

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070121\
 Data File : VX023087.D
 Acq On : 01 Jul 2021 16:03
 Operator : JC/MD
 Sample : M2892-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-6

Quant Time: Jul 02 04:30:31 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 23 19:12:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.568	168	53330	50.000	ug/l	# 0.01
34) 1,4-Difluorobenzene	6.769	114	112283	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.061	117	100940	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	37163	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	58851	60.807	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	121.620%#
35) Dibromofluoromethane	5.397	113	36311	50.925	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.860%
50) Toluene-d8	8.653	98	138700	49.634	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	99.260%
62) 4-Bromofluorobenzene	11.085	95	51082	48.980	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	97.960%
Target Compounds						
						Qvalue
11) Tert butyl alcohol	2.989	59	1068	5.295	ug/l	# 80
16) Acetone	2.392	43	2100	4.894	ug/l	89
19) Methyl tert-butyl Ether	3.123	73	24273	9.783	ug/l	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070121\
Data File : VX023087.D
Acq On : 01 Jul 2021 16:03
Operator : JC/MD
Sample : M2892-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-6

Quant Time: Jul 02 04:30:31 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260
QLast Update : Wed Jun 23 19:12:26 2021
Response via : Initial Calibration

