

Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU042319\
 Data File : VU031403.D
 Acq On : 22 Apr 2019 19:47
 Operator : JC/SP
 Sample : VSTDICV050
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 VICV53

Quant Time: Apr 23 04:06:19 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM042319WMA.M
 Quant Title : VOC Analysis
 QLast Update : Tue Apr 23 04:02:06 2019
 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	99	0.00
2 T	Dichlorodifluoromethane	0.354	0.361	-2.0	97	0.00
3 T	Chloromethane	0.452	0.445	1.5	97	0.00
4 S	Vinyl Chloride-d3	0.405	0.392	3.2	94	0.00
5 T	Vinyl chloride	0.445	0.449	-0.9	96	0.00
6 T	Bromomethane	0.260	0.271	-4.2	100	0.00
7 S	Chloroethane-d5	0.308	0.304	1.3	96	0.00
8 T	Chloroethane	0.254	0.258	-1.6	97	0.00
9 T	Trichlorofluoromethane	0.534	0.541	-1.3	95	0.00
10 T	1,1,2-Trichloro-1,2,2-trifl	0.310	0.319	-2.9	96	0.00
11 S	1,1-Dichloroethene-d2	0.705	0.698	1.0	96	0.00
12 T	1,1-Dichloroethene	0.305	0.317	-3.9	97	0.00
13 T	Acetone	0.238	0.222	6.7	90	0.00
14 T	Carbon disulfide	0.914	0.924	-1.1	97	0.00
15 T	Methyl Acetate	0.434	0.446	-2.8	100	0.00
16 T	Methylene chloride	0.347	0.352	-1.4	96	0.00
17 T	trans-1,2-Dichloroethene	0.314	0.328	-4.5	98	0.00
18 T	Methyl tert-butyl Ether	1.033	1.061	-2.7	97	0.00
19 T	1,1-Dichloroethane	0.614	0.629	-2.4	97	0.00
20 T	cis-1,2-Dichloroethene	0.356	0.365	-2.5	97	0.00
21 S	2-Butanone-d5	0.303	0.314	-3.6	99	0.00
22 T	2-Butanone	0.308	0.325	-5.5	96	0.00
23 T	Bromochloromethane	0.179	0.183	-2.2	99	0.00
24 S	Chloroform-d	0.646	0.640	0.9	97	0.00
25 T	Chloroform	0.592	0.608	-2.7	96	0.00
26 S	1,2-Dichloroethane-d4	0.387	0.379	2.1	94	0.00
27 T	1,2-Dichloroethane	0.441	0.445	-0.9	96	0.00
28 I	Chlorobenzene-d5	1.000	1.000	0.0	99	0.00
29 T	Cyclohexane	0.607	0.638	-5.1	96	0.00
30 T	1,1,1-Trichloroethane	0.533	0.552	-3.6	97	0.00
31 T	Carbon tetrachloride	0.456	0.478	-4.8	97	0.00
32 S	Benzene-d6	1.444	1.451	-0.5	97	0.00
33 T	Benzene	1.447	1.474	-1.9	96	0.00
34 T	Trichloroethene	0.383	0.376	1.8	96	0.00
35 T	Methylcyclohexane	0.561	0.576	-2.7	98	0.00
36 S	1,2-Dichloropropane-d6	0.465	0.465	0.0	96	0.00
37 T	1,2-Dichloropropane	0.385	0.404	-4.9	99	0.00
38 T	Bromodichloromethane	0.473	0.486	-2.7	96	0.00
39 T	cis-1,3-Dichloropropene	0.577	0.616	-6.8	97	0.00
40 T	4-Methyl-2-pentanone	0.586	0.620	-5.8	99	0.00
41 S	Toluene-d8	1.290	1.296	-0.5	97	0.00
42 T	Toluene	1.502	1.565	-4.2	96	0.00
43 S	trans-1,3-Dichloropropene-d	0.212	0.217	-2.4	97	0.00
44 T	trans-1,3-Dichloropropene	0.497	0.520	-4.6	97	0.00
45 T	1,1,2-Trichloroethane	0.354	0.372	-5.1	100	0.00

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	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
46 T	Tetrachloroethene	0.272	0.284	-4.4	97	0.00
47 S	2-Hexanone-d5	0.233	0.239	-2.6	95	0.00
48 T	2-Hexanone	0.457	0.491	-7.4	98	0.00
49 T	Dibromochloromethane	0.376	0.395	-5.1	99	0.00
50 T	1,2-Dibromoethane	0.372	0.399	-7.3	99	0.00
51 T	Chlorobenzene	0.932	0.962	-3.2	97	0.00
52 T	Ethylbenzene	1.609	1.684	-4.7	98	0.00
53 T	m,p-Xylene	0.587	0.614	-4.6	96	0.00
54 T	o-xylene	0.590	0.630	-6.8	98	0.00
55 T	Styrene	0.982	1.061	-8.0	99	0.00
56 T	Isopropylbenzene	1.516	1.635	-7.8	98	0.00
57 S	1,1,2,2-Tetrachloroethane-d	0.631	0.641	-1.6	97	0.00
58 T	1,1,2,2-Tetrachloroethane	0.611	0.625	-2.3	97	0.00
59	1,2,3-Trichloropropane	0.476	0.496	-4.2	97	0.00
60 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	98	0.00
61 T	Bromoform	0.615	0.623	-1.3	97	0.00
62 T	1,3-Dichlorobenzene	1.464	1.507	-2.9	97	0.00
63 T	1,4-Dichlorobenzene	1.461	1.510	-3.4	97	0.00
64 S	1,2-Dichlorobenzene-d4	1.022	1.026	-0.4	98	0.00
65 T	1,2-Dichlorobenzene	1.497	1.570	-4.9	99	0.00
66 T	1,2-Dibromo-3-chloropropane	0.277	0.301	-8.7	99	0.00
67	1,3,5-Trichlorobenzene	1.079	1.152	-6.8	98	0.00
68 T	1,2,4-trichlorobenzene	0.867	0.986	-13.7	101	0.00
69	Naphthalene	3.016	3.397	-12.6	100	0.00
70 T	1,2,3-Trichlorobenzene	0.934	1.049	-12.3	100	0.00

(#) = Out of Range

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