Enbridge Energy, Limited Partnership 1601 Pratt Avenue Marshall, Michigan49068 ENBRIDGE

October 13, 2011

Mr. Ralph Dollhopf Federal On-Scene Coordinator and Incident Commander U.S. Environmental Protection Agency 801 Garfield Avenue, #229 Traverse City, MI49686

Re: In the Matter of Enbridge Energy Partners, L.P., et al,

Docket No. CWA 1321-5-10-001

## Dear Mr. Dollhopf:

Enbridge Energy, Limited Partnership (Enbridge) is requesting the United States Environmental Protection Agency (U.S. EPA) approve the following modification to the *2011 Air Monitoring and Sampling Addendum* (dated June 21, 2011) of the *Sampling and Analysis Plan* (SAP) (dated August 2010). The proposed modifications are being requested for immediate implementation.

Enbridge is requesting an indefinite suspension of all Community Air Sampling and Monitoring activities, Work Area Monitoring, and elimination of River Opening Sampling effective immediately. The air sampling and monitoring program was designed to protect worker safety and public health during assessment and recovery operations. Work activities during these air sampling and monitoring periods included overbank surface soil oil recovery, overbank excavation oil recovery, and submerged oil recovery during a wide range of atmospheric conditions over the 2011 work periods. This request excludes Odor Response sampling activities which will continue with expedited shipping and 24-hour analytical turnaround times for samples collected.

This request for modification is supported by 2,059 Work Area and Community Air samples collected during 2010 and 1,266 Work Area and Community Air samples collected in 2011, as of September 16, 2011. In addition to the lab analyzed samples, we have collected 66,922 real-time community readings in 2010 and 18,662 in 2011, as of September 16, 2011for VOCs, H2S, SO2, and benzene (all non-detect above screening limits).

Between June 3, 2011 and September 16, 2011, a total of 10 Target Analytes have been detected above Human Health Air Screening Levels (HHASLs) in 2011 along with 22 detections of Non-Target Analytes above HHASLs, which are not related to crude oil released from Line 6B MP 608. Detections of both Target and Non-Target Analytes have been inconsistent and do not alter our conclusions regarding air sampling and monitoring results. In addition, efforts have been taken to evaluate potential lab related issues for select contaminants including isopropyl alcohol and acetone, which revealed these contaminants along with carbonyl sulfide in the laboratory carrier gas. It is important to note that these are 8-, 12-, and 24-hour samples (acute exposure) being compared to chronic exposure based screening levels (greater than one year). Therefore we also compared measures of central tendency (including median and mean values) of contaminant levels over the sampling duration in 2011, which indicated values well below the applicable HHASLs.

The attached tables and figures provide details regarding chemicals detected and comparisons with applicable HHASLs. Please refer to the attached **Table 1**: Cumulative Air Sample Summary – Target Analytes and **Table 2**: Cumulative Air Sample Summary – Non-Target

<u>Analytes</u> for a summary of analytical results from June 3, 2011 to September 16, 2011 for Target Analytes and Non-Target Analytes, respectively. In addition, see attached figures for visual depictions of sampling results compared against applicable HHASLs for each of the detected Target Analytes and specific Non-Target Analytes of interest.

Enbridge is also requesting the elimination of River Opening Sampling and Monitoring, as included in *Memorandum -RE: Modification of the 2011 Air Monitoring and Sampling Addendum to the Sampling and Analysis Plan*, dated July 6, 2011, due to the lack of applicability based on the extensive Work Area and Community Air Sampling and Monitoring conducted in 2010 and 2011 with minimal detections of Target Analytes as outlined in the above text and associated tables and figures.

Assessment of Community Air Sampling and Monitoring results lead us to conclude that our analytical data set is well established, concentrations of Target Analytes in ambient air throughout the community are stable and below levels of concern, and that work activities do not negatively influence ambient air concentrations. Additionally, as ambient air temperatures continue to decrease for the remainder of 2011, weather conditions substantiate a lesser potential for volatilization of chemicals in comparison to the higher ambient air temperatures observed during the sample collection conducted in 2011 to date. Therefore, Enbridge believes it has been demonstrated that continued air sampling and monitoring is unnecessary.

Enbridge welcomes the opportunity to present additional information regarding our sampling and monitoring strategy, a review of data collected, and a detailed assessment of results. If you have any questions regarding this request, please do not hesitate to contact me.

Sincerely,

ENBRIDGE ENERGY, LIMITED PARTNERSHIP By Enbridge Pipelines (Lakehead) L.L.C. Its General Partner

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| Target Analytes - Predominant                                 |             | Human Health Air          |                   | P    | All Sam | ple Duratio        | ons†          |     | 8-ho | ur Target Pe       | eriod         | 12-hour Target Period |      |                    |               |     | 24-ho | ur Target P        | eriod         |
|---------------------------------------------------------------|-------------|---------------------------|-------------------|------|---------|--------------------|---------------|-----|------|--------------------|---------------|-----------------------|------|--------------------|---------------|-----|-------|--------------------|---------------|
| Crude Oil VOCs Detected by TO-15<br>Analysis (including TICs) | CAS NO      | Screening Level<br>(ppbv) | Source            | N    | Ndet    | Average<br>Level** | Max<br>(ppbv) | N   | Ndet | Average<br>Level** | Max<br>(ppbv) | N                     | Ndet | Average<br>Level** | Max<br>(ppbv) | N   | Ndet  | Average<br>Level** | Max<br>(ppbv) |
| 1,2,4-Trimethylbenzene                                        | 95-63-6     | 1.5                       | EPA RfC(1)        | 1266 | 1       | 1.0                | 3.0           | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 1     | 1.0                | 3.0           |
| Benzene                                                       | 71-43-2     | 3                         | ATSDR Chr. MRL(1) | 1266 | 4       | 1.0                | 4.0           | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 4     | 1.0                | 4.0           |
| 2-Methylbutane*                                               | 78-78-4     | 6000                      | MDEQ(1)           | 1266 | 43      | 10.4               | 83.0          | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 43    | 10.4               | 83.0          |
| Cyclohexane                                                   | 110-82-7    | 2000                      | EPA RfC(1)        | 1266 | 25      | 1.0                | 63.0          | 364 | 1    | 5.0                | 63.0          | 117                   | 0    | ND                 | ND            | 785 | 24    | 1.0                | 20.0          |
| 1,3-Dimethylcyclohexane*                                      | 591-21-9    | NA                        | NA                | 1266 | 0       | ND                 | ND            | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 0     | ND                 | ND            |
| Methylcyclopentane*                                           | 96-37-7     | 200                       | MDEQ(1)           | 1266 | 2       | 2.6                | 3.7           | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 2     | 2.6                | 3.7           |
| cis-1,3-Dimethylcyclohexane*                                  | 638-04-0    | NA                        | NA                | 1266 | 0       | ND                 | ND            | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 0     | ND                 | ND            |
| Butylcyclohexane*                                             | 1678-93-9   | NA                        | NA                | 1266 | 0       | ND                 | ND            | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 0     | ND                 | ND            |
| Ethylcyclohexane*                                             | 1678-91-7   | NA                        | NA                | 1266 | 0       | ND                 | ND            | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 0     | ND                 | ND            |
| Methylcyclohexane*                                            | 108-87-2    | 4000                      | MDEQ(1)           | 1266 | 2       | 3.9                | 4.6           | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 2     | 3.9                | 4.6           |
| Propylcyclohexane*                                            | 1678-92-8   | NA                        | NA                | 1266 | 0       | ND                 | ND            | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 0     | ND                 | ND            |
| 1,1-Dimethylcyclopentane*                                     | 1638-26-2   | NA                        | NA                | 1266 | 0       | ND                 | ND            | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 0     | ND                 | ND            |
| trans-1,3-Dimethylcyclopentane*                               | 1759-58-6   | NA                        | NA                | 1266 | 0       | ND                 | ND            | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 0     | ND                 | ND            |
| trans-1,2-Dimethylcyclopentane*                               | 822-50-4    | 0.025                     | MDEQ(1)           | 1266 | 0       | ND                 | ND            | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 0     | ND                 | ND            |
| 3-Methylheptane*                                              | 589-81-1    | NA                        | NA                | 1266 | 0       | ND                 | ND            | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 0     | ND                 | ND            |
| 2,6-Dimethyl-2-Octene*                                        | 4057-42-5   | NA                        | NA                | 1266 | 0       | ND                 | ND            | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 0     | ND                 | ND            |
| Decane*                                                       | 124-18-5    | NA                        | NA                | 1266 | 3       | 4.7                | 8.9           | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 3     | 4.7                | 8.9           |
| Dodecane*                                                     | 112-40-3    | NA                        | NA                | 1266 | 4       | 2.5                | 5.6           | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 4     | 2.5                | 5.6           |
| Ethylbenzene                                                  | 100-41-4    | 60                        | ATSDR Chr. MRL(1) | 1266 | 3       | 1.0                | 3.0           | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 3     | 1.0                | 3.0           |
| 4-Ethyltoluene                                                | 622-96-8    | 70                        | MDEQ(1)           | 1266 | 1       | 1.0                | 1.0           | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 1     | 1.0                | 1.0           |
| n-Heptane                                                     | 142-82-5    | 850                       | MDEQ(1)           | 1266 | 13      | 1.0                | 14.0          | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 13    | 1.0                | 14.0          |
| n-Hexane                                                      | 110-54-3    | 200                       | EPA RfC(1)        | 1266 | 18      | 1.0                | 3.0           | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 18    | 1.0                | 3.0           |
| Naphthalene*                                                  | 91-20-3     | 1                         | ATSDR Chr. MRL(1) | 1266 | 0       | ND                 | ND            | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 0     | ND                 | ND            |
| Nonane*                                                       | 111-84-2    | 40                        | EPA RfC(1)        | 1266 | 1       | 18.0               | 18.0          | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 1     | 18.0               | 18.0          |
| 2-Methylhexane*                                               | 591-76-4    | 850                       | MDEQ(1)           | 1266 | 4       | 18.3               | 35.0          | 364 | 1    | 35.0               | 35.0          | 117                   | 0    | ND                 | ND            | 785 | 3     | 12.7               | 29.0          |
| Octane*                                                       | 111-65-9    | NA                        | NA                | 1266 | 0       | ND                 | ND            | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 0     | ND                 | ND            |
| 3-Methyloctane*                                               | 2216-33-3   | NA                        | NA                | 1266 | 0       | ND                 | ND            | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 0     | ND                 | ND            |
| 2-Methylpentane*                                              | 107-83-5    | 5000                      | MDEQ(1)           | 1266 | 1       | 1.7                | 1.7           | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 1     | 1.7                | 1.7           |
| Toluene                                                       | 108-88-3    | 80                        | ATSDR Chr. MRL(1) | 1266 | 256     | 4.5                | 270.0         | 364 | 3    | 5.0                | 54.0          | 117                   | 1    | 5.0                | 17.0          | 785 | 252   | 3.9                | 270.0         |
| 3-Methylhexane*                                               | 589-34-4    | 850                       | MDEQ(1)           | 1266 | 4       | 16.5               | 48.0          | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 4     | 16.5               | 48.0          |
| 2-Methylheptane*                                              | 592-27-8    | NA                        | NA                | 1266 | 0       | ND                 | ND            | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 0     | ND                 | ND            |
| 1,1,3- Trimethylcyclohexane*                                  | 3073-66-3   | NA                        | NA                | 1266 | 0       | ND                 | ND            | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 0     | ND                 | ND            |
| 1,3,5-Trimethylbenzene                                        | 108-67-8    | 45                        | MDEQ(1)           | 1266 | 0       | ND                 | ND            | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 0     | ND                 | ND            |
| Undecane*                                                     | 1120-21-4   | NA                        | NA                | 1266 | 1       | 1.5                | 1.5           | 364 | 0    | ND                 | ND            | 117                   | 0    | ND                 | ND            | 785 | 1     | 1.5                | 1.5           |
| m&p-Xylene                                                    | 179601-23-1 | 50                        | EPA RfC(1)        | 1266 | 6       | 2.0                | 10.0          | 364 | 0    | ND                 | ND            | 117                   |      | ND                 | ND            | 785 | 6     | 2.0                | 10.0          |
| o-Xylene                                                      | 95-47-6     | 50                        | EPA RfC(1)        | 1266 | 3       | 1.0                | 3.0           | 364 | 0    | ND                 | ND            | 117                   |      | ND                 | ND            | 785 | 3     | 1.0                | 3.0           |

<sup>\* -</sup> TIC (tentatively identified compound)

<sup>\*\* -</sup> Average Levels for compounds that were detected in <5% of samples are the median value using the detection limit for non-detects. Average Levels for compounds that were detected in >5% of samples are the mean using the detection limit for non-detects. Average Levels for tentatively identified compounds are the mean of detects.

