

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU110718\
 Data File : VU027996.D
 Acq On : 07 Nov 2018 11:30
 Operator : MD/SY
 Sample : PB114573ZHE#01
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PB114573ZHE#01

Quant Time: Nov 07 11:59:26 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	200983	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	380120	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	332922	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	137129	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	164372	55.01	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.02%	
35) Dibromofluoromethane	4.89	113	123528	51.54	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.08%	
50) Toluene-d8	7.57	98	486513	52.55	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.10%	
62) 4-Bromofluorobenzene	10.31	95	162223	47.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.00%	
Target Compounds						
16) Acetone	2.32	43	11234	6.060	ug/l	96
18) Methyl Acetate	2.62	43	6258	1.227	ug/l	94
20) Methylene Chloride	2.70	84	12431	4.592	ug/l	98
43) Isopropyl Acetate	5.55	43	245143	26.824	ug/l	100

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

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