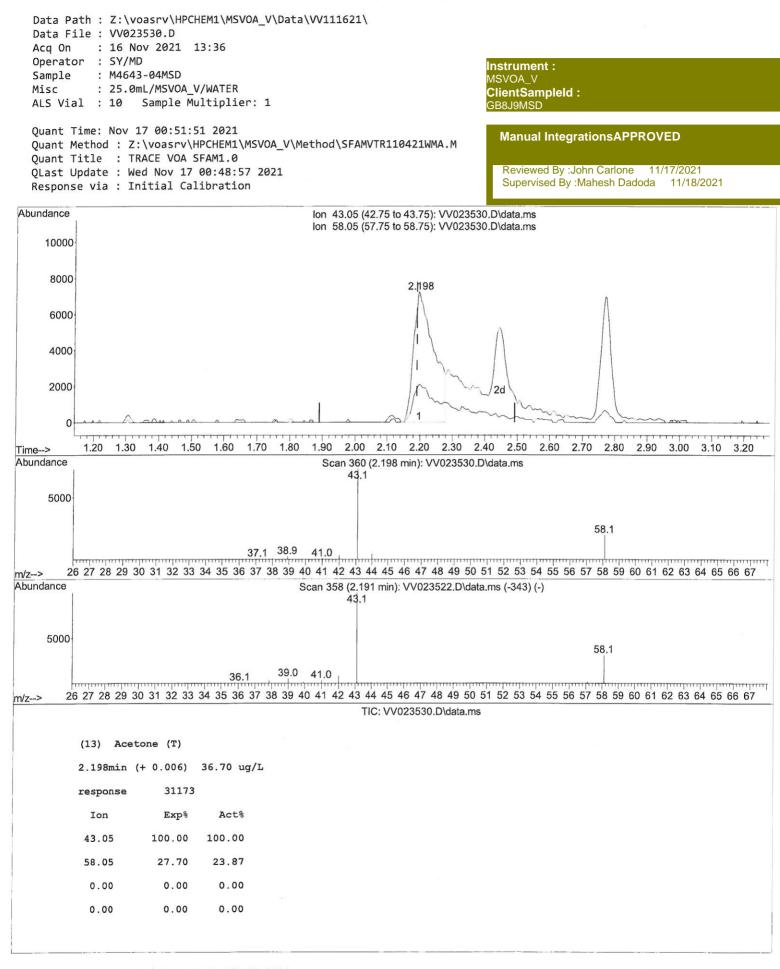
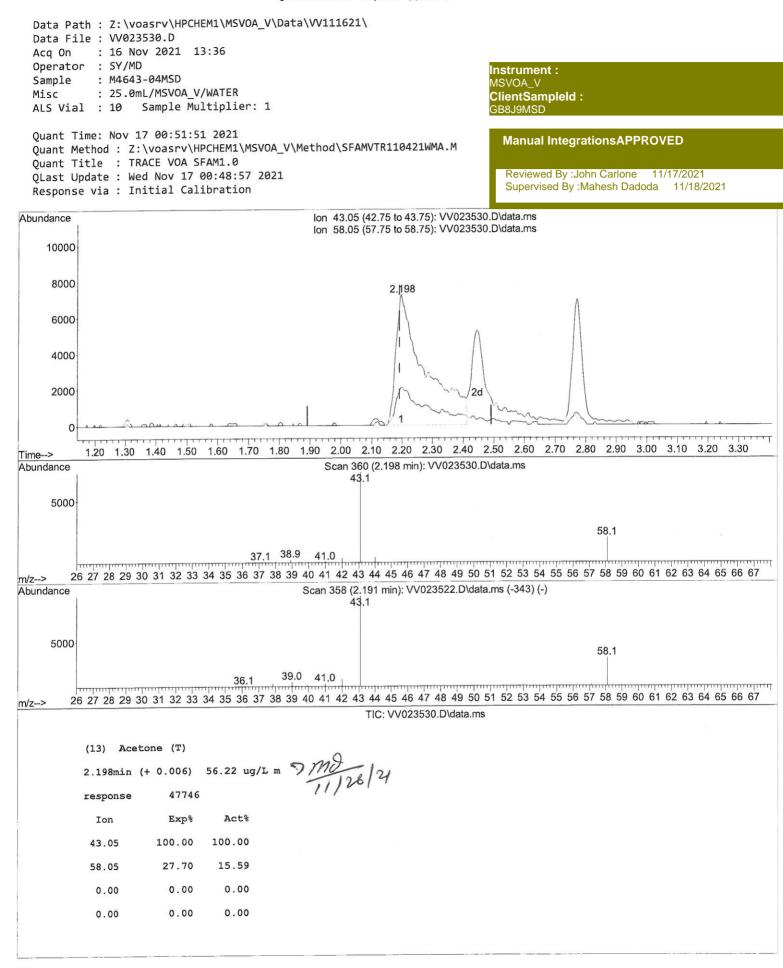