<sup>†</sup>Samples reported do not include instantaneous 'grab' samples. Samples with 8-24 hour target sample periods are included.

NA - Not Available/ND - Not Detected

 $<sup>\</sup>ensuremath{\text{\textbf{N}}}$  - Number of samples analyzed/  $\ensuremath{\text{\textbf{NDET}}}$  - Number of detections

<sup>(1) -</sup> Unless otherwise noted, the screening level was obtained from Enbridge Oil Spill Human Health Air Screening Levels, August 31, 2011.

ATSDR Chr. MRL - Agency for Toxic Substances and Disease Registry - Chronic Minimal Risk Level for Inhalation

EPA RfC - If no Chronic MRL is available, the screening level is the EPA Chronic Reference Concentration (RfC).

EPA RSL - If none of the above are available, the EPA Regional Screening Level (RSL) is used.

MDEQ - If none of the above are available, the Michigan DEQ Air Quality Division, Air Toxics Screening Level is the screening level.

|                                                                   |                        | Human Health Air |                    |      | All San | nple Duratio | nst         |     | 8-ho | ır Target Pe | riod     |     | 12-ho | ur Target P | eriod    |     | 24-ho | ur Target Pe | eriod      |
|-------------------------------------------------------------------|------------------------|------------------|--------------------|------|---------|--------------|-------------|-----|------|--------------|----------|-----|-------|-------------|----------|-----|-------|--------------|------------|
| Non-Target Analytes (including TICs)                              | CAS NO                 | Screening Level  | Source             |      |         | Average      | Max         |     |      | Average      | Max      |     |       | Average     | Max      |     |       | Average      | Max        |
| 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3                           |                        | (ppbv)           |                    | N    | Ndet    | Level**      | (ppbv)      | N   | Ndet | Level**      | (ppbv)   | N   | Ndet  | Level**     | (ppbv)   | N   | Ndet  | Level**      | (ppbv)     |
| 1,1,1-trichloroethane                                             | 71-55-6                | 1000             | EPA RfC(1)         | 1266 | 0       | ND           | 0.0         | 364 | 0    | ND           | ND       | 117 | 0     | ND          | ND       | 785 | 0     | ND           | ND         |
| 1,1,2,2-Tetrachloroethane                                         | 79-34-5                | 0.006            | EPA RSL(1)         | 1266 | 0       | ND           | 0.0         | 364 | 0    | ND           | ND       | 117 | 0     | ND          | ND       | 785 | 0     | ND           | ND         |
| 1,1,2-Trichloro-1,2,2-trifluoroethane                             | 76-13-1                | 4000             | EPA RSL(1)         | 1266 | 0       | ND           | 0.0         | 364 | 0    | ND           | ND       | 117 | 0     | ND          | ND       | 785 | 0     | ND           | ND         |
| 1,1,2-Trichloroethane                                             | 79-00-5                | 0.04             | EPA RfC(1)         | 1266 | 0       | ND           | 0.0         | 364 | 0    | ND           | ND       | 117 | 0     | ND          | ND       | 785 | 0     | ND           | ND         |
| 1,1-Dichloroethane                                                | 75-34-3                | 0.4              | EPA RSL(1)         | 1266 | 0       | ND           | 0.0         | 364 | 0    | ND           | ND       | 117 | 0     | ND          | ND       | 785 | 0     | ND           | ND         |
| 1,1-Dichloroethene                                                | 75-35-4                | 50               | EPA RfC(1)         | 1266 | 0       | ND           | 0.0         | 364 | 0    | ND           | ND       | 117 | 0     | ND          | ND       | 785 | 0     | ND           | ND         |
| 1,2,4-Trichlorobenzene                                            | 120-82-1               | 0.3              | EPA RfC(1)         | 1266 | 0       | ND           | 0.0         | 364 | 0    | ND           | ND       | 117 | 0     | ND          | ND       | 785 | 0     | ND           | ND         |
| 1,2-Dibromoethane                                                 | 106-93-4               | 1                | EPA RfC(1)         | 1266 | 0       | ND           | 0.0         | 364 | 0    | ND           | ND       | 117 | 0     | ND          | ND       | 785 | 0     | ND           | ND         |
| 1,2-Dichloro-1,1,2,2-Tetrafluoroethane                            | 76-14-2                | 9900             | MDEQ(1)            | 1266 | 0       | ND           | 0.0         | 364 | 0    | ND           | ND       | 117 | 0     | ND          | ND       | 785 | 0     | ND           | ND         |
| 1,2-Dichlorobenzene                                               | 95-50-1                | 30               | EPA RfC(1)         | 1266 | 0       | ND           | 0.0         | 364 | 0    | ND           | ND       | 117 | 0     | ND          | ND       | 785 | 0     | ND           | ND         |
| 1,2-Dichloroethane                                                | 107-06-2               | 600              | ATSDR Chr. MRL(1)  | 1266 | 2       | 1            | 3.0         | 364 | 0    | ND           | ND       | 117 | 0     | ND          | ND       | 785 | 2     | 1.0          | 3.0        |
| 1,2-Dichloropropane                                               | 78-87-5                | 1                | EPA RfC(1)         | 1266 | 0       | ND           | 0.0         | 364 | 0    | ND           | ND       | 117 | 0     | ND          | ND       | 785 | 0     | ND           | ND         |
| 1,2-PENTADIENE*                                                   | 591-95-7               | NA               | NA                 | 1266 | 1       | 1.6          | 1.6         | 364 | 0    | ND           | ND       | 117 | 0     | ND          | ND       | 785 | 1     | 1.6          | 1.6        |
| 1,3-Butadiene                                                     | 106-99-0               | 1                | EPA RfC(1)         | 1266 | 0       | ND           | 0.0         | 364 | 0    | ND           | ND       | 117 | 0     | ND          | ND       | 785 | 0     | ND           | ND         |
| 1.3-BUTADIENE. 2-METHYL-*                                         | 78-79-5                | NA               | NA                 | 1266 | 4       | 2.7          | 6.4         | 364 | 1    | 6.4          | 6.4      | 117 | 0     | ND          | ND       | 785 | 3     | 1.5          | 1.6        |
| 1,3-Dichlorobenzene                                               | 541-73-1               | 0.5              | MDEQ(1)            | 1266 |         | ND           | 0.0         | 364 | 0    | ND           | ND       | 117 | 0     | ND          | ND       | 785 | 0     | ND           | ND         |
| 1.3-PENTADIENE*                                                   | 504-60-9               | NA               | NA                 | 1266 | 2       | 8.4          | 8.9         | 364 | 0    | ND           | ND       | 117 | 0     | ND          | ND       | 785 | 2     | 8.4          | 8.9        |
| 1,3-PENTADIENE, (E)-*                                             | 2004-70-8              | NA               | NA                 | 1266 |         | 1.8          | 1.9         | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 2     | 1.8          | 1.9        |
| 1,4-Dichlorobenzene                                               | 106-46-7               | 10               | ATSDR Chr. MRL(1)  | 1266 |         | ND           | 0.0         | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 0     | ND           | ND         |
| 1,4-Dioxane                                                       | 123-91-1               | 1000             | ATSDR Chr. MRL(1)  | 1266 |         | ND           | 0.0         | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 0     | ND           | ND         |
| 1-DECENE*                                                         | 872-05-9               | NA               | NA NA              | 1266 |         | ND           | 0.0         | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 0     | ND           | ND         |
| 1-HEPTANAL, 3,5,5-TRIETHYL-*                                      | 1000160-77-0           | NA               | NA                 | 1266 |         | 14.6         | 29.0        | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 4     | 14.6         | 29.0       |
| 1-HEPTANOL, 2-PROPYL-*                                            | 10042-59-8             | NA               | NA                 | 1266 |         | 1.5          | 1.5         | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 1     | 1.5          | 1.5        |
| 1-HEXENE. 4-METHYL-*                                              | 3769-23-1              | NA               | NA                 | 1266 |         | 5.6          | 5.6         | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 1     | 5.6          | 5.6        |
| 1-METHYL-2-METHYLENECYCLOHEXANE*                                  | 2808-75-5              | NA               | NA NA              | 1266 |         | 1.2          | 1.2         | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 1     | 1.2          | 1.2        |
| 1-METHYLDECAHYDRONAPHTHALENE*                                     | 2958-75-0              | NA               | NA                 | 1266 |         | 10.9         | 26.0        | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 7     | 10.9         | 26.0       |
| 1-TRIDECENE*                                                      | 2437-56-1              | NA               | NA                 | 1266 |         | 1.3          | 1.3         | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 1     | 1.3          | 1.3        |
| 2(1H)-NAPHTHALENONE, OCTAHYDRO-4*                                 | 938-06-7               | NA               | NA                 | 1266 |         | 29           | 29.0        | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 1     | 29.0         | 29.0       |
| 2,2,4-Trimethylpentane                                            | 540-84-1               | 750              | MDEQ(1)            | 1266 |         | 1            | 1.0         | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 1     | 1.0          | 1.0        |
| 2,2,7,7-TETRAMETHYLOCTANE*                                        | 1071-31-4              | NA               | NA<br>NA           | 1266 |         | 1.2          | 1.2         | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 1     | 1.2          | 1.2        |
| 2.3.4.5-TETRAHYDROPYRIDAZINE*                                     | 694-06-4               | NA<br>NA         | NA<br>NA           | 1266 |         | 2.2          | 2.2         | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 1     | 2.2          | 2.2        |
| 2-BUTENAL*                                                        | 4170-30-3              | 3                | MDEQ(1)            | 1266 |         | 1.7          | 1.7         | 364 | 0    | ND           | ND       | 117 | -     | ND          | ND       | 785 | 1     | 1.7          | 1.7        |
| 2-BUTENAL, (E)-*                                                  | 123-73-9               | NA NA            | NA NA              | 1266 |         | 1.9          | 2.7         | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 4     | 1.9          | 2.7        |
| 2-BUTENOIC ACID, METHYL ESTER, (Z)-*                              | 4358-59-2              | NA               | NA                 | 1266 |         | 8            | 8.0         | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 1     | 8.0          | 8.0        |
| 2-Chloro-1.3-butadiene                                            | 126-99-8               | 6                | EPA RfC(1)         | 1266 |         | ND           | 0.0         | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 0     | ND           | ND         |
| 2-DECENAL, (E)-*                                                  | 3913-81-3              | NA               | NA                 | 1266 |         | ND           | 0.0         | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 0     | ND           | ND         |
| 2-HEPTANONE*                                                      | 110-43-0               | 500              | MDEQ(1)            | 1266 |         | 11           | 11.0        | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 1     | 11.0         | 11.0       |
| 2-PENTANONE*                                                      | 107-87-9               | 1500             | ( )                | 1266 |         | 23.5         | 29.0        | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 2     | 23.5         | 29.0       |
| 2-PENTANONE 2-PROPANOL, 2-METHYL-*                                | 75-65-0                | 600              | MDEQ(1)<br>MDEQ(1) | 1266 |         | 47.6         | 120.0       | 364 | 0    | ND           | ND       | 117 |       | 120.0       | 120.0    | 785 | 4     | 11.4         | 18.0       |
| ,                                                                 | 55638-50-1             | NA               | NA<br>NA           | 1266 |         | 6.4          | 6.4         | 364 | 0    | ND           | ND       | 117 |       | 120.0<br>ND | ND       | 785 | 1     | 6.4          | 6.4        |
| 3,5-DECADIENE, 2,2-DIMETHYL-, (Z*<br>3,5-OCTADIENE, 4,5-DIETHYL-* |                        | NA<br>NA         | NA<br>NA           | 1266 |         | 10           |             | 364 | 0    | ND           | ND       | 117 | -     | ND          | ND       | 785 | 1     | 10.0         |            |
|                                                                   | 67652-84-0<br>107-05-1 | 32               |                    | 1266 |         | ND           | 10.0<br>0.0 | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 0     | ND           | 10.0<br>ND |
| 3-Chloropropene                                                   |                        | -                | EPA RfC(1)         |      |         |              |             |     | 0    | ND<br>ND     | ND<br>ND | 117 |       |             | ND<br>ND |     | 1     |              |            |
| 3-HEPTANONE*                                                      | 106-35-4               | NA               | NA                 | 1266 |         | 2.4          | 2.4         | 364 | 0    |              |          |     |       | ND          |          | 785 | •     | 2.4          | 2.4        |
| 4-DECENE, 2,2-DIMETHYL-, (E)-*                                    | 55534-69-5             | NA               | NA                 | 1266 |         | ND           | 0.0         | 364 |      | ND           | ND       | 117 |       | ND          | ND       | 785 | 0     | ND<br>2.0    | ND         |
| 4-NONENE, 5-BUTYL-*                                               | 7367-38-6              | NA               | NA                 | 1266 |         | 3.9          | 3.9         | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 1     | 3.9          | 3.9        |
| 4-TRIFLUOROACETOXYTRIDECANE*                                      | 1000245-47-3           | NA               | NA                 | 1266 |         | ND           | 0.0         | 364 | 0    | ND           | ND       | 117 |       | ND          | ND       | 785 | 0     | ND           | ND         |
| 5-OCTEN-2-YN-4-OL*                                                | 1000196-87-1           | NA               | NA                 | 1266 | 1       | 11           | 11.0        | 364 | 0    | ND           | ND       | 117 | 0     | ND          | ND       | 785 | 1     | 11.0         | 11.0       |

