

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062620\
 Data File : VX016984.D
 Acq On : 26 Jun 2020 15:29
 Operator : JC/SP
 Sample : L3117-02
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KY038PS027-200624

Quant Time: Jun 29 06:14:59 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X062620W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jun 26 08:34:16 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.64	168	222754	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	399595	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	400307	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	202454	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	181464	53.95	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.90%	
35) Dibromofluoromethane	5.47	113	134959	51.14	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.28%	
50) Toluene-d8	8.70	98	501543	51.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.68%	
62) 4-Bromofluorobenzene	11.12	95	208831	55.82	ug/l	0.00
Spiked Amount 50.000			Recovery	=	111.64%	
Target Compounds						
					Qvalue	
3) Chloromethane	1.32	50	3140	1.151	ug/l	97
16) Acetone	2.42	43	3495	2.658	ug/l	98
27) cis-1,2-Dichloroethene	4.57	96	77783	25.232	ug/l	91
44) Trichloroethene	7.18	130	3778	1.246	ug/l	66
60) Dibromochloromethane	9.56	129	4164	1.266	ug/l	94
64) Tetrachloroethene	9.32	164	1446	0.534	ug/l	84
65) Chlorobenzene	10.12	112	6645	0.812	ug/l	97
71) Bromoform	10.84	173	4262	1.762	ug/l #	83

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

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