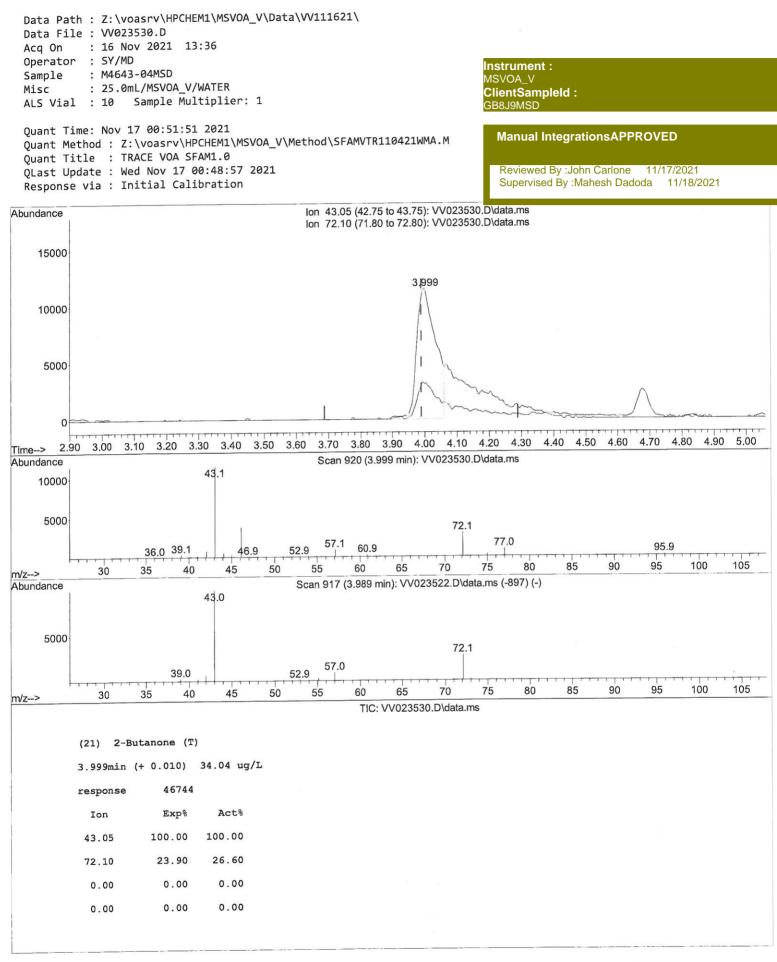


