

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX081920\
 Data File : VX017983.D
 Acq On : 19 Aug 2020 14:33
 Operator : JC/SP
 Sample : L3718-10
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB181-GW-278-280

Quant Time: Aug 29 05:49:26 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X081920W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 19 05:36:51 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.64	168	486128	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.84	114	725023	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	700229	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	336126	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.04	65	322505	45.26	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.52%	
35) Dibromofluoromethane	5.46	113	262745	48.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.32%	
50) Toluene-d8	8.70	98	872337	45.71	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.42%	
62) 4-Bromofluorobenzene	11.12	95	314522	44.62	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.24%	
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	2.36	96	5756	1.314	ug/l #	78
16) Acetone	2.42	43	14121	5.037	ug/l #	84
19) Methyl tert-butyl Ether	3.17	73	8573	0.512	ug/l #	73
24) 1,1-Dichloroethane	3.67	63	14137	1.557	ug/l #	89
27) cis-1,2-Dichloroethene	4.56	96	27716	5.216	ug/l	91
30) Chloroform	5.19	83	5671	0.571	ug/l	97
44) Trichloroethene	7.19	130	4007	0.739	ug/l	96

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

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