

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062518\
 Data File : VX002870.D
 Acq On : 25 Jun 2018 20:22
 Operator : JC/MD
 Sample : J3604-02DL 5X
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-10DL

Quant Time: Jun 27 05:47:29 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X062518W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jun 25 14:38:05 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	240311	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	328982	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	301630	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	165848	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	156052	46.98	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.96%	
35) Dibromofluoromethane	5.51	113	124153	47.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.32%	
50) Toluene-d8	8.72	98	461115	48.47	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.94%	
62) 4-Bromofluorobenzene	11.14	95	161340	43.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.94%	
Target Compounds						
27) cis-1,2-Dichloroethene	4.59	96	2894	1.00	ug/l	# 1
30) Chloroform	5.22	83	1354	0.27	ug/l	89
44) Trichloroethene	7.22	130	1317	0.42	ug/l	75
64) Tetrachloroethene	9.34	164	24205	7.22	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX062518\
Data File : VX002870.D
Acq On : 25 Jun 2018 20:22
Operator : JC/MD
Sample : J3604-02DL 5X
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10DL

Quant Time: Jun 27 05:47:29 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X062518W.M
Quant Title : SW846 8260
QLast Update : Mon Jun 25 14:38:05 2018
Response via : Initial Calibration