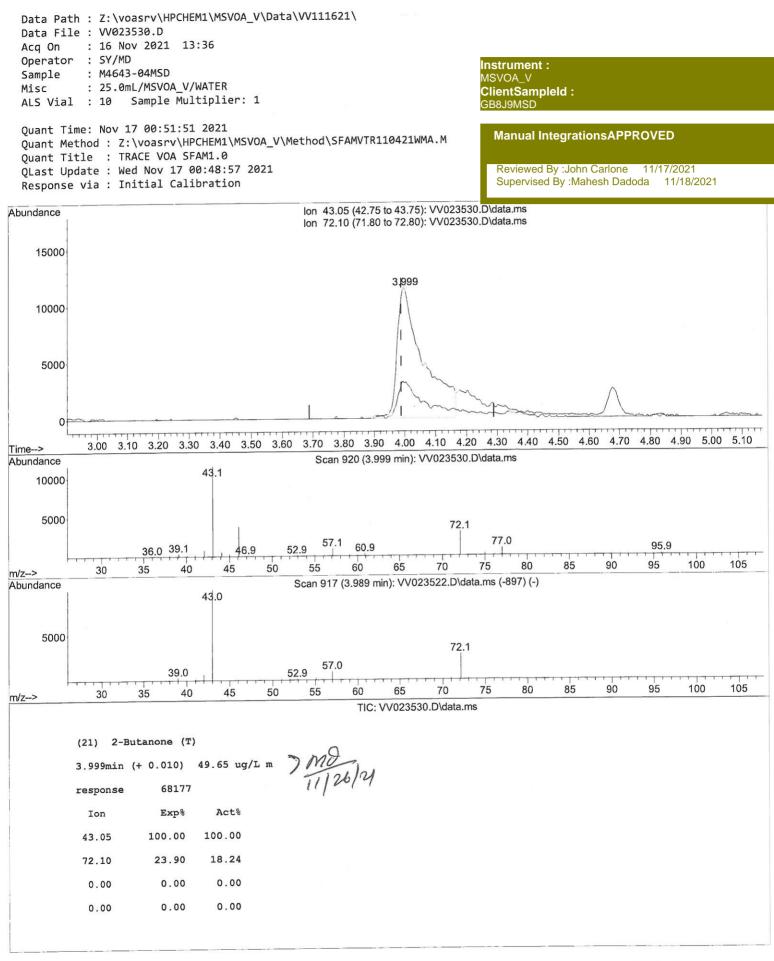
SFAMVTR110421WMA.M Wed Nov 17 01:28:26 2021







