

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100423\
 Data File : VN079409.D
 Acq On : 04 Oct 2023 17:36
 Operator : JC\MD
 Sample : PB156087TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB156087TB

Quant Time: Oct 05 02:24:32 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091123W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 12 05:05:24 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	128678	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	266579	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	244332	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	62548	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	118203	68.049	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 136.100%#	
35) Dibromofluoromethane	8.171	113	89725	50.600	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 101.200%	
50) Toluene-d8	10.571	98	302510	46.461	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 92.920%	
62) 4-Bromofluorobenzene	12.853	95	94842	45.665	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 91.340%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	22127	29.994	ug/l	100
18) Methyl Acetate	5.030	43	3679	1.213	ug/l #	59
20) Methylene Chloride	5.283	84	15634	8.708	ug/l #	89
43) Isopropyl Acetate	8.694	43	166062	37.332	ug/l #	75

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100423\
Data File : VN079409.D
Acq On : 04 Oct 2023 17:36
Operator : JC\MD
Sample : PB156087TB
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB156087TB

Quant Time: Oct 05 02:24:32 2023
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091123W.M
Quant Title : SW846 8260
QLast Update : Tue Sep 12 05:05:24 2023
Response via : Initial Calibration

