

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100418\
 Data File : VU027382.D
 Acq On : 04 Oct 2018 01:45
 Operator : MD/SY
 Sample : J5163-07
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-7

Quant Time: Oct 04 08:20:37 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	91003	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	160846	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	152839	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	77298	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	97728	64.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	128.62%	
35) Dibromofluoromethane	4.88	113	68113	61.79	ug/l	0.00
Spiked Amount 50.000			Recovery	=	123.58%	
50) Toluene-d8	7.57	98	250191	60.64	ug/l	0.00
Spiked Amount 50.000			Recovery	=	121.28%	
62) 4-Bromofluorobenzene	10.31	95	83809	54.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.32%	
Target Compounds						
11) Tert butyl alcohol	2.83	59	194063	606.445	ug/l	98
16) Acetone	2.32	43	1827	2.035	ug/l	90
19) Methyl tert-butyl Ether	3.00	73	2548069	661.557	ug/l	99
22) Diisopropyl ether	3.58	45	34589	7.455	ug/l	96

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

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