

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX030719\
 Data File : VX007908.D
 Acq On : 08 Mar 2019 02:57
 Operator : JC/SP
 Sample : K1757-08
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TP-6CD-D

Quant Time: Mar 08 08:17:51 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X030619W.M
 Quant Title : SW846 8260
 QLast Update : Thu Mar 07 07:33:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	277868	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	433948	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	425161	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	206984	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	152926	52.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.34%	
35) Dibromofluoromethane	5.50	113	141905	51.12	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.24%	
50) Toluene-d8	8.71	98	542591	53.22	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.44%	
62) 4-Bromofluorobenzene	11.13	95	199491	55.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	111.62%	
Target Compounds						
16) Acetone	2.45	43	19807	11.674	ug/l	100
18) Methyl Acetate	2.78	43	6519	1.809	ug/l	98
20) Methylene Chloride	2.86	84	3081200	1141.106	ug/l	98
25) 2-Butanone	4.70	43	5300	2.207	ug/l	99
43) Isopropyl Acetate	6.46	43	15487	2.244	ug/l	99
84) 1,2,4-Trimethylbenzene	11.80	105	12632	1.069	ug/l	97

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

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