

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072723\
 Data File : VN078671.D
 Acq On : 27 Jul 2023 19:47
 Operator : JC\MD
 Sample : 03770-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 DEL-2-RP

Quant Time: Jul 28 01:12:05 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071023W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 10 16:13:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.235	168	418386	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	781075	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	745611	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	279696	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	273772	48.810	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	97.620%
35) Dibromofluoromethane	8.177	113	269569	50.841	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.680%
50) Toluene-d8	10.576	98	957441	47.800	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	95.600%
62) 4-Bromofluorobenzene	12.859	95	320451	51.940	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	103.880%
Target Compounds						
					Qvalue	
16) Acetone	4.447	43	160739	91.999	ug/l	92
20) Methylene Chloride	5.277	84	22496	3.348	ug/l	97
25) 2-Butanone	7.500	43	81941	29.794	ug/l #	84
37) Ethyl Acetate	7.582	43	17197	2.695	ug/l #	96
43) Isopropyl Acetate	8.700	43	441597	42.028	ug/l #	72
51) 4-Methyl-2-Pentanone	10.453	43	249171	39.868	ug/l	92

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

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