

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN013125\
 Data File : VN085599.D
 Acq On : 31 Jan 2025 14:03
 Operator : JC\MD
 Sample : PB166357ZHE#04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB166357ZHE#04

Quant Time: Jan 31 22:28:18 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N011425W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jan 15 02:16:08 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	168539	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	336774	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	321227	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	117372	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	154856	56.920	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 113.840%			
35) Dibromofluoromethane	8.165	113	116930	50.047	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 100.100%			
50) Toluene-d8	10.565	98	434789	52.377	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 104.760%			
62) 4-Bromofluorobenzene	12.847	95	131532	46.321	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 92.640%			
Target Compounds						
16) Acetone	4.430	43	32865	37.387	ug/l	96
20) Methylene Chloride	5.283	84	5285	2.429	ug/l	88
30) Chloroform	7.965	83	6420	1.564	ug/l	86
43) Isopropyl Acetate	8.683	43	177079	33.390	ug/l #	84

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN013125\
Data File : VN085599.D
Acq On : 31 Jan 2025 14:03
Operator : JC\MD
Sample : PB166357ZHE#04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB166357ZHE#04

Quant Time: Jan 31 22:28:18 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N011425W.M
Quant Title : SW846 8260
QLast Update : Wed Jan 15 02:16:08 2025
Response via : Initial Calibration

