

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX110518\
 Data File : VX005778.D
 Acq On : 05 Nov 2018 15:11
 Operator : JC/MD
 Sample : J5794-04
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TP-12B-S

Quant Time: Nov 06 01:33:18 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X103118W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 31 18:23:06 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	149752	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	253572	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	224956	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	93145	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	112650	51.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.90%	
35) Dibromofluoromethane	5.51	113	87240	49.69	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.38%	
50) Toluene-d8	8.71	98	300312	48.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.54%	
62) 4-Bromofluorobenzene	11.14	95	93385	41.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	82.18%	
Target Compounds						
18) Methyl Acetate	2.78	43	3905	1.243	ug/l	95
20) Methylene Chloride	2.86	84	169763	89.916	ug/l	92
43) Isopropyl Acetate	6.46	43	113619	22.712	ug/l	98
95) Naphthalene	13.83	128	8409	1.295	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX110518\
Data File : VX005778.D
Acq On : 05 Nov 2018 15:11
Operator : JC/MD
Sample : J5794-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TP-12B-S

Quant Time: Nov 06 01:33:18 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X103118W.M
Quant Title : SW846 8260
QLast Update : Wed Oct 31 18:23:06 2018
Response via : Initial Calibration

