

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031621\
 Data File : VX021257.D
 Acq On : 16 Mar 2021 20:17
 Operator : JC/MD
 Sample : M1694-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 JG-W

Quant Time: Mar 17 02:44:02 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031621W.M
 Quant Title : SW846 8260
 QLast Update : Tue Mar 16 14:00:21 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.629	168	86244	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.829	114	137512	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.094	117	119698	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.059	152	57163	50.00	ug/l	# 0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.029	65	63745	45.06	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	90.12%		
35) Dibromofluoromethane	5.470	113	45231	48.34	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	96.68%		
50) Toluene-d8	8.694	98	168716	49.69	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	99.38%		
62) 4-Bromofluorobenzene	11.118	95	62857	47.67	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	95.34%		
Target Compounds						Qvalue
16) Acetone	2.423	43	27067	52.00	ug/l	89
25) 2-Butanone	4.635	43	43588	54.46	ug/l	91
30) Chloroform	5.176	83	112322	57.39	ug/l	99
47) Bromodichloromethane	7.876	83	6985	4.64	ug/l	93
52) Toluene	8.765	92	15077	6.39	ug/l	97
67) Ethyl Benzene	10.235	91	65649	14.43	ug/l	99
68) m/p-Xylenes	10.341	106	100059	61.54	ug/l	92
69) o-Xylene	10.682	106	49047	31.04	ug/l	93
86) p-Isopropyltoluene	12.047	119	1560287	422.36	ug/l	97

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

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