

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081222\
 Data File : VN073854.D
 Acq On : 12 Aug 2022 16:13
 Operator : JC\MD
 Sample : PB146944ZHE#01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB146944ZHE#01

Quant Time: Aug 13 05:03:29 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 10 05:09:32 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	414369	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	668467	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	671538	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.615	152	310979	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.375	65	239832	47.494	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	94.980%		
35) Dibromofluoromethane	7.951	113	226660	53.998	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	108.000%		
50) Toluene-d8	10.380	98	776861	47.032	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	94.060%		
62) 4-Bromofluorobenzene	12.674	95	294463	50.739	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	101.480%		
Target Compounds						
16) Acetone	4.228	43	71234	26.264	ug/l	96
20) Methylene Chloride	5.010	84	19328	2.303	ug/l	93
37) Ethyl Acetate	7.339	43	61579	7.816	ug/l	98
43) Isopropyl Acetate	8.498	43	250819	23.569	ug/l #	92

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081222\
Data File : VN073854.D
Acq On : 12 Aug 2022 16:13
Operator : JC\MD
Sample : PB146944ZHE#01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PB146944ZHE#01

Quant Time: Aug 13 05:03:29 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
Quant Title : SW846 8260
QLast Update : Wed Aug 10 05:09:32 2022
Response via : Initial Calibration

