Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU071918\
Data File : VU025512.D
Acq On : 19 Jul 2018 09:46
Operator : MD/SY
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTD05051

Quant Time: Jul 20 02:35:21 2018
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM071718WMA.M
Quant Title : VOC Analysis
QLast Update : Thu Jul 19 01:07:14 2018
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Difluorobenzene	1.000	1.000	0.0	95	0.00
2	T	Dichlorodifluoromethane	0.411	0.413	-0.5	103	0.00
3	T	Chloromethane	0.364	0.363	0.3	99	0.00
4	S	Vinyl Chloride-d3	0.236	0.205	13.1	83	0.00
5	T	Vinyl chloride	0.366	0.362	1.1	97	0.00
6	T	Bromomethane	0.199	0.234	-17.6	112	0.00
7	S	Chloroethane-d5	0.201	0.187	7.0	93	0.00
8	T	Chloroethane	0.219	0.222	-1.4	101	0.00
9	T	Trichlorofluoromethane	0.538	0.555	-3.2	104	0.00
10	T	1,1,2-Trichloro-1,2,2-trifl	0.292	0.317	-8.6	117	0.00
11	S	1,1-Dichloroethene-d2	0.518	0.494	4.6	93	0.00
12	T	1,1-Dichloroethene	0.271	0.274	-1.1	100	0.00
13	T	Acetone	0.219	0.260	-18.7	128	0.00
14	T	Carbon disulfide	0.876	0.844	3.7	97	0.00
15	T	Methyl Acetate	0.405	0.344	15.1	82	0.00
16	T	Methylene chloride	0.354	0.367	-3.7	101	0.00
17	T	trans-1,2-Dichloroethene	0.321	0.325	-1.2	99	0.00
18	T	Methyl tert-butyl Ether	1.136	1.104	2.8	92	0.00
19	T	1,1-Dichloroethane	0.641	0.638	0.5	96	0.00
20	T	cis-1,2-Dichloroethene	0.392	0.381	2.8	95	0.00
21	S	2-Butanone-d5	0.251	0.216	13.9	79	0.00
22	T	2-Butanone	0.302	0.297	1.7	95	0.00
23	T	Bromochloromethane	0.197	0.202	-2.5	97	0.00
24	S	Chloroform-d	0.612	0.594	2.9	92	0.00
25	T	Chloroform	0.674	0.670	0.6	95	0.00
26	S	1,2-Dichloroethane-d4	0.416	0.389	6.5	89	0.00
27	T	1,2-Dichloroethane	0.560	0.542	3.2	93	0.00
28	I	Chlorobenzene-d5	1.000	1.000	0.0	97	0.00
29	T	Cyclohexane	0.576	0.580	-0.7	103	0.00
30	T	1,1,1-Trichloroethane	0.665	0.652	2.0	97	0.00
31	T	Carbon tetrachloride	0.585	0.585	0.0	98	0.00
32	S	Benzene-d6	1.208	1.121	7.2	89	0.00
33	T	Benzene	1.479	1.452	1.8	97	0.00
34	T	Trichloroethene	0.392	0.383	2.3	99	0.00
35	T	Methylcyclohexane	0.595	0.631	-6.1	117	0.00
36	S	1,2-Dichloropropane-d6	0.400	0.382	4.5	92	0.00
37	T	1,2-Dichloropropane	0.408	0.400	2.0	95	0.00
38	T	Bromodichloromethane	0.551	0.543	1.5	96	0.00
39	T	cis-1,3-Dichloropropene	0.632	0.644	-1.9	97	0.00
40	T	4-Methyl-2-pentanone	0.592	0.501	15.4	83	0.00
41	S	Toluene-d8	1.170	1.094	6.5	91	0.00
42	T	Toluene	1.577	1.597	-1.3	101	0.00
43	S	trans-1,3-Dichloropropene-d	0.208	0.195	6.2	90	0.00
44	T	trans-1,3-Dichloropropene	0.607	0.608	-0.2	98	0.00
45	T	1,1,2-Trichloroethane	0.387	0.387	0.0	98	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU071918\
Data File : VU025512.D
Acq On : 19 Jul 2018 09:46
Operator : MD/SY
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTD05051

Quant Time: Jul 20 02:35:21 2018
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM071718WMA.M
Quant Title : VOC Analysis
QLast Update : Thu Jul 19 01:07:14 2018
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
46	T	Tetrachloroethene	0.327	0.334	-2.1	109	0.00
47	S	2-Hexanone-d5	0.207	0.169	18.4	78	0.00
48	T	2-Hexanone	0.473	0.427	9.7	89	0.00
49	T	Dibromochloromethane	0.468	0.466	0.4	97	0.00
50	T	1,2-Dibromoethane	0.427	0.409	4.2	94	0.00
51	T	Chlorobenzene	1.058	1.079	-2.0	102	0.00
52	T	Ethylbenzene	1.761	1.789	-1.6	102	0.00
53	T	m,p-Xylene	0.675	0.696	-3.1	105	0.00
54	T	o-xylene	0.681	0.698	-2.5	103	0.00
55	T	Styrene	1.124	1.172	-4.3	103	0.00
56	T	Isopropylbenzene	1.763	1.825	-3.5	105	0.00
57	S	1,1,2,2-Tetrachloroethane-d	0.594	0.562	5.4	92	0.00
58	T	1,1,2,2-Tetrachloroethane	0.666	0.617	7.4	93	0.00
59		1,2,3-Trichloropropane	0.543	0.487	10.3	90	0.00
60	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	105	0.00
61	T	Bromoform	0.705	0.658	6.7	95	0.00
62	T	1,3-Dichlorobenzene	1.657	1.666	-0.5	110	0.00
63	T	1,4-Dichlorobenzene	1.675	1.681	-0.4	111	0.00
64	S	1,2-Dichlorobenzene-d4	1.002	0.940	6.2	99	0.00
65	T	1,2-Dichlorobenzene	1.750	1.749	0.1	108	0.00
66	T	1,2-Dibromo-3-chloropropane	0.327	0.259	20.8	81	0.00
67		1,3,5-Trichlorobenzene	1.274	1.334	-4.7	122	0.00
68	T	1,2,4-trichlorobenzene	1.136	1.132	0.4	113	0.00
69		Naphthalene	3.930	3.381	14.0	89	0.00
70	T	1,2,3-Trichlorobenzene	1.225	1.187	3.1	105	0.00

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