |
| Vadose zone total porosity (unitless)                          | 0.375 | Model default                 |
| Vadose zone water-filled porosity (unitless)                   | 0.054 | Model default                 |
| Enclosed space floor thickness (cm)                            | 15    | Model default                 |
| Soil-building pressure differential (g/cm-s <sup>2</sup> )     | 40    | Model default                 |
| Enclosed space floor length (cm)                               | 1,000 | Model default                 |
| Enclosed space floor width (cm)                                | 1,000 | Model default                 |
| Modeling Enclosed space height (cm)                            | 244   | Model default                 |
| Floor-wall seam crack width (cm)                               | 0.1   | Model default                 |
| Average vapor flow rate into building, Q <sub>soil</sub> (L/m) | 5     | Model default                 |
| Indoor air exchange rate (1/hr)                                | 0.25  | Model default                 |
| Exposure duration (yrs)  | 25    | Model default (commercial)    |
| Exposure frequency (days/yr)                                   | 250   | Model default (commercial)    |
| Averaging time for carcinogens (yrs)                           | 70    | Model default (commercial)    |
| Averaging time for non-carcinogens (yrs)                       | 25    | Model default (commercial)    |



**TABLE 3**  
**MODEL ESTIMATED INDOOR AIR CONCENTRATIONS**  
**TRONOX PARCELS A/B SOIL GAS INVESTIGATION**  
**CLARK COUNTY, NEVADA**  
(Page 1 of 2)

| <b>Chemical</b>                | <b>Predicted Indoor<br/>Air Concentration<br/>(<math>\mu\text{g}/\text{m}^3</math>)<sup>1</sup></b> |
|--------------------------------|---|
| 1,1,2-Trichloroethane          | 2.1 E-4   |
| 1,1,2-Trichlorotrifluoroethane | 1.2 E-3   |
| 1,1-Dichloroethane             | 3.4 E-2   |
| 1,1-Dichloroethene             | 2.3 E-4   |
| 1,2,4-Trichlorobenzene         | 4.3 E-4   |
| 1,2,4-Trimethylbenzene         | 3.4 E-3   |
| 1,2-Dichloroethane             | 1.4 E-3   |
| 1,2-Dichloropropane            | 5.6 E-4   |
| 1,2-Dichlorotetrafluoroethane  | 2.2 E-4   |
| 1,3,5-Trimethylbenzene         | 1.9 E-3   |
| 1,3-Dichlorobenzene            | 3.9 E-4   |
| 1,4-Dichlorobenzene            | 4.4 E-2   |
| 1,4-Dioxane                    | 9.3 E-4   |
| 2-Butanone                     | 2.0 E-2   |
| 2-Hexanone                     | 1.5 E-3   |
| 4-Ethyltoluene                 | 1.8 E-3   |
| 4-Isopropyltoluene             | 3.3 E-3   |
| 4-Methyl-2-pentanone           | 9.2 E-3   |
| Acetone                        | 8.6 E-2   |
| Acrylonitrile                  | 3.3 E-4   |
| Allyl chloride                 | 3.0 E-4   |
| alpha-Methylstyrene            | 1.3 E-2   |
| Benzene                        | 5.2 E-3   |
| Bromodichloromethane           | 4.4 E-4   |
| Bromoform                      | 1.8 E-4   |
| Bromomethane                   | 1.9 E-4   |
| Carbon disulfide               | 2.1 E-2   |
| Carbon tetrachloride           | 1.3 E-2   |
| Chlorobenzene                  | 3.7 E-4   |
| Chloroethane                   | 2.1 E-2   |
| Chloroform                     | 6.6 E-1   |
| Chloromethane                  | 2.1 E-4   |
| cis-1,2-Dichloroethene         | 1.2 E-2   |
| Dibromochloromethane           | 7.9 E-5   |
| Dichlorodifluoromethane        | 4.2 E-3   |
| Ethanol                        | 5.6 E-2   |
| Ethylbenzene                   | 1.5 E-3   |
| Hexachlorobutadiene            | 2.1 E-3   |
| Isopropylbenzene               | 3.8 E-4   |
| m,p-Xylene                     | 5.7 E-3   |
| Methyl methacrylate            | 9.1 E-4   |
| Methyl tert butyl ether        | 9.4 E-3   |
| Methylene chloride             | 4.9 E-3   |
| Naphthalene                    | 3.9 E-3   |
| N-Butylbenzene                 | 8.1 E-4   |
| n-Heptane                      | 1.7 E-3   |
| n-Octane                       | 1.9 E-3   |
| N-Propylbenzene                | 7.7 E-4   |
| o-Xylene                       | 2.6 E-3   |

