

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX040720\
 Data File : VX015552.D
 Acq On : 07 Apr 2020 20:33
 Operator : JC/SP
 Sample : L2151-09
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MH-109

Quant Time: Apr 08 04:02:44 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	1072865	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	1695808	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	1650500	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	910097	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	715149	47.26	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.52%	
35) Dibromofluoromethane	5.47	113	538014	49.63	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.26%	
50) Toluene-d8	8.70	98	2042382	51.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.34%	
62) 4-Bromofluorobenzene	11.12	95	896804	55.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.58%	
Target Compounds						
16) Acetone	2.42	43	14240	2.190	ug/l	91
20) Methylene Chloride	2.83	84	19557	1.611	ug/l #	93
30) Chloroform	5.19	83	56022	2.637	ug/l	99
52) Toluene	8.77	92	13973	0.495	ug/l	91
68) m/p-Xylenes	10.34	106	14729	0.678	ug/l	89
84) 1,2,4-Trimethylbenzene	11.79	105	17553	0.338	ug/l	96

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

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