Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU101824\
Data File : VU061156.D
Acq On : 18 Oct 2024 10:26
Operator : MD/SY
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_U
LabSampleId :
VSTDCCC010

Quant Time: Oct 18 23:45:10 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U101724DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
QLast Update : Fri Oct 18 02:52:28 2024
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	Fluorobenzene	1.000	1.000	0.0	97	0.00
2	T	Dichlorodifluoromethane	0.551	0.608	-10.3	100	0.00
3	t	Chloromethane	0.524	0.574	-9.5	103	0.00
4	Rt	Vinyl Chloride	0.538	0.603	-12.1	102	0.00
5	T	Bromomethane	0.338	0.369	-9.2	98	0.00
6	T	Chloroethane	0.361	0.419	-16.1	108	0.00
7	T	Trichlorofluoromethane	0.766	0.976	-27.4	118	0.00
8		1,1,2-Trichloro-1,2,2-trifl	0.381	0.422	-10.8	99	0.00
9	Rt	1,1-Dichloroethene	0.397	0.451	-13.6	102	0.00
10	t	Iodomethane	0.472	0.620	-31.4#	116	0.00
11	t	Allyl Chloride	0.511	0.562	-10.0	98	0.00
12	t	Acrylonitrile	0.080	0.090	-12.5	104	0.00
13	T	Acetone	0.104	0.114	-9.6	93	0.00
14	T	Carbon Disulfide	1.367	1.497	-9.5	100	0.00
15	RT	Methylene Chloride	0.513	0.509	0.8	98	0.00
16	RT	trans-1,2-Dichloroethene	0.465	0.511	-9.9	99	0.00
17	t	1,1-Dichloroethane	0.825	0.921	-11.6	101	0.00
18	T	2-Butanone	0.127	0.139	-9.4	95	0.00
19		Cyclohexane	0.739	0.804	-8.8	98	0.00
20		Methylcyclohexane	0.824	0.915	-11.0	97	0.00
21	T	2,2-Dichloropropane	0.761	0.819	-7.6	98	0.00
22	RT	cis-1,2-Dichloroethene	0.507	0.564	-11.2	100	0.00
23	t	Diethyl Ether	0.308	0.338	-9.7	104	0.00
24	t	tert-Butyl Alcohol	0.032	0.041	-28.1	132	0.00
25	t	Methyl tert-Butyl Ether	1.063	1.185	-11.5	102	0.00
26	t	Bromochloromethane	0.212	0.239	-12.7	104	0.00
27	t	Chloroform	0.865	0.970	-12.1	102	0.00
28	RT	1,1,1-Trichloroethane	0.763	0.847	-11.0	99	0.00
29	T	1,1-Dichloropropene	0.674	0.736	-9.2	97	0.00
30	RT	Carbon Tetrachloride	0.662	0.743	-12.2	99	0.00
31	t	Isopropyl Ether	1.177	1.324	-12.5	101	0.00
32		Ethyl-t-butyl ether	1.210	1.368	-13.1	102	0.00
33		Tert-Amyl methyl ether	1.089	1.263	-16.0	104	0.00
34	t	Propionitrile	0.030	0.036	-20.0	101	0.00
35	RT	Benzene	1.898	2.140	-12.8	101	0.00
36	RT	1,2-Dichloroethane	0.536	0.596	-11.2	100	0.00
37	RT	Trichloroethene	0.507	0.559	-10.3	99	0.00
38	Rt	1,2-Dichloropropane	0.481	0.541	-12.5	100	0.00
39	t	Methacrylonitrile	0.109	0.130	-19.3	100	0.00
40	t	Methyl acrylate	0.222	0.246	-10.8	103	0.00
41	t	Tetrahydrofuran	0.066	0.069	-4.5	102	0.00
42	t	1-Chlorobutane	0.865	0.955	-10.4	97	0.00
