Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120921\

Data File: VV023875.D

Acq On : 09 Dec 2021 20:40

Operator : SY/MD Sample : VSTDCCC005EC

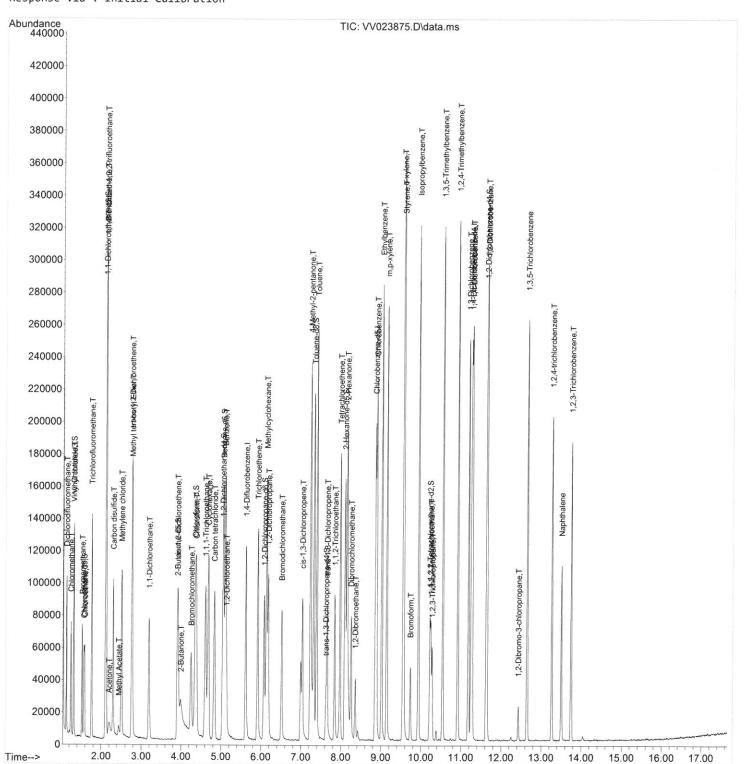
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

Quant Time: Dec 10 03:27:06 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Thu Dec 02 02:08:23 2021 Response via : Initial Calibration Instrument : MSVOA_V LabSampleId : VSTDCCC005EC

Manual IntegrationsAPPROVED



Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV120921\

Data File: VV023875.D

Acq On : 09 Dec 2021 20:40

Operator : SY/MD Sample : VSTDCCC005EC

Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

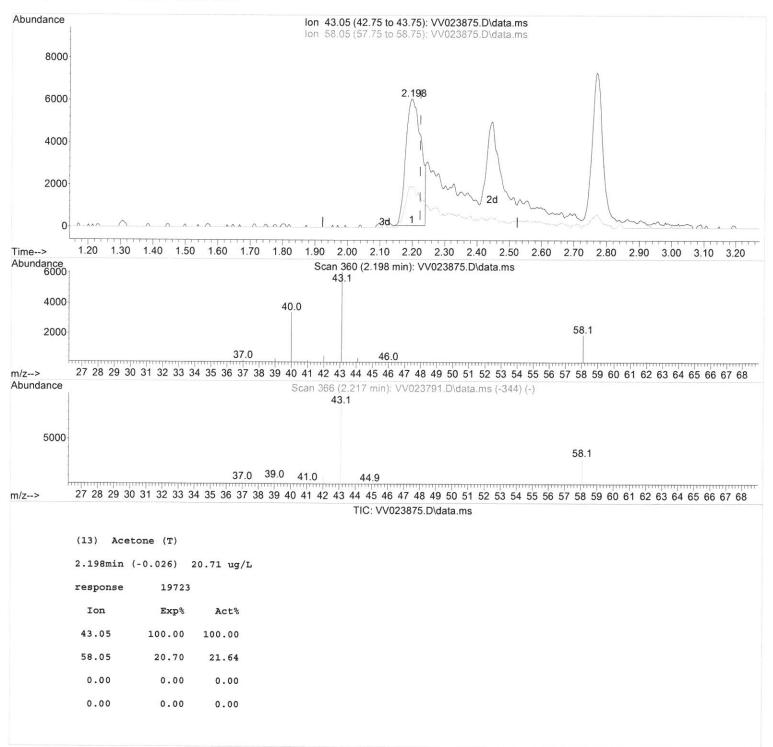
Quant Time: Dec 10 03:27:06 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Thu Dec 02 02:08:23 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED



Quantitation Report (Qedit)

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Data File: VV023875.D

Acq On : 09 Dec 2021 20:40

Operator : SY/MD Sample : VSTDCCC005EC

Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 21 Sample Multiplier: 1

Quant Time: Dec 10 03:27:06 2021

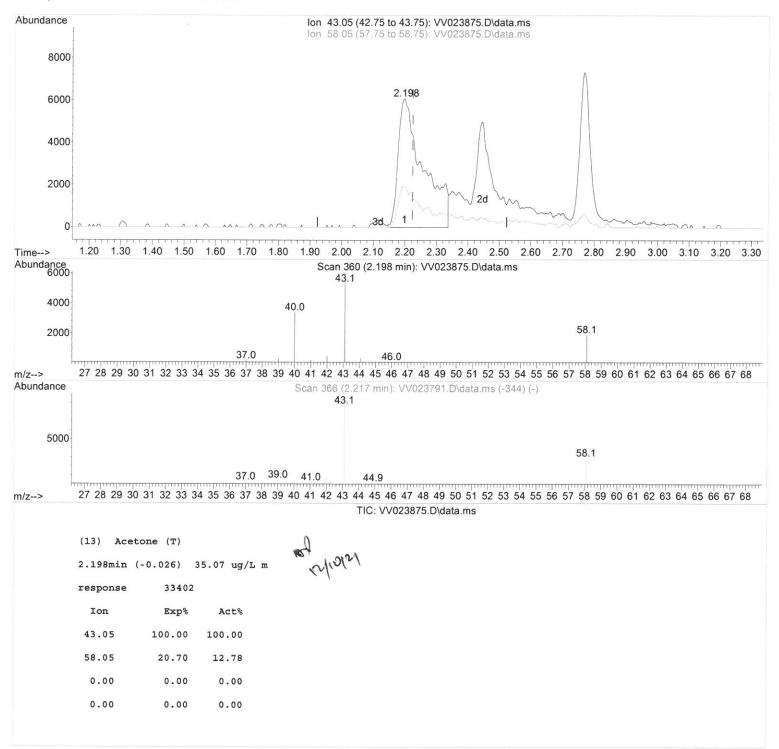
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.M

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Manual IntegrationsAPPROVED



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Acq On : 09 Dec 2021 20:40

Operator : SY/MD

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Quant Title : TRACE VOA SFAM1.0 QLast Update : Thu Dec 02 02:08:23 2021 Response via : Initial Calibration

