

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080322\
 Data File : VX030366.D
 Acq On : 03 Aug 2022 13:26
 Operator : JC/MD
 Sample : N3976-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TP1-0729

Quant Time: Aug 03 15:54:03 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 25 16:49:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	51499	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	115577	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	104855	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	54199	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	113715	109.859	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 219.720%#	
35) Dibromofluoromethane	5.397	113	77443	70.093	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 140.180%#	
50) Toluene-d8	8.653	98	194952	45.563	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 91.120%	
62) 4-Bromofluorobenzene	11.085	95	65836	41.460	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 82.920%#	
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	78395	147.931	ug/l	91
18) Methyl Acetate	2.709	43	4085	2.606	ug/l	98
20) Methylene Chloride	2.788	84	18881	19.225	ug/l	93
25) 2-Butanone	4.574	43	9704	10.897	ug/l	91
37) Ethyl Acetate	4.727	43	33671	14.954	ug/l	96
43) Isopropyl Acetate	6.348	43	131954	39.977	ug/l	95

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

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