

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU062118\
 Data File : VU024809.D
 Acq On : 21 Jun 2018 14:27
 Operator : MD/SY
 Sample : J3561-04
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 S32-EB-2018-1

Quant Time: Jun 22 04:18:36 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 13 13:55:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	195139	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	299968	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	277843	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	149527	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	139117	43.67	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.34%	
35) Dibromofluoromethane	4.89	113	115498	46.39	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.78%	
50) Toluene-d8	7.57	98	430320	47.39	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.78%	
62) 4-Bromofluorobenzene	10.31	95	159220	44.16	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.32%	
Target Compounds						
16) Acetone	2.32	43	6323	3.69	ug/l	# 88
20) Methylene Chloride	2.71	84	66533	24.77	ug/l	96
25) 2-Butanone	4.27	43	27977	12.17	ug/l	95
29) Tetrahydrofuran	4.63	42	645631	453.46	ug/l	97

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU062118\
Data File : VU024809.D
Acq On : 21 Jun 2018 14:27
Operator : MD/SY
Sample : J3561-04
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
S32-EB-2018-1

Quant Time: Jun 22 04:18:36 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260
QLast Update : Wed Jun 13 13:55:26 2018
Response via : Initial Calibration