**TABLE 3**  
**MODEL ESTIMATED INDOOR AIR CONCENTRATIONS**  
**TRONOX PARCELS A/B SOIL GAS INVESTIGATION**  
**CLARK COUNTY, NEVADA**  
**(Page 2 of 2)**

| <b>Chemical</b>        | <b>Predicted Indoor<br/>Air Concentration<br/>(<math>\mu\text{g}/\text{m}^3</math>)<sup>1</sup></b> |
|------------------------|---|
| sec-Butylbenzene       | 1.8 E-4   |
| Styrene                | 9.4 E-4   |
| t-Butyl alcohol        | 1.3 E-3   |
| tert-Butylbenzene      | 2.6 E-4   |
| Tetrachloroethene      | 2.9 E-2   |
| Toluene                | 3.2 E-2   |
| Trichloroethene        | 4.3 E-2   |
| Trichlorofluoromethane | 2.9 E-3   |
| Vinyl acetate          | 9.7 E-3   |
| Vinyl chloride         | 2.6 E-4   |

<sup>1</sup> - Calculated using the J&E Model (included on CD).

**TABLE 4**  
**SCREENING-LEVEL INDOOR AIR HEALTH RISK ASSESSMENT RESULTS**  
**TRONOX PARCELS A/B SOIL GAS INVESTIGATION**  
**CLARK COUNTY, NEVADA**  
 (Page 1 of 2)

