

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX040920\
 Data File : VX015597.D
 Acq On : 09 Apr 2020 15:24
 Operator : JC/SP
 Sample : L2178-02
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-1D

Quant Time: Apr 10 07:40:04 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	1037497	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.84	114	1608418	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	1549488	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	872608	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.04	65	684483	46.78	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.56%	
35) Dibromofluoromethane	5.47	113	508999	49.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.02%	
50) Toluene-d8	8.70	98	1951572	51.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.12%	
62) 4-Bromofluorobenzene	11.12	95	844404	54.89	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.78%	
Target Compounds						
					Qvalue	
3) Chloromethane	1.31	50	3599	0.309	ug/l #	83
16) Acetone	2.42	43	19179	3.051	ug/l #	89
27) cis-1,2-Dichloroethene	4.58	96	5817	0.463	ug/l	67
30) Chloroform	5.19	83	11052	0.538	ug/l	87
64) Tetrachloroethene	9.32	164	3739	0.320	ug/l #	93

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

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