

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030625\
 Data File : VN085910.D
 Acq On : 06 Mar 2025 19:55
 Operator : JC\MD
 Sample : PB166989ZHE#12
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB166989ZHE#12

Quant Time: Mar 07 00:45:08 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021825W.M
 Quant Title : SW846 8260
 QLast Update : Wed Feb 19 03:43:32 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	140752	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	262601	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	249216	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	91807	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	105229	57.666	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 115.340%	
35) Dibromofluoromethane	8.165	113	90798	52.841	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 105.680%	
50) Toluene-d8	10.565	98	332319	53.155	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 106.320%	
62) 4-Bromofluorobenzene	12.847	95	100628	48.852	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	= 97.700%	
Target Compounds						
						Qvalue
16) Acetone	4.424	43	32347	56.022	ug/l	98
20) Methylene Chloride	5.277	84	3845	2.071	ug/l	86
43) Isopropyl Acetate	8.688	43	65841m	18.401	ug/l	

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030625\
Data File : VN085910.D
Acq On : 06 Mar 2025 19:55
Operator : JC\MD
Sample : PB166989ZHE#12
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB166989ZHE#12

Quant Time: Mar 07 00:45:08 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021825W.M
Quant Title : SW846 8260
QLast Update : Wed Feb 19 03:43:32 2025
Response via : Initial Calibration

