

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU091019\
 Data File : VU034440.D
 Acq On : 10 Sep 2019 10:56
 Operator : JC/SP
 Sample : K4697-02
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 GW507-K2-10.0-15.0-090419

Quant Time: Sep 11 06:28:33 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	234885	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.87	114	335214	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.07	117	345224	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.46	152	203148	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	5.29	65	190851	53.70	ug/l	0.00
Spiked Amount			Recovery	=	107.40%	
35) Dibromofluoromethane	4.86	113	170133	54.54	ug/l	0.00
Spiked Amount			Recovery	=	109.08%	
50) Toluene-d8	7.55	98	593399	50.64	ug/l	0.00
Spiked Amount			Recovery	=	101.28%	
62) 4-Bromofluorobenzene	10.29	95	202736	44.81	ug/l	0.00
Spiked Amount			Recovery	=	89.62%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	2.31	43	17630	9.586	ug/l	98
30) Chloroform	4.65	83	192016	33.200	ug/l	99
47) Bromodichloromethane	6.74	83	27187	6.136	ug/l	99
60) Dibromochloromethane	8.46	129	4455	1.243	ug/l	96
64) Tetrachloroethene	8.21	164	14714	4.913	ug/l	94

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

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