

Data Path : Z:\VOASRV\HPCHEM1\MSVOA R\DATA\VR101019\  
 Data File : VR026649.D  
 Acq On : 10 Oct 2019 14:32  
 Operator : SY/MD  
 Sample : VSTD0.564  
 Misc : 25mL/MSVOA R/WATER  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 MSVOA\_R  
 ClientSampleId :  
 VSTD0.564

Manual Integrations  
 APPROVED

MMDadoda  
 10/11/2019 3:30:37 PM

Quant Time: Oct 11 06:23:18 2019  
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_R\METHODS\SOMRTR101019WMA.M  
 Quant Title : TRACE VOA SOM01.0  
 QLast Update : Fri Oct 11 05:48:01 2019  
 Response via : Initial Calibration

| Internal Standards         | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|------|-------|----------|
| 1) 1,4-Difluorobenzene     | 8.48  | 114  | 194531   | 5.00 | ug/L  | 0.00     |
| 28) Chlorobenzene-d5       | 11.30 | 117  | 163393   | 5.00 | ug/L  | 0.00     |
| 60) 1,4-Dichlorobenzene-d4 | 13.24 | 152  | 84231    | 5.00 | ug/L  | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |      |      |       |
|--------------------------------|-------|-----|--------|------|------|-------|
| 4) Vinyl Chloride-d3           | 2.20  | 65  | 7675   | 1.16 | ug/L | 0.01  |
| 7) Chloroethane-d5             | 2.64  | 69  | 6370   | 1.14 | ug/L | 0.00  |
| 11) 1,1-Dichloroethene-d2      | 3.63  | 63  | 15841m | 0.91 | ug/L | -0.02 |
| 20) 2-Butanone-d5              | 6.63  | 46  | 7218   | 5.93 | ug/L | 0.00  |
| 24) Chloroform-d               | 7.23  | 84  | 20455  | 0.85 | ug/L | 0.00  |
| 26) 1,2-Dichloroethane-d4      | 7.91  | 65  | 6813   | 0.86 | ug/L | 0.00  |
| 32) Benzene-d6                 | 7.88  | 84  | 33997  | 0.71 | ug/L | 0.00  |
| 36) 1,2-Dichloropropane-d6     | 8.91  | 67  | 8765   | 0.72 | ug/L | 0.00  |
| 41) Toluene-d8                 | 9.98  | 98  | 23600  | 0.63 | ug/L | 0.00  |
| 43) trans-1,3-Dichloropropene- | 10.24 | 79  | 2929   | 1.03 | ug/L | 0.00  |
| 46) 2-Hexanone-d5              | 10.60 | 63  | 3998   | 4.91 | ug/L | 0.00  |
| 57) 1,1,2,2-Tetrachloroethane- | 12.37 | 84  | 4753   | 0.76 | ug/L | 0.00  |
| 64) 1,2-Dichlorobenzene-d4     | 13.53 | 152 | 8454   | 0.72 | ug/L | 0.00  |

## Target Compounds

|                                |      |     |        |              | Qvalue |
|--------------------------------|------|-----|--------|--------------|--------|
| 2) Dichlorodifluoromethane     | 1.86 | 85  | 16301  | 0.676 ug/L   | 98     |
| 3) Chloromethane               | 2.07 | 50  | 8866   | 0.721 ug/L   | 93     |
| 5) Vinyl chloride              | 2.18 | 62  | 7854m  | 0.641 ug/L   |        |
| 6) Bromomethane                | 2.53 | 94  | 5417   | 0.687 ug/L   | 89     |
| 8) Chloroethane                | 2.69 | 64  | 4598   | 0.653 ug/L   | 91     |
| 9) Trichlorofluoromethane      | 2.99 | 101 | 15881  | 0.729 ug/L   | 98     |
| 10) 1,1,2-Trichloro-1,2,2-trif | 3.68 | 101 | 6504m  | 0.692 ug/L   |        |
| 12) 1,1-Dichloroethene         | 3.65 | 96  | 5757   | 0.631 ug/L # | 65     |
| 13) Acetone                    | 3.73 | 43  | 6220   | 6.779 ug/L   | 92     |
| 14) Carbon disulfide           | 3.97 | 76  | 38242  | 0.853 ug/L   | 98     |
| 15) Methyl Acetate             | 4.22 | 43  | 2150   | 0.563 ug/L   | 94     |
| 16) Methylene chloride         | 4.43 | 84  | 15533m | 0.816 ug/L   |        |
| 17) Methyl tert-butyl Ether    | 4.93 | 73  | 9951   | 0.529 ug/L # | 84     |
| 18) trans-1,2-Dichloroethene   | 4.92 | 96  | 10963  | 0.578 ug/L   | 88     |
| 19) 1,1-Dichloroethane         | 5.72 | 63  | 20162  | 0.581 ug/L   | 91     |
| 21) 2-Butanone                 | 6.72 | 43  | 7015   | 4.295 ug/L   | 89     |
| 22) cis-1,2-Dichloroethene     | 6.70 | 96  | 8888   | 0.494 ug/L   | 99     |
| 23) Bromochloromethane         | 7.10 | 128 | 3612   | 0.710 ug/L # | 47     |
| 25) Chloroform                 | 7.25 | 83  | 23024  | 0.682 ug/L   | 100    |
| 27) 1,2-Dichloroethane         | 8.01 | 62  | 8254   | 0.684 ug/L # | 91     |
| 29) 1,1,1-Trichloroethane      | 7.45 | 97  | 14164  | 0.525 ug/L   | 98     |
| 30) Cyclohexane                | 7.56 | 56  | 9560   | 0.325 ug/L # | 78     |
| 31) Carbon tetrachloride       | 7.66 | 117 | 13803  | 0.590 ug/L   | 95     |
| 33) Benzene                    | 7.93 | 78  | 36618  | 0.483 ug/L   | 100    |
| 34) Trichloroethene            | 8.72 | 95  | 10275  | 0.546 ug/L   | 86     |
| 35) Methylcyclohexane          | 8.98 | 83  | 12206  | 0.423 ug/L   | 95     |
| 37) 1,2-Dichloropropane        | 9.01 | 63  | 7720   | 0.513 ug/L   | 99     |
| 38) Bromodichloromethane       | 9.30 | 83  | 8628   | 0.530 ug/L   | 95     |
| 39) cis-1,3-Dichloropropene    | 9.73 | 75  | 5834   | 0.365 ug/L   | 95     |
| 40) 4-Methyl-2-pentanone       | 9.89 | 43  | 11942  | 3.227 ug/L   | 99     |

