

Method Path : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\

Method File : SOMVTR022820WMA.M

Title : TRACE VOA SOM01.0

Last Update : Fri Feb 28 23:00:21 2020

Response Via : Initial Calibration

Calibration Files

| | | |
|-----------------|----------------|---------------|
| 0.5 =VV014710.D | 1 =VV014711.D | 5 =VV014712.D |
| 10 =VV014713.D | 20 =VV014714.D | |

| | Compound | 0.5 | 1 | 5 | 10 | 20 | Avg | %RSD |
|----------------|---------------------------|-------|-------|-------|-------|-------|-------|-------|
| -----ISTD----- | | | | | | | | |
| 1) I | 1,4-Difluorobenzene | | | | | | | |
| 2) T | Dichlorodifluoromethane | 0.548 | 0.523 | 0.547 | 0.522 | 0.605 | 0.549 | 6.10 |
| 3) T | Chloromethane | 0.504 | 0.488 | 0.505 | 0.475 | 0.556 | 0.506 | 6.04 |
| 4) S | Vinyl Chloride-d3 | 0.351 | 0.376 | 0.402 | 0.420 | 0.409 | 0.391 | 7.08 |
| 5) T | Vinyl chloride | 0.446 | 0.443 | 0.467 | 0.445 | 0.523 | 0.465 | 7.34 |
| 6) T | Bromomethane | 0.256 | 0.248 | 0.264 | 0.251 | 0.290 | 0.262 | 6.37 |
| 7) S | Chloroethane-d5 | 0.300 | 0.285 | 0.310 | 0.319 | 0.316 | 0.306 | 4.51 |
| 8) T | Chloroethane | 0.282 | 0.239 | 0.261 | 0.244 | 0.276 | 0.260 | 7.31 |
| 9) T | Trichlorofluoromethane | 0.635 | 0.607 | 0.636 | 0.596 | 0.692 | 0.633 | 5.92 |
| 10) T | 1,1,2-Trichloro-1,2-d | 0.329 | 0.308 | 0.335 | 0.314 | 0.370 | 0.331 | 7.29 |
| 11) S | 1,1-Dichloroethene | 0.619 | 0.644 | 0.676 | 0.690 | 0.701 | 0.666 | 5.09 |
| 12) T | 1,1-Dichloroethene | 0.301 | 0.292 | 0.323 | 0.304 | 0.357 | 0.315 | 8.21 |
| 13) T | Acetone | 0.055 | 0.048 | 0.057 | 0.052 | 0.061 | 0.054 | 8.86 |
| 14) T | Carbon disulfide | 1.093 | 0.976 | 1.076 | 1.023 | 1.204 | 1.074 | 7.98 |
| 15) T | Methyl Acetate | 0.099 | 0.116 | 0.145 | 0.142 | 0.188 | 0.138 | 24.49 |
| 16) T | Methylene chloride | 0.457 | 0.374 | 0.381 | 0.353 | 0.406 | 0.394 | 10.08 |
| 17) T | Methyl tert-butyl E | 0.821 | 0.770 | 0.885 | 0.826 | 0.985 | 0.857 | 9.58 |
| 18) T | trans-1,2-Dichloroethane | 0.376 | 0.363 | 0.383 | 0.361 | 0.422 | 0.381 | 6.45 |
| 19) T | 1,1-Dichloroethane | 0.740 | 0.698 | 0.754 | 0.712 | 0.821 | 0.745 | 6.42 |
| 20) S | 2-Butanone-d5 | 0.063 | 0.066 | 0.089 | 0.092 | 0.095 | 0.081 | 18.73 |
| 21) T | 2-Butanone | 0.077 | 0.072 | 0.102 | 0.094 | 0.118 | 0.093 | 20.35 |
| 22) T | cis-1,2-Dichloroethane | 0.380 | 0.392 | 0.414 | 0.388 | 0.461 | 0.407 | 7.99 |
| 23) T | Bromochloromethane | 0.159 | 0.153 | 0.165 | 0.158 | 0.182 | 0.163 | 6.80 |
| 24) S | Chloroform-d | 0.645 | 0.651 | 0.721 | 0.753 | 0.743 | 0.703 | 7.29 |
| 25) T | Chloroform | 0.752 | 0.690 | 0.732 | 0.686 | 0.796 | 0.731 | 6.25 |
| 26) S | 1,2-Dichloroethane-d | 0.314 | 0.314 | 0.356 | 0.372 | 0.366 | 0.345 | 8.27 |
| 27) T | 1,2-Dichloroethane | 0.458 | 0.425 | 0.451 | 0.412 | 0.493 | 0.448 | 6.99 |
| 28) I | Chlorobenzene-d5 | | | | | | | |
| 29) T | 1,1,1-Trichloroethane | 0.650 | 0.635 | 0.658 | 0.622 | 0.731 | 0.659 | 6.47 |
| 30) T | Cyclohexane | 0.734 | 0.692 | 0.745 | 0.697 | 0.842 | 0.742 | 8.14 |
| 31) T | Carbon tetrachloride | 0.559 | 0.521 | 0.572 | 0.541 | 0.646 | 0.568 | 8.42 |
| 32) S | Benzene-d6 | 1.291 | 1.388 | 1.489 | 1.541 | 1.539 | 1.450 | 7.47 |
| 33) T | Benzene | 1.573 | 1.559 | 1.678 | 1.560 | 1.839 | 1.642 | 7.36 |
| 34) T | Trichloroethene | 0.446 | 0.407 | 0.420 | 0.401 | 0.476 | 0.430 | 7.19 |
| 35) T | Methylcyclohexane | 0.702 | 0.689 | 0.742 | 0.689 | 0.835 | 0.731 | 8.47 |
| 36) S | 1,2-Dichloropropane | 0.436 | 0.418 | 0.455 | 0.480 | 0.480 | 0.454 | 6.01 |
| 37) T | 1,2-Dichloropropane | 0.420 | 0.382 | 0.428 | 0.392 | 0.475 | 0.419 | 8.73 |
| 38) T | Bromodichloromethane | 0.483 | 0.483 | 0.516 | 0.483 | 0.592 | 0.512 | 9.19 |
| 39) T | cis-1,3-Dichloropropane | 0.544 | 0.554 | 0.601 | 0.604 | 0.736 | 0.608 | 12.61 |
| 40) T | 4-Methyl-2-pentanone | 0.232 | 0.226 | 0.264 | 0.241 | 0.293 | 0.251 | 10.93 |
| 41) S | Toluene-d8 | 1.225 | 1.252 | 1.361 | 1.431 | 1.431 | 1.340 | 7.29 |
| 42) T | Toluene | 1.632 | 1.622 | 1.789 | 1.674 | 1.976 | 1.739 | 8.54 |
| 43) S | trans-1,3-Dichloropropene | 0.149 | 0.158 | 0.178 | 0.193 | 0.198 | 0.175 | 12.29 |
| 44) T | trans-1,3-Dichloropropene | 0.424 | 0.415 | 0.484 | 0.461 | 0.573 | 0.471 | 13.38 |
| 45) T | 1,1,2-Trichloroethane | 0.254 | 0.259 | 0.268 | 0.254 | 0.303 | 0.268 | 7.76 |
| 46) S | 2-Hexanone-d5 | 0.058 | 0.062 | 0.074 | 0.079 | 0.082 | 0.071 | 14.60 |
| 47) T | Tetrachloroethene | 0.330 | 0.336 | 0.341 | 0.326 | 0.387 | 0.344 | 7.19 |
| 48) T | 2-Hexanone | 0.154 | 0.163 | 0.191 | 0.176 | 0.211 | 0.179 | 12.64 |
| 49) T | Dibromochloromethane | 0.268 | 0.272 | 0.307 | 0.294 | 0.373 | 0.303 | 14.00 |
| 50) T | 1,2-Dibromoethane | 0.250 | 0.246 | 0.258 | 0.238 | 0.287 | 0.256 | 7.38 |
| 51) T | Chlorobenzene | 1.072 | 1.032 | 1.105 | 1.045 | 1.251 | 1.101 | 8.05 |
| 52) T | Ethylbenzene | 1.823 | 1.817 | 2.001 | 1.918 | 2.298 | 1.971 | 10.03 |

