

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX081318\
 Data File : VX003996.D
 Acq On : 13 Aug 2018 13:29
 Operator : JC/MD
 Sample : J4439-14
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 IDW-SO3-SRI3

Manual Integrations
 APPROVED

MMDadoda
 8/14/2018 12:55:45 PM

Quant Time: Aug 14 04:08:47 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X080618W.M
 Quant Title : SW846 8260
 QLast Update : Mon Aug 13 01:34:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	289758	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	473311	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	435220	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	230438	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	213857	55.68	ug/l	0.00
Spiked Amount 50.000			Recovery	=	111.36%	
35) Dibromofluoromethane	5.51	113	177087	49.39	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.78%	
50) Toluene-d8	8.72	98	676758	48.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.06%	
62) 4-Bromofluorobenzene	11.14	95	228199	50.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.16%	
Target Compounds						
16) Acetone	2.43	43	15459	6.77	ug/l	99
18) Methyl Acetate	2.77	43	5971	1.20	ug/l	92
20) Methylene Chloride	2.86	84	10643	1.92	ug/l	96
25) 2-Butanone	4.70	43	4062m	1.18	ug/l	
43) Isopropyl Acetate	6.46	43	192538	21.94	ug/l	99
74) N-amyl acetate	6.46	43	192538	21.72	ug/l #	99

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

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