

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122821\
 Data File : VN070331.D
 Acq On : 28 Dec 2021 17:54
 Operator : JC/MD
 Sample : M5197-01
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 AUD-1230

Quant Time: Dec 28 18:22:34 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
 Quant Title : SW846 8260
 QLast Update : Mon Dec 27 18:47:12 2021
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/29/2021
 Supervised By : Mahesh Dadoda 12/30/2021

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	1185790	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2073025	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	1717564	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	586042	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	988646	57.987	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 115.980%			
35) Dibromofluoromethane	8.024	113	675579	53.862	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 107.720%			
50) Toluene-d8	10.440	98	2611239	50.962	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 101.920%			
62) 4-Bromofluorobenzene	12.731	95	788560	42.610	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 85.220%			
Target Compounds						
16) Acetone	4.272	43	35223405m	3182.537	ug/l	Qvalue
25) 2-Butanone	7.353	43	589877	42.329	ug/l	92
51) 4-Methyl-2-Pentanone	10.333	43	46670	1.752	ug/l	95
68) m/p-Xylenes	11.953	106	47028	1.935	ug/l	96

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

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