

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091318\
 Data File : VX004514.D
 Acq On : 13 Sep 2018 14:21
 Operator : JC/MD
 Sample : J4873-04
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 NF-CON-2

Quant Time: Sep 14 05:07:54 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 12 02:17:23 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	245118	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	352767	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	328386	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	185220	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	164427	52.02	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.04%	
35) Dibromofluoromethane	5.50	113	127100	39.37	ug/l	0.00
Spiked Amount 50.000			Recovery	=	78.74%	
50) Toluene-d8	8.72	98	490352	45.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.60%	
62) 4-Bromofluorobenzene	11.14	95	174310	45.11	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.22%	
Target Compounds						
16) Acetone	2.45	43	34985	22.679	ug/l	91
20) Methylene Chloride	2.86	84	64323	20.327	ug/l	96
25) 2-Butanone	4.70	43	3106	1.265	ug/l	94
95) Naphthalene	13.83	128	135943	10.751	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091318\
Data File : VX004514.D
Acq On : 13 Sep 2018 14:21
Operator : JC/MD
Sample : J4873-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
NF-CON-2

Quant Time: Sep 14 05:07:54 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260
QLast Update : Wed Sep 12 02:17:23 2018
Response via : Initial Calibration

