

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
 Data File : VU027264.D
 Acq On : 28 Sep 2018 12:22
 Operator : MD/SY
 Sample : J5038-01
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 CPMW07S-Y4S3

Quant Time: Oct 02 06:46:09 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U092218W.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 22 01:31:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	99018	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	167284	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	160710	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	76036	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	102557	51.46	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.92%	
35) Dibromofluoromethane	4.88	113	71624	49.40	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.80%	
50) Toluene-d8	7.57	98	269741	48.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.62%	
62) 4-Bromofluorobenzene	10.31	95	89382	47.14	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.28%	
Target Compounds						
40) Benzene	5.40	78	22196	3.270	ug/l	98
52) Toluene	7.64	92	29491	7.581	ug/l	96
64) Tetrachloroethene	8.23	164	3172	2.561	ug/l	92
67) Ethyl Benzene	9.25	91	8930	1.274	ug/l	98
68) m/p-Xylenes	9.38	106	10141	5.021	ug/l	100
69) o-Xylene	9.78	106	11202	4.512	ug/l	100

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

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