Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092123\
Data File : VU055451.D
Acq On : 21 Sep 2023 17:24
Operator : MD/SY
Sample : VSTDCCC005EC
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_U
LabSampleId :
VSTDCCC005EC

Quant Time: Sep 22 02:36:49 2023
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR091923WMA.M
Quant Title : TRACE VOA SFAM1.0
QLast Update : Fri Sep 22 02:32:08 2023
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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Difluorobenzene	5.000	5.000	0.0	112	0.00
2	T	Dichlorodifluoromethane	5.000	4.286	14.3	96	0.00
3	T	Chloromethane	5.000	4.484	10.3	102	0.00
4	S	Vinyl Chloride-d3	5.000	3.549	29.0	79	0.00
5	T	Vinyl chloride	5.000	4.566	8.7	104	0.00
6	T	Bromomethane	5.000	4.907	1.9	109	0.00
7	S	Chloroethane-d5	5.000	4.034	19.3	93	0.00
8	T	Chloroethane	5.000	4.683	6.3	109	0.00
9	T	Trichlorofluoromethane	5.000	4.665	6.7	110	0.00
10	T	1,1,2-Trichloro-1,2,2-trifl	5.000	4.690	6.2	108	0.00
11	S	1,1-Dichloroethene-d2	5.000	3.918	21.6	89	0.00
12	T	1,1-Dichloroethene	5.000	4.781	4.4	108	0.00
13	T	Acetone	50.000	44.795	10.4	105	0.00
14	T	Carbon disulfide	5.000	4.762	4.8	109	0.00
15	T	Methyl Acetate	5.000	5.016	-0.3	120	0.00
16	T	Methylene chloride	5.000	4.439	11.2	95	0.00
17	T	Methyl tert-butyl Ether	5.000	4.925	1.5	112	0.00
18	T	trans-1,2-Dichloroethene	5.000	4.886	2.3	110	0.00
19	T	1,1-Dichloroethane	5.000	5.008	-0.2	113	0.00
20	S	2-Butanone-d5	50.000	51.396	-2.8	113	0.00
21	T	2-Butanone	50.000	48.006	4.0	109	0.00
22	T	cis-1,2-Dichloroethene	5.000	4.714	5.7	110	0.00
23	T	Bromochloromethane	5.000	4.911	1.8	112	0.00
24	S	Chloroform-d	5.000	4.840	3.2	106	0.00
25	T	Chloroform	5.000	5.079	-1.6	114	0.00
26	S	1,2-Dichloroethane-d4	5.000	5.099	-2.0	111	0.00
27	T	1,2-Dichloroethane	5.000	5.085	-1.7	117	0.00
28	I	Chlorobenzene-d5	5.000	5.000	0.0	111	0.00
29	T	1,1,1-Trichloroethane	5.000	5.146	-2.9	113	0.00
30	T	Cyclohexane	5.000	4.785	4.3	108	0.00
31	T	Carbon tetrachloride	5.000	5.211	-4.2	113	0.00
32	S	Benzene-d6	5.000	4.665	6.7	99	0.00
33	T	Benzene	5.000	5.071	-1.4	111	0.00
34	T	Trichloroethene	5.000	5.003	-0.1	114	0.00
35	T	Methylcyclohexane	5.000	4.789	4.2	107	0.00
36	S	1,2-Dichloropropane-d6	5.000	4.820	3.6	104	0.00
37	T	1,2-Dichloropropane	5.000	4.910	1.8	112	0.00
38	T	Bromodichloromethane	5.000	5.169	-3.4	114	0.00
39	T	cis-1,3-Dichloropropene	5.000	5.264	-5.3	112	0.00
40	T	4-Methyl-2-pentanone	50.000	50.670	-1.3	113	0.00
41	S	Toluene-d8	5.000	4.657	6.9	98	0.00
42	T	Toluene	5.000	5.180	-3.6	112	0.00
43	S	trans-1,3-Dichloropropene-d	5.000	5.112	-2.2	108	0.00
44	T	trans-1,3-Dichloropropene	5.000	5.274	-5.5	115	0.00
45	T	1,1,2-Trichloroethane	5.000	5.119	-2.4	113	0.00
46	S	2-Hexanone-d5	50.000	50.728	-1.5	113	0.00
47	T	Tetrachloroethene	5.000	5.098	-2.0	111	0.00
48	T	2-Hexanone	50.000	53.189	-6.4	117	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092123\
Data File : VU055451.D
Acq On : 21 Sep 2023 17:24
Operator : MD/SY
Sample : VSTDCCC005EC
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_U
LabSampleId :
VSTDCCC005EC

Quant Time: Sep 22 02:36:49 2023
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR091923WMA.M
Quant Title : TRACE VOA SFAM1.0
QLast Update : Fri Sep 22 02:32:08 2023
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	Amount	Calc.	%Dev	Area%	Dev(min)
49	T	Dibromochloromethane	5.000	5.264	-5.3	117	0.00
50	T	1,2-Dibromoethane	5.000	5.136	-2.7	114	0.00
51	T	Chlorobenzene	5.000	5.014	-0.3	112	0.00
52	T	Ethylbenzene	5.000	5.027	-0.5	111	0.00
53	T	m,p-Xylene	5.000	5.233	-4.7	112	0.00
54	T	o-Xylene	5.000	5.128	-2.6	115	0.00
55	T	Styrene	5.000	5.337	-6.7	114	0.00
56	S	1,1,2,2-Tetrachloroethane-d	5.000	5.110	-2.2	110	0.00
57	T	1,1,2,2-Tetrachloroethane	5.000	4.971	0.6	110	0.00
58	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	110	0.00
59	T	Bromoform	5.000	5.207	-4.1	115	0.00
60		Isopropylbenzene	5.000	5.115	-2.3	112	0.00
61		1,2,3-Trichloropropane	5.000	5.017	-0.3	112	0.00
62		1,3,5-Trimethylbenzene	5.000	5.048	-1.0	113	0.00
63		1,2,4-Trimethylbenzene	5.000	5.032	-0.6	112	0.00
64	T	1,3-Dichlorobenzene	5.000	4.978	0.4	111	0.00
65	T	1,4-Dichlorobenzene	5.000	4.913	1.7	111	0.00
66	S	1,2-Dichlorobenzene-d4	5.000	4.827	3.5	106	0.00
67	T	1,2-Dichlorobenzene	5.000	5.115	-2.3	115	0.00
68	T	1,2-Dibromo-3-chloropropane	5.000	5.610	-12.2	118	0.00
69	MA	1,3,5-Trichlorobenzene	5.000	4.778	4.4	109	0.00
70	T	1,2,4-trichlorobenzene	5.000	4.867	2.7	112	0.00
71	MA	Naphthalene	5.000	4.870	2.6	116	0.00
72	T	1,2,3-Trichlorobenzene	5.000	4.892	2.2	110	0.00

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