

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082123\
 Data File : VN078974.D
 Acq On : 21 Aug 2023 14:56
 Operator : JC\MD
 Sample : 04082-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 364

Quant Time: Aug 22 00:57:36 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	188304	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	352957	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	338044	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	104797	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	147961	50.496	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	101.000%
35) Dibromofluoromethane	8.171	113	122295	51.473	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	102.940%
50) Toluene-d8	10.571	98	440969	49.985	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	99.980%
62) 4-Bromofluorobenzene	12.853	95	130041	45.097	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	90.200%
Target Compounds						
						Qvalue
16) Acetone	4.448	43	29819	24.664	ug/l	92
20) Methylene Chloride	5.289	84	9462	3.672	ug/l	90
25) 2-Butanone	7.500	43	9938	5.127	ug/l	88
43) Isopropyl Acetate	8.700	43	169947	24.045	ug/l #	74
51) 4-Methyl-2-Pentanone	10.453	43	40182	9.586	ug/l #	80

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

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