

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX051320\
 Data File : VX016176.D
 Acq On : 13 May 2020 13:58
 Operator : JC/SP
 Sample : L2597-08
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SS052MW710-200511

Quant Time: May 14 06:53:08 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X050420W.M
 Quant Title : SW846 8260
 QLast Update : Wed May 06 06:43:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.64	168	277881	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	465437	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	461349	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	231970	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	171577	46.92	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.84%	
35) Dibromofluoromethane	5.47	113	135599	45.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.98%	
50) Toluene-d8	8.70	98	573463	50.47	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.94%	
62) 4-Bromofluorobenzene	11.12	95	230016	53.23	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.46%	
Target Compounds						
					Qvalue	
3) Chloromethane	1.31	50	1086	0.318	ug/l	94
16) Acetone	2.42	43	6400	3.923	ug/l	90
27) cis-1,2-Dichloroethene	4.57	96	21810	6.198	ug/l	87
44) Trichloroethene	7.20	130	50298	10.561	ug/l	90
55) 1,1,2-Trichloroethane	9.20	97	2193	0.667	ug/l	91
64) Tetrachloroethene	9.32	164	10275	1.814	ug/l	96

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

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