

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU071218\
 Data File : VU025363.D
 Acq On : 13 Jul 2018 01:37
 Operator : MD/SY
 Sample : PB111044ZHE#01
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PB111044ZHE#01

Quant Time: Jul 13 05:00:56 2018
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jul 10 09:34:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	197671	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	302813	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	281127	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	144897	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	150665	49.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.12%	
35) Dibromofluoromethane	4.89	113	118430	46.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.72%	
50) Toluene-d8	7.57	98	434424	49.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.32%	
62) 4-Bromofluorobenzene	10.31	95	158830	45.10	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.20%	
Target Compounds						
16) Acetone	2.32	43	11045	11.778	ug/l	98
18) Methyl Acetate	2.62	43	16955	5.392	ug/l	96
20) Methylene Chloride	2.70	84	5824	3.109	ug/l	97
95) Naphthalene	13.76	128	8197	3.146	ug/l	97

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

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