

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX110920\
 Data File : VX019367.D
 Acq On : 09 Nov 2020 14:59
 Operator : JC/MD
 Sample : L4668-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-18

Quant Time: Nov 10 02:22:14 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	294199	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	517381	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	492160	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	248867	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	206219	54.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.70%	
35) Dibromofluoromethane	5.46	113	162393	49.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.94%	
50) Toluene-d8	8.70	98	636151	51.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.86%	
62) 4-Bromofluorobenzene	11.12	95	248966	53.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.32%	
Target Compounds						
3) Chloromethane	1.31	50	2589	0.945	ug/l	94
11) Tert butyl alcohol	3.01	59	2373	3.167	ug/l #	94
16) Acetone	2.42	43	6987	5.338	ug/l	99
20) Methylene Chloride	2.84	84	1158	0.352	ug/l #	94
73) Isopropylbenzene	11.00	105	3534	0.201	ug/l	98
78) n-propylbenzene	11.34	91	4201	0.209	ug/l	97

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

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