

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100418\
 Data File : VU027430.D
 Acq On : 04 Oct 2018 23:53
 Operator : MD/SY
 Sample : J5165-15
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EP1-MW-E-092718

Quant Time: Oct 05 08:38:14 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	105898	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	189388	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	187010	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	107997	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	105621	59.73	ug/l	0.00
Spiked Amount			Recovery	=	119.46%	
35) Dibromofluoromethane	4.89	113	74176	57.15	ug/l	0.00
Spiked Amount			Recovery	=	114.30%	
50) Toluene-d8	7.57	98	288878	59.47	ug/l	0.00
Spiked Amount			Recovery	=	118.94%	
62) 4-Bromofluorobenzene	10.31	95	114675	63.52	ug/l	0.00
Spiked Amount			Recovery	=	127.04%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	2.32	43	46139	44.167	ug/l	97
17) Carbon Disulfide	2.48	76	9910	2.456	ug/l	99
25) 2-Butanone	4.28	43	9071	6.368	ug/l	98
30) Chloroform	4.68	83	1097	0.445	ug/l	90
67) Ethyl Benzene	9.26	91	2726	0.391	ug/l	91
69) o-Xylene	9.78	106	8843	3.648	ug/l	97
73) Isopropylbenzene	10.17	105	51433	6.557	ug/l	100
78) n-propylbenzene	10.59	91	121182	12.857	ug/l	100
80) 1,3,5-Trimethylbenzene	10.77	105	165278	25.882	ug/l	98
83) tert-Butylbenzene	11.10	119	30299	4.908	ug/l	88
84) 1,2,4-Trimethylbenzene	11.15	105	595400	91.274	ug/l	100
85) sec-Butylbenzene	11.32	105	101522	13.052	ug/l	99
86) p-Isopropyltoluene	11.48	119	75390	11.320	ug/l	98
89) n-Butylbenzene	11.89	91	117440	17.919	ug/l #	53
95) Naphthalene	13.75	128	8202	2.631	ug/l #	80

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

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