

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX040820\
 Data File : VX015563.D
 Acq On : 08 Apr 2020 13:00
 Operator : JC/SP
 Sample : L2151-11
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CBH-140

Quant Time: Apr 09 07:02:58 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X032720W.M
 Quant Title : SW846 8260
 QLast Update : Fri Mar 27 09:35:51 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.64	168	1122833	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.84	114	1746207	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	1670151	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	868419	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.04	65	721319	45.55	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.10%	
35) Dibromofluoromethane	5.47	113	547298	49.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.06%	
50) Toluene-d8	8.70	98	2087657	50.80	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.60%	
62) 4-Bromofluorobenzene	11.12	95	900687	53.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.86%	
Target Compounds						
16) Acetone	2.42	43	28321	4.163	ug/l	95
20) Methylene Chloride	2.84	84	18412	1.449	ug/l	99
30) Chloroform	5.18	83	347968	15.651	ug/l	99
47) Bromodichloromethane	7.88	83	27272	1.513	ug/l	96
52) Toluene	8.77	92	10164	0.350	ug/l	96
68) m/p-Xylenes	10.34	106	7855	0.357	ug/l	98
84) 1,2,4-Trimethylbenzene	11.79	105	10069	0.203	ug/l	90

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

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