

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091820\
 Data File : VX018493.D
 Acq On : 18 Sep 2020 15:32
 Operator : JC/SP
 Sample : L3939-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PAV-B44-090820-GW

Quant Time: Sep 21 07:11:56 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091720W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 17 12:17:36 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	526429	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.84	114	843659	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	804438	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	427726	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	364705	48.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.94%	
35) Dibromofluoromethane	5.47	113	283181	48.55	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.10%	
50) Toluene-d8	8.70	98	1018042	48.04	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.08%	
62) 4-Bromofluorobenzene	11.12	95	413927	49.11	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.22%	
Target Compounds						
					Qvalue	
3) Chloromethane	1.31	50	3417	0.585	ug/l #	84
16) Acetone	2.42	43	28178	9.894	ug/l	93
17) Carbon Disulfide	2.56	76	7474	0.491	ug/l #	93
20) Methylene Chloride	2.84	84	4781	0.761	ug/l	91
30) Chloroform	5.18	83	165721	15.251	ug/l	100
52) Toluene	8.76	92	13728	0.930	ug/l	92

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

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