

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX090418\
 Data File : VX004352.D
 Acq On : 04 Sep 2018 17:33
 Operator : JC/MD
 Sample : J4757-02
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WC-BR

Quant Time: Sep 05 07:35:44 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X082218W.M
 Quant Title : SW846 8260
 QLast Update : Thu Aug 23 05:39:44 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	277425	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	444475	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	410410	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	220939	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	191769	53.04	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.08%	
35) Dibromofluoromethane	5.50	113	170390	51.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.06%	
50) Toluene-d8	8.72	98	634593	50.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.68%	
62) 4-Bromofluorobenzene	11.14	95	222151	47.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.40%	
Target Compounds						
18) Methyl Acetate	2.78	43	5858	1.09	ug/l	99
20) Methylene Chloride	2.86	84	145313	41.75	ug/l	99
43) Isopropyl Acetate	6.47	43	115651	15.32	ug/l	100
86) p-Isopropyltoluene	12.07	119	15534	1.21	ug/l	93
95) Naphthalene	13.83	128	97202	6.74	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX090418\
Data File : VX004352.D
Acq On : 04 Sep 2018 17:33
Operator : JC/MD
Sample : J4757-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WC-BR

Quant Time: Sep 05 07:35:44 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X082218W.M
Quant Title : SW846 8260
QLast Update : Thu Aug 23 05:39:44 2018
Response via : Initial Calibration

