

Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU110518\
 Data File : VU027967.D
 Acq On : 05 Nov 2018 19:01
 Operator : MD/SY
 Sample : J5811-04
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 DPC-91468-OW-01

Manual Integrations
 APPROVED

MMDadoda
 11/6/2018 12:06:59 PM

Quant Time: Nov 06 13:05:59 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 02 02:28:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.98	168	195713	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	371072	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.10	117	361649	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	170225	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	158043	54.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.62%	
35) Dibromofluoromethane	4.90	113	95405	40.78	ug/l	0.01
Spiked Amount 50.000			Recovery	=	81.56%	
50) Toluene-d8	7.57	98	479542	53.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.12%	
62) 4-Bromofluorobenzene	10.31	95	182319	54.69	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.38%	
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	3.05	59	4226418m	5486.630	ug/l	
18) Methyl Acetate	2.65	43	10728217	2160.252	ug/l	99
20) Methylene Chloride	2.69	84	286209	108.581	ug/l	98
37) Ethyl Acetate	4.40	43	25822938m	4534.790	ug/l	
51) 4-Methyl-2-Pentanone	7.49	43	26443303m	4438.273	ug/l	
67) Ethyl Benzene	9.26	91	5267	0.375	ug/l	97
68) m/p-Xylenes	9.38	106	6689	1.311	ug/l	91
69) o-Xylene	9.79	106	3226	0.626	ug/l	91

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

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