Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN032223\
Data File : VN077096.D
Acq On : 22 Mar 2023 10:02
Operator : JC\MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :**
MSVOA_N
**LabSampleId :**
VSTDCCC050

Quant Time: Mar 22 23:18:28 2023
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N031723W.M
Quant Title : SW846 8260
QLast Update : Mon Mar 20 08:50:30 2023
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

| | | Compound | AvgRF | CCRF | %Dev | Area% | Dev(min) |
|---|---|---|---|---|---|---|---|
| 1 | I | Pentafluorobenzene | 1.000 | 1.000 | 0.0 | 98 | 0.00 |
| 2 | T | Dichlorodifluoromethane | 0.566 | 0.343 | 39.4# | 71 | 0.00 |
| 3 | P | Chloromethane | 0.504 | 0.335 | 33.5# | 70 | 0.00 |
| 4 | C | Vinyl Chloride | 0.544 | 0.458 | 15.8# | 87 | 0.00 |
| 5 | T | Bromomethane | 0.367 | 0.343 | 6.5 | 92 | 0.00 |
| 6 | T | Chloroethane | 0.344 | 0.297 | 13.7 | 94 | 0.00 |
| 7 | T | Trichlorofluoromethane | 1.170 | 1.106 | 5.5 | 97 | 0.00 |
| 8 | T | Diethyl Ether | 0.440 | 0.421 | 4.3 | 95 | 0.00 |
| 9 | T | 1,1,2-Trichlorotrifluoroeth | 0.664 | 0.611 | 8.0 | 97 | 0.00 |
| 10 | T | Methyl Iodide | 0.747 | 0.708 | 5.2 | 92 | 0.00 |
| 11 | T | Tert butyl alcohol | 0.116 | 0.096 | 17.2 | 81 | 0.00 |
| 12 | CM | 1,1-Dichloroethene | 0.607 | 0.565 | 6.9# | 94 | 0.00 |
| 13 | T | Acrolein | 0.144 | 0.133 | 7.6 | 93 | 0.00 |
| 14 | T | Allyl chloride | 1.111 | 1.180 | -6.2 | 104 | 0.00 |
| 15 | T | Acrylonitrile | 0.391 | 0.349 | 10.7 | 88 | 0.00 |
| 16 | T | Acetone | 0.356 | 0.371 | -4.2 | 108 | 0.00 |
| 17 | T | Carbon Disulfide | 1.770 | 1.520 | 14.1 | 89 | 0.00 |
| 18 | T | Methyl Acetate | 1.342 | 1.194 | 11.0 | 90 | 0.00 |
| 19 | T | Methyl tert-butyl Ether | 2.179 | 2.142 | 1.7 | 97 | 0.00 |
| 20 | T | Methylene Chloride | 0.772 | 0.695 | 10.0 | 97 | 0.00 |
| 21 | T | trans-1,2-Dichloroethene | 0.658 | 0.612 | 7.0 | 94 | 0.00 |
| 22 | T | Diisopropyl ether | 2.435 | 2.466 | -1.3 | 98 | 0.00 |
| 23 | T | Vinyl Acetate | 1.536 | 1.536 | 0.0 | 95 | 0.00 |
| 24 | P | 1,1-Dichloroethane | 1.393 | 1.318 | 5.4 | 97 | 0.00 |
| 25 | T | 2-Butanone | 0.521 | 0.482 | 7.5 | 90 | 0.00 |
| 26 | T | 2,2-Dichloropropane | 1.048 | 1.070 | -2.1 | 100 | 0.00 |
| 27 | T | cis-1,2-Dichloroethene | 0.754 | 0.751 | 0.4 | 98 | 0.00 |
| 28 | T | Bromochloromethane | 0.605 | 0.545 | 9.9 | 96 | 0.00 |
| 29 | T | Tetrahydrofuran | 0.337 | 0.308 | 8.6 | 87 | 0.00 |
| 30 | C | Chloroform | 1.352 | 1.290 | 4.6# | 98 | 0.00 |
| 31 | T | Cyclohexane | 1.307 | 1.195 | 8.6 | 95 | 0.00 |
| 32 | T | 1,1,1-Trichloroethane | 1.156 | 1.116 | 3.5 | 97 | 0.00 |
| 33 | S | 1,2-Dichloroethane-d4 | 0.799 | 0.792 | 0.9 | 100 | 0.00 |
| 34 | I | 1,4-Difluorobenzene | 1.000 | 1.000 | 0.0 | 96 | 0.00 |
| 35 | S | Dibromofluoromethane | 0.350 | 0.337 | 3.7 | 98 | 0.00 |
| 36 | T | 1,1-Dichloropropene | 0.547 | 0.543 | 0.7 | 95 | 0.00 |
| 37 | T | Ethyl Acetate | 0.585 | 0.555 | 5.1 | 93 | 0.00 |
| 38 | T | Carbon Tetrachloride | 0.533 | 0.527 | 1.1 | 96 | 0.00 |
| 39 | T | Methylcyclohexane | 0.640 | 0.669 | -4.5 | 95 | 0.00 |
| 40 | TM | Benzene | 1.659 | 1.617 | 2.5 | 96 | 0.00 |
| 41 | T | Methacrylonitrile | 0.315 | 0.321 | -1.9 | 100 | 0.00 |
| 42 | TM | 1,2-Dichloroethane | 0.599 | 0.581 | 3.0 | 97 | 0.00 |
| 43 | T | Isopropyl Acetate | 0.914 | 0.905 | 1.0 | 92 | 0.00 |
| 44 | TM | Trichloroethene | 0.381 | 0.357 | 6.3 | 94 | 0.00 |
| 45 | C | 1,2-Dichloropropane | 0.441 | 0.441 | 0.0 | 98 | 0.00 |
| 46 | T | Dibromomethane | 0.278 | 0.268 | 3.6 | 96 | 0.00 |
| 47 | T | Bromodichloromethane | 0.583 | 0.584 | -0.2 | 99 | 0.00 |
| 48 | T | Methyl methacrylate | 0.437 | 0.452 | -3.4 | 93 | 0.00 |

