

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
 Data File : VN070446.D
 Acq On : 3 Jan 2022 18:58
 Operator : JC/MD
 Sample : M5338-08
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TW-1

Quant Time: Jan 04 01:37:01 2022
 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

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	1001647	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	1692513	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	1461367	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	494040	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	731346	50.782	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	101.560%		
35) Dibromofluoromethane	8.021	113	507068	49.516	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	99.040%		
50) Toluene-d8	10.440	98	2013854	48.140	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	96.280%		
62) 4-Bromofluorobenzene	12.731	95	706459	46.756	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	93.520%		
Target Compounds						Qvalue
16) Acetone	4.291	43	76614	8.195	ug/l	100
24) 1,1-Dichloroethane	6.385	63	28517	1.221	ug/l	97
27) cis-1,2-Dichloroethene	7.324	96	61166	4.746	ug/l	95
64) Tetrachloroethene	10.977	164	4830	0.456	ug/l	88
67) Ethyl Benzene	11.843	91	23985	0.422	ug/l	99
68) m/p-Xylenes	11.950	106	31499	1.523	ug/l	98
69) o-Xylene	12.280	106	6923	0.337	ug/l	89
84) 1,2,4-Trimethylbenzene	13.366	105	21386	0.604	ug/l	98

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

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