

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072922\
 Data File : VN073647.D
 Acq On : 29 Jul 2022 20:15
 Operator : JC\MD
 Sample : N3858-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SB-A

Quant Time: Aug 01 04:09:43 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	348954	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	572262	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	568767	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.615	152	258813	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	205241	48.263	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	96.520%
35) Dibromofluoromethane	7.951	113	190816	53.101	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	106.200%
50) Toluene-d8	10.380	98	668476	47.273	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	94.540%
62) 4-Bromofluorobenzene	12.674	95	248981	52.247	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	104.500%
Target Compounds						
					Qvalue	
16) Acetone	4.228	43	129410	56.657	ug/l	95
18) Methyl Acetate	4.793	43	7631	1.343	ug/l #	71
20) Methylene Chloride	5.022	84	22954	4.193	ug/l #	93
25) 2-Butanone	7.263	43	35940	10.132	ug/l	95
29) Tetrahydrofuran	7.622	42	11341	5.057	ug/l	98
37) Ethyl Acetate	7.345	43	53979	8.003	ug/l	99
43) Isopropyl Acetate	8.498	43	212753	23.353	ug/l #	92

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

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