

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX082522\  
 Data File : VX030768.D  
 Acq On : 25 Aug 2022 13:13  
 Operator : JC/MD  
 Sample : VSTDICV050  
 Misc : 5.0mL/MSVOA\_X/WATER  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 ClientSampleId :  
 ICVVX082522

Quant Time: Aug 26 02:39:09 2022  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X082522W.M  
 Quant Title : SW846 8260  
 QLast Update : Fri Aug 26 02:03:28 2022  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 25% Max. Rel. Area : 150%

|       | Compound                    | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|-------|-----------------------------|-------|-------|-------|-------|----------|
| 1 I   | Pentafluorobenzene          | 1.000 | 1.000 | 0.0   | 103   | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 0.463 | 0.439 | 5.2   | 103   | 0.00     |
| 3 P   | Chloromethane               | 0.432 | 0.368 | 14.8  | 102   | 0.00     |
| 4 C   | Vinyl Chloride              | 0.508 | 0.434 | 14.6# | 100   | 0.00     |
| 5 T   | Bromomethane                | 0.233 | 0.225 | 3.4   | 115   | 0.00     |
| 6 T   | Chloroethane                | 0.313 | 0.263 | 16.0  | 104   | 0.00     |
| 7 T   | Trichlorofluoromethane      | 0.765 | 0.687 | 10.2  | 103   | 0.00     |
| 8 T   | Diethyl Ether               | 0.290 | 0.265 | 8.6   | 104   | 0.00     |
| 9 T   | 1,1,2-Trichlorotrifluoroeth | 0.487 | 0.448 | 8.0   | 104   | 0.00     |
| 10 T  | Methyl Iodide               | 0.337 | 0.342 | -1.5  | 93    | 0.00     |
| 11 T  | Tert butyl alcohol          | 0.128 | 0.118 | 7.8   | 109   | 0.00     |
| 12 CM | 1,1-Dichloroethene          | 0.453 | 0.410 | 9.5#  | 105   | 0.00     |
| 13 T  | Acrolein                    | 0.097 | 0.112 | -15.5 | 107   | 0.00     |
| 14 T  | Allyl chloride              | 0.839 | 0.755 | 10.0  | 102   | 0.00     |
| 15 T  | Acrylonitrile               | 0.313 | 0.290 | 7.3   | 101   | 0.00     |
| 16 T  | Acetone                     | 0.272 | 0.255 | 6.3   | 104   | 0.00     |
| 17 T  | Carbon Disulfide            | 1.052 | 0.864 | 17.9  | 103   | 0.00     |
| 18 T  | Methyl Acetate              | 0.835 | 0.766 | 8.3   | 102   | 0.00     |
| 19 T  | Methyl tert-butyl Ether     | 1.569 | 1.509 | 3.8   | 104   | 0.00     |
| 20 T  | Methylene Chloride          | 0.559 | 0.497 | 11.1  | 105   | 0.00     |
| 21 T  | trans-1,2-Dichloroethene    | 0.487 | 0.433 | 11.1  | 103   | 0.00     |
| 22 T  | Diisopropyl ether           | 1.685 | 1.595 | 5.3   | 104   | 0.00     |
| 23 T  | Vinyl Acetate               | 1.351 | 1.302 | 3.6   | 103   | 0.00     |
| 24 P  | 1,1-Dichloroethane          | 0.955 | 0.872 | 8.7   | 103   | 0.00     |
| 25 T  | 2-Butanone                  | 0.438 | 0.410 | 6.4   | 103   | 0.00     |
| 26 T  | 2,2-Dichloropropane         | 0.692 | 0.674 | 2.6   | 104   | 0.00     |
| 27 T  | cis-1,2-Dichloroethene      | 0.601 | 0.549 | 8.7   | 104   | 0.00     |
| 28 T  | Bromochloromethane          | 0.411 | 0.360 | 12.4  | 97    | 0.00     |
| 29 T  | Tetrahydrofuran             | 0.286 | 0.262 | 8.4   | 102   | 0.00     |
| 30 C  | Chloroform                  | 1.005 | 0.910 | 9.5#  | 104   | 0.00     |
| 31 T  | Cyclohexane                 | 0.752 | 0.696 | 7.4   | 101   | 0.00     |
| 32 T  | 1,1,1-Trichloroethane       | 0.813 | 0.776 | 4.6   | 104   | 0.00     |
| 33 S  | 1,2-Dichloroethane-d4       | 0.588 | 0.540 | 8.2   | 107   | 0.00     |
| 34 I  | 1,4-Difluorobenzene         | 1.000 | 1.000 | 0.0   | 106   | 0.00     |
| 35 S  | Dibromofluoromethane        | 0.307 | 0.286 | 6.8   | 106   | 0.00     |
| 36 T  | 1,1-Dichloropropene         | 0.418 | 0.368 | 12.0  | 103   | 0.00     |
| 37 T  | Ethyl Acetate               | 0.496 | 0.447 | 9.9   | 103   | 0.00     |
| 38 T  | Carbon Tetrachloride        | 0.441 | 0.390 | 11.6  | 101   | 0.00     |
| 39 T  | Methylcyclohexane           | 0.501 | 0.459 | 8.4   | 105   | 0.00     |
| 40 TM | Benzene                     | 1.252 | 1.127 | 10.0  | 103   | 0.00     |
| 41 T  | Methacrylonitrile           | 0.277 | 0.251 | 9.4   | 100   | 0.00     |
| 42 TM | 1,2-Dichloroethane          | 0.444 | 0.402 | 9.5   | 104   | 0.00     |
| 43 T  | Isopropyl Acetate           | 0.747 | 0.709 | 5.1   | 105   | 0.00     |
| 44 TM | Trichloroethene             | 0.362 | 0.327 | 9.7   | 103   | 0.00     |
| 45 C  | 1,2-Dichloropropane         | 0.346 | 0.320 | 7.5#  | 104   | 0.00     |
| 46 T  | Dibromomethane              | 0.246 | 0.220 | 10.6  | 103   | 0.00     |
| 47 T  | Bromodichloromethane        | 0.463 | 0.439 | 5.2   | 104   | 0.00     |
| 48 T  | Methyl methacrylate         | 0.372 | 0.369 | 0.8   | 109   | 0.00     |

