

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052924\
 Data File : VN082380.D
 Acq On : 29 May 2024 19:06
 Operator : JC\MD
 Sample : P2612-09
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RBR200057

Quant Time: May 30 05:48:04 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	8.224	168	178965	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	324567	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	315985	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	136810	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	145416	51.458	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 102.920%	
35) Dibromofluoromethane	8.171	113	105239	52.709	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 105.420%	
50) Toluene-d8	10.571	98	421946	55.122	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 110.240%	
62) 4-Bromofluorobenzene	12.853	95	150503	55.086	ug/l	0.00
Spiked Amount	50.000	Range	64 - 133	Recovery	= 110.180%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	82682	71.325	ug/l	90
20) Methylene Chloride	5.283	84	12250	5.238	ug/l #	74
25) 2-Butanone	7.489	43	13015	7.405	ug/l #	85
40) Benzene	8.606	78	24230	2.427	ug/l	98
43) Isopropyl Acetate	8.688	43	265696	40.013	ug/l #	74
51) 4-Methyl-2-Pentanone	10.447	43	20874	5.533	ug/l #	69
52) Toluene	10.630	92	10497	1.752	ug/l	92

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

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