

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU070518\
 Data File : VU025212.D
 Acq On : 05 Jul 2018 13:47
 Operator : MD/SY
 Sample : J3808-02
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 WP021MW463-180627

Quant Time: Jul 06 05:46:14 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	210361	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	313221	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	288071	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	147596	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	156787	45.65	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.30%	
35) Dibromofluoromethane	4.89	113	124784	48.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.00%	
50) Toluene-d8	7.57	98	446307	47.07	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.14%	
62) 4-Bromofluorobenzene	10.31	95	164230	43.62	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.24%	
Target Compounds						
12) 1,1-Dichloroethene	2.29	96	26244	10.957	ug/l	99
16) Acetone	2.33	43	5595	3.032	ug/l	99
21) trans-1,2-Dichloroethene	2.99	96	5444	2.064	ug/l	94
24) 1,1-Dichloroethane	3.45	63	146300	28.846	ug/l	99
27) cis-1,2-Dichloroethene	4.23	96	1094028	360.712	ug/l	89
29) Tetrahydrofuran	4.65	42	3867	2.519	ug/l	94
30) Chloroform	4.68	83	1215	0.237	ug/l	95
44) Trichloroethene	6.19	130	65673	22.003	ug/l	94
55) 1,1,2-Trichloroethane	8.07	97	1425	0.520	ug/l	89

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

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