

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX040820\
 Data File : VX015573.D
 Acq On : 08 Apr 2020 16:47
 Operator : JC/SP
 Sample : L2190-17
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 M-16-200406

Quant Time: Apr 09 07:45:25 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X032720W.M
 Quant Title : SW846 8260
 QLast Update : Fri Mar 27 09:35:51 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	1067634	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.84	114	1674561	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	1605329	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	894020	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.04	65	702930	46.68	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.36%	
35) Dibromofluoromethane	5.47	113	530028	49.52	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.04%	
50) Toluene-d8	8.70	98	1974142	50.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.18%	
62) 4-Bromofluorobenzene	11.12	95	862782	53.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.74%	
Target Compounds						
3) Chloromethane	1.31	50	5471	0.457	ug/l	97
16) Acetone	2.42	43	129885	20.077	ug/l	95
27) cis-1,2-Dichloroethene	4.57	96	12862	0.994	ug/l	85
44) Trichloroethene	7.20	130	23958	1.944	ug/l	82
64) Tetrachloroethene	9.32	164	75510	6.233	ug/l	89

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX040820\
Data File : VX015573.D
Acq On : 08 Apr 2020 16:47
Operator : JC/SP
Sample : L2190-17
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
M-16-200406

Quant Time: Apr 09 07:45:25 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X032720W.M
Quant Title : SW846 8260
QLast Update : Fri Mar 27 09:35:51 2020
Response via : Initial Calibration

