

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU121718\
 Data File : VU028758.D
 Acq On : 18 Dec 2018 07:12
 Operator : MD/SY
 Sample : J6438-12 10X
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 23691

Quant Time: Dec 18 09:08:02 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	81016	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	142023	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	131481	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	52107	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	66930	54.13	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.26%	
35) Dibromofluoromethane	4.88	113	49934	55.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	111.74%	
50) Toluene-d8	7.56	98	187496	53.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.06%	
62) 4-Bromofluorobenzene	10.30	95	61654	46.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.32%	
Target Compounds						
16) Acetone	2.32	43	1826	2.263	ug/l	100
18) Methyl Acetate	2.62	43	2009	0.961	ug/l #	77
25) 2-Butanone	4.27	43	5771	5.234	ug/l	91
40) Benzene	5.39	78	2927	0.639	ug/l #	87
52) Toluene	7.64	92	1537	0.571	ug/l	95

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

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