

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021224\
 Data File : VN081012.D
 Acq On : 12 Feb 2024 18:44
 Operator : JC\MD
 Sample : P1453-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-3

Quant Time: Feb 13 03:25:35 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N020624W.M
 Quant Title : SW846 8260
 QLast Update : Wed Feb 07 06:22:23 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	591499	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	9.106	114	1105746	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	975583	50.000	ug/l	# 0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	379019	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	305259	40.910	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	81.820%
35) Dibromofluoromethane	8.171	113	281816	47.496	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	95.000%
50) Toluene-d8	10.571	98	1030328	48.008	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	96.020%
62) 4-Bromofluorobenzene	12.853	95	342078	43.876	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	87.760%
Target Compounds						
					Qvalue	
3) Chloromethane	2.360	50	17991	1.246	ug/l	98
13) Acrolein	4.177	56	9456	7.298	ug/l	94
16) Acetone	4.430	43	106438	43.357	ug/l	99
20) Methylene Chloride	5.277	84	26305	3.537	ug/l #	72
37) Ethyl Acetate	7.559	43	12295	1.841	ug/l #	82
43) Isopropyl Acetate	8.689	43	694626	56.557	ug/l #	79

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

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