

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072922\
 Data File : VN073648.D
 Acq On : 29 Jul 2022 20:40
 Operator : JC\MD
 Sample : N3920-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB-07-27-2022

Quant Time: Aug 01 04:10:08 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	361854	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	589437	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	592736	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.616	152	268005	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.375	65	214535	48.650	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	97.300%
35) Dibromofluoromethane	7.951	113	197364	53.323	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	106.640%
50) Toluene-d8	10.381	98	692391	47.538	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	95.080%
62) 4-Bromofluorobenzene	12.675	95	254721	51.952	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	103.900%
Target Compounds						
					Qvalue	
16) Acetone	4.228	43	116715	49.277	ug/l	98
18) Methyl Acetate	4.793	43	6906	1.172	ug/l	94
20) Methylene Chloride	5.016	84	21829	3.654	ug/l	97
25) 2-Butanone	7.269	43	34084	9.266	ug/l	95
29) Tetrahydrofuran	7.616	42	12002	5.161	ug/l	97
37) Ethyl Acetate	7.340	43	59245	8.528	ug/l	97
43) Isopropyl Acetate	8.498	43	228104	24.309	ug/l #	92

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

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