

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX021419\
 Data File : VX007522.D
 Acq On : 13 Feb 2019 16:04
 Operator : JC/SP
 Sample : K1460-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WC-20190208

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.605	66	74	79	rBV5	24226	57959	2.34%	0.364%
2	2.446	206	212	219	rBV	43952	75198	3.04%	0.473%
3	3.031	301	308	313	rBV2	11557	26857	1.08%	0.169%
4	5.220	653	667	676	rBV	49722	151167	6.10%	0.951%
5	5.506	702	714	729	rBV	309029	884245	35.69%	5.560%
6	5.665	729	740	753	rBV	509705	1417904	57.23%	8.916%
7	6.067	797	806	819	rBV	306014	795516	32.11%	5.002%
8	6.860	926	936	950	rBV	774598	1790720	72.28%	11.260%
9	7.896	1099	1106	1115	rVB2	24467	50001	2.02%	0.314%
10	8.713	1232	1240	1247	rBV	1609594	2477609	100.00%	15.580%
11	8.780	1247	1251	1257	rVB	86453	137076	5.53%	0.862%
12	9.335	1335	1342	1348	rBV	39272	63243	2.55%	0.398%
13	9.585	1375	1383	1389	rBV2	19741	33424	1.35%	0.210%
14	10.109	1463	1469	1477	rBV	1612221	2192255	88.48%	13.785%
15	10.353	1505	1509	1515	rVB	22162	33940	1.37%	0.213%
16	10.823	1582	1586	1597	rBV	81431	112302	4.53%	0.706%
17	11.140	1631	1638	1644	rBV	1329357	1766030	71.28%	11.105%
18	11.518	1695	1700	1706	rVB8	13613	24857	1.00%	0.156%
19	12.078	1786	1792	1799	rBV	1760283	2166321	87.44%	13.622%
20	12.158	1799	1805	1810	rVB9	15195	26247	1.06%	0.165%
21	12.267	1819	1823	1831	rBV2	28305	48639	1.96%	0.306%
22	12.908	1923	1928	1933	rBV3	21171	29762	1.20%	0.187%
23	13.151	1965	1968	1972	rVB2	33633	43038	1.74%	0.271%
24	13.474	2017	2021	2028	rBV	143311	193820	7.82%	1.219%
25	13.535	2028	2031	2035	rBV2	31963	46128	1.86%	0.290%
26	13.645	2044	2049	2059	rBV2	294270	418736	16.90%	2.633%
27	13.804	2072	2075	2078	rVB	30458	36243	1.46%	0.228%
28	13.913	2089	2093	2095	rBV	43918	54931	2.22%	0.345%
29	13.944	2095	2098	2108	rVB2	82519	126540	5.11%	0.796%
30	14.249	2144	2148	2162	rBV	253070	391789	15.81%	2.464%
31	14.505	2185	2190	2193	rBV4	26121	41445	1.67%	0.261%
32	14.541	2193	2196	2205	rVB2	27182	42737	1.72%	0.269%
33	14.651	2211	2214	2216	rBV2	23489	27235	1.10%	0.171%
34	14.682	2216	2219	2226	rVB3	38974	59243	2.39%	0.373%

