

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX061920\
 Data File : VX016905.D
 Acq On : 20 Jun 2020 03:59
 Operator : JC/SP
 Sample : L3063-03
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TW-3

Manual Integrations
 APPROVED

MMDadoda
 6/22/2020 12:22:51 PM

Quant Time: Jun 22 07:11:52 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061820W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jun 19 12:07:48 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	304526	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	492194	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	475600	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	287196	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	172352	47.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.12%	
35) Dibromofluoromethane	5.47	113	159486	50.65	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.30%	
50) Toluene-d8	8.70	98	567394	45.04	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.08%	
62) 4-Bromofluorobenzene	11.12	95	240578	49.85	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.70%	
Target Compounds						
8) Diethyl Ether	2.17	74	2380	1.025	ug/l	90
16) Acetone	2.42	43	8144	6.125	ug/l	99
17) Carbon Disulfide	2.55	76	7515	0.894	ug/l	98
20) Methylene Chloride	2.84	84	3652	1.145	ug/l	95
68) m/p-Xylenes	10.35	106	2187	0.346	ug/l	98
78) n-propylbenzene	11.34	91	4313	0.209	ug/l	98
84) 1,2,4-Trimethylbenzene	11.79	105	7676	0.467	ug/l	99
89) n-Butylbenzene	12.37	91	3715m	0.258	ug/l	

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

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