

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX061418\
 Data File : VX002586.D
 Acq On : 15 Jun 2018 07:21
 Operator : JC/MD
 Sample : J3239-14
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 AU-06-061318-C

Quant Time: Jun 15 08:13:47 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jun 05 03:22:35 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	226630	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	330163	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	308717	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.09	152	179558	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	156107	48.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.06%	
35) Dibromofluoromethane	5.51	113	117797	45.37	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.74%	
50) Toluene-d8	8.72	98	478143	48.05	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.10%	
62) 4-Bromofluorobenzene	11.14	95	170526	44.44	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.88%	
Target Compounds						
18) Methyl Acetate	2.78	43	21909	4.280	ug/l	95
20) Methylene Chloride	2.86	84	8917	0.758	ug/l	96
25) 2-Butanone	4.71	43	6409	2.884	ug/l	99
95) Naphthalene	13.84	128	181279	13.870	ug/l	100

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061418\
Data File : VX002586.D
Acq On : 15 Jun 2018 07:21
Operator : JC/MD
Sample : J3239-14
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AU-06-061318-C

Quant Time: Jun 15 08:13:47 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260
QLast Update : Tue Jun 05 03:22:35 2018
Response via : Initial Calibration

