

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071124\
 Data File : VN082881.D
 Acq On : 11 Jul 2024 17:09
 Operator : JC\MD
 Sample : P3206-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 626-A

Quant Time: Jul 12 01:09:34 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N070824W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 08 13:46:04 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	117304	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	237337	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	231650	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	104192	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	86588	52.657	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 105.320%		
35) Dibromofluoromethane	8.171	113	70155	46.066	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 92.140%		
50) Toluene-d8	10.571	98	268627	47.757	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 95.520%		
62) 4-Bromofluorobenzene	12.853	95	103187	51.070	ug/l	0.00
Spiked Amount	50.000	Range 64 - 133	Recovery	= 102.140%		
Target Compounds						
						Qvalue
11) Tert butyl alcohol	5.530	59	8333	30.130	ug/l #	90
13) Acrolein	4.195	56	7390	37.010	ug/l	96
16) Acetone	4.430	43	158086	273.062	ug/l #	79
25) 2-Butanone	7.489	43	32720	37.430	ug/l	91
40) Benzene	8.612	78	4949	0.766	ug/l	96
51) 4-Methyl-2-Pentanone	10.447	43	6476	3.056	ug/l #	78
52) Toluene	10.630	92	2423	0.601	ug/l	89

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

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