

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX071618\
 Data File : VX003404.D
 Acq On : 16 Jul 2018 21:53
 Operator : JC/MD
 Sample : PB111123ZHE#12
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB111123ZHE#12

Quant Time: Jul 17 02:28:49 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071618W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 16 13:15:28 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	269802	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	390631	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	348588	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	203829	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	167416	47.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.76%	
35) Dibromofluoromethane	5.51	113	151680	47.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.98%	
50) Toluene-d8	8.72	98	585585	50.21	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.42%	
62) 4-Bromofluorobenzene	11.14	95	195785	47.02	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.04%	
Target Compounds						
16) Acetone	2.45	43	26114	17.77	ug/l #	87
18) Methyl Acetate	2.78	43	9490	1.98	ug/l #	90
20) Methylene Chloride	2.86	84	30471	8.85	ug/l	85
43) Isopropyl Acetate	6.48	43	128438	19.09	ug/l #	93
74) N-amyl acetate	6.48	43	128438	20.46	ug/l #	92

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

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