

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU070718\
 Data File : VU025249.D
 Acq On : 06 Jul 2018 18:53
 Operator : MD/SY
 Sample : J3808-11
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 DUP-0473-180628

Manual Integrations
 APPROVED

MMDadoda
 7/9/2018 12:58:00 PM

Quant Time: Jul 09 10:38:22 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 13 13:55:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	200584	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	302319	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	279147	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	147911	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	153635	46.92	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.84%	
35) Dibromofluoromethane	4.89	113	120074	47.86	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.72%	
50) Toluene-d8	7.57	98	431930	47.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.40%	
62) 4-Bromofluorobenzene	10.31	95	156793	43.15	ug/l	0.00
Spiked Amount 50.000			Recovery	=	86.30%	
Target Compounds						
27) cis-1,2-Dichloroethene	4.23	96	5789	2.002	ug/l	51
44) Trichloroethene	6.19	130	6645	2.307	ug/l	92
65) Chlorobenzene	9.12	112	21607	3.030	ug/l	96
73) Isopropylbenzene	10.17	105	12610	1.110	ug/l	94
78) n-propylbenzene	10.59	91	19791	1.479	ug/l	98
83) tert-Butylbenzene	11.10	119	56563	5.974	ug/l	91
85) sec-Butylbenzene	11.33	105	88596	7.476	ug/l	96
89) n-Butylbenzene	11.90	91	8610m	0.914	ug/l	

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

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