

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU121318\
 Data File : VU028649.D
 Acq On : 12 Dec 2018 13:57
 Operator : MD/SY
 Sample : J6351-03
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 KY038PS035-181210

Quant Time: Dec 13 03:04:10 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U120618W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 07 00:39:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.98	168	105320	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.88	114	188760	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	169288	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	67857	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	80999	50.40	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.80%	
35) Dibromofluoromethane	4.88	113	62458	52.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.16%	
50) Toluene-d8	7.57	98	241121	51.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.62%	
62) 4-Bromofluorobenzene	10.30	95	80660	45.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.86%	
Target Compounds						
16) Acetone	2.32	43	42269	40.302	ug/l	100
18) Methyl Acetate	2.62	43	3124	1.149	ug/l	94
25) 2-Butanone	4.28	43	8122	5.667	ug/l	97
27) cis-1,2-Dichloroethene	4.23	96	1694	1.134	ug/l	88

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

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