

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX052820\
 Data File : VX016486.D
 Acq On : 28 May 2020 15:32
 Operator : JC/SP
 Sample : L2782-04
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SOURCE-WATER

Quant Time: May 29 09:13:44 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	286772	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	467526	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	456077	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	222654	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	170835	47.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.94%	
35) Dibromofluoromethane	5.46	113	135537	47.08	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.16%	
50) Toluene-d8	8.70	98	574516	51.27	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.54%	
62) 4-Bromofluorobenzene	11.12	95	225734	52.14	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.28%	
Target Compounds						
16) Acetone	2.42	43	9925	6.361	ug/l	96
17) Carbon Disulfide	2.55	76	11564	1.411	ug/l #	91
47) Bromodichloromethane	7.88	83	2388	0.512	ug/l #	95
60) Dibromochloromethane	9.57	129	4500	1.210	ug/l	100
65) Chlorobenzene	10.12	112	2367	0.255	ug/l	94
71) Bromoform	10.84	173	4757	1.586	ug/l #	96

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

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