| <b>Chemical</b>             | <b>Non-Cancer Reference Concentration (mg/kg-d)</b> | <b>Unit Risk Factor (mg/kg-d)<sup>-1</sup></b> | <b>J&amp;E Predicted Conc.<sup>a</sup></b> | <b>Non-Cancer Hazard Index</b> | <b>Incremental Lifetime Cancer Risk</b> |
|-----------------------------|---|--|--|--------------------------------|---|
| 1,1,2-Trichloroethane       | 1.4 E-2   | 1.6 E-5  | 2.1 E-4                                    | 0.00001                        | 8 E-10                                  |
| 1,1,2-Trichlorotrifluoroeth | 3.0 E+1   | NA   | 1.2 E-3                                    | 0.00000003                     | NA                                      |
| 1,1-Dichloroethane          | 7.0 E-1   | NA   | 3.4 E-2                                    | 0.00003                        | NA                                      |
| 1,1-Dichloroethene          | 2.0 E-1   | NA   | 2.3 E-4                                    | 0.0000008                      | NA                                      |
| 1,2,4-Trichlorobenzene      | 4.0 E-3   | NA   | 4.3 E-4                                    | 0.00007                        | NA                                      |
| 1,2,4-Trimethylbenzene      | 7.0 E-3   | NA   | 3.4 E-3                                    | 0.0003                         | NA                                      |
| 1,2-Dichloroethane          | 4.9 E-3   | 2.6 E-5  | 1.4 E-3                                    | 0.0002                         | 9 E-9                                   |
| 1,2-Dichloropropane         | 4.0 E-3   | 1.9 E-5  | 5.6 E-4                                    | 0.0001                         | 3 E-9                                   |
| 1,2-Dichlorotetrafluoroeth  | NA  | NA   | 2.2 E-4                                    | NA                             | NA                                      |
| 1,3,5-Trimethylbenzene      | 6.0 E-3   | NA   | 1.9 E-3                                    | 0.0002                         | NA                                      |
| 1,3-Dichlorobenzene         | 8.0 E-3   | NA   | 3.9 E-4                                    | 0.00003                        | NA                                      |
| 1,4-Dichlorobenzene         | 8.0 E-1   | 6.9 E-6  | 4.4 E-2                                    | 0.00004                        | 7 E-8                                   |
| 1,4-Dioxane                 | NA  | 3.1 E-6  | 9.3 E-4                                    | NA                             | 7 E-10                                  |
| 2-Butanone                  | 5.0 E+0   | NA   | 2.0 E-2                                    | 0.000003                       | NA                                      |
| 2-Hexanone                  | NA  | NA   | 1.5 E-3                                    | NA                             | NA                                      |
| 4-Ethyltoluene              | NA  | NA   | 1.8 E-3                                    | NA                             | NA                                      |
| 4-Isopropyltoluene          | NA  | NA   | 3.3 E-3                                    | NA                             | NA                                      |
| 4-Methyl-2-pentanone        | 3.0 E+0   | NA   | 9.2 E-3                                    | 0.000002                       | NA                                      |
| Acetone                     | 3.2 E+0   | NA   | 8.6 E-2                                    | 0.00002                        | NA                                      |
| Acrylonitrile               | 2.0 E-3   | 6.8 E-5  | 3.3 E-4                                    | 0.0001                         | 6 E-9                                   |
| Allyl chloride              | 1.0 E-3   | NA   | 3.0 E-4                                    | 0.0002                         | NA                                      |
| alpha-Methylstyrene         | 4.0 E-2   | NA   | 1.3 E-2                                    | 0.0002                         | NA                                      |
| Benzene                     | 3.0 E-2   | 7.8 E-6  | 5.2 E-3                                    | 0.0001                         | 1 E-8                                   |
| Bromodichloromethane        | 7.0 E-2   | 1.8 E-5  | 4.4 E-4                                    | 0.000004                       | 2 E-9                                   |
| Bromoform                   | 7.0 E-2   | 1.1 E-6  | 1.8 E-4                                    | 0.000002                       | 5 E-11                                  |
| Bromomethane                | 5.0 E-3   | NA   | 1.9 E-4                                    | 0.00003                        | NA                                      |
| Carbon disulfide            | 7.0 E-1   | NA   | 2.1 E-2                                    | 0.00002                        | NA                                      |
| Carbon tetrachloride        | NA  | 1.5 E-5  | 1.3 E-2                                    | NA                             | 5 E-8                                   |
| Chlorobenzene               | 5.0 E-2   | NA   | 3.7 E-4                                    | 0.000005                       | NA                                      |
| Chloroethane                | 1.0 E+1   | 8.3 E-7  | 2.1 E-2                                    | 0.000001                       | 4 E-9                                   |
| Chloroform                  | 4.5 E-2   | 2.3 E-5  | 6.6 E-1                                    | 0.01                           | 4 E-6                                   |
| Chloromethane               | 9.0 E-2   | NA   | 2.1 E-4                                    | 0.000002                       | NA                                      |
| cis-1,2-Dichloroethene      | 3.5 E-2   | NA   | 1.2 E-2                                    | 0.0002                         | NA                                      |
| Dibromochloromethane        | 7.0 E-2   | 2.4 E-5  | 7.9 E-5                                    | 0.0000008                      | 5 E-10                                  |
| Dichlorodifluoromethane     | 2.0 E-1   | NA   | 4.2 E-3                                    | 0.00001                        | NA                                      |
| Ethanol                     | NA  | NA   | 5.6 E-2                                    | NA                             | NA                                      |
| Ethylbenzene                | 1.0 E+0   | NA   | 1.5 E-3                                    | 0.000001                       | NA                                      |
| Hexachlorobutadiene         | NA  | 2.2 E-5  | 2.1 E-3                                    | NA                             | 1 E-8                                   |
| Isopropylbenzene            | 4.0 E-1   | NA   | 3.8 E-4                                    | 0.0000006                      | NA                                      |
| m,p-Xylene                  | 1.0 E-1   | NA   | 5.7 E-3                                    | 0.00004                        | NA                                      |
| Methyl methacrylate         | 7.0 E-1   | NA   | 9.1 E-4                                    | 0.0000009                      | NA                                      |
| Methyl tert butyl ether     | 3.0 E+0   | NA   | 9.4 E-3                                    | 0.000002                       | NA                                      |
| Methylene chloride          | NA  | 4.7 E-7  | 4.9 E-3                                    | NA                             | 6 E-10                                  |
| Naphthalene                 | 3.0 E-3   | NA   | 3.9 E-3                                    | 0.0009                         | NA                                      |
| N-Butylbenzene              | 1.4 E-1   | NA   | 8.1 E-4                                    | 0.000004                       | NA                                      |
| n-Heptane                   | NA  | NA   | 1.7 E-3                                    | NA                             | NA                                      |
| n-Octane                    | NA  | NA   | 1.9 E-3                                    | NA                             | NA                                      |
| N-Propylbenzene             | 1.4 E-1   | NA   | 7.7 E-4                                    | 0.000004                       | NA                                      |
| o-Xylene                    | 1.0 E-1   | NA   | 2.6 E-3                                    | 0.00002                        | NA                                      |