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 Sample : VSTDICV050  
 Misc : 5.0mL/MSVOA\_X/WATER  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 ClientSampleId :  
 ICSVVX082522

Quant Time: Aug 26 02:39:09 2022  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X082522W.M  
 Quant Title : SW846 8260  
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Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 25% Max. Rel. Area : 150%

|       | Compound                    | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|-------|-----------------------------|-------|-------|------|-------|----------|
| 49 T  | 1,4-Dioxane                 | 0.009 | 0.009 | 0.0  | 107   | 0.00     |
| 50 S  | Toluene-d8                  | 1.071 | 0.994 | 7.2  | 105   | 0.00     |
| 51 T  | 4-Methyl-2-Pentanone        | 0.504 | 0.463 | 8.1  | 101   | 0.00     |
| 52 CM | Toluene                     | 0.781 | 0.715 | 8.5# | 102   | 0.00     |
| 53 T  | t-1,3-Dichloropropene       | 0.448 | 0.438 | 2.2  | 104   | 0.00     |
| 54 T  | cis-1,3-Dichloropropene     | 0.504 | 0.484 | 4.0  | 103   | 0.00     |
| 55 T  | 1,1,2-Trichloroethane       | 0.347 | 0.324 | 6.6  | 103   | 0.00     |
| 56 T  | Ethyl methacrylate          | 0.498 | 0.494 | 0.8  | 104   | 0.00     |
| 57 T  | 1,3-Dichloropropane         | 0.560 | 0.520 | 7.1  | 104   | 0.00     |
| 58 T  | 2-Chloroethyl Vinyl ether   | 0.266 | 0.269 | -1.1 | 105   | 0.00     |
| 59 T  | 2-Hexanone                  | 0.378 | 0.358 | 5.3  | 102   | 0.00     |
| 60 T  | Dibromochloromethane        | 0.362 | 0.343 | 5.2  | 102   | 0.00     |
| 61 T  | 1,2-Dibromoethane           | 0.352 | 0.332 | 5.7  | 103   | 0.00     |
| 62 S  | 4-Bromofluorobenzene        | 0.399 | 0.382 | 4.3  | 105   | 0.00     |
| 63 I  | Chlorobenzene-d5            | 1.000 | 1.000 | 0.0  | 105   | 0.00     |
| 64 T  | Tetrachloroethene           | 0.394 | 0.356 | 9.6  | 101   | 0.00     |
| 65 PM | Chlorobenzene               | 0.967 | 0.881 | 8.9  | 102   | 0.00     |
| 66 T  | 1,1,1,2-Tetrachloroethane   | 0.363 | 0.343 | 5.5  | 101   | 0.00     |
| 67 C  | Ethyl Benzene               | 1.671 | 1.595 | 4.5# | 104   | 0.00     |
| 68 T  | m/p-Xylenes                 | 0.640 | 0.612 | 4.4  | 104   | 0.00     |
| 69 T  | o-Xylene                    | 0.634 | 0.605 | 4.6  | 103   | 0.00     |
| 70 T  | Styrene                     | 1.051 | 1.038 | 1.2  | 103   | 0.00     |
| 71 P  | Bromoform                   | 0.300 | 0.285 | 5.0  | 102   | 0.00     |
