

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX121218\
 Data File : VX006474.D
 Acq On : 13 Dec 2018 05:06
 Operator : JC/MD
 Sample : J6344-03
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RBR-200027

Quant Time: Dec 13 07:45:29 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112918W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 30 03:55:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	673931	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	1045410	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	962904	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	457429	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	354555	51.54	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.08%	
35) Dibromofluoromethane	5.51	113	313970	46.39	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.78%	
50) Toluene-d8	8.71	98	1240220	50.37	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.74%	
62) 4-Bromofluorobenzene	11.13	95	454262	48.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.92%	
Target Compounds						
16) Acetone	2.45	43	22756	7.048	ug/l #	86
18) Methyl Acetate	2.78	43	16423	1.521	ug/l	93
20) Methylene Chloride	2.86	84	11945	1.738	ug/l #	88
43) Isopropyl Acetate	6.46	43	30787	1.873	ug/l #	94

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

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