

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102124\
 Data File : VN084442.D
 Acq On : 21 Oct 2024 17:40
 Operator : JC\MD
 Sample : P4438-07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BP-VPB-190-GW-198-200

Quant Time: Oct 22 01:38:44 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N093024W.M
 Quant Title : SW846 8260
 QLast Update : Tue Oct 01 07:11:01 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	167949	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	288550	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	244682	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	95771	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	118526	47.598	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	95.200%
35) Dibromofluoromethane	8.165	113	97274	51.135	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	102.260%
50) Toluene-d8	10.565	98	345899	49.423	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	98.840%
62) 4-Bromofluorobenzene	12.847	95	114060	44.734	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	89.460%
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	2703	2.527	ug/l	100
30) Chloroform	7.959	83	1413	0.358	ug/l #	73
64) Tetrachloroethene	11.106	164	872	0.512	ug/l #	73

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

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