

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092818\
 Data File : VU027265.D
 Acq On : 28 Sep 2018 12:45
 Operator : MD/SY
 Sample : J5038-02
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 FD1-Y4S3

Quant Time: Oct 02 13:15:08 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	94978	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	162433	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	155671	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	73662	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	99357	51.98	ug/l	0.00
Spiked Amount	50.000		Recovery	=	103.96%	
35) Dibromofluoromethane	4.89	113	69023	49.02	ug/l	0.00
Spiked Amount	50.000		Recovery	=	98.04%	
50) Toluene-d8	7.57	98	260074	48.47	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.94%	
62) 4-Bromofluorobenzene	10.31	95	87138	47.33	ug/l	0.00
Spiked Amount	50.000		Recovery	=	94.66%	
Target Compounds						
					Qvalue	
40) Benzene	5.39	78	20446	3.102	ug/l	100
52) Toluene	7.64	92	27739	7.343	ug/l	97
64) Tetrachloroethene	8.23	164	3108	2.591	ug/l	92
67) Ethyl Benzene	9.25	91	8520	1.255	ug/l	97
68) m/p-Xylenes	9.38	106	9650	4.959	ug/l	99
69) o-Xylene	9.78	106	10338	4.299	ug/l	96

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

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