

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN020524\
 Data File : VN080906.D
 Acq On : 05 Feb 2024 19:02
 Operator : JC\MD
 Sample : P1366-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 EB-2-6-8

Quant Time: Feb 06 04:47:18 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N012924W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jan 29 12:54:23 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	79127	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	191463	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	214097	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	83046	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	64998	47.483	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	94.960%
35) Dibromofluoromethane	8.165	113	53172	45.393	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	90.780%
50) Toluene-d8	10.571	98	217398	47.418	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	94.840%
62) 4-Bromofluorobenzene	12.853	95	85794	53.700	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	107.400%
Target Compounds						
					Qvalue	
16) Acetone	4.430	43	50760	102.245	ug/l	95
18) Methyl Acetate	5.030	43	4213	2.435	ug/l #	69
20) Methylene Chloride	5.265	84	7865	6.035	ug/l	90
25) 2-Butanone	7.494	43	8032	8.518	ug/l	96
37) Ethyl Acetate	7.565	43	4292	1.915	ug/l #	90
43) Isopropyl Acetate	8.694	43	121751	29.837	ug/l #	89
51) 4-Methyl-2-Pentanone	10.447	43	12580	5.515	ug/l #	73

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

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