

Method Path : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\
 Method File : SOMULM041719WMA.M
 Title : VOC Analysis
 Last Update : Wed Apr 17 04:13:40 2019
 Response Via : Initial Calibration

Calibration Files

5 =VU031213.D 10 =VU031214.D 50 =VU031215.D
 100 =VU031216.D 200 =VU031217.D

| | Compound | 5 | 10 | 50 | 100 | 200 | Avg | %RSD |
|----------------|---------------------|-------|-------|-------|-------|-------|-------|-------|
| -----ISTD----- | | | | | | | | |
| 1) I | 1,4-Difluorobenzene | | | | | | | |
| 2) T | Dichlorodifluoromet | 0.524 | 0.475 | 0.436 | 0.438 | 0.434 | 0.462 | 8.44 |
| 3) T | Chloromethane | 0.652 | 0.569 | 0.507 | 0.510 | 0.510 | 0.550 | 11.43 |
| 4) S | Vinyl Chloride-d3 | 0.478 | 0.501 | 0.431 | 0.455 | 0.437 | 0.460 | 6.35 |
| 5) T | Vinyl chloride | 0.626 | 0.609 | 0.497 | 0.514 | 0.489 | 0.547 | 11.93 |
| 6) T | Bromomethane | 0.413 | 0.381 | 0.340 | 0.330 | 0.334 | 0.359 | 10.10 |
| 7) S | Chloroethane-d5 | 0.407 | 0.414 | 0.401 | 0.391 | 0.386 | 0.400 | 2.82 |
| 8) T | Chloroethane | 0.394 | 0.374 | 0.335 | 0.325 | 0.331 | 0.352 | 8.67 |
| 9) T | Trichlorofluorometh | 0.896 | 0.790 | 0.713 | 0.707 | 0.701 | 0.762 | 10.96 |
| 10) T | 1,1,2-Trichloro-1,2 | 0.448 | 0.410 | 0.368 | 0.395 | 0.371 | 0.398 | 8.15 |
| 11) S | 1,1-Dichloroethene- | 0.996 | 0.926 | 0.842 | 0.914 | 0.884 | 0.912 | 6.24 |
| 12) T | 1,1-Dichloroethene | 0.439 | 0.405 | 0.358 | 0.385 | 0.367 | 0.391 | 8.29 |
| 13) T | Acetone | 0.279 | 0.283 | 0.239 | 0.271 | 0.218 | 0.258 | 10.96 |
| 14) T | Carbon disulfide | 1.179 | 1.051 | 0.983 | 1.028 | 1.052 | 1.059 | 6.90 |
| 15) T | Methyl Acetate | 0.477 | 0.441 | 0.425 | 0.435 | 0.467 | 0.449 | 4.88 |
| 16) T | Methylene chloride | 0.427 | 0.381 | 0.362 | 0.356 | 0.380 | 0.381 | 7.28 |
| 17) T | trans-1,2-Dichloroe | 0.416 | 0.342 | 0.319 | 0.349 | 0.354 | 0.356 | 10.14 |
| 18) T | Methyl tert-butyl E | 1.210 | 1.056 | 1.045 | 1.134 | 1.162 | 1.121 | 6.28 |
| 19) T | 1,1-Dichloroethane | 0.812 | 0.704 | 0.653 | 0.681 | 0.688 | 0.708 | 8.65 |
| 20) T | cis-1,2-Dichloroeth | 0.450 | 0.420 | 0.414 | 0.392 | 0.396 | 0.414 | 5.53 |
| 21) S | 2-Butanone-d5 | 0.313 | 0.328 | 0.354 | 0.337 | 0.346 | 0.335 | 4.73 |
| 22) T | 2-Butanone | 0.367 | 0.345 | 0.374 | 0.339 | 0.344 | 0.354 | 4.42 |
| 23) T | Bromochloromethane | 0.228 | 0.206 | 0.204 | 0.194 | 0.195 | 0.205 | 6.71 |
| 24) S | Chloroform-d | 0.787 | 0.783 | 0.783 | 0.742 | 0.747 | 0.768 | 2.83 |
| 25) T | Chloroform | 0.774 | 0.761 | 0.737 | 0.679 | 0.679 | 0.726 | 6.21 |
| 26) S | 1,2-Dichloroethane- | 0.486 | 0.472 | 0.482 | 0.459 | 0.466 | 0.473 | 2.36 |
| 27) T | 1,2-Dichloroethane | 0.587 | 0.580 | 0.547 | 0.519 | 0.527 | 0.552 | 5.51 |
| -----ISTD----- | | | | | | | | |
| 28) I | Chlorobenzene-d5 | | | | | | | |
| 29) T | Cyclohexane | 0.743 | 0.631 | 0.728 | 0.681 | 0.733 | 0.703 | 6.66 |
| 30) T | 1,1,1-Trichloroetha | 0.753 | 0.659 | 0.645 | 0.607 | 0.645 | 0.662 | 8.24 |
| 31) T | Carbon tetrachlorid | 0.633 | 0.566 | 0.559 | 0.537 | 0.567 | 0.572 | 6.24 |
| 32) S | Benzene-d6 | 1.603 | 1.501 | 1.587 | 1.553 | 1.658 | 1.581 | 3.70 |
| 33) T | Benzene | 1.809 | 1.675 | 1.660 | 1.589 | 1.676 | 1.682 | 4.72 |
| 34) T | Trichloroethene | 0.479 | 0.424 | 0.402 | 0.400 | 0.399 | 0.421 | 8.15 |
| 35) T | Methylcyclohexane | 0.648 | 0.595 | 0.597 | 0.629 | 0.664 | 0.626 | 4.88 |
| 36) S | 1,2-Dichloropropane | 0.513 | 0.509 | 0.503 | 0.508 | 0.532 | 0.513 | 2.14 |
| 37) T | 1,2-Dichloropropane | 0.481 | 0.453 | 0.423 | 0.426 | 0.444 | 0.445 | 5.27 |
| 38) T | Bromodichloromethan | 0.628 | 0.555 | 0.546 | 0.537 | 0.564 | 0.566 | 6.36 |
| 39) T | cis-1,3-Dichloropro | 0.674 | 0.623 | 0.640 | 0.672 | 0.728 | 0.667 | 6.03 |
| 40) T | 4-Methyl-2-pentanon | 0.617 | 0.594 | 0.629 | 0.631 | 0.670 | 0.628 | 4.35 |
| 41) S | Toluene-d8 | 1.273 | 1.327 | 1.406 | 1.433 | 1.516 | 1.391 | 6.78 |
| 42) T | Toluene | 1.693 | 1.584 | 1.640 | 1.676 | 1.772 | 1.673 | 4.14 |
| 43) S | trans-1,3-Dichlorop | 0.233 | 0.224 | 0.244 | 0.251 | 0.274 | 0.245 | 7.79 |
| 44) T | trans-1,3-Dichlorop | 0.578 | 0.525 | 0.572 | 0.584 | 0.638 | 0.579 | 6.89 |
| 45) T | 1,1,2-Trichloroetha | 0.429 | 0.400 | 0.388 | 0.383 | 0.403 | 0.401 | 4.41 |
| 46) T | Tetrachloroethene | 0.336 | 0.313 | 0.305 | 0.305 | 0.323 | 0.316 | 4.26 |
| 47) S | 2-Hexanone-d5 | 0.208 | 0.205 | 0.247 | 0.264 | 0.288 | 0.242 | 14.81 |
| 48) T | 2-Hexanone | 0.497 | 0.460 | 0.506 | 0.505 | 0.531 | 0.500 | 5.13 |
| 49) T | Dibromochloromethan | 0.452 | 0.411 | 0.421 | 0.427 | 0.455 | 0.433 | 4.50 |
| 50) T | 1,2-Dibromoethane | 0.430 | 0.425 | 0.412 | 0.414 | 0.437 | 0.423 | 2.46 |
| 51) T | Chlorobenzene | 1.142 | 1.017 | 1.015 | 1.007 | 1.047 | 1.046 | 5.35 |
| 52) T | Ethylbenzene | 1.832 | 1.720 | 1.765 | 1.790 | 1.999 | 1.821 | 5.90 |