|                                        |              | Human Health Air |                   |      | All San | nple Duratio | ons†   |     | 8-hou | ır Target Pe | riod   |          | 12-ho | ur Target P | eriod  |     | 24-ho | ur Target P | eriod  |
|----------------------------------------|--------------|------------------|-------------------|------|---------|--------------|--------|-----|-------|--------------|--------|----------|-------|-------------|--------|-----|-------|-------------|--------|
| Non-Target Analytes (including TICs)   | CAS NO       | Screening Level  | Source            |      |         | Average      | Max    |     |       | Average      | Max    | <u> </u> |       | Average     | Max    |     |       | Average     | Max    |
|                                        |              | (ppbv)           |                   | N    | Ndet    | Level**      | (ppbv) | N   | Ndet  | Level**      | (ppbv) | N        | Ndet  | Level**     | (ppbv) | N   | Ndet  | Level**     | (ppbv) |
| 7-METHYLBICYCLO[4.2.0]OCTANE*          | 1000210-90-2 | NA               | NA                | 1266 | 1       | 4.5          | 4.5    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 1     | 4.5         | 4.5    |
| 7-NORBORNYL T-BUTYL ETHER*             | 3391-07-9    | NA               | NA                | 1266 | 1       | 14           | 14.0   | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 1     | 14.0        | 14.0   |
| ACETALDEHYDE*                          | 75-07-0      | 5                | EPA RfC(1)        | 1266 | 26      | 6.9          | 15.0   | 364 | 5     | 7.7          | 10.0   | 117      | 3     | 8.2         | 11.0   | 785 | 18    | 6.5         | 15.0   |
| ACETIC ACID, 1,1-DIMETHYLETHYL ESTER*  | 540-88-5     | 2000             | MDEQ(1)           | 1266 | 1       | 2.2          | 2.2    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 1     | 2.2         | 2.2    |
| ACETIC ACID, 2-ETHYLHEXYL ESTER*       | 103-09-3     | 2                | MDEQ(1)           | 1266 | 1       | 6.2          | 6.2    | 364 | 0     | ND           | ND     | 117      | 1     | 6.2         | 6.2    | 785 | 0     | ND          | ND     |
| ACETOACETIC ACID, 1-THIO-, S-ALLYL*    | 15780-65-1   | NA               | NA                | 1266 | 3       | 1.6          | 1.8    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 3     | 1.6         | 1.8    |
| Acetone                                | 67-64-1      | 13000            | EPA RSL(1)        | 1266 | 1096    | 8.1          | 300.0  | 364 | 339   | 9.9          | 120.0  | 117      | 116   | 17.3        | 300.0  | 785 | 641   | 5.9         | 89.0   |
| ACETONITRILE*                          | 75-05-8      | 36               | EPA RfC(1)        | 1266 | 62      | 7.7          | 47.0   | 364 | 5     | 8.3          | 12.0   | 117      | 2     | 5.9         | 6.4    | 785 | 55    | 7.7         | 47.0   |
| Acrylonitrile                          | 107-13-1     | 1                | EPA RfC(1)        | 1266 | 0       | ND           | 0.0    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 0     | ND          | ND     |
| ALPHA-PINENE*                          | 80-56-8      | 200              | MDEQ(1)           | 1266 | 2       | 7.1          | 13.0   | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 2     | 7.1         | 13.0   |
| BENZENE, 1,3,5-TRICHLORO-*             | 108-70-3     | NA               | NA                | 1266 | 0       | ND           | 0.0    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 0     | ND          | ND     |
| BENZENE, 1,3-DIETHYL-*                 | 141-93-5     | NA               | NA                | 1266 | 1       | 1.9          | 1.9    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 1     | 1.9         | 1.9    |
| BENZENE, 1-METHYL-4-(1-METHYLETHYL)*   | 99-87-6      | 1.8              | MDEQ(1)           | 1266 | 1       | 2.7          | 2.7    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 1     | 2.7         | 2.7    |
| Benzyl chloride                        | 100-44-7     | 0.2              | EPA RfC(1)        | 1266 | 0       | ND           | 0.0    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 0     | ND          | ND     |
| BETA-PINENE*                           | 127-91-3     | 200              | MDEQ(1)           | 1266 | 0       | ND           | 0.0    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 0     | ND          | ND     |
| Bromodichloromethane                   | 75-27-4      | 0.01             | EPA RSL(1)        | 1266 | 0       | ND           | 0.0    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 0     | ND          | ND     |
| Bromoform                              | 75-25-2      | 0.2              | EPA RSL(1)        | 1266 | 0       | ND           | 0.0    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 0     | ND          | ND     |
| BROMOMETHANE                           | 74-83-9      | 5                | ATSDR Chr. MRL(1) | 1266 | 10      | 1            | 23.0   | 364 | 1     | 5.0          | 15.0   | 117      | 3     | 5.0         | 23.0   | 785 | 6     | 1.0         | 7.1    |
| BUTANAMIDE, 3,3-DIMETHYL-*             | 926-04-5     | NA               | NA                | 1266 | 1       | 5.6          | 5.6    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 1     | 5.6         | 5.6    |
| BUTANE*                                | 106-97-8     | 10000            | MDEQ(1)           | 1266 | 7       | 4.8          | 10.0   | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 7     | 4.8         | 10.0   |
| BUTANOIC ACID, 2-METHYLPROPYL ESTER*   | 539-90-2     | NA               | NA                | 1266 |         | 1.9          | 1.9    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 1     | 1.9         | 1.9    |
| BUTANOIC ACID, BUTYL ESTER*            | 109-21-7     | NA               | NA                | 1266 | 1       | 14           | 14.0   | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 1     | 14.0        | 14.0   |
| BUTANOIC ACID, ETHYL ESTER*            | 105-54-4     | NA               | NA                | 1266 | 1       | 17           | 17.0   | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 1     | 17.0        | 17.0   |
| BUTANOIC ACID, METHYL ESTER*           | 623-42-7     | NA               | NA                | 1266 | 1       | 4.8          | 4.8    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 1     | 4.8         | 4.8    |
| BUTANOIC ACID, PROPYL ESTER*           | 105-66-8     | NA               | NA                | 1266 | 1       | 16           | 16.0   | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 1     | 16.0        | 16.0   |
| CARBON DISULFIDE                       | 75-15-0      | 200              | EPA RfC(1)        | 1266 | 60      | 2            | 120.0  | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 60    | 3.1         | 120.0  |
| Carbon tetrachloride                   | 56-23-5      | 30               | ATSDR Chr. MRL(1) | 1266 | 0       | ND           | 0.0    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 0     | ND          | ND     |
| CARBONYL SULFIDE*                      | 463-58-1     | 4                | MDEQ(1)           | 1266 | 5       | 2.3          | 5.2    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 5     | 2.3         | 5.2    |
| Chlorobenzene                          | 108-90-7     | 10               | EPA RfC(1)        | 1266 | 0       | ND           | 0.0    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 0     | ND          | ND     |
| Chloroethane                           | 75-00-3      | 4000             | EPA RfC(1)        | 1266 | 0       | ND           | 0.0    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 0     | ND          | ND     |
| Chloroform                             | 67-66-3      | 20               | ATSDR Chr. MRL(1) | 1266 | 0       | ND           | 0.0    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 0     | ND          | ND     |
| Chloromethane                          | 74-87-3      | 50               | ATSDR Chr. MRL(1) | 1266 | 0       | ND           | 0.0    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 0     | ND          | ND     |
| cis-1,2-dichloroethene                 | 156-59-2     | 9                | MDEQ(1)           | 1266 | 0       | ND           | 0.0    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 0     | ND          | ND     |
| cis-1,3-Dichloropropene                | 10061-02-6   | 4                | MDEQ(1)           | 1266 | 0       | ND           | 0.0    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 0     | ND          | ND     |
| Cumene                                 | 98-82-8      | 80               | EPA RfC(1)        | 1266 | 0       | ND           | 0.0    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 0     | ND          | ND     |
| CYCLODECENE, 1-METHYL-*                | 66633-38-3   | NA               | NA                | 1266 | 2       | 4.7          | 5.2    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 2     | 4.7         | 5.2    |
| CYCLODODECANE*                         | 294-62-2     | NA               | NA                | 1266 | 1       | 1.7          | 1.7    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 1     | 1.7         | 1.7    |
| CYCLODODECENE*                         | 1501-82-2    | NA               | NA                | 1266 | 1       | 8.6          | 8.6    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 1     | 8.6         | 8.6    |
| CYCLODODECENE, (E)-*                   | 1486-75-5    | NA               | NA                | 1266 | 1       | 7.7          | 7.7    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 1     | 7.7         | 7.7    |
| CYCLOHEXANE, 1,2,3-TRIMETHYL-*         | 7667-55-2    | NA               | NA                | 1266 | 1       | 3.8          | 3.8    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 1     | 3.8         | 3.8    |
| CYCLOHEXANE, 2-BUTYL-1,1,3-TRIMETHYL-* | 54676-39-0   | NA               | NA                | 1266 |         | 22.8         | 48.0   | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 4     | 22.8        | 48.0   |
| CYCLOHEXANOL, 3-(3,3-DIMETHYLBUTYL)-*  | 40564-98-5   | NA               | NA                | 1266 |         | 10.2         | 15.0   | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 3     | 10.2        | 15.0   |
| CYCLOHEXANONE, 2,3-DIMETHYL-*          | 13395-76-1   | NA               | NA                | 1266 | 1       | 26           | 26.0   | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 1     | 26.0        | 26.0   |
| CYCLOPENTANE*                          | 287-92-3     | 6000             | MDEQ(1)           | 1266 |         | 3.4          | 5.1    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 2     | 3.4         | 5.1    |
| CYCLOPENTANE, 1,1'-ETHYLIDENEBIS-*     | 4413-21-2    | NA               | NA                | 1266 |         | 9.7          | 9.7    | 364 | 0     | ND           | ND     | 117      |       | ND          | ND     | 785 | 1     | 9.7         | 9.7    |
| CYCLOPENTANE, 1-HYDROXYMETHYL-*        | 1000156-73-8 | NA               | NA                | 1266 |         | 6.1          | 6.1    | 364 | 0     | ND           | ND     | 117      | 0     | ND          | ND     | 785 | 1     | 6.1         | 6.1    |
| CYCLOPENTANE, 1-PENTYL-2-PROPYL-*      | 62199-51-3   | NA               | NA                | 1266 |         | 1.1          | 1.1    | 364 | 0     | ND           | ND     | 117      |       | ND          | ND     | 785 | 1     | 1.1         | 1.1    |

