

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072522\
 Data File : VN073579.D
 Acq On : 25 Jul 2022 15:08
 Operator : JC\MD
 Sample : N3771-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-10F

Quant Time: Jul 26 02:17:17 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062822W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 29 01:54:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	250972	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	410364	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	381008	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.615	152	183067	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.375	65	158848	43.844	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	87.680%
35) Dibromofluoromethane	7.957	113	136924	50.151	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	100.300%
50) Toluene-d8	10.380	98	438615	45.233	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	90.460%
62) 4-Bromofluorobenzene	12.674	95	180029	48.311	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	96.620%
Target Compounds						
					Qvalue	
16) Acetone	4.234	43	5861	2.882	ug/l	97
30) Chloroform	7.751	83	37640	6.218	ug/l	96
44) Trichloroethene	9.145	130	2638	0.800	ug/l	80
47) Bromodichloromethane	9.704	83	5691	1.310	ug/l	92
60) Dibromochloromethane	11.174	129	893	0.284	ug/l	98
64) Tetrachloroethene	10.910	164	773	0.248	ug/l #	48

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

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