SFAMVTR110421WMA.M Wed Nov 17 01:27:24 2021



SFAMVTR110421WMA.M Wed Nov 17 01:27:39 2021

1

| Data Path : Z:\voasrv\HPCHEM1 | MSVOA_V\Da | ta\VV1 | 11621\ | | | |
|--|-------------------|--------------|-------------------|---|----------|---|
| Data File : VV023530.D | | | | | | |
| Acq On : 16 Nov 2021 13:30 | 5 | | | | | |
| Operator : SY/MD | | Instrument : | | | | |
| Sample : M4643-04MSD | TED | MSVOA_V | | | | |
| Misc : 25.0mL/MSVOA_V/WA ALS Vial : 10 Sample Multip | | | | | | ClientSampleId : |
| ALS VIAL . 10 Sample Harel | | | | | | GB8J9MSD |
| Quant Time: Nov 17 00:51:51 20 | 921 | | | | | |
| Quant Method : Z:\voasrv\HPCH | EM1\MSVOA_V | \Metho | d\SFAMVTR11 | 10421WMA | . M | Manual IntegrationsAPPROVED |
| Quant Title : TRACE VOA SFAM | 1.0 | | | | | |
| QLast Update : Wed Nov 17 00:4 | 48:57 2021 | | | | | Reviewed By :John Carlone 11/17/2021 |
| Response via : Initial Calibra | ation | | | | | Supervised By :Mahesh Dadoda 11/18/2021 |
| | | 0.7 | D | Conc Uni | to Dou(A | 110 |
| Compound | R.I. | QION | Response (| LOUC OUT | LS Dev(| 111) |
| Internal Standards | | | | | | |
| 1) 1,4-Difluorobenzene | 5.619 | 114 | 128790 | 5.000 | ug/L | 0.00 |
| 28) Chlorobenzene-d5 | 8.854 | | 126220 | 5.000 | | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 11.249 | | 68679 | 5.000 | ug/L | 0.00 |
| 56) 1,4 516110 555111 | | | | | 170 | |
| System Monitoring Compounds | | | | | | |
| 4) Vinyl Chloride-d3 | 1.307 | 65 | 34818 | 4.315 | ug/L | 0.00 |
| Spiked Amount 5.000 | Range 40 | - 130 | Recovery | | 86.400% | |
| Chloroethane-d5 | 1.568 | 69 | 30147 | 4.585 | | 0.00 |
| Spiked Amount 5.000 | Range 65 | | Recovery | | 91.600% | 0.00 |
| 11) 1,1-Dichloroethene-d2 | 2.111 | 63 | 66679 Decever | 4.415 | | 0.00 |
| Spiked Amount 5.000 | Range 60 | - 125 46 | Recovery 61533 | y = 44.268 | 88.200% | 0.00 |
| 20) 2-Butanone-d5 | 3.908 Range 40 | | Recovery | | 88.540% | 0.00 |
| Spiked Amount 50.000 24) Chloroform-d | 4.349 | 84 | 75506 | 4.391 | | 0.00 |
| Spiked Amount 5.000 | Range 70 | | Recovery | | 87.800% | |
| 26) 1,2-Dichloroethane-d4 | 5.034 | 65 | 35697 | 4.617 | | 0.00 |
| Spiked Amount 5.000 | Range 70 | - 130 | Recovery | y = | 92.400% | |
| 32) Benzene-d6 | 5.050 | 84 | 148190 | 4.576 | | 0.00 |
| Spiked Amount 5.000 | Range 70 | - 125 | Recovery | | 91.600% | |
| 36) 1,2-Dichloropropane-d6 | 6.072 | 67 | 43254 | 4.537 | | 0.00 |
| Spiked Amount 5.000 | Range 60 | | Recovery | The second se | 90.800% | 0.00 |
| 41) Toluene-d8 | 7.317 | 98 120 | 144583 Bacovon | 4.764 v = | 95.200% | 0.00 |
| Spiked Amount 5.000 | Range 70 | - 150 | Recovery 17390 | 4.811 | | 0.00 |
| <pre>43) trans-1,3-Dichloroprop. Spiked Amount 5.000</pre> | Range 55 | | Recovery | | 96.200% | |
| 46) 2-Hexanone-d5 | 8.092 | 63 | 62401 | 46.917 | | 0.00 |
| Spiked Amount 50.000 | Range 45 | | Recovery | | 93.840% | |
| 56) 1,1,2,2-Tetrachloroeth. | | 84 | 31782 | 4.636 | ug/L | 0.00 |
| Spiked Amount 5.000 | Range 65 | | Recovery | | 92.800% | |
| 66) 1,2-Dichlorobenzene-d4 | 11.625 | 152 | 54458 | 4.762 | | 0.00 |
| Spiked Amount 5.000 | Range 80 | - 120 | Recovery | y = | 95.200% | |
| | | | | | Qva. | |
| Target Compounds | 1 1 20 | ØE | 51907 | 4.133 | 100 C | 99 |
| 2) Dichlorodifluoromethane | 1.130 1.240 | 85 50 | 46857 | 4.388 | | 97 |
| 3) Chloromethane 5) Vinyl chloride | 1.311 | 62 | 48303 | 4.530 | | 98 |
| 6) Bromomethane | 1.523 | 94 | 28272 | 4.148 | | 97 |
| 8) Chloroethane | 1.584 | 64 | 28384 | 4.612 | ug/L | 96 |
| 9) Trichlorofluoromethane | 1.754 | 101 | 74863 | 4.672 | ug/L | 99 |
| 10) 1,1,2-Trichloro-1,2,2 | 2.118 | 101 | 38391 | 4.759 | | 96 |
| 12) 1,1-Dichloroethene | 2.118 | 96 | 35066 | 4.566 | | 95 |
| 13) Acetone | 2.198 | 43 | 47746m | 56.216 | | and mile |
| 14) Carbon disulfide | 2.294 | 76 | 118627 | 4.093 | | 99 11/26/21 |
| 15) Methyl Acetate | 2.446 | 43 | 10212 41731 | 4.248 3.723 | | 96 |
| 16) Methylene chloride | 2.507 2.770 | 84 73 | 80728 | 4.775 | | 94 |
| 17) Methyl tert-butyl Ether 18) trans-1,2-Dichloroethen | e 2.761 | 96 | 42905 | 4.544 | | 97 |
| 19) 1,1-Dichloroethane | 3.191 | 63 | 72881 | 4.572 | | 96 |
| 21) 2-Butanone | 3.999 | 43 | 68177m | 49.649 | | 7 MD |
| 22) cis-1,2-Dichloroethene | 3.912 | 96 | 41919 | | ug/L # | 93 MD 84 11/26/21 |
| 23) Bromochloromethane | 4.249 | 128 | 19720 | 4.707 | ug/L # | 84 [/] / |
| errorate S zoency personality | | | | | | |

SFAMVTR110421WMA.M Wed Nov 17 01:28:25 2021

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111621\ Data File : VV023530.D Acg On : 16 Nov 2021 13:36 Operator : SY/MD Sample : M4643-04MSD : 25.0mL/MSVOA_V/WATER Misc ALS Vial : 10 Sample Multiplier: 1

Quant Time: Nov 17 00:51:51 2021 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M Quant Title : TRACE VOA SFAM1.0 QLast Update : Wed Nov 17 00:48:57 2021

