

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX121119\
 Data File : VX013964.D
 Acq On : 12 Dec 2019 00:30
 Operator : JC/SP
 Sample : K6249-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SA-PZ190I-20191208

Quant Time: Dec 12 06:42:44 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 26 15:53:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.65	168	534500	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	822950	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	720168	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	330437	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	6.05	65	302485	48.75	ug/l	0.00
Spiked Amount			Recovery	=	97.50%	
35) Dibromofluoromethane	5.49	113	243739	47.91	ug/l	0.00
Spiked Amount			Recovery	=	95.82%	
50) Toluene-d8	8.71	98	973507	49.11	ug/l	0.00
Spiked Amount			Recovery	=	98.22%	
62) 4-Bromofluorobenzene	11.14	95	330740	46.27	ug/l	0.00
Spiked Amount			Recovery	=	92.54%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
5) Bromomethane	1.63	94	3114	0.617	ug/l	98
6) Chloroethane	1.70	64	108701	25.424	ug/l	99
12) 1,1-Dichloroethene	2.34	96	40893	8.039	ug/l	94
16) Acetone	2.45	43	8938	2.918	ug/l	99
24) 1,1-Dichloroethane	3.68	63	1741023	179.176	ug/l	99
32) 1,1,1-Trichloroethane	5.48	97	47204	5.732	ug/l	98
40) Benzene	6.13	78	11992	0.527	ug/l	98
65) Chlorobenzene	10.14	112	9929	0.664	ug/l #	86
83) tert-Butylbenzene	11.76	119	9266	0.480	ug/l	86
85) sec-Butylbenzene	11.94	105	29440	1.335	ug/l	82
87) 1,3-Dichlorobenzene	12.03	146	11082	1.015	ug/l	93
88) 1,4-Dichlorobenzene	12.09	146	45451	4.088	ug/l	94
91) 1,2-Dichlorobenzene	12.39	146	16056	1.489	ug/l	96
93) 1,2,4-Trichlorobenzene	13.64	180	3899	0.519	ug/l #	85

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

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