

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100419\
 Data File : VX012811.D
 Acq On : 04 Oct 2019 15:01
 Operator : JC/SP
 Sample : K5189-13
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SA-MW128D-20190930

Quant Time: Oct 05 05:29:10 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092319W.M
 Quant Title : SW846 8260
 QLast Update : Mon Sep 23 12:27:01 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.65	168	183093	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	307821	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	257961	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	95782	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.05	65	120437	45.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.34%	
35) Dibromofluoromethane	5.49	113	93488	51.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.00%	
50) Toluene-d8	8.71	98	360293	51.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.06%	
62) 4-Bromofluorobenzene	11.14	95	121393	43.73	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.46%	
Target Compounds						
					Qvalue	
3) Chloromethane	1.32	50	1904	0.805	ug/l #	88
16) Acetone	2.43	43	7530	5.332	ug/l	90
19) Methyl tert-butyl Ether	3.19	73	2556	0.343	ug/l #	86
24) 1,1-Dichloroethane	3.69	63	12904	3.146	ug/l	96

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

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