

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100118\
 Data File : VX004896.D
 Acq On : 02 Oct 2018 03:31
 Operator : JC/MD
 Sample : J5041-04
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 S13-0247

Manual Integrations
 APPROVED

MMDadoda
 10/4/2018 7:07:55 PM

Quant Time: Oct 02 05:49:02 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 26 14:17:02 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	238430	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	374987	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	368729	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	213104	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	172761	50.72	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.44%	
35) Dibromofluoromethane	5.50	113	145592	45.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.62%	
50) Toluene-d8	8.72	98	544799	51.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.12%	
62) 4-Bromofluorobenzene	11.14	95	208570	54.80	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.60%	
Target Compounds						
					Qvalue	
3) Chloromethane	1.32	50	3213	0.815	ug/l	94
16) Acetone	2.45	43	8680	4.365	ug/l	99
31) Cyclohexane	5.57	56	437610	100.980	ug/l	93
39) Methylcyclohexane	7.46	83	457288	103.281	ug/l	94
40) Benzene	6.15	78	3584	0.264	ug/l	92
52) Toluene	8.79	92	2753	0.369	ug/l	95
73) Isopropylbenzene	11.02	105	49178	3.661	ug/l	98
78) n-propylbenzene	11.36	91	27737	1.800	ug/l	96
83) tert-Butylbenzene	11.77	119	10700	0.968	ug/l	88
85) sec-Butylbenzene	11.94	105	39276	2.916	ug/l	87
89) n-Butylbenzene	12.38	91	4903m	0.442	ug/l	
95) Naphthalene	13.83	128	41395	3.021	ug/l #	89

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

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