

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX103120\
 Data File : VX019194.D
 Acq On : 29 Oct 2020 17:40
 Operator : JC/MD
 Sample : L4470-02RE
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-TT-MW179D-GW-20201015RE

Quant Time: Oct 30 04:17:28 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102720W.M
 Quant Title : SW846 8260
 QLast Update : Tue Oct 27 13:55:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	294054	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	510298	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	483345	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	250213	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	197741	52.14	ug/l	0.00
Spiked Amount			Recovery	=	104.28%	
35) Dibromofluoromethane	5.46	113	157191	49.04	ug/l	0.00
Spiked Amount			Recovery	=	98.08%	
50) Toluene-d8	8.70	98	624565	51.69	ug/l	0.00
Spiked Amount			Recovery	=	103.38%	
62) 4-Bromofluorobenzene	11.12	95	244420	53.41	ug/l	0.00
Spiked Amount			Recovery	=	106.82%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Chloromethane	1.31	50	1777	0.649	ug/l	97
12) 1,1-Dichloroethene	2.36	96	1261	0.442	ug/l	81
16) Acetone	2.42	43	3315	2.534	ug/l	97
17) Carbon Disulfide	2.55	76	1186	0.904	ug/l #	95
19) Methyl tert-butyl Ether	3.17	73	8358	0.846	ug/l	95
21) trans-1,2-Dichloroethene	3.14	96	1097	0.334	ug/l	96
27) cis-1,2-Dichloroethene	4.56	96	135608	36.998	ug/l	92
44) Trichloroethene	7.19	130	510947	133.533	ug/l	97
64) Tetrachloroethene	9.32	164	58019	15.755	ug/l	94

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

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