

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX011522\
 Data File : VX026468.D
 Acq On : 14 Jan 2022 21:22
 Operator : JC/MD
 Sample : PB142010ZHE#07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB142010ZHE#07

Quant Time: Jan 17 00:42:17 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jan 11 14:45:13 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	128280	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	215480	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	188213	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	79181	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	89725	50.070	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 100.140%	
35) Dibromofluoromethane	5.391	113	68945	49.547	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 99.100%	
50) Toluene-d8	8.653	98	249892	49.835	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 99.680%	
62) 4-Bromofluorobenzene	11.079	95	90383	50.740	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 101.480%	
Target Compounds						Qvalue
16) Acetone	2.386	43	5750	6.292	ug/l	98
20) Methylene Chloride	2.794	84	2974	1.970	ug/l	96
43) Isopropyl Acetate	6.348	43	12284	3.422	ug/l	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX011522\
Data File : VX026468.D
Acq On : 14 Jan 2022 21:22
Operator : JC/MD
Sample : PB142010ZHE#07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PB142010ZHE#07

Quant Time: Jan 17 00:42:17 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.M
Quant Title : SW846 8260
QLast Update : Tue Jan 11 14:45:13 2022
Response via : Initial Calibration