Instrument : MSVOA_V ClientSampleId : GB8J9MSD

Manual IntegrationsAPPROVED

Reviewed By : John Carlone 11/17/2021 Supervised By :Mahesh Dadoda 11/18/2021

Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc Units Dev | v(Min) |
|---|--------|------|----------|----------------|--------|
| 25) Chloroform | 4.375 | 83 | 82189 | 4.837 ug/L | 99 |
| 27) 1,2-Dichloroethane | 5.134 | 62 | 41784 | 4.623 ug/L | 96 |
| 29) 1,1,1-Trichloroethane | 4.609 | 97 | 71833 | 4.686 ug/L | 99 |
| 30) Cyclohexane | 4.680 | 56 | 61006 | 4.441 ug/L | 97 |
| 31) Carbon tetrachloride | 4.828 | 117 | 65189 | 4.735 ug/L | 97 |
| 33) Benzene | 5.101 | 78 | 167599 | 4.751 ug/L | 100 |
| 34) Trichloroethene | 5.915 | 95 | 43851 | 4.674 ug/L | 99 |
| 35) Methylcyclohexane | 6.130 | 83 | 66840 | 4.514 ug/L | 95 |
| 37) 1,2-Dichloropropane | 6.175 | 63 | 38801 | 4.711 ug/L | 98 |
| 38) Bromodichloromethane | 6.510 | 83 | 51360 | 4.653 ug/L | 99 |
| 39) cis-1,3-Dichloropropene | 7.031 | 75 | 56413 | 4.762 ug/L | 96 |
| 40) 4-Methyl-2-pentanone | 7.230 | 43 | 204618 | 53.568 ug/L | 97 |
| 42) Toluene | 7.387 | 91 | 187330 | 4.965 ug/L | 98 |
| 44) trans-1,3-Dichloropropene | 7.651 | 75 | 47673 | 4.850 ug/L | 97 |
| 45) 1,1,2-Trichloroethane | 7.841 | 97 | 28087 | 4.746 ug/L | 98 |
| 47) Tetrachloroethene | 7.976 | 164 | 75994 | 9.347 ug/L | 98 |
| 48) 2-Hexanone | 8.143 | 43 | 148221 | 55.377 ug/L | 97 |
| 49) Dibromochloromethane | 8.246 | 129 | 36687 | 4.893 ug/L | 99 |
| 50) 1,2-Dibromoethane | 8.355 | 107 | 26503 | 4.833 ug/L | 97 |
| 51) Chlorobenzene | 8.883 | 112 | 117059 | 4.667 ug/L | 97 |
| 52) Ethylbenzene | 9.011 | 91 | 188176 | 4.728 ug/L | 98 |
| 53) m,p-xylene | 9.140 | 106 | 74624 | 4.778 ug/L | 96 |
| 54) o-xylene | 9.545 | 106 | 71549 | 4.883 ug/L | 98 |
| 55) Styrene | 9.561 | 104 | 123071 | 4.903 ug/L | 98 |
| 57) 1,1,2,2-Tetrachloroethane | 10.243 | 83 | 30030 | 4.632 ug/L | 97 |
| 59) Bromoform | 9.731 | 173 | 20259 | 4.939 ug/L | 99 |
| 60) Isopropylbenzene | 9.931 | 105 | 192837 | 4.893 ug/L | 100 |
| 61) 1,2,3-Trichloropropane | 10.275 | 75 | 22397 | 4.909 ug/L | 97 |
| 62) 1,3,5-Trimethylbenzene | 10.538 | 105 | 160272 | 4.905 ug/L | 99 |
| 63) 1,2,4-Trimethylbenzene | 10.915 | 105 | 161645 | 4.970 ug/L | 98 |
| 64) 1,3-Dichlorobenzene | 11.182 | 146 | 97596 | 4.847 ug/L | 99 |
| 65) 1,4-Dichlorobenzene | 11.272 | 146 | 96693 | 4.702 ug/L | 99 |
| 67) 1,2-Dichlorobenzene | 11.641 | 146 | 89086 | 4.944 ug/L | 99 |
| 68) 1,2-Dibromo-3-chloropr | 12.429 | 75 | 4370 | 4.496 ug/L # | |
| 69) 1,3,5-Trichlorobenzene | 12.644 | 180 | 74156 | 4.703 ug/L | 100 |
| 70) 1,2,4-trichlorobenzene | 13.262 | 180 | 54923 | 4.350 ug/L | 99 |
| 71) Naphthalene | 13.503 | 128 | 78651 | 4.225 ug/L | 99 |
| 72) 1,2,3-Trichlorobenzene | 13.744 | 180 | 50514 | 4.573 ug/L | 98 |
| | | | | | |

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