

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX030819\
 Data File : VX007952.D
 Acq On : 09 Mar 2019 00:55
 Operator : JC/SP
 Sample : K1793-04
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 DUP01-20190305

Quant Time: Mar 11 12:10:06 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	268763	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	419069	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	395107	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	191832	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	148880	52.51	ug/l	0.00
Spiked Amount						
			Recovery	=	105.02%	
35) Dibromofluoromethane	5.51	113	136861	51.06	ug/l	0.00
Spiked Amount						
			Recovery	=	102.12%	
50) Toluene-d8	8.71	98	506986	51.50	ug/l	0.00
Spiked Amount						
			Recovery	=	103.00%	
62) 4-Bromofluorobenzene	11.13	95	186167	53.94	ug/l	0.00
Spiked Amount						
			Recovery	=	107.88%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
9) 1,1,2-Trichlorotrifluoroet	2.39	101	9301	3.921	ug/l	98
12) 1,1-Dichloroethene	2.38	96	1723	0.753	ug/l	96
16) Acetone	2.45	43	2518	1.534	ug/l #	86
18) Methyl Acetate	2.78	43	3039	0.872	ug/l	99
19) Methyl tert-butyl Ether	3.21	73	4216	0.536	ug/l #	90
24) 1,1-Dichloroethane	3.70	63	2072	0.437	ug/l #	94
27) cis-1,2-Dichloroethene	4.60	96	13105	4.552	ug/l	88
30) Chloroform	5.22	83	7994	1.783	ug/l	98
44) Trichloroethene	7.21	130	504168	161.948	ug/l	99
49) 1,4-Dioxane	7.76	88	1540	21.396	ug/l	94
64) Tetrachloroethene	9.34	164	37147	11.345	ug/l	99

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

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