Method Path : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\
 Method File : SOMULM041719WMA.M
 Title : VOC Analysis
 Last Update : Wed Apr 17 04:13:40 2019
 Response Via : Initial Calibration

Calibration Files

5 =VU031213.D 10 =VU031214.D 50 =VU031215.D
 100 =VU031216.D 200 =VU031217.D

| | Compound | 5 | 10 | 50 | 100 | 200 | Avg | %RSD |
|-------|-----------------------|----------------|-------|-------|-------|-------|-------|-------|
| 53) T | m,p-Xylene | 0.561 | 0.633 | 0.641 | 0.655 | 0.724 | 0.643 | 9.04 |
| 54) T | o-xylene | 0.621 | 0.603 | 0.642 | 0.642 | 0.747 | 0.651 | 8.60 |
| 55) T | Styrene | 1.042 | 0.920 | 1.111 | 1.121 | 1.298 | 1.099 | 12.52 |
| 56) T | Isopropylbenzene | 1.637 | 1.528 | 1.666 | 1.620 | 1.890 | 1.668 | 8.05 |
| 57) S | 1,1,2,2-Tetrachloro | 0.765 | 0.694 | 0.689 | 0.683 | 0.730 | 0.712 | 4.90 |
| 58) T | 1,1,2,2-Tetrachloro | 0.760 | 0.693 | 0.667 | 0.641 | 0.690 | 0.690 | 6.43 |
| 59) T | 1,2,3-Trichloroprop | 0.588 | 0.529 | 0.515 | 0.484 | 0.532 | 0.530 | 7.08 |
| 60) I | 1,4-Dichlorobenzene-d | -----ISTD----- | | | | | | |
| 61) T | Bromoform | 0.750 | 0.727 | 0.664 | 0.651 | 0.670 | 0.692 | 6.31 |
| 62) T | 1,3-Dichlorobenzene | 1.888 | 1.598 | 1.538 | 1.671 | 1.567 | 1.652 | 8.52 |
| 63) T | 1,4-Dichlorobenzene | 1.888 | 1.648 | 1.578 | 1.596 | 1.574 | 1.657 | 8.00 |
| 64) S | 1,2-Dichlorobenzene | 1.220 | 1.123 | 1.066 | 1.241 | 1.083 | 1.147 | 6.96 |
| 65) T | 1,2-Dichlorobenzene | 1.987 | 1.713 | 1.608 | 1.811 | 1.571 | 1.738 | 9.66 |
| 66) T | 1,2-Dibromo-3-chlor | 0.382 | 0.307 | 0.302 | 0.363 | 0.304 | 0.332 | 11.48 |
| 67) T | 1,3,5-Trichlorobenz | 1.306 | 1.187 | 1.179 | 1.291 | 1.220 | 1.237 | 4.74 |
| 68) T | 1,2,4-trichlorobenz | 0.924 | 0.902 | 1.002 | 1.084 | 1.099 | 1.002 | 8.96 |
| 69) T | Naphthalene | 2.835 | 2.442 | 3.280 | 3.566 | 3.598 | 3.144 | 15.83 |
| 70) T | 1,2,3-Trichlorobenz | 1.108 | 1.025 | 1.065 | 1.124 | 1.120 | 1.088 | 3.89 |

(#) = Out of Range