

Data Path : Z:\voasrv\HPCHEM1\MSVOA W\Data\VW072318\
 Data File : VW004153.D
 Acq On : 23 Jul 2018 23:25
 Operator : SY/AP
 Sample : VSTDCCC050
 Misc : 5.00G/5ML/MSVOA W/SOIL
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 MSVOA_W
 LabSampleId :
 VSTDCCC050

Quant Time: Jul 24 06:37:56 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W071618S.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 16 15:06:12 2018
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	Pentafluorobenzene	1.000	1.000	0.0	90	0.00
2 T	Dichlorodifluoromethane	0.281	0.239	14.9	80	0.00
3 P	Chloromethane	0.370	0.377	-1.9	96	0.00
4 C	Vinyl Chloride	0.508	0.513	-1.0#	92	0.00
5 T	Bromomethane	0.318	0.306	3.8	93	0.00
6 T	Chloroethane	0.289	0.284	1.7	92	0.00
7 T	Trichlorofluoromethane	0.213	0.208	2.3	93	0.00
8 T	Diethyl Ether	0.248	0.279	-12.5	104	0.00
9 T	1,1,2-Trichlorotrifluoroeth	0.492	0.516	-4.9	97	0.00
10 T	Methyl Iodide	0.707	0.771	-9.1	99	0.00
11 T	Tert butyl alcohol	0.039	0.045	-15.4	105	0.00
12 CM	1,1-Dichloroethene	0.486	0.516	-6.2#	98	0.00
13 T	Acrolein	0.040	0.055	-37.5#	138	0.00
14 T	Allyl chloride	0.802	0.910	-13.5	106	0.00
15 T	Acrylonitrile	0.127	0.155	-22.0#	111	0.00
16 T	Acetone	0.134	0.124	7.5	87	0.00
17 T	Carbon Disulfide	1.517	1.615	-6.5	97	0.00
18 T	Methyl Acetate	0.322	0.417	-29.5#	116	0.00
19 T	Methyl tert-butyl Ether	0.931	1.101	-18.3	105	0.00
20 T	Methylene Chloride	0.833	0.593	28.8#	92	0.00
21 T	trans-1,2-Dichloroethene	0.533	0.588	-10.3	101	0.00
22 T	Diisopropyl ether	1.621	1.925	-18.8	105	0.00
23 T	Vinyl Acetate	1.027	1.195	-16.4	101	0.00
24 P	1,1-Dichloroethane	0.989	1.117	-12.9	103	0.00
25 T	2-Butanone	0.181	0.206	-13.8	103	0.00
26 T	2,2-Dichloropropane	0.568	0.524	7.7	88	0.00
27 T	cis-1,2-Dichloroethene	0.576	0.634	-10.1	100	0.00
28 T	Bromochloromethane	0.405	0.416	-2.7	97	0.00
29 T	Tetrahydrofuran	0.113	0.138	-22.1#	109	0.00
30 C	Chloroform	0.984	1.104	-12.2#	104	0.00
31 T	Cyclohexane	0.927	0.961	-3.7	98	0.00
32 T	1,1,1-Trichloroethane	0.768	0.831	-8.2	100	0.00
33 S	1,2-Dichloroethane-d4	0.559	0.662	-18.4	101	0.00
34 I	1,4-Difluorobenzene	1.000	1.000	0.0	94	0.00
35 S	Dibromofluoromethane	0.321	0.366	-14.0	100	0.00
36 T	1,1-Dichloropropene	0.514	0.545	-6.0	99	0.00
37 T	Ethyl Acetate	0.248	0.294	-18.5	110	0.00
38 T	Carbon Tetrachloride	0.481	0.501	-4.2	98	0.00
39 T	Methylcyclohexane	0.602	0.627	-4.2	94	0.00
40 TM	Benzene	1.427	1.567	-9.8	103	0.00
41 T	Methacrylonitrile	0.137	0.172	-25.5#	122	0.00
42 TM	1,2-Dichloroethane	0.435	0.484	-11.3	104	0.00
43 T	Isopropyl Acetate	0.464	0.537	-15.7	106	0.00
44 TM	Trichloroethene	0.370	0.396	-7.0	101	0.00
45 C	1,2-Dichloropropane	0.364	0.413	-13.5#	106	0.00