Data Path : Z:\VOASRV\HPCHEM1\MSVOA R\DATA\VR101019\  
 Data File : VR026649.D  
 Acq On : 10 Oct 2019 14:32  
 Operator : SY/MD  
 Sample : VSTD0.564  
 Misc : 25mL/MSVOA R/WATER  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 MSVOA\_R  
 ClientSampleId :  
 VSTD0.564

Manual Integrations  
 APPROVED

MMDadoda  
 10/11/2019 3:30:37 PM

Quant Time: Oct 11 06:23:18 2019  
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_R\METHODS\SOMRTR101019WMA.M  
 Quant Title : TRACE VOA SOM01.0  
 QLast Update : Fri Oct 11 05:48:01 2019  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| -----                          |       |      |          |       |        |          |
| 42) Toluene                    | 10.05 | 91   | 27911    | 0.419 | ug/L   | 81       |
| 44) trans-1,3-Dichloropropene  | 10.28 | 75   | 4852     | 0.465 | ug/L   | 100      |
| 45) 1,1,2-Trichloroethane      | 10.46 | 97   | 3475     | 0.533 | ug/L # | 86       |
| 47) Tetrachloroethene          | 10.53 | 164  | 6798     | 0.596 | ug/L # | 83       |
| 48) 2-Hexanone                 | 10.65 | 43   | 9192     | 3.892 | ug/L   | 92       |
| 49) Dibromochloromethane       | 10.80 | 129  | 4042     | 0.569 | ug/L   | 81       |
| 50) 1,2-Dibromoethane          | 10.91 | 107  | 2831     | 0.501 | ug/L # | 72       |
| 51) Chlorobenzene              | 11.33 | 112  | 19204    | 0.498 | ug/L   | 92       |
| 52) Ethylbenzene               | 11.41 | 91   | 30876    | 0.412 | ug/L   | 96       |
| 53) m,p-Xylene                 | 11.52 | 106  | 10570    | 0.391 | ug/L   | 97       |
| 54) o-Xylene                   | 11.84 | 106  | 9254     | 0.349 | ug/L   | 89       |
| 55) Styrene                    | 11.86 | 104  | 13015    | 0.343 | ug/L   | 100      |
| 56) Isopropylbenzene           | 12.15 | 105  | 26476    | 0.353 | ug/L   | 92       |
| 58) 1,1,2,2-Tetrachloroethane  | 12.40 | 83   | 4504     | 0.594 | ug/L # | 96       |
| 59) 1,2,3-Trichloropropane     | 12.45 | 75   | 3315     | 0.615 | ug/L # | 66       |
| 61) Bromoform                  | 12.02 | 173  | 1855     | 0.551 | ug/L # | 69       |
| 62) 1,3-Dichlorobenzene        | 13.18 | 146  | 13578    | 0.445 | ug/L   | 95       |
| 63) 1,4-Dichlorobenzene        | 13.26 | 146  | 16770    | 0.550 | ug/L   | 98       |
| 65) 1,2-Dichlorobenzene        | 13.54 | 146  | 12204    | 0.475 | ug/L   | 92       |
| 66) 1,2-Dibromo-3-chloropropan | 14.16 | 75   | 764      | 0.587 | ug/L # | 77       |
| 67) 1,3,5-Trichlorobenzene     | 14.31 | 180  | 10419    | 0.427 | ug/L   | 98       |
| 68) 1,2,4-trichlorobenzene     | 14.80 | 180  | 6255     | 0.365 | ug/L   | 87       |
| 69) Naphthalene                | 15.01 | 128  | 5591     | 0.250 | ug/L # | 82       |
| 70) 1,2,3-Trichlorobenzene     | 15.20 | 180  | 7378     | 0.474 | ug/L   | 89       |
| -----                          |       |      |          |       |        |          |

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

Data Path : Z:\VOASRV\HPCHEM1\MSVOA R\DATA\VR101019\  
Data File : VR026649.D  
Acq On : 10 Oct 2019 14:32  
Operator : SY/MD  
Sample : VSTD0.564  
Misc : 25mL/MSVOA R/WATER  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
MSVOA\_R  
ClientSampleId :  
VSTD0.564

Manual Integrations  
APPROVED

MMDadoda  
10/11/2019 3:30:37 PM

Quant Time: Oct 11 06:23:18 2019  
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_R\METHODS\SOMRTR101019WMA.M  
Quant Title : TRACE VOA SOM01.0  
QLast Update : Fri Oct 11 05:48:01 2019  
Response via : Initial Calibration

