

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101121\
 Data File : VX024747.D
 Acq On : 12 Oct 2021 06:31
 Operator : JC/MD
 Sample : M4158-13
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 T8-3-TOP-GRAB

Quant Time: Oct 12 08:12:03 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092321W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 23 18:24:34 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	126603	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	222264	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	220793	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	90693	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	97556	54.674	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 109.340%	
35) Dibromofluoromethane	5.391	113	75074	50.844	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 101.680%	
50) Toluene-d8	8.653	98	281493	51.797	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 103.600%	
62) 4-Bromofluorobenzene	11.085	95	101489	47.497	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 95.000%	
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	21815	19.632	ug/l #	86
20) Methylene Chloride	2.794	84	6839	4.072	ug/l #	83
43) Isopropyl Acetate	6.348	43	16199	4.005	ug/l	93

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101121\
Data File : VX024747.D
Acq On : 12 Oct 2021 06:31
Operator : JC/MD
Sample : M4158-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T8-3-TOP-GRAB

Quant Time: Oct 12 08:12:03 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092321W.M
Quant Title : SW846 8260
QLast Update : Thu Sep 23 18:24:34 2021
Response via : Initial Calibration

