

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX080320\
 Data File : VX017750.D
 Acq On : 03 Aug 2020 16:54
 Operator : JC/SP
 Sample : L3526-12
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB179-GW-298-300

Manual Integrations
 APPROVED

MMDadoda
 8/4/2020 9:51:29 AM

Quant Time: Aug 04 11:54:39 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X072320W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 24 05:48:09 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	591629	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.84	114	874116	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	789640	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	408117	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.04	65	383466	50.46	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.92%	
35) Dibromofluoromethane	5.47	113	296490	50.72	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.44%	
50) Toluene-d8	8.70	98	997103	47.28	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.56%	
62) 4-Bromofluorobenzene	11.12	95	378460	44.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.68%	
Target Compounds						
16) Acetone	2.42	43	14012	4.189	ug/l	92
19) Methyl tert-butyl Ether	3.17	73	4459	0.211	ug/l	97
27) cis-1,2-Dichloroethene	4.57	96	96745	13.606	ug/l	88
30) Chloroform	5.18	83	3885m	0.324	ug/l	
44) Trichloroethene	7.19	130	878788	125.110	ug/l	95
64) Tetrachloroethene	9.32	164	76064	10.733	ug/l	98

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

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