

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU090719\
 Data File : VU034352.D
 Acq On : 07 Sep 2019 00:18
 Operator : JC/SP
 Sample : K4580-40
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 T12-A2-(3.5)

Quant Time: Sep 07 05:07:09 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	303256	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.87	114	469536	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.07	117	466676	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.46	152	275582	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.29	65	243051	52.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.94%	
35) Dibromofluoromethane	4.86	113	211877	48.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.00%	
50) Toluene-d8	7.55	98	816265	49.73	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.46%	
62) 4-Bromofluorobenzene	10.29	95	292177	46.10	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.20%	
Target Compounds						
16) Acetone	2.31	43	13811	5.817	ug/l	97
20) Methylene Chloride	2.69	84	9347	2.020	ug/l	96
43) Isopropyl Acetate	5.53	43	63430	6.754	ug/l	99
95) Naphthalene	13.74	128	18297	4.883	ug/l #	92

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU090719\
Data File : VU034352.D
Acq On : 07 Sep 2019 00:18
Operator : JC/SP
Sample : K4580-40
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
T12-A2-(3.5)

Quant Time: Sep 07 05:07:09 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U090619W.M
Quant Title : SW846 8260
QLast Update : Fri Sep 06 05:10:13 2019
Response via : Initial Calibration

