

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX101018\
 Data File : VX005241.D
 Acq On : 11 Oct 2018 08:32
 Operator : JC/MD
 Sample : J5364-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KY038PS035-181008

Quant Time: Oct 12 10:08:46 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 03 04:00:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	172800	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	265054	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	257880	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	138860	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	132363	53.62	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.24%	
35) Dibromofluoromethane	5.51	113	113447	50.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.00%	
50) Toluene-d8	8.72	98	377953	50.61	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.22%	
62) 4-Bromofluorobenzene	11.14	95	130986	48.69	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.38%	
Target Compounds						
16) Acetone	2.45	43	22540	15.64	ug/l #	85
21) trans-1,2-Dichloroethene	3.17	96	1729	0.84	ug/l #	80
27) cis-1,2-Dichloroethene	4.60	96	55789	24.37	ug/l	93
44) Trichloroethene	7.21	130	6381	2.71	ug/l	97
64) Tetrachloroethene	9.34	164	10164	4.00	ug/l	96
65) Chlorobenzene	10.14	112	1677	0.25	ug/l	99

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

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