

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX111318\
 Data File : VX005959.D
 Acq On : 13 Nov 2018 19:54
 Operator : JC/MD
 Sample : J5931-03
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KY038PS035-181112

Quant Time: Nov 14 04:37:23 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 07 14:40:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	93193	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	133479	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	121665	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	59338	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	59315	55.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.72%	
35) Dibromofluoromethane	5.50	113	48865	53.82	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.64%	
50) Toluene-d8	8.71	98	160590	53.21	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.42%	
62) 4-Bromofluorobenzene	11.14	95	54216	49.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.92%	
Target Compounds						
16) Acetone	2.45	43	9550	16.555	ug/l	97
27) cis-1,2-Dichloroethene	4.59	96	23289	24.276	ug/l	91
44) Trichloroethene	7.21	130	1713	1.798	ug/l	97
64) Tetrachloroethene	9.34	164	2253	2.175	ug/l #	85
65) Chlorobenzene	10.14	112	1739	0.713	ug/l	97

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

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