

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062823\
 Data File : VN078437.D
 Acq On : 28 Jun 2023 13:26
 Operator : JC\MD
 Sample : 03428-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PNNL-APM-0621-1

Quant Time: Jun 29 06:09:05 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060623W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 07 06:01:44 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	456561	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	831750	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	791397	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	237441	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	308076	50.609	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 101.220%		
35) Dibromofluoromethane	8.177	113	287725	54.710	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 109.420%		
50) Toluene-d8	10.571	98	986859	48.676	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 97.360%		
62) 4-Bromofluorobenzene	12.853	95	321159	50.955	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 101.920%		
Target Compounds						
						Qvalue
16) Acetone	4.442	43	102909	58.732	ug/l	97
20) Methylene Chloride	5.289	84	29529	5.100	ug/l	85
25) 2-Butanone	7.500	43	23410	8.688	ug/l	97
37) Ethyl Acetate	7.571	43	56347	9.562	ug/l	95
43) Isopropyl Acetate	8.700	43	264530	23.472	ug/l #	85
51) 4-Methyl-2-Pentanone	10.453	43	166654	28.081	ug/l	89

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

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