

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121818\
 Data File : VX006623.D
 Acq On : 19 Dec 2018 03:25
 Operator : JC/MD
 Sample : J6408-17
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SA.PZ163I-20181208

Quant Time: Jan 02 05:38:35 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
 Quant Title : SW846 8260
 QLast Update : Tue Dec 18 13:23:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	635579	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	1014519	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	935163	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	444824	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	355820	51.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.88%	
35) Dibromofluoromethane	5.50	113	301254	46.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.92%	
50) Toluene-d8	8.71	98	1214184	50.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.40%	
62) 4-Bromofluorobenzene	11.13	95	445726	50.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.64%	
Target Compounds						
12) 1,1-Dichloroethene	2.37	96	5375	0.947	ug/l	85
16) Acetone	2.45	43	5243	1.666	ug/l	90
24) 1,1-Dichloroethane	3.70	63	144523	13.845	ug/l	99
64) Tetrachloroethene	9.34	164	3459	0.467	ug/l	95
85) sec-Butylbenzene	11.94	105	22761	0.740	ug/l	84

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

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