

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
 Data File : VU027299.D
 Acq On : 29 Sep 2018 02:42
 Operator : MD/SY
 Sample : J4776-06
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EO-01-092718-A

Quant Time: Oct 01 02:13:55 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	108093	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	184596	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	179506	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	85823	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	107266	49.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.62%	
35) Dibromofluoromethane	4.88	113	74970	46.85	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.70%	
50) Toluene-d8	7.57	98	292221	47.92	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.84%	
62) 4-Bromofluorobenzene	10.31	95	99388	47.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.00%	
Target Compounds						
16) Acetone	2.32	43	36168	15.752	ug/l	96
20) Methylene Chloride	2.70	84	43530	22.342	ug/l	98
43) Isopropyl Acetate	5.55	43	107079	22.015	ug/l	99
95) Naphthalene	13.75	128	25303	5.130	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092818\
Data File : VU027299.D
Acq On : 29 Sep 2018 02:42
Operator : MD/SY
Sample : J4776-06
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EO-01-092718-A

Quant Time: Oct 01 02:13:55 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

