

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092218\
 Data File : VU026999.D
 Acq On : 22 Sep 2018 06:08
 Operator : MD/SY
 Sample : J4934-17
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 SA-PZ145D-20180912

Quant Time: Sep 24 03:20:12 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 22 01:31:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	74135	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	148564	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	151829	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	73637	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	90280	60.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	121.02%	
35) Dibromofluoromethane	4.88	113	67573	52.47	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.94%	
50) Toluene-d8	7.57	98	237708	48.44	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.88%	
62) 4-Bromofluorobenzene	10.31	95	83903	49.82	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.64%	
Target Compounds						
24) 1,1-Dichloroethane	3.45	63	35936	14.689	ug/l	Qvalue 99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092218\
Data File : VU026999.D
Acq On : 22 Sep 2018 06:08
Operator : MD/SY
Sample : J4934-17
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 39 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
SA-PZ145D-20180912

Quant Time: Sep 24 03:20:12 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.M
Quant Title : SW846 8260
QLast Update : Sat Sep 22 01:31:57 2018
Response via : Initial Calibration

