

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120321\
 Data File : VN069746.D
 Acq On : 4 Dec 2021 3:57
 Operator : JC/MD
 Sample : M4919-06
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FDR-MW-4-20211202-A

Quant Time: Dec 04 05:41:55 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	1451264	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2554442	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	2637022	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	1043167	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	1106132	49.364	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	98.720%		
35) Dibromofluoromethane	8.021	113	774847	48.426	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	96.860%		
50) Toluene-d8	10.441	98	3270628	50.907	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	101.820%		
62) 4-Bromofluorobenzene	12.731	95	1238833	53.225	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	106.460%		
Target Compounds					Qvalue	
16) Acetone	4.299	43	25467	1.889	ug/l	99
17) Carbon Disulfide	4.537	76	26105	0.731	ug/l	98
95) Naphthalene	15.509	128	21153	4.367	ug/l	97

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

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