

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW053023\  
 Data File : VW026085.D  
 Acq On : 30 May 2023 12:58  
 Operator : SY/MD  
 Sample : VSTDCCC025  
 Misc : 5.00g/10mL/MSVOA\_W/SOIL  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_W  
 LabSampleId :  
 VSTDCCC025

Quant Time: May 31 01:52:05 2023  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM053023SMA.M  
 Quant Title : SFAM01.0  
 QLast Update : Wed May 31 01:45:52 2023  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 25% Max. Rel. Area : 200%

|      | Compound                    | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|-----------------------------|-------|-------|-------|-------|----------|
| 1 I  | 1,4-Difluorobenzene         | 1.000 | 1.000 | 0.0   | 99    | 0.00     |
| 2 T  | Dichlorodifluoromethane     | 0.306 | 0.315 | -2.9  | 98    | 0.00     |
| 3 T  | Chloromethane               | 0.393 | 0.442 | -12.5 | 106   | 0.00     |
| 4 S  | Vinyl Chloride-d3           | 0.352 | 0.401 | -13.9 | 98    | 0.00     |
| 5 T  | Vinyl chloride              | 0.420 | 0.459 | -9.3  | 104   | 0.00     |
| 6 T  | Bromomethane                | 0.229 | 0.212 | 7.4   | 98    | 0.00     |
| 7 S  | Chloroethane-d5             | 0.263 | 0.264 | -0.4  | 92    | 0.00     |
| 8 T  | Chloroethane                | 0.227 | 0.226 | 0.4   | 101   | 0.00     |
| 9 T  | Trichlorofluoromethane      | 0.270 | 0.275 | -1.9  | 112   | 0.00     |
| 10 T | 1,1,2-Trichloro-1,2,2-trifl | 0.341 | 0.355 | -4.1  | 106   | 0.00     |
| 11 S | 1,1-Dichloroethene-d2       | 0.173 | 0.188 | -8.7  | 97    | 0.00     |
| 12 T | 1,1-Dichloroethene          | 0.318 | 0.333 | -4.7  | 104   | 0.00     |
| 13 T | Acetone                     | 0.079 | 0.073 | 7.6   | 97    | 0.00     |
| 14 T | Carbon disulfide            | 1.124 | 1.189 | -5.8  | 106   | 0.00     |
| 15 T | Methyl Acetate              | 0.178 | 0.187 | -5.1  | 108   | 0.00     |
| 16 T | Methylene chloride          | 0.399 | 0.382 | 4.3   | 108   | 0.00     |
| 17 T | trans-1,2-Dichloroethene    | 0.357 | 0.370 | -3.6  | 108   | 0.00     |
| 18 T | Methyl tert-butyl Ether     | 0.563 | 0.582 | -3.4  | 106   | 0.00     |
| 19 T | 1,1-Dichloroethane          | 0.727 | 0.745 | -2.5  | 108   | 0.00     |
| 20 T | cis-1,2-Dichloroethene      | 0.378 | 0.397 | -5.0  | 109   | 0.00     |
| 21 S | 2-Butanone-d5               | 0.102 | 0.108 | -5.9  | 99    | 0.00     |
| 22 T | 2-Butanone                  | 0.113 | 0.116 | -2.7  | 105   | 0.00     |
| 23 T | Bromochloromethane          | 0.161 | 0.161 | 0.0   | 106   | 0.00     |
| 24 S | Chloroform-d                | 0.732 | 0.765 | -4.5  | 101   | 0.00     |
| 25 T | Chloroform                  | 0.677 | 0.691 | -2.1  | 107   | 0.00     |
| 26 S | 1,2-Dichloroethane-d4       | 0.400 | 0.412 | -3.0  | 100   | 0.00     |
| 27 T | 1,2-Dichloroethane          | 0.467 | 0.475 | -1.7  | 106   | 0.00     |
| 28 I | Chlorobenzene-d5            | 1.000 | 1.000 | 0.0   | 92    | 0.00     |
| 29 T | Cyclohexane                 | 0.772 | 0.860 | -11.4 | 105   | 0.00     |
| 30 T | 1,1,1-Trichloroethane       | 0.607 | 0.651 | -7.2  | 105   | 0.00     |
| 31 T | Carbon tetrachloride        | 0.554 | 0.598 | -7.9  | 107   | 0.00     |
| 32 S | Benzene-d6                  | 1.682 | 1.880 | -11.8 | 101   | 0.00     |
| 33 T | Benzene                     | 1.713 | 1.885 | -10.0 | 108   | 0.00     |
| 34 T | Trichloroethene             | 0.439 | 0.478 | -8.9  | 108   | 0.00     |
| 35 T | Methylcyclohexane           | 0.776 | 0.818 | -5.4  | 100   | 0.00     |
| 36 S | 1,2-Dichloropropane-d6      | 0.533 | 0.552 | -3.6  | 95    | 0.00     |
| 37 T | 1,2-Dichloropropane         | 0.464 | 0.481 | -3.7  | 104   | 0.00     |
| 38 T | Bromodichloromethane        | 0.539 | 0.545 | -1.1  | 100   | 0.00     |
| 39 T | cis-1,3-Dichloropropene     | 0.682 | 0.703 | -3.1  | 100   | 0.00     |
| 40 T | 4-Methyl-2-pentanone        | 0.273 | 0.291 | -6.6  | 99    | 0.00     |
| 41 S | Toluene-d8                  | 1.492 | 1.561 | -4.6  | 96    | 0.00     |
| 42 T | Toluene                     | 1.752 | 1.874 | -7.0  | 104   | 0.00     |
| 43 S | trans-1,3-Dichloropropene-d | 0.211 | 0.215 | -1.9  | 96    | 0.00     |
| 44 T | trans-1,3-Dichloropropene   | 0.568 | 0.593 | -4.4  | 101   | 0.00     |
| 45 T | 1,1,2-Trichloroethane       | 0.306 | 0.320 | -4.6  | 104   | 0.00     |
| 46 T | Tetrachloroethene           | 0.310 | 0.341 | -10.0 | 109   | 0.00     |
| 47 S | 2-Hexanone-d5               | 0.084 | 0.095 | -13.1 | 99    | 0.00     |
| 48 T | 2-Hexanone                  | 0.193 | 0.201 | -4.1  | 98    | 0.00     |

