

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX020519\
 Data File : VX007384.D
 Acq On : 05 Feb 2019 18:31
 Operator : JC/SP
 Sample : K1380-02
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KY038PS027-190204

Quant Time: Feb 06 01:36:47 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
 Quant Title : SW846 8260
 QLast Update : Sat Feb 02 01:32:36 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	439684	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	680191	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	634203	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	317726	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	238565	51.12	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.24%	
35) Dibromofluoromethane	5.51	113	209975	47.44	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.88%	
50) Toluene-d8	8.71	98	811496	49.42	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.84%	
62) 4-Bromofluorobenzene	11.14	95	287784	48.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.00%	
Target Compounds						
					Qvalue	
16) Acetone	2.48	43	13135	6.144	ug/l #	86
18) Methyl Acetate	2.79	43	95096	17.463	ug/l	93
27) cis-1,2-Dichloroethene	4.59	96	191898	39.287	ug/l	92
44) Trichloroethene	7.21	130	31139	5.751	ug/l	98
64) Tetrachloroethene	9.33	164	33762	5.648	ug/l	92
65) Chlorobenzene	10.14	112	9049	0.654	ug/l	95

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

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