

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU110718\
 Data File : VU027999.D
 Acq On : 07 Nov 2018 12:39
 Operator : MD/SY
 Sample : PB114573ZHE#04
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PB114573ZHE#04

Quant Time: Nov 07 14:21:36 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 02 02:28:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	199010	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	378701	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	336959	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	145932	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	166556	56.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	112.58%	
35) Dibromofluoromethane	4.89	113	125737	52.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.32%	
50) Toluene-d8	7.57	98	491828	53.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.64%	
62) 4-Bromofluorobenzene	10.31	95	164820	48.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.90%	
Target Compounds						
16) Acetone	2.32	43	11388	6.203	ug/l	95
18) Methyl Acetate	2.62	43	5786	1.146	ug/l	99
20) Methylene Chloride	2.70	84	12465	4.651	ug/l	98
43) Isopropyl Acetate	5.55	43	241963	26.575	ug/l	99

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

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