

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010623\
 Data File : VN076141.D
 Acq On : 06 Jan 2023 18:12
 Operator : JC\MD
 Sample : 01061-11 50X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TOTE-3

Quant Time: Jan 06 22:57:40 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
 Quant Title : SW846 8260
 QLast Update : Wed Dec 14 13:03:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	321723	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	529971	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	460006	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	210850	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	175360	50.102	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 100.200%	
35) Dibromofluoromethane	8.177	113	158435	48.273	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 96.540%	
50) Toluene-d8	10.571	98	608025	49.549	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 99.100%	
62) 4-Bromofluorobenzene	12.853	95	196246	48.285	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 96.560%	
Target Compounds						
					Qvalue	
16) Acetone	4.465	43	6259	5.188	ug/l	91
30) Chloroform	7.977	83	2967	0.488	ug/l	95
52) Toluene	10.635	92	3921	0.448	ug/l	93
67) Ethyl Benzene	11.965	91	9525	0.598	ug/l	96
68) m/p-Xylenes	12.071	106	12855	2.038	ug/l	99
69) o-Xylene	12.406	106	7845	1.257	ug/l	96
78) n-propylbenzene	13.041	91	6294	0.375	ug/l	97
80) 1,3,5-Trimethylbenzene	13.170	105	10103	0.792	ug/l	98
84) 1,2,4-Trimethylbenzene	13.488	105	39997	3.162	ug/l	100

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

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