

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU061318\
 Data File : VU024510.D
 Acq On : 13 Jun 2018 16:13
 Operator : MD/SY
 Sample : J3425-01 200X
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PGWPAL-AQ060818

Quant Time: Jun 13 17:20:56 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 13 13:55:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	184317	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	310944	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	301677	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	176430	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.32	65	147683	49.08	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.16%	
35) Dibromofluoromethane	4.89	113	126462	49.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.00%	
50) Toluene-d8	7.57	98	462150	49.10	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.20%	
62) 4-Bromofluorobenzene	10.31	95	178310	47.71	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.42%	
Target Compounds						
16) Acetone	2.32	43	385540	238.467	ug/l	97
19) Methyl tert-butyl Ether	3.00	73	41506	5.261	ug/l	99
40) Benzene	5.40	78	10396	0.999	ug/l	98
52) Toluene	7.64	92	10383	1.550	ug/l	99
68) m/p-Xylenes	9.38	106	6380	1.230	ug/l	95
84) 1,2,4-Trimethylbenzene	11.15	105	11847	0.990	ug/l	98
95) Naphthalene	13.75	128	14756	1.801	ug/l	98

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

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