

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091420\
 Data File : VX018385.D
 Acq On : 14 Sep 2020 15:50
 Operator : JC/SP
 Sample : L3952-02
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 40614

Quant Time: Sep 16 05:07:31 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X082120W.M
 Quant Title : SW846 8260
 QLast Update : Fri Aug 21 03:03:56 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.64	168	441899	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	701454	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	665484	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	322802	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	6.04	65	297824	48.55	ug/l	0.00
Spiked Amount			Recovery	=	97.10%	
35) Dibromofluoromethane	5.46	113	239040	50.04	ug/l	0.00
Spiked Amount			Recovery	=	100.08%	
50) Toluene-d8	8.70	98	861246	48.40	ug/l	0.00
Spiked Amount			Recovery	=	96.80%	
62) 4-Bromofluorobenzene	11.12	95	307910	43.98	ug/l	0.00
Spiked Amount			Recovery	=	87.96%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	2.42	43	2448736	1024.383	ug/l	96
25) 2-Butanone	4.64	43	49103	12.499	ug/l #	89
40) Benzene	6.12	78	39998	2.104	ug/l	97
51) 4-Methyl-2-Pentanone	8.62	43	14010	1.869	ug/l	98
52) Toluene	8.77	92	68973	5.722	ug/l	98
59) 2-Hexanone	9.47	43	2935	0.515	ug/l	88
67) Ethyl Benzene	10.24	91	11090	0.491	ug/l	98
68) m/p-Xylenes	10.34	106	21952	2.602	ug/l	99
69) o-Xylene	10.68	106	11187	1.368	ug/l	90
84) 1,2,4-Trimethylbenzene	11.79	105	11233	0.628	ug/l	74
86) p-Isopropyltoluene	12.05	119	93842	5.203	ug/l	95
89) n-Butylbenzene	12.37	91	5227	0.318	ug/l	89
95) Naphthalene	13.82	128	11559	0.541	ug/l #	87

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

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