Evaluate Continuing Calibration Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV100418\
Data File : VV008055.D
Acq On : 04 Oct 2018 11:17
Operator : SY/MD
Sample : VSTDCCC005
Misc : 25.0 mL/MSVOA_V/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument : MSVOA_V
LabSampleId : VSTD00570

Quant Time: Oct 05 05:18:15 2018
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR092718WMA.M
Quant Title : TRACE VOA SOM01.0
QLast Update : Thu Oct 04 07:54:45 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	102	0.00
2	T	Dichlorodifluoromethane	0.419	0.425	-1.4	104	0.00
3	T	Chloromethane	0.373	0.350	6.2	96	0.00
4	S	Vinyl Chloride-d3	0.268	0.231	13.8	89	0.00
5	T	Vinyl chloride	0.399	0.376	5.8	96	0.00
6	T	Bromomethane	0.246	0.231	6.1	96	0.00
7	S	Chloroethane-d5	0.232	0.205	11.6	92	0.00
8	T	Chloroethane	0.241	0.229	5.0	98	0.00
9	T	Trichlorofluoromethane	0.652	0.639	2.0	100	0.00
10	T	1,1,2-Trichloro-1,2,2-trifl	0.370	0.375	-1.4	104	0.00
11	S	1,1-Dichloroethene-d2	0.605	0.534	11.7	93	0.00
12	T	1,1-Dichloroethene	0.331	0.317	4.2	98	0.00
13	T	Acetone	0.059	0.053	10.2	99	0.00
14	T	Carbon disulfide	0.896	0.775	13.5	93	0.00
15	T	Methyl Acetate	0.150	0.133	11.3	104	0.00
16	T	Methylene chloride	0.367	0.341	7.1	102	0.00
17	T	Methyl tert-butyl Ether	0.817	0.780	4.5	101	0.00
18	T	trans-1,2-Dichloroethene	0.364	0.344	5.5	99	0.00
19	T	1,1-Dichloroethane	0.622	0.637	-2.4	103	0.00
20	S	2-Butanone-d5	0.081	0.080	1.2	108	0.00
21	T	2-Butanone	0.094	0.088	6.4	103	0.00
22	T	cis-1,2-Dichloroethene	0.369	0.388	-5.1	108	0.00
23	T	Bromochloromethane	0.160	0.175	-9.4	112	0.00
24	S	Chloroform-d	0.611	0.596	2.5	101	0.00
25	T	Chloroform	0.672	0.703	-4.6	107	0.00
26	S	1,2-Dichloroethane-d4	0.336	0.307	8.6	97	0.00
27	T	1,2-Dichloroethane	0.433	0.415	4.2	98	0.00
28	I	Chlorobenzene-d5	1.000	1.000	0.0	101	0.00
29	T	1,1,1-Trichloroethane	0.601	0.644	-7.2	110	0.00
30	T	Cyclohexane	0.565	0.542	4.1	99	0.00
31	T	Carbon tetrachloride	0.530	0.576	-8.7	111	0.00
32	S	Benzene-d6	1.194	1.151	3.6	100	0.00
33	T	Benzene	1.432	1.477	-3.1	106	0.00
34	T	Trichloroethene	0.410	0.412	-0.5	105	0.00
35	T	Methylcyclohexane	0.615	0.599	2.6	101	0.00
36	S	1,2-Dichloropropane-d6	0.378	0.363	4.0	101	0.00
37	T	1,2-Dichloropropane	0.366	0.386	-5.5	111	0.00
38	T	Bromodichloromethane	0.437	0.477	-9.2	113	0.00
39	T	cis-1,3-Dichloropropene	0.489	0.517	-5.7	111	0.00
40	T	4-Methyl-2-pentanone	0.230	0.224	2.6	104	0.00
41	S	Toluene-d8	1.215	1.151	5.3	97	0.00
42	T	Toluene	1.559	1.637	-5.0	107	0.00
43	S	trans-1,3-Dichloropropene-d	0.146	0.143	2.1	106	0.00
44	T	trans-1,3-Dichloropropene	0.407	0.430	-5.7	109	0.00
45	T	1,1,2-Trichloroethane	0.264	0.267	-1.1	108	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV100418\
Data File : VV008055.D
Acq On : 04 Oct 2018 11:17
Operator : SY/MD
Sample : VSTDCCC005
Misc : 25.0 mL/MSVOA_V/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument : MSVOA_V
LabSampleId : VSTD00570

Quant Time: Oct 05 05:18:15 2018
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR092718WMA.M
Quant Title : TRACE VOA SOM01.0
QLast Update : Thu Oct 04 07:54:45 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	0.064	0.061	4.7	101	0.00
47	T	Tetrachloroethene	0.380	0.391	-2.9	107	0.00
48	T	2-Hexanone	0.160	0.158	1.3	103	0.00
49	T	Dibromochloromethane	0.301	0.338	-12.3	117	0.00
50	T	1,2-Dibromoethane	0.241	0.249	-3.3	105	0.00
51	T	Chlorobenzene	1.077	1.116	-3.6	108	0.00
52	T	Ethylbenzene	1.739	1.780	-2.4	105	0.00
53	T	m,p-xylene	0.662	0.676	-2.1	105	0.00
54	T	o-xylene	0.640	0.654	-2.2	105	0.00
55	T	Styrene	1.093	1.146	-4.8	106	0.00
56	T	Isopropylbenzene	1.733	1.808	-4.3	106	0.00
57	S	1,1,2,2-Tetrachloroethane-d	0.292	0.278	4.8	101	0.00
58	T	1,1,2,2-Tetrachloroethane	0.310	0.312	-0.6	106	0.00
59		1,2,3-Trichloropropane	0.228	0.231	-1.3	107	0.00
60	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	99	0.00
61	T	Bromoform	0.300	0.347	-15.7	124	0.00
62	T	1,3-Dichlorobenzene	1.651	1.701	-3.0	107	0.00
63	T	1,4-Dichlorobenzene	1.716	1.739	-1.3	106	0.00
64	S	1,2-Dichlorobenzene-d4	0.968	0.860	11.2	97	0.00
65	T	1,2-Dichlorobenzene	1.556	1.647	-5.8	110	0.00
66	T	1,2-Dibromo-3-chloropropane	0.079	0.085	-7.6	109	0.00
67		1,3,5-Trichlorobenzene	1.413	1.483	-5.0	109	0.00
68	T	1,2,4-trichlorobenzene	1.195	1.126	5.8	101	0.00
69		Naphthalene	1.671	1.374	17.8	86	0.00
70	T	1,2,3-Trichlorobenzene	1.082	1.066	1.5	102	0.00

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