

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX020122\
 Data File : VX026719.D
 Acq On : 01 Feb 2022 19:57
 Operator : JC/MD
 Sample : PB142403ZHE#12
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB142403ZHE#12

Quant Time: Feb 02 04:38:00 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	135553	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	229349	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	206262	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	89300	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	85485	45.145	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	90.280%
35) Dibromofluoromethane	5.391	113	70661	47.710	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	95.420%
50) Toluene-d8	8.653	98	269169	50.433	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	100.860%
62) 4-Bromofluorobenzene	11.079	95	98009	51.694	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	103.380%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	8263	8.557	ug/l	99
18) Methyl Acetate	2.709	43	2040	1.063	ug/l	91
20) Methylene Chloride	2.794	84	3027	1.897	ug/l	97
43) Isopropyl Acetate	6.342	43	21894	5.731	ug/l	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX020122\
Data File : VX026719.D
Acq On : 01 Feb 2022 19:57
Operator : JC/MD
Sample : PB142403ZHE#12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PB142403ZHE#12

Quant Time: Feb 02 04:38:00 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

