

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX081718\
 Data File : VX004140.D
 Acq On : 17 Aug 2018 18:49
 Operator : JC/MD
 Sample : PB112172TB
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB112172TB

Quant Time: Aug 18 03:59:30 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X081418W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 14 13:27:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	253682	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	412377	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	385307	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	207534	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	188937	52.86	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.72%	
35) Dibromofluoromethane	5.51	113	162391	52.62	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.24%	
50) Toluene-d8	8.72	98	599618	47.92	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.84%	
62) 4-Bromofluorobenzene	11.14	95	203596	44.86	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.72%	
Target Compounds						
3) Chloromethane	1.32	50	5592	1.85	ug/l	Qvalue 100
16) Acetone	2.44	43	16740	9.01	ug/l	95
18) Methyl Acetate	2.78	43	6303	1.53	ug/l	94
20) Methylene Chloride	2.86	84	14858	4.34	ug/l	95
43) Isopropyl Acetate	6.47	43	195041	25.89	ug/l	99
74) N-amyl acetate	6.47	43	195041	29.53	ug/l #	99

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

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