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 LabSampleId :  
 VSTDCCC025

Quant Time: May 31 01:52:05 2023  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM053023SMA.M  
 Quant Title : SFAM01.0  
 QLast Update : Wed May 31 01:45:52 2023  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 25% Max. Rel. Area : 200%

|      | Compound                    | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|-----------------------------|-------|-------|------|-------|----------|
| 49 T | Dibromochloromethane        | 0.321 | 0.313 | 2.5  | 99    | 0.00     |
| 50 T | 1,2-Dibromoethane           | 0.283 | 0.276 | 2.5  | 98    | 0.00     |
| 51 T | Chlorobenzene               | 1.098 | 1.120 | -2.0 | 102   | 0.00     |
| 52 T | Ethylbenzene                | 1.976 | 2.094 | -6.0 | 103   | 0.00     |
| 53 T | m,p-Xylene                  | 0.735 | 0.791 | -7.6 | 105   | 0.00     |
| 54 T | o-Xylene                    | 0.694 | 0.763 | -9.9 | 109   | 0.00     |
| 55 T | Styrene                     | 1.190 | 1.308 | -9.9 | 106   | 0.00     |
| 56 S | 1,1,2,2-Tetrachloroethane-d | 0.388 | 0.423 | -9.0 | 102   | 0.00     |
| 57 T | 1,1,2,2-Tetrachloroethane   | 0.370 | 0.394 | -6.5 | 106   | 0.00     |
| 58 I | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0  | 95    | 0.00     |
| 59 T | Bromoform                   | 0.375 | 0.369 | 1.6  | 107   | 0.00     |
| 60   | Isopropylbenzene            | 3.927 | 4.110 | -4.7 | 107   | 0.00     |
| 61   | 1,2,3-Trichloropropane      | 0.560 | 0.563 | -0.5 | 107   | 0.00     |
| 62   | 1,3,5-Trimethylbenzene      | 3.122 | 3.331 | -6.7 | 109   | 0.00     |
| 63   | 1,2,4-Trimethylbenzene      | 2.859 | 3.045 | -6.5 | 108   | 0.00     |
| 64 T | 1,3-Dichlorobenzene         | 1.651 | 1.705 | -3.3 | 109   | 0.00     |
| 65 T | 1,4-Dichlorobenzene         | 1.662 | 1.638 | 1.4  | 105   | 0.00     |
| 66 S | 1,2-Dichlorobenzene-d4      | 0.957 | 0.959 | -0.2 | 98    | 0.00     |
| 67 T | 1,2-Dichlorobenzene         | 1.483 | 1.466 | 1.1  | 106   | 0.00     |
| 68 T | 1,2-Dibromo-3-chloropropane | 0.130 | 0.123 | 5.4  | 103   | 0.00     |
| 69   | 1,3,5-Trichlorobenzene      | 1.134 | 1.080 | 4.8  | 106   | 0.00     |
| 70 T | 1,2,4-trichlorobenzene      | 0.936 | 0.927 | 1.0  | 108   | 0.00     |
| 71   | Naphthalene                 | 1.668 | 1.727 | -3.5 | 112   | 0.00     |
| 72 T | 1,2,3-Trichlorobenzene      | 0.818 | 0.803 | 1.8  | 111   | 0.00     |

(#) = Out of Range

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