

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU091019\
 Data File : VU034442.D
 Acq On : 10 Sep 2019 11:42
 Operator : JC/SP
 Sample : K4698-01
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 GW507-K2-35.0-40.0-090419

Quant Time: Sep 11 06:32:53 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	239786	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.87	114	344296	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.07	117	348908	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.46	152	205504	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.29	65	204098	56.26	ug/l	0.00
Spiked Amount 50.000			Recovery	=	112.52%	
35) Dibromofluoromethane	4.86	113	179078	55.90	ug/l	0.00
Spiked Amount 50.000			Recovery	=	111.80%	
50) Toluene-d8	7.55	98	639074	53.10	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.20%	
62) 4-Bromofluorobenzene	10.29	95	217185	46.74	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.48%	
Target Compounds						
16) Acetone	2.31	43	8814	4.695	ug/l	98
30) Chloroform	4.66	83	7745	1.312	ug/l	97
89) n-Butylbenzene	11.87	91	3284	0.310	ug/l	96
95) Naphthalene	13.74	128	12054	4.776	ug/l #	87

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

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