

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121721\
 Data File : VN070164.D
 Acq On : 17 Dec 2021 15:19
 Operator : JC/MD
 Sample : M5131-01DL 20X
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 211215028-VOA-02DL

Quant Time: Dec 18 05:45:37 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 03 08:46:47 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	1086815	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	1821627	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	1616350	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	567258	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	739899	44.093	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	88.180%		
35) Dibromofluoromethane	8.021	113	485770	42.573	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	85.140%		
50) Toluene-d8	10.443	98	1874336	40.910	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	81.820%		
62) 4-Bromofluorobenzene	12.733	95	609818	36.740	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	73.480%		
Target Compounds					Qvalue	
11) Tert butyl alcohol	5.387	59	446379	140.729	ug/l	98
16) Acetone	4.291	43	407647	40.384	ug/l	95
17) Carbon Disulfide	4.524	76	15162	0.567	ug/l	98
25) 2-Butanone	7.340	43	720631	52.543	ug/l	91
29) Tetrahydrofuran	7.697	42	125077	15.019	ug/l	87
51) 4-Methyl-2-Pentanone	10.333	43	32158	1.298	ug/l	93
52) Toluene	10.507	92	9625	0.297	ug/l	99
67) Ethyl Benzene	11.846	91	15197	0.255	ug/l	93
68) m/p-Xylenes	11.956	106	8493	0.392	ug/l	100
95) Naphthalene	15.509	128	83588	6.760	ug/l	99

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

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