

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX121918\
 Data File : VX006655.D
 Acq On : 19 Dec 2018 23:24
 Operator : JC/MD
 Sample : J6194-08
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OK-01-121818-A

Quant Time: Dec 20 04:58:08 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
 Quant Title : SW846 8260
 QLast Update : Wed Dec 19 15:20:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	633766	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	1005821	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	959280	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	461525	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	348178	50.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.94%	
35) Dibromofluoromethane	5.50	113	299039	47.01	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.02%	
50) Toluene-d8	8.71	98	1222944	51.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.00%	
62) 4-Bromofluorobenzene	11.13	95	439636	50.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.12%	
Target Compounds						
16) Acetone	2.45	43	31933	10.176	ug/l	95
18) Methyl Acetate	2.78	43	24915	2.700	ug/l	96
20) Methylene Chloride	2.86	84	186862	28.796	ug/l	99
43) Isopropyl Acetate	6.46	43	27114	2.055	ug/l #	94

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX121918\
Data File : VX006655.D
Acq On : 19 Dec 2018 23:24
Operator : JC/MD
Sample : J6194-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OK-01-121818-A

Quant Time: Dec 20 04:58:08 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260
QLast Update : Wed Dec 19 15:20:58 2018
Response via : Initial Calibration

