

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100218\
 Data File : VU027338.D
 Acq On : 02 Oct 2018 10:41
 Operator : MD/SY
 Sample : PB113364TB
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PB113364TB

Quant Time: Oct 03 02:16:24 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
 Quant Title : SW846 8260
 QLast Update : Tue Oct 02 02:04:34 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	97940	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	167810	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	162265	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	74500	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	99430	60.80	ug/l	0.00
Spiked Amount 50.000			Recovery	=	121.60%	
35) Dibromofluoromethane	4.88	113	69575	60.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	121.00%	
50) Toluene-d8	7.57	98	264368	61.42	ug/l	0.00
Spiked Amount 50.000			Recovery	=	122.84%	
62) 4-Bromofluorobenzene	10.31	95	86874	54.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.62%	
Target Compounds						
16) Acetone	2.32	43	10047	10.399	ug/l	98
18) Methyl Acetate	2.62	43	3027	1.373	ug/l	96
20) Methylene Chloride	2.70	84	2477	1.841	ug/l	91
43) Isopropyl Acetate	5.55	43	107077	29.278	ug/l	100

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

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