43	T	Dibromomethane	0.249	0.275	-10.4	101	0.00
44	T	Bromodichloromethane	0.606	0.683	-12.7	101	0.00
45	T	4-Methyl-2-Pentanone	0.250	0.275	-10.0	99	0.00
46	t	t-1,4-Dichloro-2-butene	0.100	0.108	-8.0	90	0.00
47	t	Methyl methacrylate	0.217	0.250	-15.2	101	0.00
48	t	Ethyl methacrylate	0.462	0.528	-14.3	101	0.00
49	Rt	Toluene	1.199	1.354	-12.9	100	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
50	T	t-1,3-Dichloropropene	0.584	0.671	-14.9	100	0.00
51	T	cis-1,3-Dichloropropene	0.717	0.821	-14.5	100	0.00
52	RT	1,1,2-Trichloroethane	0.336	0.376	-11.9	102	0.00
53	t	1,3-Dichloropropane	0.591	0.657	-11.2	100	0.00
54	t	2-Hexanone	0.206	0.220	-6.8	92	0.00
55	t	Dibromochloromethane	0.397	0.458	-15.4	101	0.00
56	T	1,2-Dibromoethane	0.323	0.360	-11.5	100	0.00
57	S	4-Bromofluorobenzene	0.531	0.544	-2.4	95	0.00
58	RT	Tetrachloroethene	0.430	0.472	-9.8	99	0.00
59	Rt	Chlorobenzene	1.292	1.439	-11.4	99	0.00
60	T	1,1,1,2-Tetrachloroethane	0.436	0.501	-14.9	102	0.00
61	t	Pentachloroethane	0.333	0.395	-18.6	101	0.00
62	t	Hexachloroethane	0.344	0.410	-19.2	101	0.00
63	Rt	Ethyl Benzene	2.314	2.593	-12.1	99	0.00
64	RT	m/p-Xylenes	0.881	0.990	-12.4	98	0.00
65	RT	o-Xylene	0.867	0.978	-12.8	99	0.00
66	RT	Styrene	1.349	1.564	-15.9	100	0.00
67	t	Bromoform	0.208	0.246	-18.3	101	0.00
68	S	1,2-Dichlorobenzene-d4	0.524	0.556	-6.1	98	0.00
69	T	Isopropylbenzene	2.064	2.321	-12.5	98	0.00
70	T	1,1,2,2-Tetrachloroethane	0.408	0.467	-14.5	103	0.00
71	T	1,2,3-Trichloropropane	0.333	0.358	-7.5	99	0.00
72	t	Bromobenzene	0.496	0.549	-10.7	98	0.00
73	t	n-propylbenzene	0.602	0.694	-15.3	99	0.00
74	t	2-Chlorotoluene	0.521	0.588	-12.9	99	0.00
75	t	1,3,5-Trimethylbenzene	1.861	2.128	-14.3	99	0.00
76	t	4-Chlorotoluene	0.529	0.600	-13.4	98	0.00
77	t	tert-Butylbenzene	1.839	2.103	-14.4	99	0.00
78	t	1,2,4-Trimethylbenzene	1.878	2.151	-14.5	99	0.00
79	t	sec-Butylbenzene	2.395	2.746	-14.7	98	0.00
80		Nitrobenzene	0.009	0.011	-22.2	95	0.00
81	t	p-Isopropyltoluene	1.983	2.284	-15.2	99	0.00
82	t	1,3-Dichlorobenzene	0.952	1.084	-13.9	99	0.00
83	Rt	1,4-Dichlorobenzene	0.949	1.088	-14.6	99	0.00
84	t	n-Butylbenzene	1.840	2.162	-17.5	98	0.00
85	Rt	1,2-Dichlorobenzene	0.890	1.014	-13.9	100	0.00
86	t	1,2-Dibromo-3-Chloropropane	0.061	0.073	-19.7	98	0.00
87	Rt	1,2,4-Trichlorobenzene	0.509	0.620	-21.8	100	0.00
88	t	Hexachlorobutadiene	0.288	0.331	-14.9	99	0.00
89	t	Naphthalene	0.898	1.099	-22.4	103	0.00
90	t	1,2,3-Trichlorobenzene	0.457	0.544	-19.0	99	0.00

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