

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100418\
 Data File : VU027429.D
 Acq On : 04 Oct 2018 23:30
 Operator : MD/SY
 Sample : J5165-14
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EP1-MW-092718FD-W

Manual Integrations
 APPROVED

MMDadoda
 10/8/2018 1:52:38 PM

Quant Time: Oct 05 08:32:42 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
 Quant Title : SW846 8260
 QLast Update : Tue Oct 02 02:04:34 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	107532	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	191980	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	188211	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	108122	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	5.31	65	106345	59.23	ug/l	0.00
Spiked Amount 50.000			Recovery	=	118.46%	
35) Dibromofluoromethane	4.89	113	76124	57.86	ug/l	0.00
Spiked Amount 50.000			Recovery	=	115.72%	
50) Toluene-d8	7.57	98	295607	60.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	120.06%	
62) 4-Bromofluorobenzene	10.31	95	115148	62.92	ug/l	0.00
Spiked Amount 50.000			Recovery	=	125.84%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	2.32	43	63914	60.253	ug/l	99
17) Carbon Disulfide	2.48	76	16632	4.060	ug/l	100
25) 2-Butanone	4.28	43	10458	7.230	ug/l	94
30) Chloroform	4.68	83	4704	1.879	ug/l	89
73) Isopropylbenzene	10.17	105	6019	0.766	ug/l	100
78) n-propylbenzene	10.59	91	9462	1.003	ug/l	94
83) tert-Butylbenzene	11.10	119	22888	3.703	ug/l	90
84) 1,2,4-Trimethylbenzene	11.15	105	173822	26.616	ug/l	100
85) sec-Butylbenzene	11.32	105	22005	2.826	ug/l	98
86) p-Isopropyltoluene	11.48	119	12583	1.887	ug/l	97
89) n-Butylbenzene	11.89	91	5620m	2.713	ug/l	

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

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