

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010821\
 Data File : VN065447.D
 Acq On : 8 Jan 2021 16:39
 Operator : JC/MD
 Sample : M1022-15DL 10X
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleID :
 41095DL

Quant Time: Jan 08 22:49:08 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122220W.M
 Quant Title : SW846 8260
 QLast Update : Tue Dec 22 15:42:18 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	7.627	168	289275	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.550	114	473194	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.380	117	507974	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.318	152	234500	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.991	65	162408	49.73	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	99.46%		
35) Dibromofluoromethane	7.553	113	143486	50.26	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	100.52%		
50) Toluene-d8	10.058	98	591975	55.33	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	110.66%		
62) 4-Bromofluorobenzene	12.376	95	246973	57.51	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	115.02%		
Target Compounds						Qvalue
11) Tert butyl alcohol	4.788	59	44736	83.72	ug/l	99
16) Acetone	3.792	43	60354	34.78	ug/l	99
18) Methyl Acetate	4.293	43	7302	2.49	ug/l #	72
25) 2-Butanone	6.804	43	192546	105.02	ug/l	99
37) Ethyl Acetate	6.897	43	23227m	5.45	ug/l	
40) Benzene	8.004	78	14201	1.11	ug/l #	84
49) 1,4-Dioxane	9.177	88	684	11.83	ug/l #	26
51) 4-Methyl-2-Pentanone	9.952	43	7290	1.76	ug/l #	75
52) Toluene	10.126	92	8865	1.10	ug/l	99
68) m/p-Xylenes	11.589	106	5282	0.75	ug/l	98
69) o-Xylene	11.920	106	3872	0.57	ug/l	100
70) Styrene	11.939	104	8743	0.77	ug/l	97
95) Naphthalene	15.103	128	160905	15.72	ug/l	100

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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