

Method Path : Z:\VOASRV\HPCHEM1\MSVOA U\METHOD\  
 Method File : SOMULM051419WMA.M  
 Title : VOC Analysis  
 Last Update : Wed May 15 04:26:34 2019  
 Response Via : Initial Calibration

## Calibration Files

5 =VU031982.D 10 =VU031983.D 50 =VU031988.D  
 100 =VU031985.D 200 =VU031986.D

|       | Compound             | 5              | 10    | 50    | 100   | 200   | Avg   | %RSD  |
|-------|----------------------|----------------|-------|-------|-------|-------|-------|-------|
| ----- |                      |                |       |       |       |       |       |       |
| 1) I  | 1,4-Difluorobenzene  | -----ISTD----- |       |       |       |       |       |       |
| 2) T  | Dichlorodifluoromet  | 0.530          | 0.510 | 0.439 | 0.416 | 0.406 | 0.460 | 12.22 |
| 3) T  | Chloromethane        | 0.593          | 0.576 | 0.494 | 0.458 | 0.450 | 0.514 | 12.94 |
| 4) S  | Vinyl Chloride-d3    | 0.482          | 0.452 | 0.429 | 0.423 | 0.408 | 0.439 | 6.54  |
| 5) T  | Vinyl chloride       | 0.616          | 0.580 | 0.515 | 0.495 | 0.487 | 0.539 | 10.49 |
| 6) T  | Bromomethane         | 0.329          | 0.310 | 0.270 | 0.264 | 0.243 | 0.283 | 12.39 |
| 7) S  | Chloroethane-d5      | 0.378          | 0.374 | 0.354 | 0.343 | 0.327 | 0.355 | 6.04  |
| 8) T  | Chloroethane         | 0.375          | 0.350 | 0.300 | 0.290 | 0.279 | 0.319 | 13.09 |
| 9) T  | Trichlorofluorometh  | 0.713          | 0.668 | 0.574 | 0.552 | 0.539 | 0.610 | 12.64 |
| 10) T | 1,1,2-Trichloro-1,2  | 0.441          | 0.399 | 0.354 | 0.337 | 0.328 | 0.372 | 12.73 |
| 11) S | 1,1-Dichloroethene-  | 0.851          | 0.835 | 0.765 | 0.746 | 0.735 | 0.786 | 6.76  |
| 12) T | 1,1-Dichloroethene   | 0.415          | 0.395 | 0.354 | 0.346 | 0.336 | 0.369 | 9.29  |
| 13) T | Acetone              | 0.374          | 0.377 | 0.303 | 0.274 | 0.252 | 0.316 | 18.08 |
| 14) T | Carbon disulfide     | 1.340          | 1.302 | 1.128 | 1.105 | 1.089 | 1.193 | 9.97  |
| 15) T | Methyl Acetate       | 0.544          | 0.510 | 0.461 | 0.450 | 0.433 | 0.480 | 9.62  |
| 16) T | Methylene chloride   | 0.511          | 0.501 | 0.428 | 0.410 | 0.402 | 0.450 | 11.44 |
| 17) T | trans-1,2-Dichloroe  | 0.430          | 0.396 | 0.371 | 0.363 | 0.360 | 0.384 | 7.62  |
| 18) T | Methyl tert-butyl E  | 1.208          | 1.216 | 1.155 | 1.160 | 1.163 | 1.180 | 2.46  |
| 19) T | 1,1-Dichloroethane   | 0.838          | 0.817 | 0.723 | 0.702 | 0.692 | 0.754 | 9.04  |
| 20) T | cis-1,2-Dichloroeth  | 0.452          | 0.436 | 0.417 | 0.412 | 0.408 | 0.425 | 4.39  |
| 21) S | 2-Butanone-d5        | 0.282          | 0.297 | 0.316 | 0.321 | 0.320 | 0.307 | 5.53  |
| 22) T | 2-Butanone           | 0.355          | 0.370 | 0.376 | 0.366 | 0.355 | 0.364 | 2.58  |
| 23) T | Bromochloromethane   | 0.228          | 0.228 | 0.202 | 0.193 | 0.190 | 0.208 | 8.98  |
| 24) S | Chloroform-d         | 0.769          | 0.765 | 0.726 | 0.713 | 0.700 | 0.734 | 4.25  |