Instrument : MSVOA_V LabSampleId : VSTDCCC005EC

Manual IntegrationsAPPROVED

Reviewed By :John Carlone 12/10/2021 Supervised By: Mahesh Dadoda 12/10/2021

| Compound | R.T. | QIon | Response Conc Units Dev(Mir | 1) |
|--|-------------------|-------------|--|----------|
| Internal Standards | | | | |
| 1,4-Difluorobenzene | 5.619 | 114 | 107257 5.000 ug/L 0 | 0.00 |
| 28) Chlorobenzene-d5 | 8.854 | 117 | <u> </u> | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 11.249 | 152 | 60474 5.000 ug/L 0 | 9.00 |
| System Monitoring Compounds | | | | |
| 4) Vinyl Chloride-d3 | 1.307 | 65 | 37917 4.306 ug/L 0. | .00 |
| Spiked Amount 5.000 | Range 40 | | Recovery = 86.200% | |
| 7) Chloroethane-d5 | 1.568 | 69 | | .00 |
| Spiked Amount 5.000 | Range 65 | | Recovery = 90.800% | 200 |
| 11) 1,1-Dichloroethene-d2 | 2.111 | 63 | 5. | .00 |
| Spiked Amount 5.000 | Range 60 | | Recovery = 88.400% | |
| 20) 2-Butanone-d5 | 3.905 | 46 | <u>.</u> | 00 |
| Spiked Amount 50.000 | O | - 130 | Recovery = 90.160% | 2.2 |
| 24) Chloroform-d | 4.352 | 84 | 9. | 00 |
| Spiked Amount 5.000 | 0 | - 125 | Recovery = 105.400% | |
| 26) 1,2-Dichloroethane-d4 | 5.037 | 65 | 0, | 99 |
| Spiked Amount 5.000 | Range 70 | | Recovery = 102.800% | |
| 32) Benzene-d6 | 5.053 | 84 | 142540 4.991 ug/L 0. | 99 |
| Spiked Amount 5.000 | Range 70 | | Recovery = 99.800% | 00 |
| 36) 1,2-Dichloropropane-d6 | 6.072 | 67 | 42657 5.328 ug/L 0. | 99 |
| Spiked Amount 5.000 41) Toluene-d8 | Range 60 | | Recovery = 106.600% | ~~ |
| | 7.317 | 98 | 134656 5.046 ug/L 0. | 90 |
| Spiked Amount 5.000 43) trans-1,3-Dichloroprop. | | - 130 | Recovery = 101.000% | 00 |
| Spiked Amount 5.000 | | 79 | 17523 5.430 ug/L 0. | 90 |
| 46) 2-Hexanone-d5 | • | - 130 | Recovery = 108.600% | 00 |
| Spiked Amount 50.000 | 8.092 Range 45 | 63 - 130 | 77072 71.880 ug/L 0. Recovery = 143.760%# | 99 |
| 56) 1,1,2,2-Tetrachloroeth. | | 84 | | aa |
| Spiked Amount 5.000 | Range 65 | | 33558 5.825 ug/L 0.0 Recovery = 116.400% | 99 |
| 66) 1,2-Dichlorobenzene-d4 | 11.625 | 152 | 55169 5.160 ug/L 0.0 | aa |
| Spiked Amount 5.000 | Range 80 | | Recovery = 103.200% | 00 |
| Target Compounds | | | Ovalue | |
| 2) Dichlorodifluoromethane | 1.130 | 85 | c | 99 |
| 3) Chloromethane | 1.240 | 50 | G. | 97 |
| 5) Vinyl chloride | 1.310 | 62 | O. | 98 |
| 6) Bromomethane | 1.523 | 94 | | 99 |
| 8) Chloroethane | 1.587 | 64 | | 96 |
| 9) Trichlorofluoromethane | 1.754 | 101 | | 00 |
| 10) 1,1,2-Trichloro-1,2,2 | | 101 | | 98 |
| 12) 1,1-Dichloroethene | 2.121 | 96 | | 98 |
| 13) Acetone | 2.198 | 43 | 33402m /35.069 ug/L | |
| 14) Carbon disulfide | 2.298 | 76 | | 99 |
| 15) Methyl Acetate | 2.445 | 43 | | 98 |
| 16) Methylene chloride | 2.507 | 84 | | 97 |
| 17) Methyl tert-butyl Ether | 2.770 | 73 | | 99 |
| 18) trans-1,2-Dichloroethene | 2.764 | 96 | | 99 |
| 19) 1,1-Dichloroethane | 3.191 | 63 | 0. | 98 |
| / -j- brenzen de chane | 2 005 | 43 | | 96 |
| 21) 2-Butanone | 3.995 | 43 | 303/3 32.333 ug/L | |
| | 3.995 | 96 | . | 98 |
| 21) 2-Butanone | | | 42797 5.452 ug/L | 98 94 |

MUSIN

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120921\

Data File : VV023875.D

Acq On : 09 Dec 2021 20:40 Operator : SY/MD

Sample : VSTDCCC005EC
Misc : 25.0mL/MSVOA_V/WATER ALS Vial : 21 Sample Multiplier: 1

Quant Time: Dec 10 03:27:06 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.M