|                                                               |              | Human Health Air |            |      | All San  | nple Duration | nst    | I   | 8-hou  | ır Target Pe | riod       |     | 12-ho  | ur Target P | erind  | 1        | 24-ho   | ur Target P | eriod  |
|---------------------------------------------------------------|--------------|------------------|------------|------|----------|---------------|--------|-----|--------|--------------|------------|-----|--------|-------------|--------|----------|---------|-------------|--------|
| Non-Target Analytes (including TICs)                          | CAS NO       | Screening Level  | Source     |      | All Sall | Average       | Max    |     | 6-1100 | Average      | Max        |     | 12-110 | Average     | Max    | <u> </u> | 24-1100 | Average     | Max    |
| ger/, (                                                       |              | (ppbv)           | 554.55     | N    | Ndet     | Level**       | (ppbv) | N   | Ndet   | Level**      | (ppbv)     | N   | Ndet   | Level**     | (ppbv) | N        | Ndet    | Level**     | (ppbv) |
| CYCLOPENTANE, PROPYL-*                                        | 2040-96-2    | NA               | NA         | 1266 | 1        | 6.1           | 6.1    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 6.1         | 6.1    |
| DECANAL*                                                      | 112-31-2     | NA               | NA         | 1266 | 2        | 7.9           | 8.7    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 2       | 7.9         | 8.7    |
| DECANE, 2,2,4-TRIMETHYL-*                                     | 62237-98-3   | NA               | NA         | 1266 | 1        | 1.1           | 1.1    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 1.1         | 1.1    |
| DECANE, 2,2,6-TRIMETHYL-*                                     | 62237-97-2   | NA               | NA         | 1266 | 1        | 4.3           | 4.3    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 4.3         | 4.3    |
| DECANE, 2,2,8-TRIMETHYL-*                                     | 62238-01-1   | NA               | NA         | 1266 | 1        | 14            | 14.0   | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 14.0        | 14.0   |
| DECANE, 2,2-DIMETHYL-*                                        | 17302-37-3   | NA               | NA         | 1266 | 1        | 1.3           | 1.3    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 1.3         | 1.3    |
| DECANE, 2,3,4-TRIMETHYL-*                                     | 62238-15-7   | NA               | NA         | 1266 | 1        | 1.5           | 1.5    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 1.5         | 1.5    |
| DECANE, 2,3,5-TRIMETHYL-*                                     | 62238-11-3   | NA               | NA         | 1266 | 1        | 6.1           | 6.1    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 6.1         | 6.1    |
| DECANE, 2,6,7-TRIMETHYL-*                                     | 62108-25-2   | NA               | NA         | 1266 | 1        | 6.6           | 6.6    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 6.6         | 6.6    |
| DIACETYL*                                                     | 431-03-8     | NA               | NA         | 1266 | 1        | 1.8           | 1.8    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 1.8         | 1.8    |
| Dibromochloromethane                                          | 124-48-1     | 0.01             | EPA RSL(1) | 1266 |          | ND            | 0.0    | 364 | 0      | ND           | ND         | 117 |        | ND          | ND     | 785      | 0       | ND          | ND     |
| DICHLOROACETIC ACID, 4-TRIDECYL ESTER*                        | 1000280-64-3 | NA               | NA NA      | 1266 |          | ND            | 0.0    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 0       | ND          | ND     |
| Dichlorodifluoromethane                                       | 75-71-8      | 20               | EPA RfC(1) | 1266 |          | ND            | 0.0    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 0       | ND          | ND     |
| DIMETHYL ETHER*                                               | 115-10-6     | 35               | MDEQ(1)    | 1266 |          | ND            | 0.0    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 0       | ND          | ND     |
| DISULFIDE. DIMETHYL*                                          | 624-92-0     | 7.3              | MDEQ(1)    | 1266 |          | 2.4           | 2.4    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 2.4         | 2.4    |
| DISULFIDE, FLUOROMETHYL METHYL*                               | 60307-49-5   | NA               | NA<br>NA   | 1266 |          | 2.4           | 2.0    | 364 | 0      | ND           | ND         | 117 | -      | ND          | ND     | 785      | 1       | 2.0         | 2.0    |
| DODECANE. 2.2.11.11-TETRAMETHYL-*                             | 127204-12-0  | NA<br>NA         | NA         | 1266 |          | 6.6           | 6.6    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 6.6         | 6.6    |
| DODECANE, 2,2,11,11-1ETRAMETITE  DODECANE, 2,6,11-TRIMETHYL-* | 31295-56-4   | NA               | NA         | 1266 |          | 1.6           | 1.6    | 364 | 0      | ND           | ND         | 117 | -      | ND          | ND     | 785      | 1       | 1.6         | 1.6    |
|                                                               | 6044-71-9    | NA<br>NA         | NA<br>NA   | 1266 |          | 1.7           | 1.7    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       |             | 1.7    |
| DODECANE, 6-METHYL-*                                          |              |                  |            |      |          |               |        |     |        |              |            |     |        |             |        |          |         | 1.7         |        |
| ETHANE, 1,1-DIFLUORO-*                                        | 75-37-6      | 14800            | EPA RfC(1) | 1266 |          | 6             | 12.0   | 364 | 0      | ND           | ND<br>70.0 | 117 |        | ND          | ND     | 785      | 9       | 6.0         | 12.0   |
| ETHYL ACETATE                                                 | 141-78-6     | 890              | MDEQ(1)    | 1266 |          | 3             | 76.0   | 364 | 2      | 5.0          | 76.0       | 117 | 0      | ND          | ND     | 785      | 65      | 1.6         | 33.0   |
| ETHYL ALCOHOL*                                                | 64-17-5      | 10000            | MDEQ(1)    | 1266 |          | 68.9          | 220.0  | 364 | 1      | 7.3          | 7.3        | 117 |        | 15.0        | 15.0   | 785      | 27      | 73.2        | 220.0  |
| EUCALYPTOL*                                                   | 470-82-6     | NA               | NA         | 1266 |          | 6.9           | 6.9    | 364 | 0      | ND           | ND         | 117 | 1      | 6.9         | 6.9    | 785      | 0       | ND          | ND     |
| FURAN, 2,3-DIHYDRO-3-METHYL-*                                 | 1708-27-6    | NA               | NA         | 1266 |          | 1.5           | 1.5    | 364 | 0      | ND           | ND         | 117 |        | ND          | ND     | 785      | 1       | 1.5         | 1.5    |
| HENEICOSANE*                                                  | 629-94-7     | NA               | NA         | 1266 |          | ND            | 0.0    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 0       | ND          | ND     |
| HEPTADECANE, 7-METHYL-*                                       | 20959-33-5   | NA               | NA         | 1266 |          | 1.4           | 1.4    | 364 | 0      | ND           | ND         | 117 |        | ND          | ND     | 785      | 1       | 1.4         | 1.4    |
| HEPTANE, 2,2,3,4,6,6-HEXAMETHYL-*                             | 62108-32-1   | NA               | NA         | 1266 |          | 3.9           | 6.5    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 2       | 3.9         | 6.5    |
| HEPTANE, 2,2,4,6,6-PENTAMETHYL-*                              | 13475-82-6   | NA               | NA         | 1266 |          | 3.1           | 3.9    | 364 | 0      | ND           | ND         | 117 |        | ND          | ND     | 785      | 2       | 3.1         | 3.9    |
| HEPTANE, 2,2,4-TRIMETHYL-*                                    | 14720-74-2   | NA               | NA         | 1266 | 1        | 13            | 13.0   | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 13.0        | 13.0   |
| HEPTANE, 2,2-DIMETHYL-*                                       | 1071-26-7    | NA               | NA         | 1266 | 1        | 57            | 57.0   | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 57.0        | 57.0   |
| HEPTANE, 4-ETHYL-2,2,6,6-TETRAMER*                            | 62108-31-0   | NA               | NA         | 1266 | 1        | 6.4           | 6.4    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 6.4         | 6.4    |
| Hexachlorobutadiene                                           | 87-68-3      | 0.01             | EPA RSL(1) | 1266 | 0        | ND            | 0.0    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 0       | ND          | ND     |
| HEXADECANE*                                                   | 544-76-3     | NA               | NA         | 1266 | 1        | 1.5           | 1.5    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 1.5         | 1.5    |
| HEXADECANE, 2,6,10,14-TETRAMETHYL-*                           | 638-36-8     | NA               | NA         | 1266 | 1        | 29            | 29.0   | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 29.0        | 29.0   |
| HEXANAL*                                                      | 66-25-1      | 0.49             | MDEQ(1)    | 1266 | 7        | 3.9           | 9.2    | 364 | 1      | 6.0          | 6.0        | 117 | 0      | ND          | ND     | 785      | 6       | 3.6         | 9.2    |
| HEXANE, 2,2,5,5-TETRAMETHYL-*                                 | 1071-81-4    | NA               | NA         | 1266 | 1        | 1.2           | 1.2    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 1.2         | 1.2    |
| HEXANE, 2,2,5-TRIMETHYL-*                                     | 3522-94-9    | NA               | NA         | 1266 | 2        | 3.5           | 5.2    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 2       | 3.5         | 5.2    |
| HEXANE, 2,3,4-TRIMETHYL-*                                     | 921-47-1     | NA               | NA         | 1266 | 1        | 15            | 15.0   | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 15.0        | 15.0   |
| HEXANE, 2,4-DIMETHYL-*                                        | 589-43-5     | NA               | NA         | 1266 | 3        | 5             | 8.8    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 3       | 5.0         | 8.8    |
| HEXANOIC ACID, METHYL ESTER*                                  | 106-70-7     | NA               | NA         | 1266 | 1        | 2.4           | 2.4    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 2.4         | 2.4    |
| HEXATRIACONTANE*                                              | 630-06-8     | NA               | NA         | 1266 | 1        | 5.2           | 5.2    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 5.2         | 5.2    |
| HEXYL N-VALERATE*                                             | 1117-59-5    | NA               | NA         | 1266 | 1        | 2.2           | 2.2    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 1       | 2.2         | 2.2    |
| ISOBUTANE*                                                    | 75-28-5      | 10000            | MDEQ(1)    | 1266 | 37       | 11.7          | 38.0   | 364 | 10     | 13.2         | 38.0       | 117 | 0      | ND          | ND     | 785      | 27      | 11.1        | 38.0   |
| Isopropyl Alcohol                                             | 67-63-0      | 3000             | EPA RfC(1) | 1266 |          | 11.8          | 530.0  | 364 | 274    | 11.9         | 150.0      | 117 | 113    | 22.8        | 130.0  | 785      | 711     | 10.1        | 530.0  |
| LIMONENE*                                                     | 138-86-3     | NA               | NA         | 1266 |          | 7.9           | 21.0   | 364 | 1      | 21.0         | 21.0       | 117 | 0      | ND          | ND     | 785      | 5       | 5.3         | 8.1    |
| METHYL 3-BUTENOATE*                                           | 3724-55-8    | NA               | NA         | 1266 |          | 5.1           | 8.8    | 364 | 0      | ND           | ND         | 117 | 0      | ND          | ND     | 785      | 2       | 5.1         | 8.8    |
| Methyl butyl ketone                                           | 591-78-6     | 10               | EPA RfC(1) | 1266 |          | ND            | 0.0    | 364 |        | ND           | ND         | 117 |        | ND          | ND     | 785      | 0       | ND          | ND     |

