

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081022\
 Data File : VN073800.D
 Acq On : 10 Aug 2022 16:10
 Operator : JC\MD
 Sample : N4111-03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BP-MW158S1-GW-225-227

Quant Time: Aug 10 22:54:50 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 10 05:09:32 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.022	168	380845	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	597099	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	565431	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.616	152	262147	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.375	65	237685	51.212	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 102.420%	
35) Dibromofluoromethane	7.957	113	221208	58.998	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 118.000%	
50) Toluene-d8	10.380	98	708693	48.033	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 96.060%	
62) 4-Bromofluorobenzene	12.674	95	257124	49.662	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 99.320%	
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	4.116	96	2408	0.666	ug/l #	69
16) Acetone	4.234	43	12062	4.839	ug/l #	84
24) 1,1-Dichloroethane	6.310	63	21768	2.839	ug/l	95
44) Trichloroethene	9.151	130	2470	0.561	ug/l	90
64) Tetrachloroethene	10.910	164	1448	0.322	ug/l #	75

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

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