

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN110723\
 Data File : VN079969.D
 Acq On : 07 Nov 2023 16:58
 Operator : JC\MD
 Sample : 05107-25DL 10X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 1049DL

Quant Time: Nov 08 04:31:02 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	8.229	168	282441	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	582166	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	604686	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	267681	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	241526	56.129	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 112.260%			
35) Dibromofluoromethane	8.171	113	201319	48.860	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 97.720%			
50) Toluene-d8	10.570	98	740173	48.797	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 97.600%			
62) 4-Bromofluorobenzene	12.853	95	279099	55.480	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 110.960%			
Target Compounds					Qvalue	
16) Acetone	4.441	43	22431	14.840	ug/l	96
20) Methylene Chloride	5.288	84	4445	1.165	ug/l #	88
40) Benzene	8.612	78	1105194	62.721	ug/l	100
43) Isopropyl Acetate	8.700	43	30180	3.190	ug/l #	76
51) 4-Methyl-2-Pentanone	10.453	43	96653	18.144	ug/l	87
52) Toluene	10.635	92	459781	44.013	ug/l	99
67) Ethyl Benzene	11.970	91	48222	2.110	ug/l	99
68) m/p-Xylenes	12.070	106	90456	10.626	ug/l	96
69) o-Xylene	12.400	106	65808	8.028	ug/l	90
80) 1,3,5-Trimethylbenzene	13.176	105	18167	1.149	ug/l	99
84) 1,2,4-Trimethylbenzene	13.488	105	54332	3.494	ug/l	87
95) Naphthalene	15.647	128	1112580	78.994	ug/l	99

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

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