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 Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
46 T	Dibromomethane	0.189	0.213	-12.7	105	0.00
47 T	Bromodichloromethane	0.466	0.516	-10.7	103	0.00
48 T	Methyl methacrylate	0.224	0.265	-18.3	106	0.00
49 T	1,4-Dioxane	0.003	0.004	-33.3#	115	0.00
50 S	Toluene-d8	1.245	1.417	-13.8	98	0.00
51 T	4-Methyl-2-Pentanone	0.247	0.297	-20.2#	108	0.00
52 CM	Toluene	0.883	0.970	-9.9#	102	0.00
53 T	t-1,3-Dichloropropene	0.487	0.521	-7.0	98	0.00
54 T	cis-1,3-Dichloropropene	0.557	0.595	-6.8	98	0.00
55 T	1,1,2-Trichloroethane	0.266	0.302	-13.5	106	0.00
56 T	Ethyl methacrylate	0.344	0.414	-20.3#	105	0.00
57 T	1,3-Dichloropropane	0.462	0.527	-14.1	105	0.00
58 T	2-Chloroethyl Vinyl ether	0.158	0.182	-15.2	103	0.00
59 T	2-Hexanone	0.174	0.208	-19.5	107	0.00
60 T	Dibromochloromethane	0.310	0.352	-13.5	105	0.00
61 T	1,2-Dibromoethane	0.253	0.292	-15.4	107	0.00
62 S	4-Bromofluorobenzene	0.449	0.518	-15.4	99	0.00
63 I	Chlorobenzene-d5	1.000	1.000	0.0	96	0.00
64 T	Tetrachloroethene	0.343	0.399	-16.3	115	0.00
65 PM	Chlorobenzene	1.063	1.136	-6.9	103	0.00
66 T	1,1,1,2-Tetrachloroethane	0.369	0.402	-8.9	103	0.00
67 C	Ethyl Benzene	1.868	2.020	-8.1#	102	0.00
68 T	m/p-Xylenes	0.714	0.772	-8.1	102	0.00
69 T	o-Xylene	0.662	0.719	-8.6	101	0.00
70 T	Styrene	1.106	1.245	-12.6	104	0.00
71 P	Bromoform	0.209	0.232	-11.0	103	0.00
72 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	95	0.00
73 T	Isopropylbenzene	3.533	3.736	-5.7	100	0.00
74 T	N-amyl acetate	0.921	1.042	-13.1	104	0.00
75 P	1,1,2,2-Tetrachloroethane	0.683	0.760	-11.3	106	0.00
76 T	1,2,3-Trichloropropane	0.496	0.527	-6.3	106	0.00
77 T	Bromobenzene	0.845	0.895	-5.9	102	0.00
78 T	n-propylbenzene	4.335	4.643	-7.1	101	0.00
79 T	2-Chlorotoluene	2.445	2.610	-6.7	103	0.00
80 T	1,3,5-Trimethylbenzene	2.973	3.201	-7.7	101	0.00
81 T	trans-1,4-Dichloro-2-butene	0.224	0.227	-1.3	95	0.00
82 T	4-Chlorotoluene	2.624	2.780	-5.9	102	0.00
83 T	tert-Butylbenzene	2.496	2.679	-7.3	99	0.00
84 T	1,2,4-Trimethylbenzene	3.059	3.304	-8.0	101	0.00
85 T	sec-Butylbenzene	3.704	3.946	-6.5	100	0.00
86 T	p-Isopropyltoluene	3.249	3.467	-6.7	99	0.00
87 T	1,3-Dichlorobenzene	1.689	1.776	-5.2	101	0.00
88 T	1,4-Dichlorobenzene	1.693	1.751	-3.4	102	0.00
89 T	n-Butylbenzene	3.194	3.363	-5.3	98	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
90 T	Hexachloroethane	0.599	0.628	-4.8	100	0.00
91 T	1,2-Dichlorobenzene	1.504	1.607	-6.8	103	0.00
92 T	1,2-Dibromo-3-Chloropropane	0.122	0.135	-10.7	107	0.00
93 T	1,2,4-Trichlorobenzene	1.057	1.109	-4.9	99	0.00
94 T	Hexachlorobutadiene	0.607	0.631	-4.0	101	0.00
95 T	Naphthalene	1.950	2.226	-14.2	104	0.00
96 T	1,2,3-Trichlorobenzene	0.935	1.023	-9.4	103	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 6