

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060324\
 Data File : VN082444.D
 Acq On : 03 Jun 2024 19:16
 Operator : JC\MD
 Sample : P2668-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-TP-2

Quant Time: Jun 04 00:34:38 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N051524W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 16 05:21:06 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	156086	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	286821	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	287833	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	117945	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	133512	54.171	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 108.340%	
35) Dibromofluoromethane	8.171	113	96894	54.916	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 109.840%	
50) Toluene-d8	10.571	98	380133	56.195	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 112.380%	
62) 4-Bromofluorobenzene	12.853	95	126216	52.276	ug/l	0.00
Spiked Amount	50.000	Range	64 - 133	Recovery	= 104.560%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	82369	81.470	ug/l	91
20) Methylene Chloride	5.277	84	14027	6.876	ug/l #	80
25) 2-Butanone	7.488	43	13166	8.589	ug/l #	78
43) Isopropyl Acetate	8.694	43	218387	37.216	ug/l #	76
51) 4-Methyl-2-Pentanone	10.453	43	18562	5.567	ug/l #	69

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

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