

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX030819\
 Data File : VX007949.D
 Acq On : 08 Mar 2019 23:35
 Operator : JC/SP
 Sample : K1793-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RE131D1-20190305

Quant Time: Mar 11 12:07:30 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X030619W.M
 Quant Title : SW846 8260
 QLast Update : Thu Mar 07 07:33:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	269054	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	421076	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	399392	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	193933	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	6.07	65	146239	51.52	ug/l	0.00
Spiked Amount			Recovery	=	103.04%	
35) Dibromofluoromethane	5.51	113	135763	50.40	ug/l	0.00
Spiked Amount			Recovery	=	100.80%	
50) Toluene-d8	8.71	98	511289	51.69	ug/l	0.00
Spiked Amount			Recovery	=	103.38%	
62) 4-Bromofluorobenzene	11.13	95	190095	54.81	ug/l	0.00
Spiked Amount			Recovery	=	109.62%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
9) 1,1,2-Trichlorotrifluoroet	2.39	101	8018	3.376	ug/l	96
12) 1,1-Dichloroethene	2.39	96	1599	0.699	ug/l	97
16) Acetone	2.45	43	3039	1.850	ug/l	96
18) Methyl Acetate	2.78	43	7532	2.158	ug/l	100
19) Methyl tert-butyl Ether	3.21	73	4252	0.540	ug/l	93
24) 1,1-Dichloroethane	3.71	63	2224	0.469	ug/l #	85
27) cis-1,2-Dichloroethene	4.60	96	12910	4.479	ug/l	92
30) Chloroform	5.22	83	8012	1.785	ug/l	100
44) Trichloroethene	7.21	130	504858	161.396	ug/l	98
49) 1,4-Dioxane	7.75	88	1817	25.124	ug/l	98
64) Tetrachloroethene	9.34	164	37286	11.266	ug/l	97

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

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