

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082823\
 Data File : VN079073.D
 Acq On : 28 Aug 2023 17:44
 Operator : JC\MD
 Sample : 04188-11
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-4

Quant Time: Aug 29 04:40:47 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081523W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 16 02:42:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	209148	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	401073	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	391051	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	123022	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.589	65	161954	49.763	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	99.520%
35) Dibromofluoromethane	8.171	113	137718	51.010	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	102.020%
50) Toluene-d8	10.571	98	493519	49.231	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	98.460%
62) 4-Bromofluorobenzene	12.853	95	158036	48.231	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	96.460%
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	31906	23.760	ug/l	97
20) Methylene Chloride	5.283	84	11026	3.853	ug/l #	80
25) 2-Butanone	7.495	43	11098	5.155	ug/l #	62
43) Isopropyl Acetate	8.694	43	150918	18.791	ug/l #	74
51) 4-Methyl-2-Pentanone	10.453	43	32449	6.812	ug/l #	81

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

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