

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX092019\
 Data File : VX012566.D
 Acq On : 20 Sep 2019 18:42
 Operator : JC/SP
 Sample : K4883-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW5-R2-1

Quant Time: Sep 21 03:34:00 2019
 Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091719W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 17 15:27:10 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.65	168	172787	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	275199	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	227806	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	88914	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	117023	47.26	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.52%	
35) Dibromofluoromethane	5.48	113	83826	50.33	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.66%	
50) Toluene-d8	8.71	98	312024	50.61	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.22%	
62) 4-Bromofluorobenzene	11.14	95	111484	43.55	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.10%	
Target Compounds						
16) Acetone	2.43	43	24193	13.862	ug/l	97
27) cis-1,2-Dichloroethene	4.59	96	24326	14.435	ug/l	86
44) Trichloroethene	7.21	130	5615	4.013	ug/l	99
64) Tetrachloroethene	9.33	164	332711	318.721	ug/l	93

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

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