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ALS Vial : 1 Sample Multiplier: 1

**Instrument :**
MSVOA_N
**LabSampleId :**
VSTDCCC050

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Max. RRF Dev : 25% Max. Rel. Area : 150%

| | | Compound | AvgRF | CCRF | %Dev | Area% | Dev(min) |
|---|---|---|---|---|---|---|---|
| 49 | T | 1,4-Dioxane | 0.007 | 0.006 | 14.3 | 81 | 0.00 |
| 50 | S | Toluene-d8 | 1.251 | 1.286 | -2.8 | 98 | 0.00 |
| 51 | T | 4-Methyl-2-Pentanone | 0.569 | 0.547 | 3.9 | 89 | 0.00 |
| 52 | CM | Toluene | 0.990 | 0.994 | -0.4# | 95 | 0.00 |
| 53 | T | t-1,3-Dichloropropene | 0.571 | 0.617 | -8.1 | 96 | 0.00 |
| 54 | T | cis-1,3-Dichloropropene | 0.644 | 0.682 | -5.9 | 98 | 0.00 |
| 55 | T | 1,1,2-Trichloroethane | 0.391 | 0.387 | 1.0 | 97 | 0.00 |
| 56 | T | Ethyl methacrylate | 0.587 | 0.622 | -6.0 | 93 | 0.00 |
| 57 | T | 1,3-Dichloropropane | 0.715 | 0.700 | 2.1 | 96 | 0.00 |
| 58 | T | 2-Chloroethyl Vinyl ether | 0.250 | 0.233 | 6.8 | 83 | 0.00 |
| 59 | T | 2-Hexanone | 0.413 | 0.405 | 1.9 | 90 | 0.00 |
| 60 | T | Dibromochloromethane | 0.392 | 0.396 | -1.0 | 94 | 0.00 |
| 61 | T | 1,2-Dibromoethane | 0.387 | 0.378 | 2.3 | 95 | 0.00 |
| 62 | S | 4-Bromofluorobenzene | 0.442 | 0.483 | -9.3 | 99 | 0.00 |
| 63 | I | Chlorobenzene-d5 | 1.000 | 1.000 | 0.0 | 97 | 0.00 |
| 64 | T | Tetrachloroethene | 0.336 | 0.307 | 8.6 | 92 | 0.00 |
| 65 | PM | Chlorobenzene | 1.150 | 1.111 | 3.4 | 96 | 0.00 |
| 66 | T | 1,1,1,2-Tetrachloroethane | 0.426 | 0.412 | 3.3 | 97 | 0.00 |
| 67 | C | Ethyl Benzene | 2.090 | 2.164 | -3.5# | 97 | 0.00 |
| 68 | T | m/p-Xylenes | 0.775 | 0.803 | -3.6 | 97 | 0.00 |
| 69 | T | o-Xylene | 0.770 | 0.788 | -2.3 | 96 | 0.00 |
| 70 | T | Styrene | 1.199 | 1.297 | -8.2 | 96 | 0.00 |
| 71 | P | Bromoform | 0.288 | 0.286 | 0.7 | 93 | 0.00 |
