

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX072320\
 Data File : VX017513.D
 Acq On : 24 Jul 2020 00:20
 Operator : JC/SP
 Sample : PB130403ZHE#03
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB130403ZHE#03

Quant Time: Jul 24 07:08:26 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X072320W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 24 05:48:09 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.64	168	511445	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	826905	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	867964	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	439766	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.04	65	357109	54.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.72%	
35) Dibromofluoromethane	5.47	113	291490	52.71	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.42%	
50) Toluene-d8	8.70	98	1000338	50.14	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.28%	
62) 4-Bromofluorobenzene	11.12	95	443012	55.49	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.98%	
Target Compounds						
16) Acetone	2.42	43	67370	23.299	ug/l	99
18) Methyl Acetate	2.75	43	8139	1.195	ug/l #	80
20) Methylene Chloride	2.84	84	51331	8.356	ug/l	90
43) Isopropyl Acetate	6.42	43	178222	12.448	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX072320\
Data File : VX017513.D
Acq On : 24 Jul 2020 00:20
Operator : JC/SP
Sample : PB130403ZHE#03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 38 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PB130403ZHE#03

Quant Time: Jul 24 07:08:26 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X072320W.M
Quant Title : SW846 8260
QLast Update : Fri Jul 24 05:48:09 2020
Response via : Initial Calibration