|                                                    |              | Human Health Air |                   | All Sample Durations† |      |                    |               |     | 8-hou | ır Target Pe       | riod          |     | 12-ho | ur Target Po       | eriod         |     | 24-ho   | ur Target Pe       | eriod         |
|----------------------------------------------------|--------------|------------------|-------------------|-----------------------|------|--------------------|---------------|-----|-------|--------------------|---------------|-----|-------|--------------------|---------------|-----|---------|--------------------|---------------|
| Non-Target Analytes (including TICs)               | CAS NO       | Screening Level  | Source            | N                     | Ndet | Average<br>Level** | Max<br>(ppbv) | N   | Ndet  | Average<br>Level** | Max<br>(ppbv) | N   | Ndet  | Average<br>Level** | Max<br>(ppbv) |     | Ndet    | Average<br>Level** | Max<br>(ppbv) |
| METHYL ETHYL KETONE                                | 78-93-3      | 2000             | EPA RfC(1)        | 1266                  | 48   | 1                  | 120.0         | 364 | 0     | ND                 | ND            | 117 | 1     | 5.0                | 30.0          | 785 | 47      | 1.7                | 120.0         |
| Methyl isobutyl ketone                             | 108-10-1     | 700              | EPA RfC(1)        | 1266                  | 0    | ND                 | 0.0           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 0       | ND                 | ND            |
| Methyl tert-butyl ether                            | 1634-04-4    | 700              | ATSDR Chr. MRL(1) | 1266                  | 0    | ND                 | 0.0           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 0       | ND                 | ND            |
| METHYLENE CHLORIDE                                 | 75-09-2      | 300              | ATSDR Chr. MRL(1) | 1266                  | 18   | 1                  | 9.0           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 18      | 1.0                | 9.0           |
| NAPHTHALENE, 2-(1,1-DIMETHYLETHY*                  | 54934-96-2   | NA               | NA                | 1266                  | 1    | 7.8                | 7.8           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 7.8                | 7.8           |
| NAPHTHALENE, DECAHYDRO-1,6-DIMET*                  | 1750-51-2    | NA               | NA                | 1266                  | 1    | 7.2                | 7.2           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 7.2                | 7.2           |
| NAPHTHALENE, DECAHYDRO-2,6-DIMETHYL-*              | 1618-22-0    | NA               | NA                | 1266                  | 2    | 6.8                | 6.9           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 2       | 6.8                | 6.9           |
| NAPHTHALENE, DECAHYDRO-2-METHYL-*                  | 2958-76-1    | NA               | NA                | 1266                  | 4    | 10                 | 23.0          | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 4       | 10.0               | 23.0          |
| NONANAL*                                           | 124-19-6     | NA               | NA                | 1266                  | 17   | 3.5                | 13.0          | 364 | 1     | 13.0               | 13.0          | 117 | 1     | 9.1                | 9.1           | 785 | 15      | 2.4                | 9.1           |
| NONANE, 2,2,3-TRIMETHYL-*                          | 55499-04-2   | NA               | NA                | 1266                  | 1    | 1.1                | 1.1           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 1.1                | 1.1           |
| NONANE, 2-METHYL-*                                 | 871-83-0     | NA               | NA                | 1266                  | 1    | 16                 | 16.0          | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 16.0               | 16.0          |
| NONANE, 3-METHYL-5-PROPYL-*                        | 31081-18-2   | NA               | NA                | 1266                  | 2    | 4.4                | 5.6           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 2       | 4.4                | 5.6           |
| OCTADECANE, 2,6-DIMETHYL-*                         | 75163-97-2   | NA               | NA                | 1266                  | 1    | 7.8                | 7.8           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 7.8                | 7.8           |
| OCTANAL*                                           | 124-13-0     | NA               | NA                | 1266                  | 1    | 1.1                | 1.1           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 1.1                | 1.1           |
| OCTANE, 2,2-DIMETHYL-*                             | 15869-87-1   | NA               | NA                | 1266                  | 1    | 2.5                | 2.5           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 2.5                | 2.5           |
| OCTANE, 2,3.3-TRIMETHYL-*                          | 62016-30-2   | NA               | NA                | 1266                  | 1    | 5.7                | 5.7           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 5.7                | 5.7           |
| OCTANE. 2.6-DIMETHYL-*                             | 2051-30-1    | NA               | NA                | 1266                  | 3    | 12.3               | 25.0          | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 3       | 12.3               | 25.0          |
| OCTANE, 2,7-DIMETHYL-*                             | 1072-16-8    | NA               | NA                | 1266                  | 1    | 4.8                | 4.8           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 4.8                | 4.8           |
| OCTANE, 4-ETHYL-*                                  | 15869-86-0   | NA               | NA                | 1266                  | 1    | 14                 | 14.0          | 364 | 0     | ND                 | ND            | 117 | _     | ND                 | ND            | 785 | 1       | 14.0               | 14.0          |
| o-Veratramide*                                     | 1521-39-7    | NA<br>NA         | NA<br>NA          | 1266                  | 1    | 24                 | 24.0          | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 24.0               | 24.0          |
| PENTADECANE, 2,6,10,14-TETRAMETHYL-*               | 1921-70-6    | NA<br>NA         | NA<br>NA          | 1266                  | 1    | 1.1                | 1.1           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 1.1                | 1.1           |
| PENTANE*                                           | 109-66-0     | 300              | EPA RfC(1)        | 1266                  | 59   | 21.2               | 180.0         | 364 | 12    | 10.5               | 51.0          | 117 | 0     | ND                 | ND            | 785 | 47      | 23.9               | 180.0         |
| PENTANE. 2.3-DIMETHYL-*                            | 565-59-3     | 850              | MDEQ(1)           | 1266                  | 2    | 9.2                | 16.0          | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 2       | 9.2                | 16.0          |
| PENTANE, 3-ETHYL-2,2-DIMETHYL-*                    | 16747-32-3   | NA               | NA NA             | 1266                  | 1    | 4.6                | 4.6           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 4.6                | 4.6           |
| PENTANE, 3-ETHTL-2,2-DIMETHTL- PENTANE, 3-METHYL-* | 96-14-0      | 1000             | MDEQ(1)           | 1266                  | 1    | 1.4                | 1.4           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 1.4                | 1.4           |
| PROPANE*                                           | 74-98-6      | NA               | NA<br>NA          | 1266                  | 10   | 4.4                | 25.0          | 364 | 1     | 7.0                | 7.0           | 117 | 0     | ND                 | ND            | 785 | 9       | 4.1                | 25.0          |
| PROPANOIC ACID, 2-OXO-*                            | 127-17-3     | NA<br>NA         | NA<br>NA          | 1266                  | 10   | 5.5                | 5.5           | 364 | 0     | ND                 | ND            | 117 | Ŭ     | ND                 | ND            | 785 | 1       | 5.5                | 5.5           |
| PROPYLENE                                          |              |                  |                   | 1266                  | •    | 1                  |               | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 |         | 1.0                |               |
| STYRENE                                            | 115-07-1     | 2000             | EPA RfC(1)        |                       | 18   | •                  | 1.0           | 364 |       |                    |               | 117 | _     |                    |               | 785 | 18<br>2 |                    | 1.0           |
|                                                    | 100-42-5     | 200              | ATSDR Chr. MRL(1) | 1266                  | 2    | 1                  | 5.0           |     | 0     | ND                 | ND            |     |       | ND                 | ND            |     |         | 1.0                | 5.0           |
| SULFUROUS ACID, BUTYL HEPTADECYL*                  | 1000309-18-4 | NA<br>40         | NA                | 1266                  | 1    | 6.1                | 6.1           | 364 | -     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 6.1                | 6.1           |
| TETRACHLOROETHYLENE                                | 127-18-4     | 40               | ATSDR Chr. MRL(1) | 1266                  | 1    | 1                  | 2.0           | 364 | 0     | ND                 | ND            | 117 |       | ND                 | ND            | 785 | 1       | 1.0                | 2.0           |
| TETRADECANE*                                       | 629-59-4     | NA               | NA                | 1266                  | 1    | 1.3                | 1.3           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 1.3                | 1.3           |
| TETRADECANE, 2,2-DIMETHYL-*                        | 59222-86-5   | NA               | NA<br>NA          | 1266                  | 1    | 34                 | 34.0          | 364 | 0     | ND<br>5.0          | ND            | 117 | -     | ND                 | ND            | 785 | 1       | 34.0               | 34.0          |
| TETRAHYDROFURAN                                    | 109-99-9     | 6                | MDEQ(1)           | 1266                  | 4    | 1                  | 44.0          | 364 | 1     | 5.0                | 44.0          | 117 | 0     | ND                 | ND            | 785 | 3       | 1.0                | 3.0           |
| TRANS,TRANS-1,8-DIMETHYLSPIRO[4*                   | 1000111-72-8 | NA               | NA                | 1266                  | 1    | 6.3                | 6.3           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 6.3                | 6.3           |
| trans-1,2-Dichloroethene                           | 156-60-5     | 15               | EPA RfC(1)        | 1266                  | 0    | ND                 | 0.0           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 0       | ND                 | ND            |
| trans-1,3-Dichloropropene                          | 10061-01-5   | 0.5              | MDEQ(1)           | 1266                  | 0    | ND                 | 0.0           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 0       | ND                 | ND            |
| TRANS-DECALIN, 2-METHYL-*                          | 1000152-47-3 | NA               | NA                | 1266                  | 1    | 11                 | 11.0          | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 11.0               | 11.0          |
| Trichloroethene                                    | 79-01-6      | 2                | EPA RfC(1)        | 1266                  | 1    | 1                  | 1.0           | 364 | 0     | ND                 | ND            | 117 |       | ND                 | ND            | 785 | 1       | 1.0                | 1.0           |
| Trichlorofluoromethane                             | 75-69-4      | 130              | EPA RfC(1)        | 1266                  | 0    | ND                 | 0.0           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 0       | ND                 | ND            |
| TRICOSYL PENTAFLUOROPROPIONATE*                    | 1000351-81-0 | NA               | NA                | 1266                  | 1    | 1.9                | 1.9           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 1.9                | 1.9           |
| TRIDECANE*                                         | 629-50-5     | NA               | NA                | 1266                  | 1    | 10                 | 10.0          | 364 | 0     | ND                 | ND            | 117 | _     | ND                 | ND            | 785 | 1       | 10.0               | 10.0          |
| TRIDECANE, 4,8-DIMETHYL-*                          | 55030-62-1   | NA               | NA                | 1266                  | 1    | 4.8                | 4.8           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 4.8                | 4.8           |
| TRIDECANE, 6-METHYL-*                              | 13287-21-3   | NA               | NA                | 1266                  | 1    | 45                 | 45.0          | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 45.0               | 45.0          |
| TRIDECYL HEPTAFLUOROBUTYRATE*                      | 1000351-82-8 | NA               | NA                | 1266                  | 0    | ND                 | 0.0           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 0       | ND                 | ND            |
| TRIETHYLAMINE*                                     | 121-44-8     | 2                | EPA RfC(1)        | 1266                  | 2    | 12.9               | 18.0          | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 2       | 12.9               | 18.0          |
| TRISULFIDE, DIMETHYL*                              | 3658-80-8    | NA               | NA                | 1266                  | 1    | 5.1                | 5.1           | 364 | 0     | ND                 | ND            | 117 | 0     | ND                 | ND            | 785 | 1       | 5.1                | 5.1           |