| 25) T | Chloroform           | 0.837          | 0.813 | 0.709 | 0.692 | 0.677 | 0.745 | 9.89  |
| 26) S | 1,2-Dichloroethane-  | 0.501          | 0.481 | 0.452 | 0.444 | 0.434 | 0.463 | 5.98  |
| 27) T | 1,2-Dichloroethane   | 0.631          | 0.625 | 0.553 | 0.543 | 0.534 | 0.577 | 8.14  |
| ----- |                      |                |       |       |       |       |       |       |
| 28) I | Chlorobenzene-d5     | -----ISTD----- |       |       |       |       |       |       |
| 29) T | Cyclohexane          | 0.560          | 0.595 | 0.622 | 0.649 | 0.639 | 0.613 | 5.88  |
| 30) T | 1,1,1-Trichloroetha  | 0.650          | 0.625 | 0.546 | 0.547 | 0.531 | 0.580 | 9.25  |
| 31) T | Carbon tetrachlorid  | 0.536          | 0.520 | 0.459 | 0.455 | 0.449 | 0.484 | 8.41  |
| 32) S | Benzene-d6           | 1.454          | 1.452 | 1.461 | 1.464 | 1.407 | 1.448 | 1.60  |
| 33) T | Benzene              | 1.758          | 1.735 | 1.598 | 1.575 | 1.523 | 1.638 | 6.30  |
| 34) T | Trichloroethene      | 0.444          | 0.429 | 0.384 | 0.388 | 0.383 | 0.406 | 7.10  |
| 35) T | Methylcyclohexane    | 0.632          | 0.586 | 0.618 | 0.638 | 0.636 | 0.622 | 3.51  |
| 36) S | 1,2-Dichloropropane  | 0.491          | 0.489 | 0.466 | 0.476 | 0.466 | 0.478 | 2.55  |
| 37) T | 1,2-Dichloropropane  | 0.469          | 0.470 | 0.425 | 0.426 | 0.410 | 0.440 | 6.34  |
| 38) T | Bromodichloromethan  | 0.588          | 0.561 | 0.509 | 0.519 | 0.511 | 0.538 | 6.61  |
| 39) T | cis-1,3-Dichloropro  | 0.581          | 0.599 | 0.628 | 0.650 | 0.664 | 0.625 | 5.52  |
| 40) T | 4-Methyl-2-pentanone | 0.540          | 0.580 | 0.600 | 0.625 | 0.639 | 0.597 | 6.51  |
| 41) S | Toluene-d8           | 1.223          | 1.265 | 1.319 | 1.364 | 1.314 | 1.297 | 4.18  |
| 42) T | Toluene              | 1.643          | 1.716 | 1.691 | 1.690 | 1.663 | 1.681 | 1.69  |
| 43) S | trans-1,3-Dichlorop  | 0.193          | 0.196 | 0.211 | 0.225 | 0.229 | 0.211 | 7.74  |
| 44) T | trans-1,3-Dichlorop  | 0.522          | 0.529 | 0.545 | 0.563 | 0.583 | 0.548 | 4.60  |
| 45) T | 1,1,2-Trichloroetha  | 0.434          | 0.434 | 0.381 | 0.386 | 0.382 | 0.403 | 6.93  |
| 46) T | Tetrachloroethene    | 0.312          | 0.300 | 0.274 | 0.275 | 0.270 | 0.286 | 6.53  |
| 47) S | 2-Hexanone-d5        | 0.140          | 0.166 | 0.203 | 0.225 | 0.235 | 0.194 | 20.65 |
| 48) T | 2-Hexanone           | 0.405          | 0.487 | 0.493 | 0.517 | 0.520 | 0.484 | 9.60  |
| 49) T | Dibromochloromethan  | 0.425          | 0.417 | 0.385 | 0.393 | 0.403 | 0.405 | 4.04  |
| 50) T | 1,2-Dibromoethane    | 0.462          | 0.440 | 0.406 | 0.414 | 0.412 | 0.427 | 5.51  |
| 51) T | Chlorobenzene        | 1.165          | 1.116 | 1.015 | 1.017 | 1.023 | 1.067 | 6.50  |
| 52) T | Ethylbenzene         | 1.696          | 1.788 | 1.790 | 1.855 | 1.864 | 1.799 | 3.75  |

