

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX071720\
 Data File : VX017417.D
 Acq On : 17 Jul 2020 14:17
 Operator : JC/SP
 Sample : PB130262ZHE#09
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB130262ZHE#09

Quant Time: Jul 18 07:27:51 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jul 14 02:59:51 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	246534	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	446047	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	506769	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	252705	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.04	65	179841	58.12	ug/l	0.00
Spiked Amount 50.000			Recovery	=	116.24%	
35) Dibromofluoromethane	5.47	113	154675	51.48	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.96%	
50) Toluene-d8	8.70	98	584419	53.92	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.84%	
62) 4-Bromofluorobenzene	11.12	95	239377	55.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	111.20%	
Target Compounds						
16) Acetone	2.42	43	32492	29.434	ug/l	98
18) Methyl Acetate	2.75	43	3496	1.282	ug/l #	87
20) Methylene Chloride	2.84	84	36882	13.694	ug/l	96
43) Isopropyl Acetate	6.41	43	86474	12.754	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX071720\
Data File : VX017417.D
Acq On : 17 Jul 2020 14:17
Operator : JC/SP
Sample : PB130262ZHE#09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PB130262ZHE#09

Quant Time: Jul 18 07:27:51 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
Quant Title : SW846 8260
QLast Update : Tue Jul 14 02:59:51 2020
Response via : Initial Calibration