| 72 | I | 1,4-Dichlorobenzene-d4 | 1.000 | 1.000 | 0.0 | 99 | 0.00 |
| 73 | T | Isopropylbenzene | 4.481 | 4.427 | 1.2 | 96 | 0.00 |
| 74 | T | N-amyl acetate | 1.939 | 1.872 | 3.5 | 92 | 0.00 |
| 75 | P | 1,1,2,2-Tetrachloroethane | 1.623 | 1.358 | 16.3 | 93 | 0.00 |
| 76 | T | 1,2,3-Trichloropropane | 1.293 | 1.040 | 19.6 | 79 | 0.00 |
| 77 | T | Bromobenzene | 1.004 | 0.931 | 7.3 | 98 | 0.00 |
| 78 | T | n-propylbenzene | 5.173 | 5.359 | -3.6 | 98 | 0.00 |
| 79 | T | 2-Chlorotoluene | 3.296 | 3.209 | 2.6 | 97 | 0.00 |
| 80 | T | 1,3,5-Trimethylbenzene | 3.741 | 3.796 | -1.5 | 97 | 0.00 |
| 81 | T | trans-1,4-Dichloro-2-butene | 0.470 | 0.435 | 7.4 | 92 | 0.00 |
| 82 | T | 4-Chlorotoluene | 3.134 | 3.168 | -1.1 | 100 | 0.00 |
| 83 | T | tert-Butylbenzene | 3.141 | 3.155 | -0.4 | 96 | 0.00 |
| 84 | T | 1,2,4-Trimethylbenzene | 3.733 | 3.817 | -2.3 | 98 | 0.00 |
| 85 | T | sec-Butylbenzene | 4.468 | 4.697 | -5.1 | 97 | 0.00 |
| 86 | T | p-Isopropyltoluene | 3.588 | 3.812 | -6.2 | 98 | 0.00 |
| 87 | T | 1,3-Dichlorobenzene | 1.897 | 1.795 | 5.4 | 99 | 0.00 |
| 88 | T | 1,4-Dichlorobenzene | 1.887 | 1.760 | 6.7 | 97 | 0.00 |
| 89 | T | n-Butylbenzene | 3.148 | 3.506 | -11.4 | 100 | 0.00 |
| 90 | T | Hexachloroethane | 0.730 | 0.693 | 5.1 | 99 | 0.00 |
| 91 | T | 1,2-Dichlorobenzene | 1.891 | 1.735 | 8.2 | 95 | 0.00 |
| 92 | T | 1,2-Dibromo-3-Chloropropane | 0.301 | 0.255 | 15.3 | 83 | 0.00 |
| 93 | T | 1,2,4-Trichlorobenzene | 0.855 | 0.898 | -5.0 | 93 | 0.00 |
| 94 | T | Hexachlorobutadiene | 0.471 | 0.450 | 4.5 | 96 | 0.00 |
| 95 | T | Naphthalene | 2.729 | 2.856 | -4.7 | 86 | 0.00 |

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|---|---|---|---|---|---|---|
| 96 T | 1,2,3-Trichlorobenzene | 0.858 | 0.858 | 0.0 | 92 | 0.00 |

(#) = Out of Range SPCC's out = 0 CCC's out = 5