

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX092420\
 Data File : VX018580.D
 Acq On : 24 Sep 2020 18:09
 Operator : JC/SP
 Sample : L4118-16
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW202D1-20200919

Quant Time: Sep 25 02:26:48 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091720W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 17 12:17:36 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	435623	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.84	114	695306	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	675977	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	329167	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	313671	50.90	ug/l	0.00
Spiked Amount			Recovery	=	101.80%	
35) Dibromofluoromethane	5.46	113	244300	50.82	ug/l	0.00
Spiked Amount			Recovery	=	101.64%	
50) Toluene-d8	8.70	98	856369	49.03	ug/l	0.00
Spiked Amount			Recovery	=	98.06%	
62) 4-Bromofluorobenzene	11.12	95	321054	46.21	ug/l	0.00
Spiked Amount			Recovery	=	92.42%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Chloromethane	1.31	50	6217	1.286	ug/l #	82
16) Acetone	2.44	43	12459	5.286	ug/l	89
19) Methyl tert-butyl Ether	3.17	73	88033	5.944	ug/l	96
20) Methylene Chloride	2.84	84	3566	0.686	ug/l #	70
21) trans-1,2-Dichloroethene	3.14	96	3897	0.795	ug/l #	65
22) Diisopropyl ether	3.82	45	31256	2.034	ug/l #	82
24) 1,1-Dichloroethane	3.67	63	29661	3.347	ug/l #	93
27) cis-1,2-Dichloroethene	4.57	96	165902	30.189	ug/l	87
30) Chloroform	5.18	83	13899	1.546	ug/l	89
42) 1,2-Dichloroethane	6.17	62	10566	1.404	ug/l	90
44) Trichloroethene	7.19	130	97656	17.442	ug/l	91
64) Tetrachloroethene	9.32	164	107739	20.245	ug/l	94

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

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