

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062920\
 Data File : VX017016.D
 Acq On : 29 Jun 2020 18:55
 Operator : JC/SP
 Sample : L3020-15
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-26

Quant Time: Jul 01 05:24:22 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X062920W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jun 29 17:10:38 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.64	168	277089	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	479611	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	480823	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	249698	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	6.03	65	178527	52.26	ug/l	0.00
Spiked Amount			Recovery	=	104.52%	
35) Dibromofluoromethane	5.47	113	153696	50.66	ug/l	0.00
Spiked Amount			Recovery	=	101.32%	
50) Toluene-d8	8.70	98	579211	51.63	ug/l	0.00
Spiked Amount			Recovery	=	103.26%	
62) 4-Bromofluorobenzene	11.12	95	244219	55.69	ug/l	0.00
Spiked Amount			Recovery	=	111.38%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Chloromethane	1.32	50	6453	1.989	ug/l	97
16) Acetone	2.42	43	8199	6.416	ug/l	95
17) Carbon Disulfide	2.55	76	4664	0.632	ug/l #	86
30) Chloroform	5.17	83	5951	1.088	ug/l #	83
47) Bromodichloromethane	7.88	83	5309	1.103	ug/l	98
60) Dibromochloromethane	9.57	129	4708	1.193	ug/l	99
64) Tetrachloroethene	9.32	164	6079	1.567	ug/l	95
67) Ethyl Benzene	10.23	91	4883	0.295	ug/l	94
68) m/p-Xylenes	10.34	106	5621	0.883	ug/l	98
69) o-Xylene	10.68	106	2063	0.341	ug/l	90
87) 1,3-Dichlorobenzene	12.01	146	1945	0.238	ug/l	94
89) n-Butylbenzene	12.37	91	3667	0.282	ug/l	93
93) 1,2,4-Trichlorobenzene	13.63	180	2219	0.426	ug/l	98
94) Hexachlorobutadiene	13.76	225	1185	1.434	ug/l	88
96) 1,2,3-Trichlorobenzene	14.00	180	1845	0.366	ug/l	91

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

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