Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120518\
Data File : VV008852.D
Acq On : 05 Dec 2018 20:23
Operator : SY/MD
Sample : VSTDCCC005EC
Misc : 25.0 mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument : MSVOA_V
LabSampleId : VSTD00576

Quant Time: Dec 06 05:00:38 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	91	0.00
2	T	Dichlorodifluoromethane	0.475	0.467	1.7	91	0.00
3	T	Chloromethane	0.315	0.312	1.0	91	0.00
4	S	Vinyl Chloride-d3	0.267	0.243	9.0	87	0.00
5	T	Vinyl chloride	0.321	0.321	0.0	92	0.00
6	T	Bromomethane	0.188	0.141	25.0	69	0.00
7	S	Chloroethane-d5	0.197	0.184	6.6	87	0.00
8	T	Chloroethane	0.180	0.170	5.6	90	0.00
9	T	Trichlorofluoromethane	0.517	0.529	-2.3	94	0.00
10	T	1,1,2-Trichloro-1,2,2-trifl	0.281	0.289	-2.8	95	0.00
11	S	1,1-Dichloroethene-d2	0.514	0.491	4.5	90	0.00
12	T	1,1-Dichloroethene	0.245	0.250	-2.0	93	0.00
13	T	Acetone	0.037	0.035	5.4	83	-0.02
14	T	Carbon disulfide	0.737	0.723	1.9	91	0.00
15	T	Methyl Acetate	0.081	0.097	-19.8	109	0.00
16	T	Methylene chloride	0.265	0.262	1.1	97	0.00
17	T	Methyl tert-butyl Ether	0.602	0.604	-0.3	93	0.00
18	T	trans-1,2-Dichloroethene	0.268	0.267	0.4	95	0.00
19	T	1,1-Dichloroethane	0.585	0.577	1.4	89	0.00
20	S	2-Butanone-d5	0.066	0.067	-1.5	92	-0.02
21	T	2-Butanone	0.069	0.069	0.0	87	-0.02
22	T	cis-1,2-Dichloroethene	0.367	0.366	0.3	90	0.00
23	T	Bromochloromethane	0.155	0.159	-2.6	94	0.00
24	S	Chloroform-d	0.619	0.603	2.6	94	0.00
25	T	Chloroform	0.634	0.626	1.3	90	0.00
26	S	1,2-Dichloroethane-d4	0.293	0.293	0.0	94	0.00
27	T	1,2-Dichloroethane	0.362	0.371	-2.5	91	0.00
28	I	Chlorobenzene-d5	1.000	1.000	0.0	92	0.00
29	T	1,1,1-Trichloroethane	0.592	0.578	2.4	90	0.00
30	T	Cyclohexane	0.569	0.560	1.6	91	0.00
31	T	Carbon tetrachloride	0.533	0.520	2.4	92	0.00
32	S	Benzene-d6	1.307	1.236	5.4	91	0.00
33	T	Benzene	1.422	1.406	1.1	92	0.00
34	T	Trichloroethene	0.413	0.404	2.2	92	0.00
35	T	Methylcyclohexane	0.654	0.616	5.8	88	0.00
36	S	1,2-Dichloropropane-d6	0.384	0.370	3.6	93	0.00
37	T	1,2-Dichloropropane	0.348	0.345	0.9	89	0.00
38	T	Bromodichloromethane	0.429	0.409	4.7	90	0.00
39	T	cis-1,3-Dichloropropene	0.488	0.464	4.9	88	0.00
40	T	4-Methyl-2-pentanone	0.177	0.177	0.0	91	0.00
41	S	Toluene-d8	1.237	1.174	5.1	90	0.00
42	T	Toluene	1.534	1.526	0.5	92	0.00
43	S	trans-1,3-Dichloropropene-d	0.145	0.130	10.3	87	0.00
44	T	trans-1,3-Dichloropropene	0.379	0.364	4.0	86	0.00
45	T	1,1,2-Trichloroethane	0.245	0.246	-0.4	93	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120518\
Data File : VV008852.D
Acq On : 05 Dec 2018 20:23
Operator : SY/MD
Sample : VSTDCCC005EC
Misc : 25.0 mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument : MSVOA_V
LabSampleId : VSTD00576

Quant Time: Dec 06 05:00:38 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	94	0.00
47	T	Tetrachloroethene	0.332	0.329	0.9	92	0.00
48	T	2-Hexanone	0.125	0.122	2.4	88	0.00
49	T	Dibromochloromethane	0.286	0.284	0.7	91	0.00
50	T	1,2-Dibromoethane	0.229	0.226	1.3	90	0.00
51	T	Chlorobenzene	1.010	0.998	1.2	93	0.00
52	T	Ethylbenzene	1.675	1.648	1.6	91	0.00
53	T	m,p-xylene	0.640	0.639	0.2	92	0.00
54	T	o-xylene	0.610	0.606	0.7	90	0.00
55	T	Styrene	1.010	1.020	-1.0	90	0.00
56	T	Isopropylbenzene	1.651	1.669	-1.1	93	0.00
57	S	1,1,2,2-Tetrachloroethane-d	0.257	0.245	4.7	92	0.00
58	T	1,1,2,2-Tetrachloroethane	0.247	0.256	-3.6	92	0.00
59		1,2,3-Trichloropropane	0.208	0.202	2.9	89	0.00
60	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	92	0.00
61	T	Bromoform	0.297	0.289	2.7	90	0.00
62	T	1,3-Dichlorobenzene	1.617	1.569	3.0	91	0.00
63	T	1,4-Dichlorobenzene	1.653	1.599	3.3	91	0.00
64	S	1,2-Dichlorobenzene-d4	0.921	0.891	3.3	93	0.00
65	T	1,2-Dichlorobenzene	1.493	1.483	0.7	91	0.00
66	T	1,2-Dibromo-3-chloropropane	0.075	0.078	-4.0	93	0.00
67		1,3,5-Trichlorobenzene	1.203	1.198	0.4	91	0.00
68	T	1,2,4-trichlorobenzene	0.894	0.875	2.1	86	0.00
69		Naphthalene	1.298	1.209	6.9	82	0.00
70	T	1,2,3-Trichlorobenzene	0.837	0.843	-0.7	89	0.00

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