

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX061220\
 Data File : VX016726.D
 Acq On : 12 Jun 2020 14:24
 Operator : JC/SP
 Sample : L3001-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SOURCE-WATER

Quant Time: Jun 15 07:51:32 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X052720W.M
 Quant Title : SW846 8260
 QLast Update : Wed May 27 16:31:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	225016	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	354542	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	365431	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	186400	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	133755	47.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.74%	
35) Dibromofluoromethane	5.47	113	111128	50.90	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.80%	
50) Toluene-d8	8.70	98	436957	51.42	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.84%	
62) 4-Bromofluorobenzene	11.12	95	176780	53.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.68%	
Target Compounds						
16) Acetone	2.41	43	8824	7.207	ug/l	93
25) 2-Butanone	4.64	43	2260	1.089	ug/l	95
30) Chloroform	5.18	83	31930	7.340	ug/l	96
47) Bromodichloromethane	7.88	83	47848	13.519	ug/l	99
60) Dibromochloromethane	9.57	129	65881	23.357	ug/l	97
65) Chlorobenzene	10.12	112	3667	0.492	ug/l	99
71) Bromoform	10.84	173	29293	12.188	ug/l #	96

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

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