

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX012819\
 Data File : VX007266.D
 Acq On : 28 Jan 2019 17:45
 Operator : JC/SP
 Sample : K1230-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MH-6

Quant Time: Jan 29 09:51:23 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X010819W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jan 08 14:06:59 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	534035	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	812742	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	776338	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	377309	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	286048	51.25	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.50%	
35) Dibromofluoromethane	5.51	113	258945	49.75	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.50%	
50) Toluene-d8	8.71	98	982115	51.55	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.10%	
62) 4-Bromofluorobenzene	11.13	95	340666	51.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.58%	
Target Compounds						
8) Diethyl Ether	2.19	74	11004	3.117	ug/l	96
16) Acetone	2.45	43	30873	11.189	ug/l	99
30) Chloroform	5.22	83	5798	0.626	ug/l #	80
65) Chlorobenzene	10.13	112	16438	0.971	ug/l	100
68) m/p-Xylenes	10.36	106	28909	2.625	ug/l	97
84) 1,2,4-Trimethylbenzene	11.80	105	14906	0.667	ug/l	99

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

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