

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX070620\
 Data File : VX017169.D
 Acq On : 06 Jul 2020 15:40
 Operator : JC/SP
 Sample : L3159-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 50BR-20200629

Quant Time: Jul 07 07:02:09 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X062920W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jun 29 17:10:38 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	273545	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	441326	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	459237	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	270020	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	167946	49.80	ug/l	0.00
Spiked Amount	50.000		Recovery	=	99.60%	
35) Dibromofluoromethane	5.47	113	143168	51.28	ug/l	0.00
Spiked Amount	50.000		Recovery	=	102.56%	
50) Toluene-d8	8.70	98	533514	51.68	ug/l	0.00
Spiked Amount	50.000		Recovery	=	103.36%	
62) 4-Bromofluorobenzene	11.12	95	237726	58.91	ug/l	0.00
Spiked Amount	50.000		Recovery	=	117.82%	
Target Compounds						
16) Acetone	2.42	43	8707	6.902	ug/l	95
17) Carbon Disulfide	2.56	76	2247	0.320	ug/l #	93
27) cis-1,2-Dichloroethene	4.56	96	10071	2.959	ug/l	93
44) Trichloroethene	7.20	130	11454	3.097	ug/l	86
79) 2-Chlorotoluene	11.40	91	18834	1.556	ug/l	98
82) 4-Chlorotoluene	11.49	91	10606	0.733	ug/l	96

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX070620\
Data File : VX017169.D
Acq On : 06 Jul 2020 15:40
Operator : JC/SP
Sample : L3159-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
50BR-20200629

Quant Time: Jul 07 07:02:09 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X062920W.M
Quant Title : SW846 8260
QLast Update : Mon Jun 29 17:10:38 2020
Response via : Initial Calibration

