

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU101018\
 Data File : VU027540.D
 Acq On : 09 Oct 2018 22:42
 Operator : MD/SY
 Sample : J5365-01
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 CY-1

Quant Time: Oct 10 07:24:15 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	116010	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	193394	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	167048	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	71945	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	96649	49.89	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.78%	
35) Dibromofluoromethane	4.89	113	66887	50.47	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.94%	
50) Toluene-d8	7.57	98	247611	49.92	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.84%	
62) 4-Bromofluorobenzene	10.31	95	79574	43.16	ug/l	0.00
Spiked Amount 50.000			Recovery	=	86.32%	
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.83	59	27918	68.437	ug/l	100
16) Acetone	2.32	43	230764	201.646	ug/l	100
25) 2-Butanone	4.27	43	60207	38.580	ug/l	98
51) 4-Methyl-2-Pentanone	7.47	43	21173	7.731	ug/l	94
52) Toluene	7.64	92	3569	0.986	ug/l	91
59) 2-Hexanone	8.37	43	4931	2.321	ug/l	88

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

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