

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX121918\
 Data File : VX006641.D
 Acq On : 19 Dec 2018 17:11
 Operator : JC/MD
 Sample : J6486-03
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 S72-EB-2018-2

Quant Time: Dec 20 04:24:30 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	620186	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	990579	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	903675	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	426880	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	343886	51.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.90%	
35) Dibromofluoromethane	5.50	113	294142	46.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.92%	
50) Toluene-d8	8.71	98	1172414	49.64	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.28%	
62) 4-Bromofluorobenzene	11.13	95	426851	49.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.70%	
Target Compounds						
16) Acetone	2.45	43	5596	1.822	ug/l	89
20) Methylene Chloride	2.86	84	139231	21.926	ug/l	99
25) 2-Butanone	4.71	43	5747	1.330	ug/l	97
29) Tetrahydrofuran	5.15	42	211498	74.340	ug/l	99

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

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