|                                      |            | Human Health Air          |                   |      |      | nple Duratio       | ns†           | 8-hour Target Period |      |                    |               |     | 12-hοι | ur Target Pe       | eriod         |     | 24-hou | our Target Period  |               |
|--------------------------------------|------------|---------------------------|-------------------|------|------|--------------------|---------------|----------------------|------|--------------------|---------------|-----|--------|--------------------|---------------|-----|--------|--------------------|---------------|
| Non-Target Analytes (including TICs) | CAS NO     | Screening Level<br>(ppbv) | Source            | N    | Ndet | Average<br>Level** | Max<br>(ppbv) | N                    | Ndet | Average<br>Level** | Max<br>(ppbv) | N   | Ndet   | Average<br>Level** | Max<br>(ppbv) | N   | Ndet   | Average<br>Level** | Max<br>(ppbv) |
| UNDECANE, 2,2-DIMETHYL-*             | 17312-64-0 | NA                        | NA                | 1266 | 1    | 2.7                | 2.7           | 364                  | 0    | ND                 | ND            | 117 | 0      | ND                 | ND            | 785 | 1      | 2.7                | 2.7           |
| UNDECANE, 2,6-DIMETHYL-*             | 17301-23-4 | NA                        | NA                | 1266 | 6    | 13.4               | 52.0          | 364                  | 0    | ND                 | ND            | 117 | 0      | ND                 | ND            | 785 | 6      | 13.4               | 52.0          |
| UNDECANE, 3,6-DIMETHYL-*             | 17301-28-9 | NA                        | NA                | 1266 | 2    | 8.7                | 16.0          | 364                  | 0    | ND                 | ND            | 117 | 0      | ND                 | ND            | 785 | 2      | 8.7                | 16.0          |
| UNDECANE, 3,9-DIMETHYL-*             | 17301-31-4 | NA                        | NA                | 1266 | 1    | 4.9                | 4.9           | 364                  | 0    | ND                 | ND            | 117 | 0      | ND                 | ND            | 785 | 1      | 4.9                | 4.9           |
| UNDECANE, 5-METHYL-*                 | 1632-70-8  | NA                        | NA                | 1266 | 1    | 1.2                | 1.2           | 364                  | 0    | ND                 | ND            | 117 | 0      | ND                 | ND            | 785 | 1      | 1.2                | 1.2           |
| UNKNOWN*                             | 78-99-9    | NA                        | NA                | 1266 | 1    | 1.6                | 1.6           | 364                  | 0    | ND                 | ND            | 117 | 0      | ND                 | ND            | 785 | 1      | 1.6                | 1.6           |
| UNKNOWN*                             | 288-47-1   | NA                        | NA                | 1266 | 1    | 1.2                | 1.2           | 364                  | 0    | ND                 | ND            | 117 | 0      | ND                 | ND            | 785 | 1      | 1.2                | 1.2           |
| UNKNOWN*                             | 62-55-5    | NA                        | NA                | 1266 | 2    | 1.4                | 1.4           | 364                  | 0    | ND                 | ND            | 117 | 0      | ND                 | ND            | 785 | 2      | 1.4                | 1.4           |
| Vinyl acetate                        | 108-05-4   | 10                        | ATSDR Int. MRL(1) | 1266 | 3    | 1                  | 2.0           | 364                  | 0    | ND                 | ND            | 117 | 0      | ND                 | ND            | 785 | 3      | 1.0                | 2.0           |
| Vinyl bromide                        | 593-60-2   | 1                         | EPA RfC(1)        | 1266 | 0    | ND                 | 0.0           | 364                  | 0    | ND                 | ND            | 117 | 0      | ND                 | ND            | 785 | 0      | ND                 | ND            |
| Vinyl chloride                       | 75-01-4    | 30                        | ATSDR Int. MRL(1) | 1266 | 0    | ND                 | 0.0           | 364                  | 0    | ND                 | ND            | 117 | 0      | ND                 | ND            | 785 | 0      | ND                 | ND            |

NA - Not Available/ND - Not Detected

N - Number of samples analyzed/NDET - Number of detections

(¹) - Unless otherwise noted, the screening level was obtained from Enbridge Oil Spill Human Health Air Screening Levels, August 31, 2011.

ATSDR Chr. MRL - Agency for Toxic Substances and Disease Registry - Chronic Minimal Risk Level for Inhalation

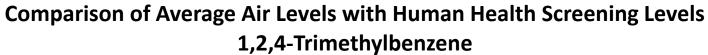
EPA RfC - If no Chronic MRL is available, the screening level is the EPA Chronic Reference Concentration (RfC). For 2 chemicals, vinyl acetate and vinyl chloride, the ATSDR Intermediate (Int.) MRL was selected. These were developed more recently than the EPA RfC and were lower than the EPA RfC.

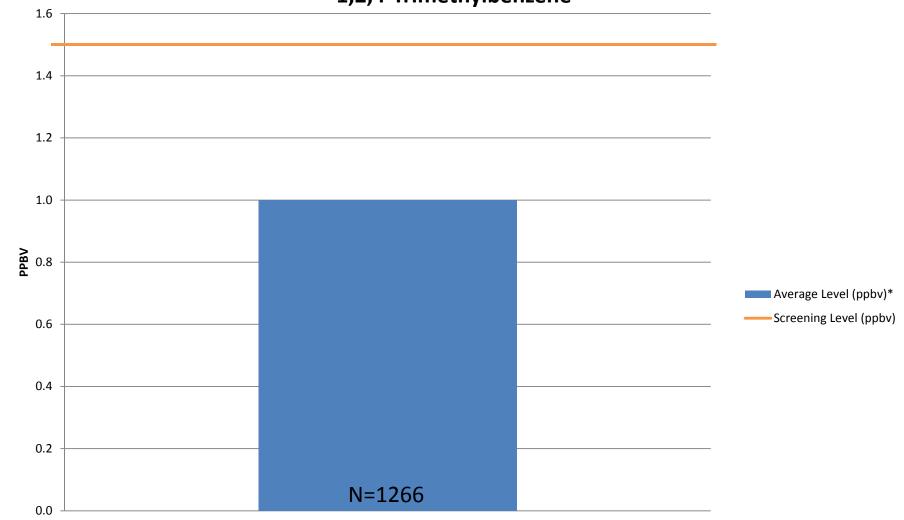
EPA RSL - If none of the above are available, the EPA Regional Screening Level (RSL) is used.

MDEQ - If none of the above are available, the Michigan DEQ Air Quality Division, Air Toxics Screening Level is the screening level.

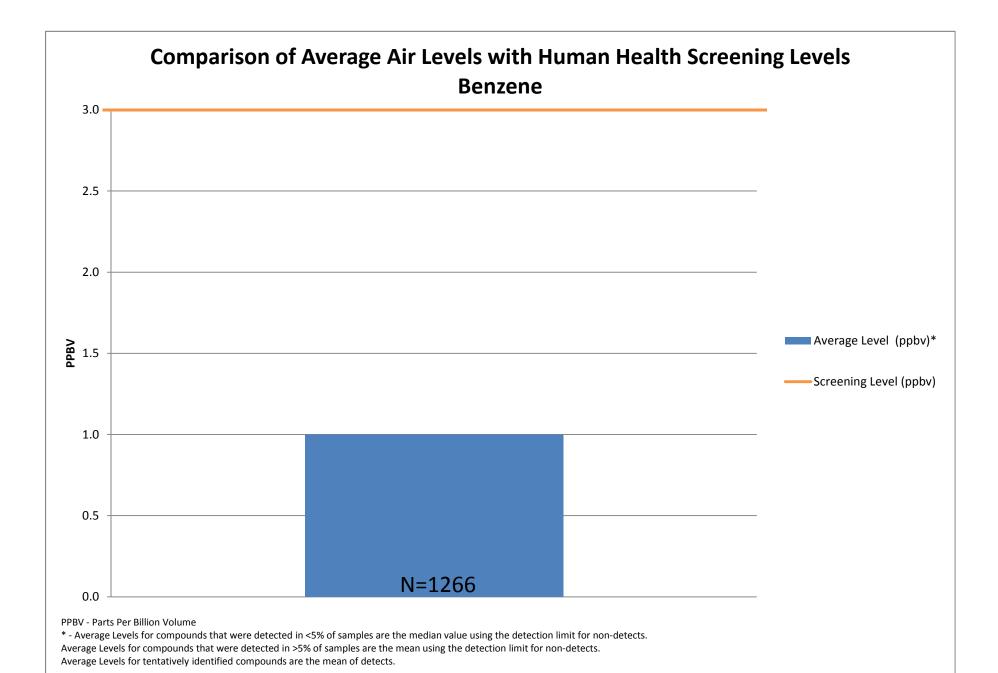
<sup>\* -</sup> TIC (tentatively identified compound)

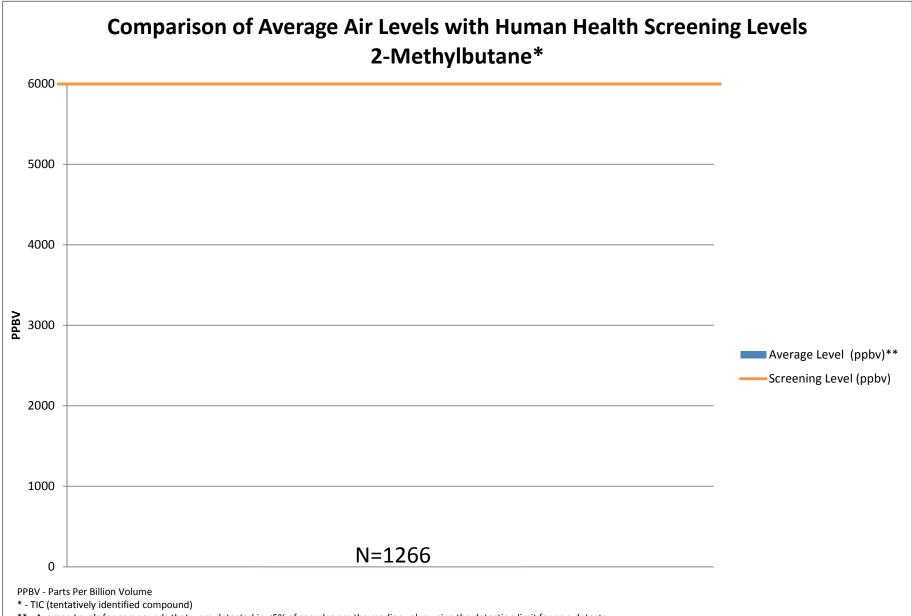
<sup>\*\* -</sup> Average Levels for compounds that were detected in <5% of samples are the median value using the detection limit for non-detects. Average Levels for compounds that were detected in >5% of samples are the mean using the detection limit for non-detects. Average Levels for tentatively identified compounds are the mean of detects. †Samples reported do not include instantaneous 'grab' samples. Samples with 8-24 hour target sample periods are included.





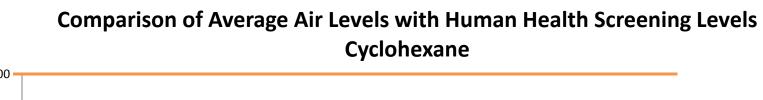
<sup>\* -</sup> Average Levels for compounds that were detected in <5% of samples are the median value using the detection limit for non-detects. Average Levels for compounds that were detected in >5% of samples are the mean using the detection limit for non-detects. Average Levels for tentatively identified compounds are the mean of detects.

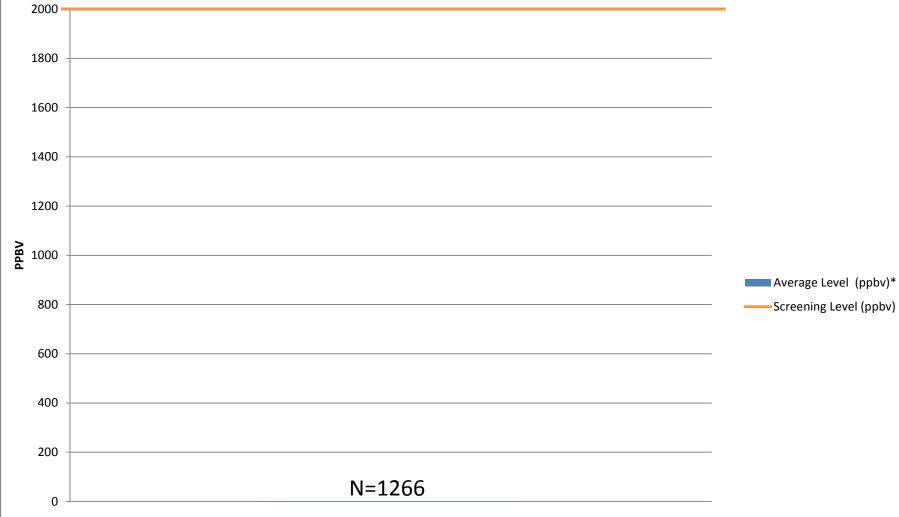




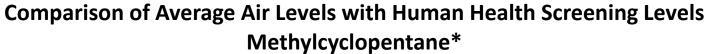
<sup>\*\* -</sup> Average Levels for compounds that were detected in <5% of samples are the median value using the detection limit for non-detects.