Quant Title : TRACE VOA SFAM1.0

QLast Update : Thu Dec 02 02:08:23 2021 Response via : Initial Calibration

Instrument : MSVOA_V LabSampleId : VSTDCCC005EC

Manual IntegrationsAPPROVED

| Compound | R.T. | QIon | Response | Conc Units Dev | (Min) |
|---------------------------------|--------|------|----------|----------------|-------|
| 27) 1,2-Dichloroethane | 5.133 | 62 | 43552 | 5.343 ug/L | 97 |
| 29) 1,1,1-Trichloroethane | 4.609 | 97 | 75949 | 5.544 ug/L | 99 |
| 30) Cyclohexane | 4.680 | 56 | 60135 | 5.260 ug/L | 99 |
| 31) Carbon tetrachloride | 4.828 | 117 | 68090 | 5.426 ug/L | 98 |
| 33) Benzene | 5.101 | 78 | 161785 | 5.413 ug/L | 100 |
| 34) Trichloroethene | 5.915 | 95 | 43414 | 5.422 ug/L | 97 |
| 35) Methylcyclohexane | 6.130 | 83 | 66672 | 5.341 ug/L | 98 |
| 37) 1,2-Dichloropropane | 6.175 | 63 | 37597 | 5.292 ug/L | 99 |
| 38) Bromodichloromethane | 6.510 | 83 | 53024 | 5.499 ug/L | 97 |
| 39) cis-1,3-Dichloropropene | 7.027 | 75 | 53805 | 5.322 ug/L | 100 |
| 40) 4-Methyl-2-pentanone | 7.227 | 43 | 193646 | 57.099 ug/L | 98 |
| 42) Toluene | 7.387 | 91 | 188518 | 5.816 ug/L | 98 |
| 44) trans-1,3-Dichloropropene | 7.651 | 75 | 46848 | 5.513 ug/L | 90 |
| 45) 1,1,2-Trichloroethane | 7.841 | 97 | 28035 | 5.699 ug/L | 98 |
| 47) Tetrachloroethene | 7.976 | 164 | 39893 | 5.474 ug/L | 96 |
| 48) 2-Hexanone | 8.143 | 43 | 139976 | 55.842 ug/L | 98 |
| 49) Dibromochloromethane | 8.246 | 129 | 38715 | 5.743 ug/L | 96 |
| 50) 1,2-Dibromoethane | 8.352 | 107 | 27140 | 5.658 ug/L | 97 |
| 51) Chlorobenzene | 8.882 | 112 | 121035 | 5.632 ug/L | 99 |
| 52) Ethylbenzene | 9.011 | 91 | 194303 | 5.735 ug/L | 98 |
| 53) m,p-xylene | 9.140 | 106 | 77633 | 5.757 ug/L | 100 |
| 54) o-xylene | 9.542 | 106 | 74589 | 5.816 ug/L | 98 |
| 55) Styrene | 9.561 | 104 | 129862 | 6.010 ug/L | 99 |
| 57) 1,1,2,2-Tetrachloroethane | 10.243 | 83 | 30555 | 5.583 ug/L | 99 |
| 59) Bromoform | 9.731 | 173 | 20877 | 5.227 ug/L | 98 |
| <pre>60) Isopropylbenzene</pre> | 9.931 | 105 | 204101 | 5.652 ug/L | 100 |
| 61) 1,2,3-Trichloropropane | 10.275 | 75 | 22355 | 5.212 ug/L | 96 |
| 62) 1,3,5-Trimethylbenzene | 10.538 | 105 | 168542 | 5.609 ug/L | 99 |
| 63) 1,2,4-Trimethylbenzene | 10.915 | 105 | 170366 | 5.734 ug/L | 99 |
| 64) 1,3-Dichlorobenzene | 11.181 | 146 | 103000 | 5.579 ug/L | 98 |
| 65) 1,4-Dichlorobenzene | 11.271 | 146 | 101366 | 5.463 ug/L | 100 |
| 67) 1,2-Dichlorobenzene | 11.641 | 146 | 92683 | 5.485 ug/L | 98 |
| 68) 1,2-Dibromo-3-chloropr | 12.429 | 75 | 5076 | 5.960 ug/L | 95 |
| 69) 1,3,5-Trichlorobenzene | 12.644 | 180 | 81066 | 5.624 ug/L | 100 |
| 70) 1,2,4-trichlorobenzene | 13.262 | 180 | 61827 | 5.528 ug/L | 98 |
| 71) Naphthalene | 13.503 | 128 | 86880 | 5.772 ug/L | 99 |
| 72) 1,2,3-Trichlorobenzene | 13.744 | 180 | 55670 | 5.745 ug/L | 99 |
| | | | | | |

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