

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX110720\
 Data File : VX019340.D
 Acq On : 06 Nov 2020 14:36
 Operator : JC/MD
 Sample : L4662-05
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB178-GW-278-280

Quant Time: Nov 07 02:01:59 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102720W.M
 Quant Title : SW846 8260
 QLast Update : Tue Oct 27 13:55:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	296282	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	524309	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	490030	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	246842	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	206956	54.16	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.32%	
35) Dibromofluoromethane	5.46	113	161925	49.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.34%	
50) Toluene-d8	8.70	98	634802	51.13	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.26%	
62) 4-Bromofluorobenzene	11.12	95	245654	52.25	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.50%	
Target Compounds						
16) Acetone	2.42	43	4341	3.293	ug/l	90
19) Methyl tert-butyl Ether	3.16	73	3923	0.394	ug/l #	79
30) Chloroform	5.18	83	2780	0.482	ug/l	94
44) Trichloroethene	7.19	130	1496	0.381	ug/l	94
64) Tetrachloroethene	9.32	164	912	0.244	ug/l	89

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

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