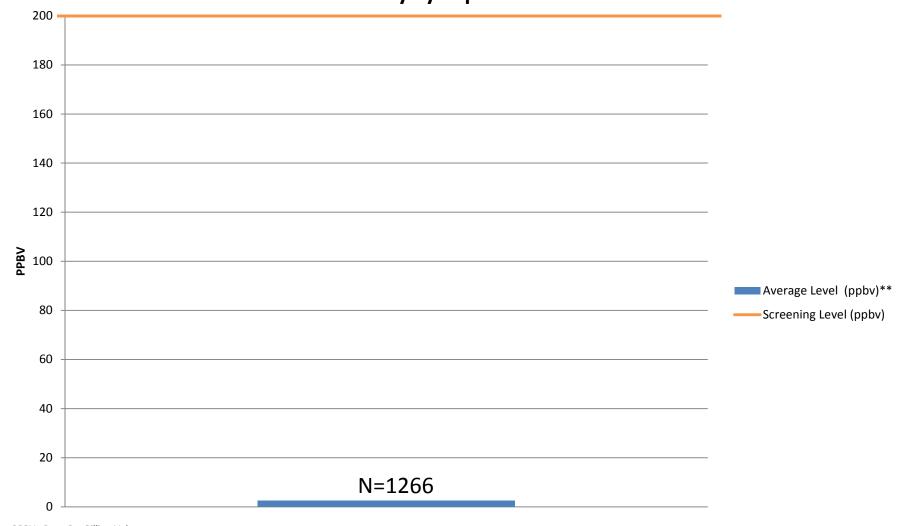
Average Levels for compounds that were detected in >5% of samples are the mean using the detection limit for non-detects.





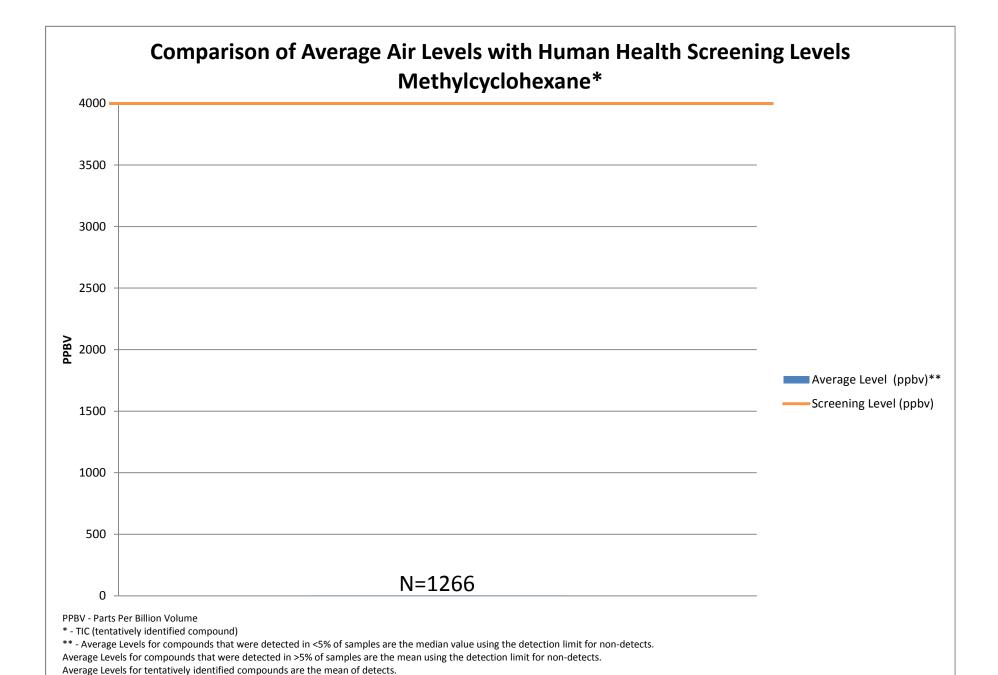
\* - Average Levels for compounds that were detected in <5% of samples are the median value using the detection limit for non-detects. Average Levels for compounds that were detected in >5% of samples are the mean using the detection limit for non-detects. Average Levels for tentatively identified compounds are the mean of detects.

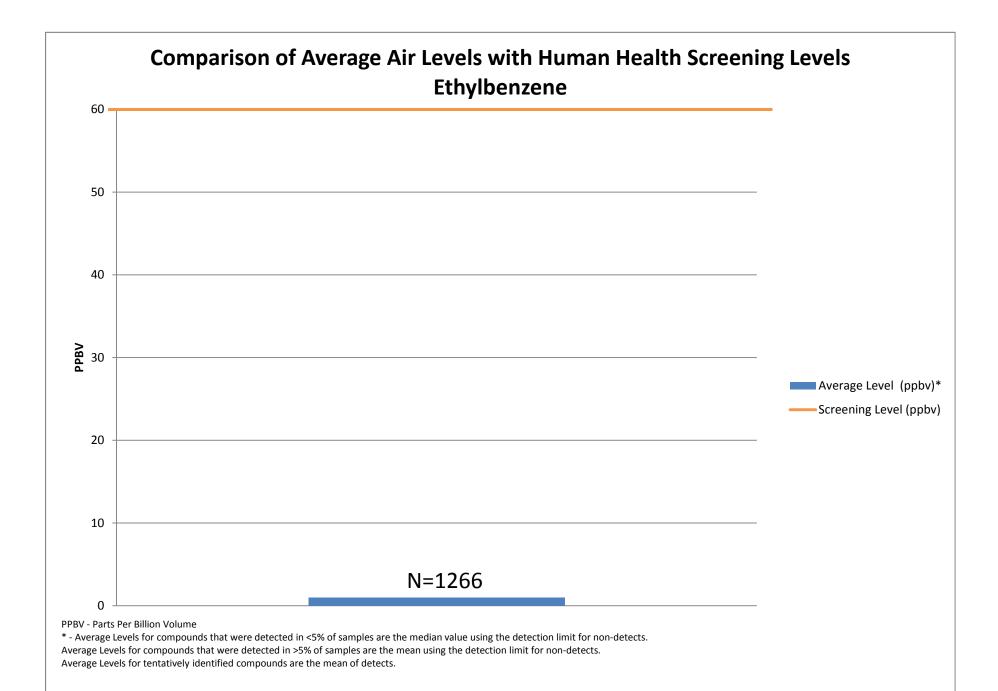


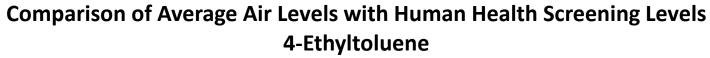


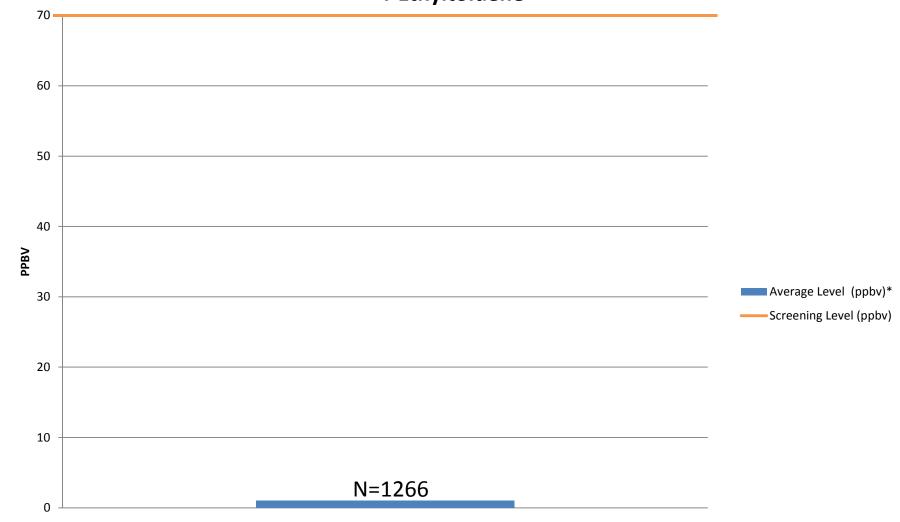
\*\* - Average Levels for compounds that were detected in <5% of samples are the median value using the detection limit for non-detects. Average Levels for compounds that were detected in >5% of samples are the mean using the detection limit for non-detects. Average Levels for tentatively identified compounds are the mean of detects.

<sup>\* -</sup> TIC (tentatively identified compound)