**TABLE 4**  
**SCREENING-LEVEL INDOOR AIR HEALTH RISK ASSESSMENT RESULTS**  
**TRONOX PARCELS A/B SOIL GAS INVESTIGATION**  
**CLARK COUNTY, NEVADA**  
 (Page 2 of 2)

| <b>Chemical</b>        | <b>Non-Cancer Reference Concentration (mg/kg-d)</b> | <b>Unit Risk Factor (mg/kg-d)<sup>-1</sup></b> | <b>J&amp;E Predicted Conc.<sup>a</sup></b> | <b>Non-Cancer Hazard Index</b> | <b>Incremental Lifetime Cancer Risk</b> |
|------------------------|---|--|--|--------------------------------|---|
| sec-Butylbenzene       | 1.4 E-1   | NA   | 1.8 E-4                                    | 0.0000009                      | NA                                      |
| Styrene                | 1.0 E+0   | NA   | 9.4 E-4                                    | 0.0000006                      | NA                                      |
| t-Butyl alcohol        | NA  | NA   | 1.3 E-3                                    | NA                             | NA                                      |
| tert-Butylbenzene      | 1.4 E-1   | NA   | 2.6 E-4                                    | 0.000001                       | NA                                      |
| Tetrachloroethene      | 6.0 E-1   | 5.9 E-6  | 2.9 E-2                                    | 0.00003                        | 4 E-8                                   |
| Toluene                | 5.0 E+0   | NA   | 3.2 E-2                                    | 0.000004                       | NA                                      |
| Trichloroethene        | 4.0 E-2   | 1.1 E-4  | 4.3 E-2                                    | 0.000004                       | NA                                      |
| Trichlorofluoromethane | 7.0 E-1   | NA   | 2.9 E-3                                    | 0.000003                       | NA                                      |
| Vinyl acetate          | 2.0 E-1   | NA   | 9.7 E-3                                    | 0.00003                        | NA                                      |
| Vinyl chloride         | 1.0 E-1   | 4.4 E-6  | 2.6 E-4                                    | 0.000002                       | 3 E-10                                  |
| <b>Total</b>           |   |  |  | 0.01                           | 4 E-6                                   |

<sup>a</sup>From Table 3; concentration is in  $\mu\text{g}/\text{m}^3$ .

NA - Toxicity criteria has not been established.

ATTACHMENT A

SCREENING-LEVEL INDOOR AIR HEALTH RISK ASSESSMENT  
CALCULATION SPREADSHEETS (ON CD)