

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120518\
 Data File : VV008840.D
 Acq On : 05 Dec 2018 12:14
 Operator : SY/MD
 Sample : VSTDICV005
 Misc : 25.0 mL/MSVOA_V/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 VICV69

Quant Time: Dec 06 04:54:09 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR120518WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Thu Dec 06 04:51:41 2018
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Difluorobenzene	1.000	1.000	0.0	110	0.00
2 T	Dichlorodifluoromethane	0.475	0.473	0.4	112	0.00
3 T	Chloromethane	0.315	0.314	0.3	111	0.00
4 S	Vinyl Chloride-d3	0.267	0.234	12.4	102	0.00
5 T	Vinyl chloride	0.321	0.322	-0.3	112	0.00
6 T	Bromomethane	0.188	0.198	-5.3	118	0.00
7 S	Chloroethane-d5	0.197	0.179	9.1	103	0.00
8 T	Chloroethane	0.180	0.182	-1.1	116	0.00
9 T	Trichlorofluoromethane	0.517	0.532	-2.9	114	0.00
10 T	1,1,2-Trichloro-1,2,2-trifl	0.281	0.288	-2.5	114	0.00
11 S	1,1-Dichloroethene-d2	0.514	0.490	4.7	109	0.00
12 T	1,1-Dichloroethene	0.245	0.245	0.0	111	0.00
13 T	Acetone	0.037	0.038	-2.7	109	0.00
14 T	Carbon disulfide	0.737	0.766	-3.9	116	0.00
15 T	Methyl Acetate	0.081	0.084	-3.7	115	0.00
16 T	Methylene chloride	0.265	0.258	2.6	116	0.00
17 T	Methyl tert-butyl Ether	0.602	0.605	-0.5	113	0.00
18 T	trans-1,2-Dichloroethene	0.268	0.269	-0.4	116	0.00
19 T	1,1-Dichloroethane	0.585	0.613	-4.8	115	0.00
20 S	2-Butanone-d5	0.066	0.064	3.0	106	0.00
21 T	2-Butanone	0.069	0.072	-4.3	110	0.00
22 T	cis-1,2-Dichloroethene	0.367	0.388	-5.7	116	0.00
23 T	Bromochloromethane	0.155	0.160	-3.2	114	0.00
24 S	Chloroform-d	0.619	0.578	6.6	109	0.00
25 T	Chloroform	0.634	0.637	-0.5	111	0.00
26 S	1,2-Dichloroethane-d4	0.293	0.279	4.8	109	0.00
27 T	1,2-Dichloroethane	0.362	0.374	-3.3	111	0.00
28 I	Chlorobenzene-d5	1.000	1.000	0.0	110	0.00
29 T	1,1,1-Trichloroethane	0.592	0.608	-2.7	113	0.00
30 T	Cyclohexane	0.569	0.611	-7.4	118	0.00
31 T	Carbon tetrachloride	0.533	0.547	-2.6	115	0.00
32 S	Benzene-d6	1.307	1.218	6.8	107	0.00
33 T	Benzene	1.422	1.498	-5.3	117	0.00
34 T	Trichloroethene	0.413	0.425	-2.9	115	0.00
35 T	Methylcyclohexane	0.654	0.675	-3.2	115	0.00
36 S	1,2-Dichloropropane-d6	0.384	0.359	6.5	108	0.00
37 T	1,2-Dichloropropane	0.348	0.370	-6.3	115	0.00
38 T	Bromodichloromethane	0.429	0.445	-3.7	117	0.00
39 T	cis-1,3-Dichloropropene	0.488	0.514	-5.3	116	0.00
40 T	4-Methyl-2-pentanone	0.177	0.188	-6.2	116	0.00
41 S	Toluene-d8	1.237	1.168	5.6	107	0.00
42 T	Toluene	1.534	1.599	-4.2	115	0.00
43 S	trans-1,3-Dichloropropene-d	0.145	0.139	4.1	112	0.00
44 T	trans-1,3-Dichloropropene	0.379	0.406	-7.1	115	0.00
45 T	1,1,2-Trichloroethane	0.245	0.245	0.0	111	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120518\
 Data File : VV008840.D
 Acq On : 05 Dec 2018 12:14
 Operator : SY/MD
 Sample : VSTDICV005
 Misc : 25.0 mL/MSVOA_V/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 VICV69

Quant Time: Dec 06 04:54:09 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR120518WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Thu Dec 06 04:51:41 2018
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
46 S	2-Hexanone-d5	0.055	0.056	-1.8	113	0.00
47 T	Tetrachloroethene	0.332	0.348	-4.8	117	0.00
48 T	2-Hexanone	0.125	0.135	-8.0	116	0.00
49 T	Dibromochloromethane	0.286	0.302	-5.6	115	0.00
50 T	1,2-Dibromoethane	0.229	0.245	-7.0	116	0.00
51 T	Chlorobenzene	1.010	1.054	-4.4	117	0.00
52 T	Ethylbenzene	1.675	1.776	-6.0	118	0.00
53 T	m,p-xylene	0.640	0.682	-6.6	117	0.00
54 T	o-xylene	0.610	0.646	-5.9	115	0.00
55 T	Styrene	1.010	1.091	-8.0	116	0.00
56 T	Isopropylbenzene	1.651	1.768	-7.1	117	0.00
57 S	1,1,2,2-Tetrachloroethane-d	0.257	0.240	6.6	108	0.00
58 T	1,1,2,2-Tetrachloroethane	0.247	0.268	-8.5	115	0.00
59	1,2,3-Trichloropropane	0.208	0.211	-1.4	112	0.00
60 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	114	0.00
61 T	Bromoform	0.297	0.295	0.7	114	0.00
62 T	1,3-Dichlorobenzene	1.617	1.642	-1.5	118	0.00
63 T	1,4-Dichlorobenzene	1.653	1.697	-2.7	118	0.00
64 S	1,2-Dichlorobenzene-d4	0.921	0.852	7.5	109	0.00
65 T	1,2-Dichlorobenzene	1.493	1.544	-3.4	117	0.00
66 T	1,2-Dibromo-3-chloropropane	0.075	0.080	-6.7	116	0.00
67	1,3,5-Trichlorobenzene	1.203	1.294	-7.6	121	0.00
68 T	1,2,4-trichlorobenzene	0.894	1.007	-12.6	122	0.00
69	Naphthalene	1.298	1.552	-19.6	130	0.00
70 T	1,2,3-Trichlorobenzene	0.837	0.952	-13.7	124	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0