

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080621\
 Data File : VX023589.D
 Acq On : 06 Aug 2021 12:24
 Operator : JC/MD
 Sample : M3298-03 20X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 1917

Quant Time: Aug 07 04:42:49 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 23 06:05:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	160587	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	270911	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	242430	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	93095	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	107539	53.68	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 107.36%	
35) Dibromofluoromethane	5.397	113	88096	52.01	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 104.02%	
50) Toluene-d8	8.653	98	327774	52.67	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 105.34%	
62) 4-Bromofluorobenzene	11.085	95	105171	49.70	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 99.40%	
Target Compounds						
11) Tert butyl alcohol	2.977	59	5041	11.02	ug/l	96
16) Acetone	2.386	43	2217	2.26	ug/l #	73
17) Carbon Disulfide	2.514	76	1142	0.29	ug/l #	84

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080621\
Data File : VX023589.D
Acq On : 06 Aug 2021 12:24
Operator : JC/MD
Sample : M3298-03 20X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
1917

Quant Time: Aug 07 04:42:49 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072221W.M
Quant Title : SW846 8260
QLast Update : Fri Jul 23 06:05:38 2021
Response via : Initial Calibration

