

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052224\
 Data File : VN082291.D
 Acq On : 22 May 2024 14:17
 Operator : JC\MD
 Sample : PB161037ZHE#04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB161037ZHE#04

Quant Time: May 23 01:52:56 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N051524W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 16 05:21:06 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	194892	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	345471	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	329651	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	143938	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	159876	51.952	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 103.900%		
35) Dibromofluoromethane	8.171	113	109710	51.624	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 103.240%		
50) Toluene-d8	10.571	98	440049	54.009	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 108.020%		
62) 4-Bromofluorobenzene	12.853	95	156878	53.945	ug/l	0.00
Spiked Amount	50.000	Range 64 - 133	Recovery	= 107.880%		
Target Compounds						
						Qvalue
16) Acetone	4.436	43	22663	17.952	ug/l	99
20) Methylene Chloride	5.277	84	11264	4.422	ug/l #	82
43) Isopropyl Acetate	8.694	43	291991	41.312	ug/l #	72

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052224\
Data File : VN082291.D
Acq On : 22 May 2024 14:17
Operator : JC\MD
Sample : PB161037ZHE#04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB161037ZHE#04

Quant Time: May 23 01:52:56 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N051524W.M
Quant Title : SW846 8260
QLast Update : Thu May 16 05:21:06 2024
Response via : Initial Calibration

