

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022324\
 Data File : VN081172.D
 Acq On : 23 Feb 2024 20:33
 Operator : JC\MD
 Sample : P1574-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SB-2

Quant Time: Feb 23 22:05:21 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021624W.M
 Quant Title : SW846 8260
 QLast Update : Fri Feb 16 23:45:20 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	281692	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	541928	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	492503	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	189865	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	181160	52.468	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 104.940%		
35) Dibromofluoromethane	8.165	113	157459	51.383	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 102.760%		
50) Toluene-d8	10.571	98	505395	46.434	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 92.860%		
62) 4-Bromofluorobenzene	12.853	95	185054	45.249	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 90.500%		
Target Compounds						
						Qvalue
16) Acetone	4.436	43	66330	48.823	ug/l	99
18) Methyl Acetate	5.030	43	6042	1.823	ug/l #	84
20) Methylene Chloride	5.277	84	19288	4.645	ug/l #	94
37) Ethyl Acetate	7.559	43	8123	1.714	ug/l #	93
43) Isopropyl Acetate	8.694	43	475397	57.368	ug/l #	79
51) 4-Methyl-2-Pentanone	10.447	43	23918	5.017	ug/l #	76

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

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