

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX102519\
 Data File : VX013206.D
 Acq On : 24 Oct 2019 12:50
 Operator : JC/SP
 Sample : K5531-03
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-TT-MW203D1-GW-20191022

Quant Time: Oct 25 03:34:12 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102419W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 24 03:56:38 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.65	168	79507	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	119420	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	120389	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	75751	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.05	65	88934	53.48	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.96%	
35) Dibromofluoromethane	5.48	113	68023	51.08	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.16%	
50) Toluene-d8	8.71	98	246585	49.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.92%	
62) 4-Bromofluorobenzene	11.14	95	91894	47.04	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.08%	
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	2.37	96	7993	5.902	ug/l	95
16) Acetone	2.43	43	1661	2.120	ug/l	100
24) 1,1-Dichloroethane	3.69	63	48227	18.806	ug/l	97
27) cis-1,2-Dichloroethene	4.59	96	2850	1.718	ug/l	82
30) Chloroform	5.19	83	2536	0.929	ug/l #	74
32) 1,1,1-Trichloroethane	5.48	97	6201	2.554	ug/l	97
44) Trichloroethene	7.21	130	10209	5.997	ug/l	95

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

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