

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN110923\
 Data File : VN080040.D
 Acq On : 09 Nov 2023 22:33
 Operator : JC\MD
 Sample : 05318-01 50X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 KEARNY-DRUM

Quant Time: Nov 10 00:44:03 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110123W.M
 Quant Title : SW846 8260
 QLast Update : Thu Nov 02 05:49:08 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	279045	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	601593	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	623247	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	269292	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	255976	60.211	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	120.420%
35) Dibromofluoromethane	8.171	113	212443	49.895	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.780%
50) Toluene-d8	10.571	98	744215	47.479	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	94.960%
62) 4-Bromofluorobenzene	12.853	95	269750	51.890	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	103.780%
Target Compounds						
					Qvalue	
16) Acetone	4.447	43	8250	5.525	ug/l #	71
17) Carbon Disulfide	4.712	76	21380	2.468	ug/l	99
18) Methyl Acetate	5.030	43	7780	1.089	ug/l #	80
40) Benzene	8.606	78	27497	1.510	ug/l	98
52) Toluene	10.635	92	13798	1.278	ug/l	100

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

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