

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042121\
 Data File : VX021791.D
 Acq On : 21 Apr 2021 14:41
 Operator : JC/MD
 Sample : M2122-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB-RW8-GW-148-150

Quant Time: Apr 22 08:57:44 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042021W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 20 13:46:22 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.629	168	158662	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.830	114	272115	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.098	117	248261	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.061	152	121247	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.031	65	93336	51.86	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	103.72%
35) Dibromofluoromethane	5.464	113	86755	49.85	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.70%
50) Toluene-d8	8.696	98	329055	49.18	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	98.36%
62) 4-Bromofluorobenzene	11.116	95	120901	48.39	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	96.78%
Target Compounds						
16) Acetone	2.422	43	2481	4.01	ug/l	96
17) Carbon Disulfide	2.556	76	1828	0.44	ug/l #	93
19) Methyl tert-butyl Ether	3.166	73	1534	0.30	ug/l #	89

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042121\
Data File : VX021791.D
Acq On : 21 Apr 2021 14:41
Operator : JC/MD
Sample : M2122-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BP-VPB-RW8-GW-148-150

Quant Time: Apr 22 08:57:44 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042021W.M
Quant Title : SW846 8260
QLast Update : Tue Apr 20 13:46:22 2021
Response via : Initial Calibration

