

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX063020\
 Data File : VX017046.D
 Acq On : 30 Jun 2020 15:39
 Operator : JC/SP
 Sample : L3147-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 02BR-20200626

Quant Time: Jul 01 09:06:35 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	238063	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.84	114	402358	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	414665	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	219106	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	155203	52.88	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.76%	
35) Dibromofluoromethane	5.46	113	134289	52.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.52%	
50) Toluene-d8	8.70	98	494054	52.49	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.98%	
62) 4-Bromofluorobenzene	11.12	95	203688	55.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.72%	
Target Compounds						
3) Chloromethane	1.32	50	1775	0.637	ug/l	91
4) Vinyl Chloride	1.39	62	5186	2.005	ug/l	98
16) Acetone	2.42	43	5700	5.192	ug/l	98
27) cis-1,2-Dichloroethene	4.57	96	97842	33.037	ug/l	89
44) Trichloroethene	7.19	130	214608	69.727	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX063020\
Data File : VX017046.D
Acq On : 30 Jun 2020 15:39
Operator : JC/SP
Sample : L3147-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
02BR-20200626

Quant Time: Jul 01 09:06:35 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

