

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX101818\
 Data File : VX005431.D
 Acq On : 18 Oct 2018 15:43
 Operator : JC/MD
 Sample : J5588-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MH-6

Quant Time: Oct 19 04:46:44 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X101218W.M
 Quant Title : SW846 8260
 QLast Update : Fri Oct 12 11:28:43 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	317989	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	448625	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	391311	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	190944	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	175380	49.69	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.38%	
35) Dibromofluoromethane	5.50	113	138464	46.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.20%	
50) Toluene-d8	8.72	98	524817	51.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.90%	
62) 4-Bromofluorobenzene	11.14	95	174804	45.62	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.24%	
Target Compounds						
8) Diethyl Ether	2.19	74	4445	2.137	ug/l	95
11) Tert butyl alcohol	3.07	59	1581	1.588	ug/l #	62
16) Acetone	2.45	43	13328	6.532	ug/l	99
29) Tetrahydrofuran	5.17	42	4262	2.238	ug/l	93
30) Chloroform	5.21	83	5844	1.070	ug/l	96
47) Bromodichloromethane	7.90	83	1371	0.351	ug/l #	92
65) Chlorobenzene	10.14	112	4022	0.483	ug/l	95
68) m/p-Xylenes	10.36	106	5300	1.009	ug/l	100
84) 1,2,4-Trimethylbenzene	11.80	105	3784	0.378	ug/l	92

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

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