

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU090719\
 Data File : VU034331.D
 Acq On : 06 Sep 2019 15:25
 Operator : JC/SP
 Sample : K4700-02
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 SINK

Quant Time: Sep 06 17:22:29 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U090619W.M
 Quant Title : SW846 8260
 QLast Update : Fri Sep 06 05:10:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.96	168	309347	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.87	114	465753	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.07	117	445922	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.46	152	278659	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.29	65	232934	49.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.54%	
35) Dibromofluoromethane	4.86	113	204905	47.28	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.56%	
50) Toluene-d8	7.55	98	784149	48.16	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.32%	
62) 4-Bromofluorobenzene	10.29	95	283114	45.04	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.08%	
Target Compounds						
16) Acetone	2.31	43	20491	8.460	ug/l	99
30) Chloroform	4.65	83	285907	37.534	ug/l	98
47) Bromodichloromethane	6.74	83	59246	9.623	ug/l	99
60) Dibromochloromethane	8.46	129	9669	1.942	ug/l	98
95) Naphthalene	13.73	128	140787	10.945	ug/l	98

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

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