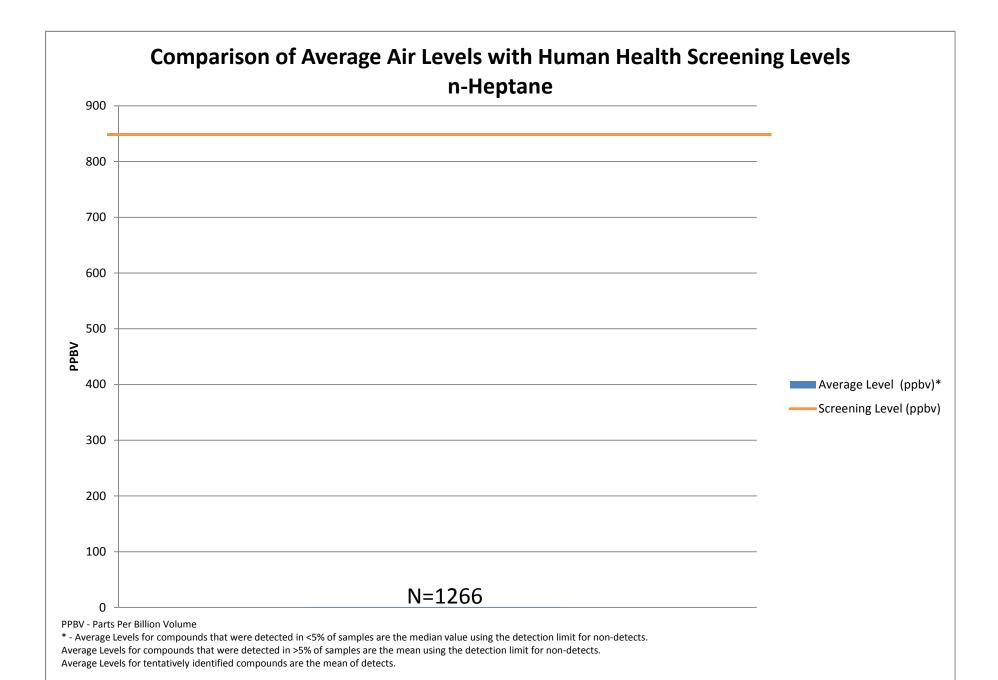


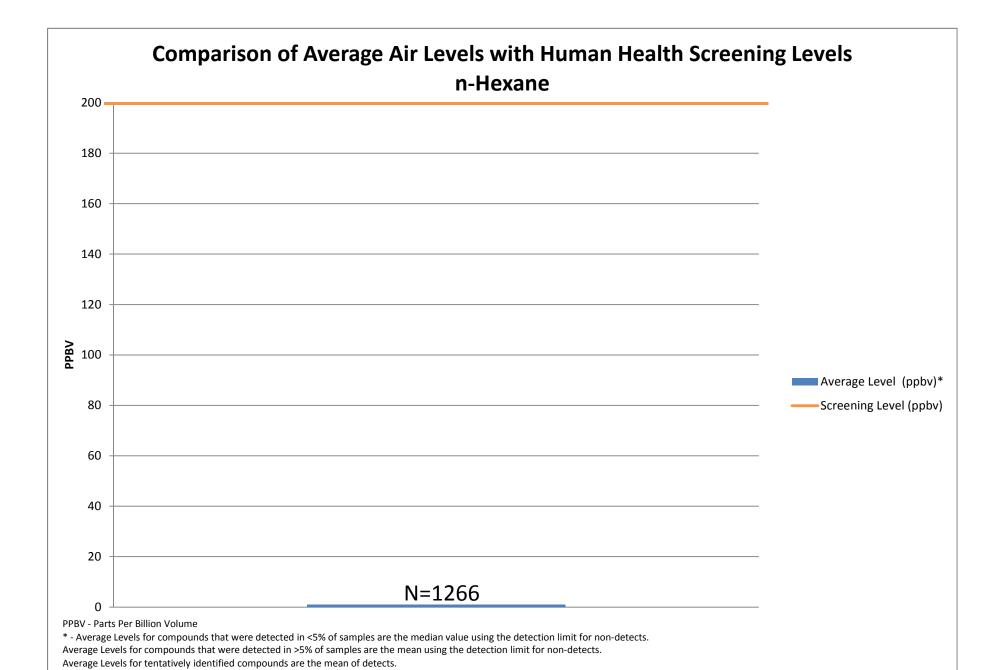


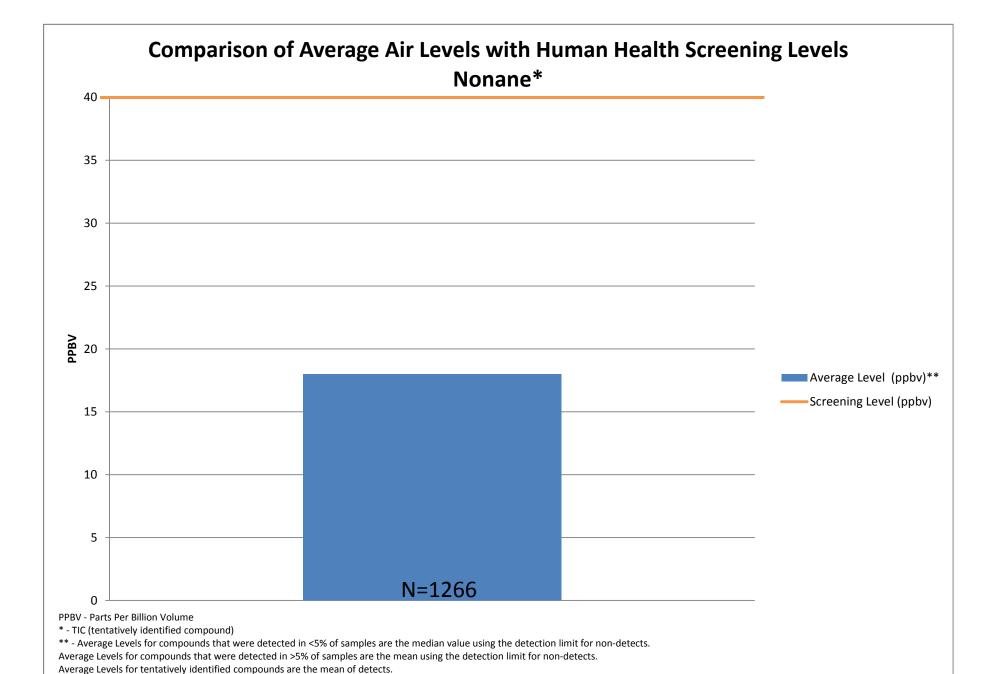


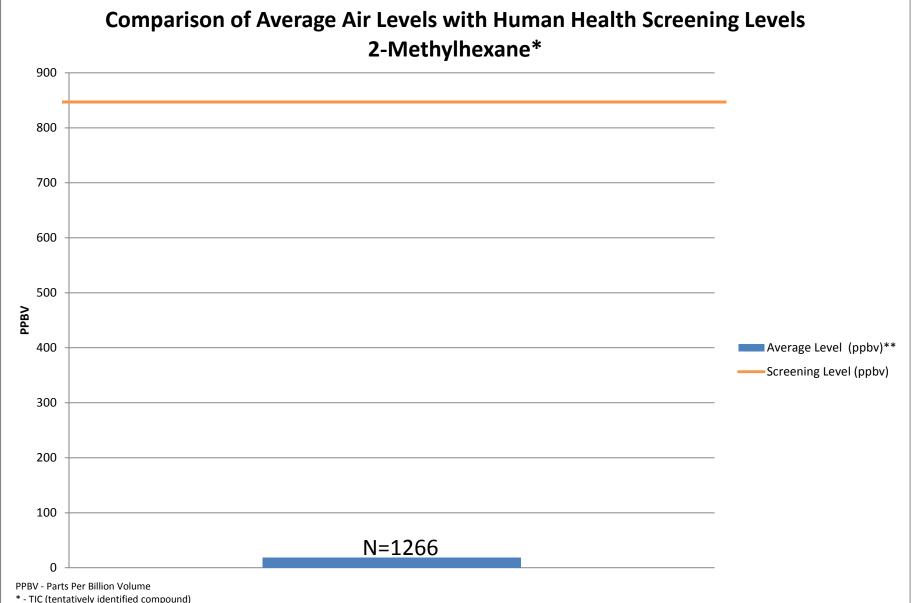


\* - Average Levels for compounds that were detected in <5% of samples are the median value using the detection limit for non-detects. Average Levels for compounds that were detected in >5% of samples are the mean using the detection limit for non-detects. Average Levels for tentatively identified compounds are the mean of detects.



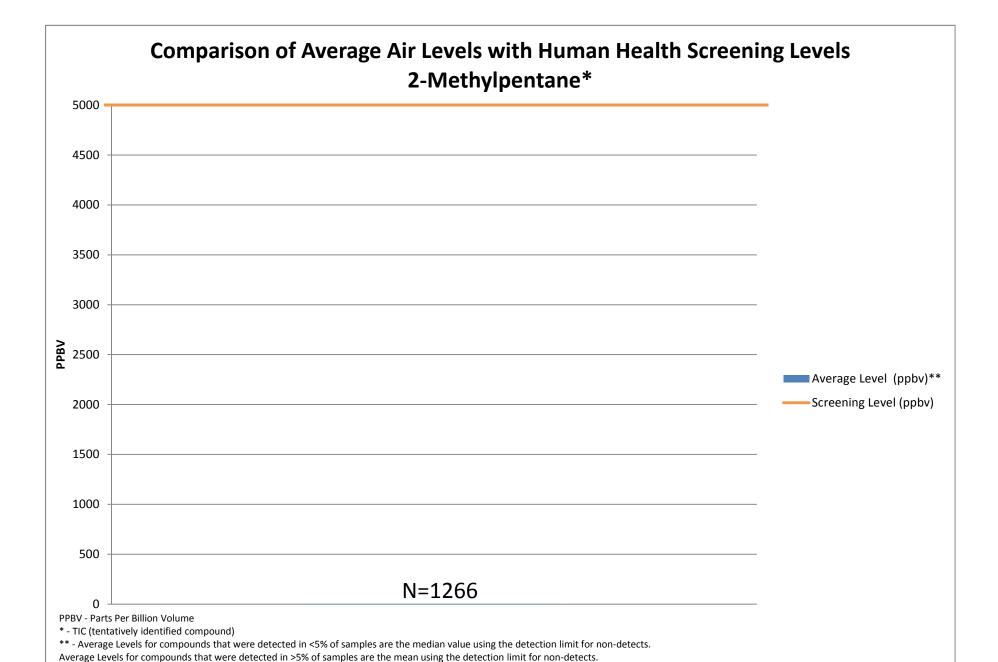


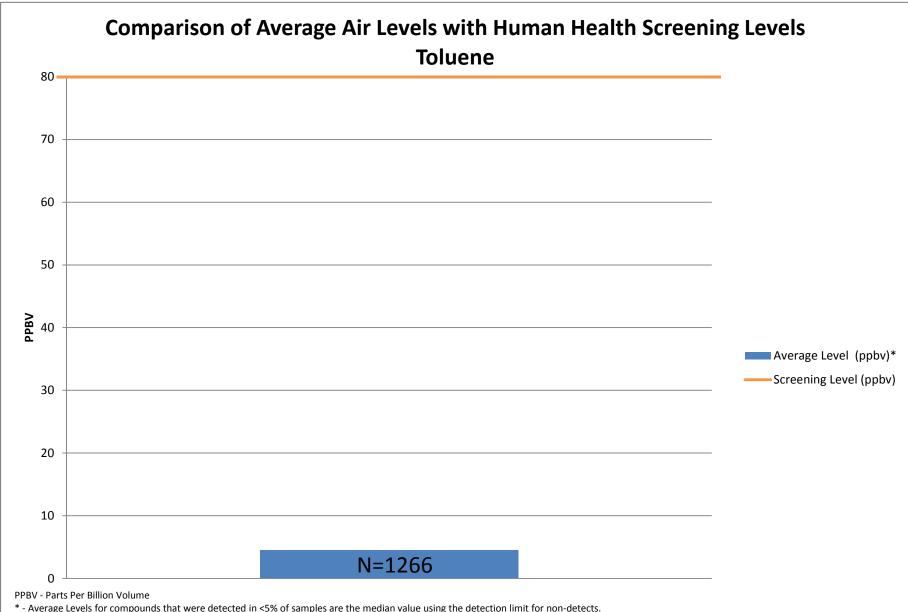




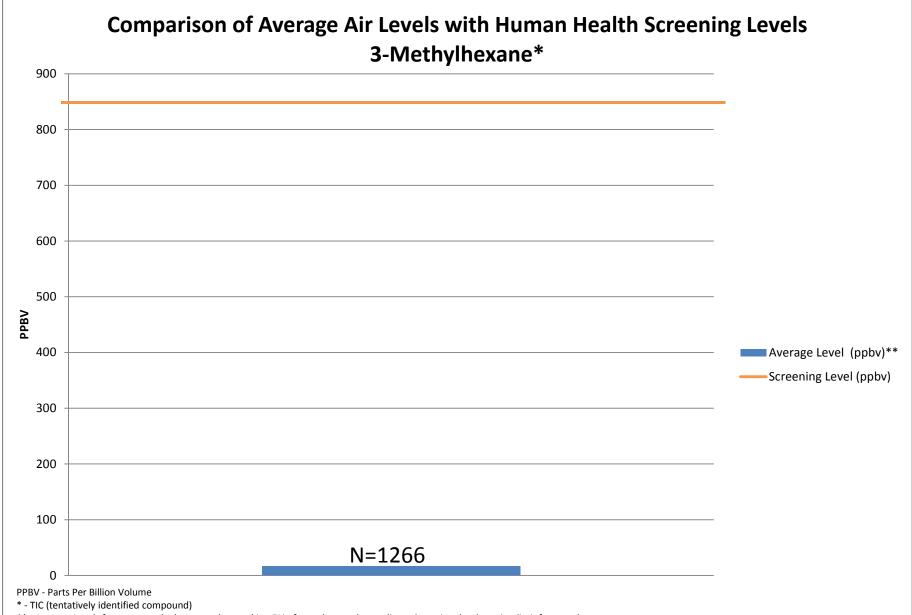
Average Levels for compounds that were detected in >5% of samples are the mean using the detection limit for non-detects.

<sup>\*\* -</sup> Average Levels for compounds that were detected in <5% of samples are the median value using the detection limit for non-detects.





<sup>\* -</sup> Average Levels for compounds that were detected in <5% of samples are the median value using the detection limit for non-detects. Average Levels for compounds that were detected in >5% of samples are the mean using the detection limit for non-detects. Average Levels for tentatively identified compounds are the mean of detects.



<sup>\*\* -</sup> Average Levels for compounds that were detected in <5% of samples are the median value using the detection limit for non-detects.

 $Average\ Levels\ for\ compounds\ that\ were\ detected\ in\ >5\%\ of\ samples\ are\ the\ mean\ using\ the\ detection\ limit\ for\ non-detects.$ 

