

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX102318\
 Data File : VX005524.D
 Acq On : 23 Oct 2018 16:40
 Operator : JC/MD
 Sample : J5639-13
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BS-DRUM

Quant Time: Oct 24 02:28:56 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	5.67	168	239641	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	340362	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	289034	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	139043	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	140461	52.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.62%	
35) Dibromofluoromethane	5.50	113	117951	52.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.64%	
50) Toluene-d8	8.71	98	383864	49.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.20%	
62) 4-Bromofluorobenzene	11.14	95	127782	43.95	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.90%	
Target Compounds						
					Qvalue	
16) Acetone	2.45	43	15270	9.930	ug/l #	87
17) Carbon Disulfide	2.57	76	3174	0.499	ug/l	97
18) Methyl Acetate	2.78	43	5499	1.257	ug/l	98
30) Chloroform	5.21	83	1567	0.381	ug/l	87
51) 4-Methyl-2-Pentanone	8.65	43	2675	0.684	ug/l	93

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

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