

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN020524\
 Data File : VN080908.D
 Acq On : 05 Feb 2024 19:50
 Operator : JC\MD
 Sample : P1366-06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 EB-1-4-6

Quant Time: Feb 06 04:48:05 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	77404	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	188057	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	204636	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	81796	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	65509	48.921	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	97.840%
35) Dibromofluoromethane	8.171	113	49342	42.886	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	85.780%
50) Toluene-d8	10.565	98	216374	48.050	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	96.100%
62) 4-Bromofluorobenzene	12.853	95	84370	53.765	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	107.540%
Target Compounds						
					Qvalue	
16) Acetone	4.430	43	13141	27.059	ug/l #	86
18) Methyl Acetate	5.036	43	4385	2.591	ug/l #	66
20) Methylene Chloride	5.265	84	6264	4.914	ug/l #	90
25) 2-Butanone	7.489	43	6924	7.506	ug/l	95
37) Ethyl Acetate	7.565	43	4434	2.014	ug/l #	93
43) Isopropyl Acetate	8.694	43	126048	31.449	ug/l #	87

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

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