Method Path : Z:\VOASRV\HPCHEM1\MSVOA U\METHOD\  
 Method File : SOMULM051419WMA.M  
 Title : VOC Analysis  
 Last Update : Wed May 15 04:26:34 2019  
 Response Via : Initial Calibration

## Calibration Files

5 =VU031982.D 10 =VU031983.D 50 =VU031988.D  
 100 =VU031985.D 200 =VU031986.D

|       | Compound              | 5              | 10    | 50    | 100   | 200   | Avg   | %RSD  |
|-------|-----------------------|----------------|-------|-------|-------|-------|-------|-------|
| 53) T | m,p-Xylene            | 0.593          | 0.626 | 0.678 | 0.692 | 0.693 | 0.656 | 6.83  |
| 54) T | o-xylene              | 0.574          | 0.624 | 0.665 | 0.685 | 0.692 | 0.648 | 7.57  |
| 55) T | Styrene               | 0.910          | 1.075 | 1.152 | 1.197 | 1.229 | 1.112 | 11.43 |
| 56) T | Isopropylbenzene      | 1.470          | 1.596 | 1.680 | 1.755 | 1.801 | 1.660 | 7.93  |
| 57) S | 1,1,2,2-Tetrachloro   | 0.710          | 0.693 | 0.690 | 0.717 | 0.724 | 0.707 | 2.09  |
| 58) T | 1,1,2,2-Tetrachloro   | 0.763          | 0.772 | 0.685 | 0.702 | 0.718 | 0.728 | 5.23  |
| 59)   | 1,2,3-Trichloroprop   | 0.619          | 0.596 | 0.542 | 0.554 | 0.558 | 0.574 | 5.63  |
| 60) I | 1,4-Dichlorobenzene-d | -----ISTD----- |       |       |       |       |       |       |
| 61) T | Bromoform             | 0.713          | 0.744 | 0.628 | 0.619 | 0.633 | 0.667 | 8.53  |
| 62) T | 1,3-Dichlorobenzene   | 1.682          | 1.712 | 1.597 | 1.589 | 1.600 | 1.636 | 3.48  |
| 63) T | 1,4-Dichlorobenzene   | 1.937          | 1.829 | 1.672 | 1.605 | 1.609 | 1.730 | 8.47  |
| 64) S | 1,2-Dichlorobenzene   | 1.079          | 1.076 | 1.061 | 1.035 | 1.011 | 1.052 | 2.76  |
| 65) T | 1,2-Dichlorobenzene   | 1.818          | 1.885 | 1.695 | 1.650 | 1.620 | 1.734 | 6.55  |
| 66) T | 1,2-Dibromo-3-chlor   | 0.344          | 0.365 | 0.321 | 0.326 | 0.318 | 0.335 | 5.84  |
| 67)   | 1,3,5-Trichlorobenz   | 1.151          | 1.212 | 1.167 | 1.173 | 1.162 | 1.173 | 1.99  |
| 68) T | 1,2,4-trichlorobenz   | 0.745          | 0.797 | 0.947 | 1.012 | 1.018 | 0.904 | 13.93 |
| 69)   | Naphthalene           | 1.890          | 2.457 | 3.351 | 3.685 | 3.659 | 3.009 | 26.55 |
| 70) T | 1,2,3-Trichlorobenz   | 0.792          | 0.970 | 1.057 | 1.066 | 1.034 | 0.984 | 11.54 |

(#) = Out of Range