

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100418\
 Data File : VU027435.D
 Acq On : 05 Oct 2018 01:50
 Operator : MD/SY
 Sample : J5165-09
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EP-MW-H-092718

Quant Time: Oct 05 08:50:24 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	117737	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	202102	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	198070	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	113140	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	110097	56.00	ug/l	0.00
Spiked Amount			Recovery	=	112.00%	
35) Dibromofluoromethane	4.89	113	79213	57.19	ug/l	0.00
Spiked Amount			Recovery	=	114.38%	
50) Toluene-d8	7.57	98	310473	59.89	ug/l	0.00
Spiked Amount			Recovery	=	119.78%	
62) 4-Bromofluorobenzene	10.31	95	121849	63.25	ug/l	0.00
Spiked Amount			Recovery	=	126.50%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	2.32	43	82676	71.184	ug/l	98
17) Carbon Disulfide	2.48	76	13202	2.943	ug/l	99
25) 2-Butanone	4.28	43	21541	13.601	ug/l	98
30) Chloroform	4.69	83	1620	0.591	ug/l	90
69) o-Xylene	9.78	106	11404	4.441	ug/l	96
73) Isopropylbenzene	10.17	105	265850	32.353	ug/l	99
78) n-propylbenzene	10.59	91	569557	57.682	ug/l	100
80) 1,3,5-Trimethylbenzene	10.77	105	257417	38.478	ug/l	99
83) tert-Butylbenzene	11.10	119	59680	9.228	ug/l	89
84) 1,2,4-Trimethylbenzene	11.15	105	5761278	843.052	ug/l	99
85) sec-Butylbenzene	11.32	105	262023	32.154	ug/l	100
86) p-Isopropyltoluene	11.48	119	281568	40.356	ug/l	100
89) n-Butylbenzene	11.89	91	216948	30.108	ug/l #	72
95) Naphthalene	13.74	128	39104	6.080	ug/l #	96

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

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