

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072922\
 Data File : VN073646.D
 Acq On : 29 Jul 2022 19:51
 Operator : JC\MD
 Sample : N3858-06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SB-B

Quant Time: Aug 01 04:09:21 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 29 02:20:25 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	340166	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	567387	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	553162	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.616	152	239800	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	204748	49.391	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	98.780%	
35) Dibromofluoromethane	7.951	113	188618	52.940	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	105.880%	
50) Toluene-d8	10.381	98	659459	47.036	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	94.080%	
62) 4-Bromofluorobenzene	12.674	95	232709	49.732	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	99.460%	
Target Compounds						
						Qvalue
16) Acetone	4.228	43	116792	52.454	ug/l	91
18) Methyl Acetate	4.787	43	8166	1.474	ug/l	93
20) Methylene Chloride	5.010	84	23376	4.483	ug/l	96
25) 2-Butanone	7.269	43	35002	10.123	ug/l	94
29) Tetrahydrofuran	7.622	42	12019	5.498	ug/l	97
37) Ethyl Acetate	7.340	43	55219	8.257	ug/l	99
43) Isopropyl Acetate	8.498	43	212983	23.579	ug/l #	92

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

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