| 72 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0  | 100   | 0.00     |
| 73 T  | Isopropylbenzene            | 3.351 | 3.224 | 3.8  | 103   | 0.00     |
| 74 T  | N-amyl acetate              | 1.305 | 1.338 | -2.5 | 103   | 0.00     |
| 75 P  | 1,1,2,2-Tetrachloroethane   | 1.146 | 1.067 | 6.9  | 102   | 0.00     |
| 76 T  | 1,2,3-Trichloropropane      | 1.095 | 1.017 | 7.1  | 89    | 0.00     |
| 77 T  | Bromobenzene                | 0.826 | 0.781 | 5.4  | 105   | 0.00     |
| 78 T  | n-propylbenzene             | 3.854 | 3.805 | 1.3  | 104   | 0.00     |
| 79 T  | 2-Chlorotoluene             | 2.361 | 2.228 | 5.6  | 104   | 0.00     |
| 80 T  | 1,3,5-Trimethylbenzene      | 2.804 | 2.765 | 1.4  | 105   | 0.00     |
| 81 T  | trans-1,4-Dichloro-2-butene | 0.340 | 0.350 | -2.9 | 104   | 0.00     |
| 82 T  | 4-Chlorotoluene             | 2.646 | 2.652 | -0.2 | 106   | 0.00     |
| 83 T  | tert-Butylbenzene           | 2.699 | 2.746 | -1.7 | 105   | 0.00     |
| 84 T  | 1,2,4-Trimethylbenzene      | 2.716 | 2.739 | -0.8 | 103   | 0.00     |
| 85 T  | sec-Butylbenzene            | 3.517 | 3.481 | 1.0  | 101   | 0.00     |
| 86 T  | p-Isopropyltoluene          | 2.892 | 3.041 | -5.2 | 107   | 0.00     |
| 87 T  | 1,3-Dichlorobenzene         | 1.557 | 1.490 | 4.3  | 96    | 0.00     |
| 88 T  | 1,4-Dichlorobenzene         | 1.598 | 1.542 | 3.5  | 100   | 0.00     |
| 89 T  | n-Butylbenzene              | 2.678 | 2.730 | -1.9 | 99    | 0.00     |
| 90 T  | Hexachloroethane            | 0.561 | 0.553 | 1.4  | 102   | 0.00     |
| 91 T  | 1,2-Dichlorobenzene         | 1.639 | 1.485 | 9.4  | 95    | 0.00     |
| 92 T  | 1,2-Dibromo-3-Chloropropane | 0.295 | 0.288 | 2.4  | 101   | 0.00     |
| 93 T  | 1,2,4-Trichlorobenzene      | 1.025 | 1.116 | -8.9 | 107   | 0.00     |
| 94 T  | Hexachlorobutadiene         | 0.427 | 0.453 | -6.1 | 107   | 0.00     |
| 95 T  | Naphthalene                 | 3.569 | 3.681 | -3.1 | 97    | 0.00     |

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MSVOA\_X  
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X082522W.M  
Quant Title : SW846 8260  
QLast Update : Fri Aug 26 02:03:28 2022  
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min  
Max. RRF Dev : 25% Max. Rel. Area : 150%

|      | Compound               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 96 T | 1,2,3-Trichlorobenzene | 1.084 | 1.141 | -5.3 | 102   | 0.00     |

(#) = Out of Range

SPCC's out = 0 CCC's out = 6