

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX102519\
 Data File : VX013204.D
 Acq On : 24 Oct 2019 12:04
 Operator : JC/SP
 Sample : K5531-02
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-TT-MW203D2-GW-20191021

Quant Time: Oct 26 02:06:46 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102419W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 24 03:56:38 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.65	168	77787	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	113933	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	113266	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	72595	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	83750	51.48	ug/l	0.00
Spiked Amount			Recovery	=	102.96%	
35) Dibromofluoromethane	5.49	113	63542	50.01	ug/l	0.00
Spiked Amount			Recovery	=	100.02%	
50) Toluene-d8	8.71	98	237073	50.35	ug/l	0.00
Spiked Amount			Recovery	=	100.70%	
62) 4-Bromofluorobenzene	11.14	95	89296	47.92	ug/l	0.00
Spiked Amount			Recovery	=	95.84%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Chloromethane	1.32	50	858	1.628	ug/l #	84
16) Acetone	2.44	43	2199	2.869	ug/l	90
21) trans-1,2-Dichloroethene	3.17	96	373	0.744	ug/l #	80
24) 1,1-Dichloroethane	3.69	63	4208	1.677	ug/l	99
27) cis-1,2-Dichloroethene	4.59	96	2621	1.615	ug/l	83
30) Chloroform	5.20	83	965	0.361	ug/l	85
42) 1,2-Dichloroethane	6.18	62	2674	1.239	ug/l	99
44) Trichloroethene	7.21	130	1243	0.765	ug/l	83
52) Toluene	8.78	92	958	0.253	ug/l	97
65) Chlorobenzene	10.13	112	1191	0.300	ug/l #	80
84) 1,2,4-Trimethylbenzene	11.80	105	1387	0.251	ug/l	100
87) 1,3-Dichlorobenzene	12.03	146	1263	0.442	ug/l	93
91) 1,2-Dichlorobenzene	12.39	146	988	0.352	ug/l	91
95) Naphthalene	13.83	128	3722	0.686	ug/l	96

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

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