

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022125\
 Data File : VN085807.D
 Acq On : 21 Feb 2025 17:03
 Operator : JC\MD
 Sample : PB166809TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB166809TB

Quant Time: Feb 22 02:37:41 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	8.218	168	230909	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	437050	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	428742	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	151786	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	173290	57.886	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	115.780%
35) Dibromofluoromethane	8.165	113	151293	52.903	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	105.800%
50) Toluene-d8	10.565	98	564793	54.281	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	108.560%
62) 4-Bromofluorobenzene	12.847	95	167508	48.862	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	97.720%
Target Compounds						
						Qvalue
16) Acetone	4.424	43	21831	23.047	ug/l	97
20) Methylene Chloride	5.277	84	8202	2.693	ug/l	87
30) Chloroform	7.959	83	8023	1.503	ug/l	96
43) Isopropyl Acetate	8.683	43	164323	27.869	ug/l #	91

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022125\
Data File : VN085807.D
Acq On : 21 Feb 2025 17:03
Operator : JC\MD
Sample : PB166809TB
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB166809TB

Quant Time: Feb 22 02:37:41 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

