

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX081319\
 Data File : VX011541.D
 Acq On : 13 Aug 2019 17:14
 Operator : JC/SP
 Sample : K4336-02
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KY038PS027-190812

Quant Time: Aug 14 02:23:17 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X080119W.M
 Quant Title : SW846 8260
 QLast Update : Fri Aug 02 01:55:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	203937	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	304929	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	273257	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	122214	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	114427	55.72	ug/l	0.00
Spiked Amount 50.000			Recovery	=	111.44%	
35) Dibromofluoromethane	5.49	113	93613	47.02	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.04%	
50) Toluene-d8	8.71	98	356032	47.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.76%	
62) 4-Bromofluorobenzene	11.13	95	121595	45.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.64%	
Target Compounds						
					Qvalue	
3) Chloromethane	1.32	50	1245	0.704	ug/l	99
16) Acetone	2.43	43	3785	3.872	ug/l #	86
27) cis-1,2-Dichloroethene	4.59	96	80763	35.723	ug/l	93
44) Trichloroethene	7.21	130	5418	2.253	ug/l	90
64) Tetrachloroethene	9.33	164	2996	1.291	ug/l	90
65) Chlorobenzene	10.13	112	6998	1.194	ug/l	98

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

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