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Response Via : Initial Calibration

Calibration Files

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| 0.5 | =VV014710.D | 1 | =VV014711.D | 5 | =VV014712.D |
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| | Compound | 0.5 | 1 | 5 | 10 | 20 | Avg | %RSD |
|-----|-------------------------|----------------|-------|-------|-------|-------|-------|-------|
| 53) | T m,p-xylene | 0.654 | 0.645 | 0.740 | 0.707 | 0.862 | 0.722 | 12.11 |
| 54) | T o-xylene | 0.633 | 0.644 | 0.720 | 0.684 | 0.833 | 0.703 | 11.47 |
| 55) | T Styrene | 1.032 | 1.036 | 1.229 | 1.178 | 1.427 | 1.180 | 13.81 |
| 56) | T Isopropylbenzene | 1.669 | 1.712 | 1.951 | 1.886 | 2.272 | 1.898 | 12.63 |
| 57) | S 1,1,2,2-Tetrachloro | 0.316 | 0.295 | 0.326 | 0.335 | 0.350 | 0.325 | 6.39 |
| 58) | T 1,1,2,2-Tetrachloro | 0.303 | 0.323 | 0.338 | 0.322 | 0.388 | 0.335 | 9.59 |
| 59) | T 1,2,3-Trichloroprop | 0.243 | 0.227 | 0.247 | 0.229 | 0.279 | 0.245 | 8.55 |
| 60) | I 1,4-Dichlorobenzene-d | -----ISTD----- | | | | | | |
| 61) | T Bromoform | 0.239 | 0.255 | 0.301 | 0.294 | 0.383 | 0.294 | 19.04 |
| 62) | T 1,3-Dichlorobenzene | 1.581 | 1.615 | 1.747 | 1.632 | 1.972 | 1.709 | 9.32 |
| 63) | T 1,4-Dichlorobenzene | 1.766 | 1.638 | 1.736 | 1.637 | 1.955 | 1.746 | 7.46 |
| 64) | S 1,2-Dichlorobenzene | 0.895 | 0.860 | 0.928 | 0.972 | 0.987 | 0.928 | 5.66 |
| 65) | T 1,2-Dichlorobenzene | 1.576 | 1.475 | 1.580 | 1.492 | 1.773 | 1.579 | 7.49 |
| 66) | T 1,2-Dibromo-3-chlor | 0.113 | 0.095 | 0.100 | 0.093 | 0.123 | 0.105 | 12.08 |
| 67) | T 1,3,5-Trichlorobenz | 1.264 | 1.305 | 1.417 | 1.332 | 1.617 | 1.387 | 10.10 |
| 68) | T 1,2,4-trichlorobenz | 1.117 | 1.082 | 1.196 | 1.151 | 1.411 | 1.191 | 10.90 |
| 69) | Naphthalene | 1.703 | 1.694 | 1.999 | 1.900 | 2.401 | 1.939 | 14.90 |
| 70) | T 1,2,3-Trichlorobenz | 1.001 | 0.977 | 1.077 | 1.018 | 1.243 | 1.063 | 10.06 |

(#) = Out of Range