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Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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35	14.974	2262	2267	2273	rBV	37723	59708	2.41%	0.375%
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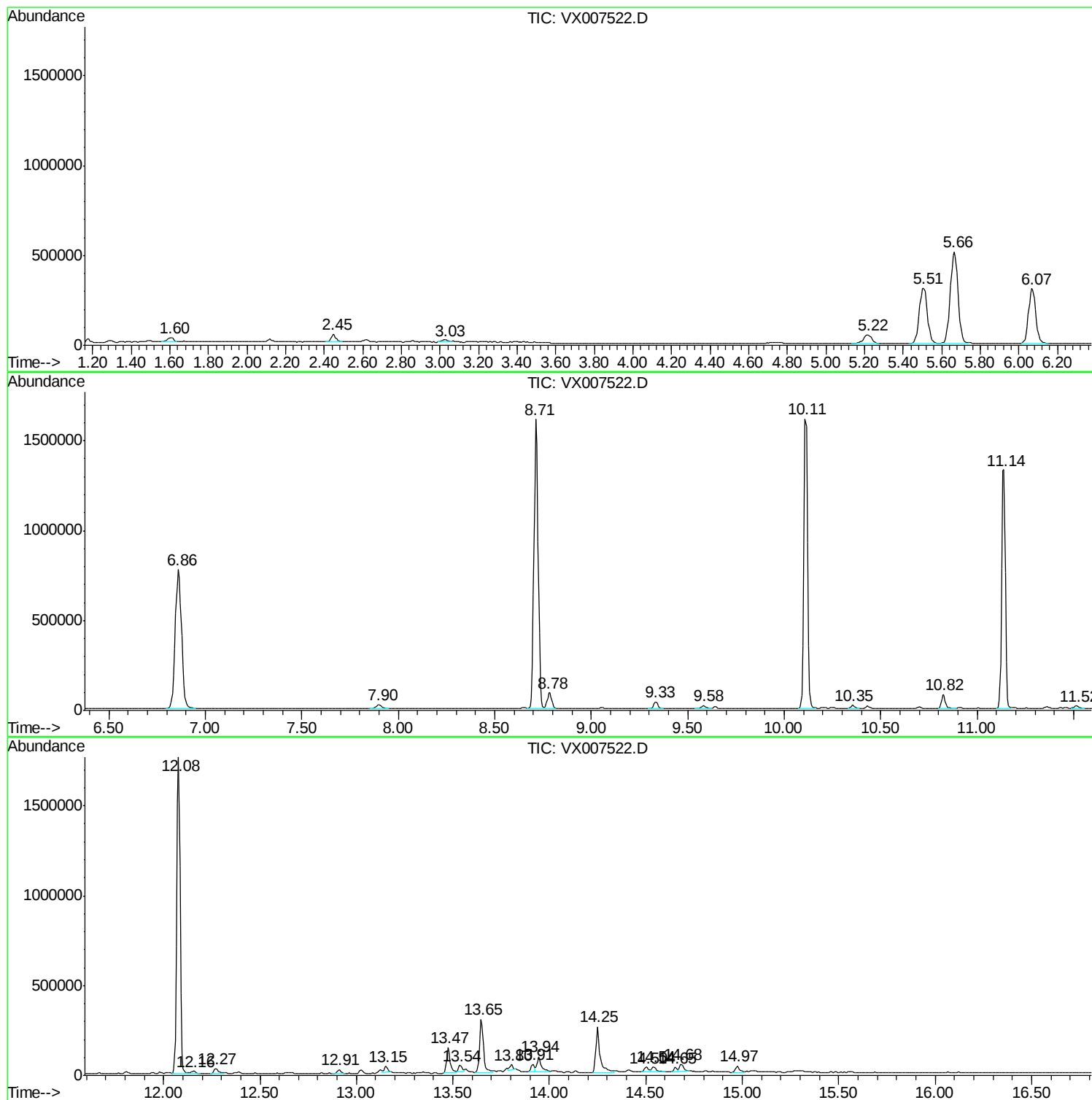
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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Sample : K1460-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

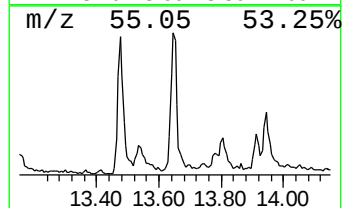
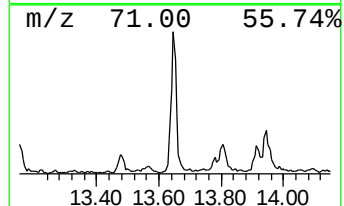
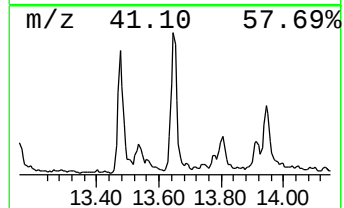
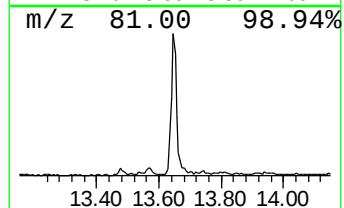
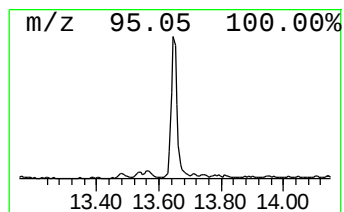
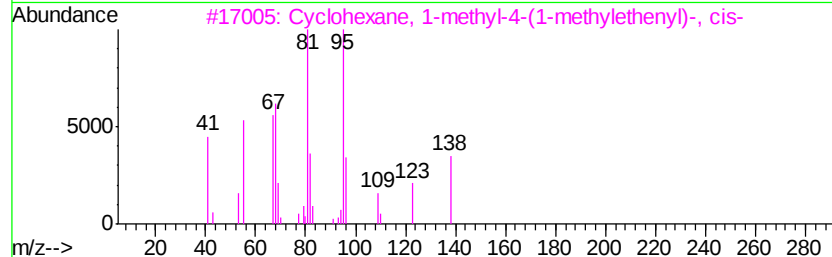
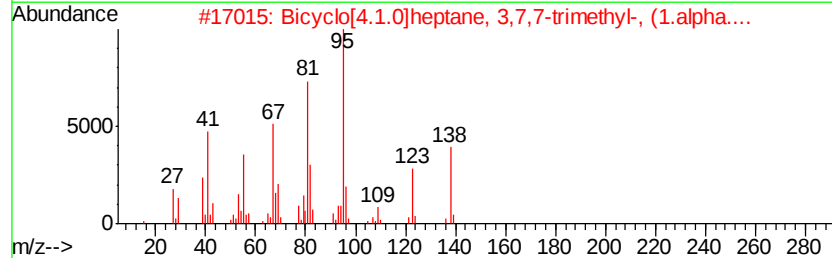
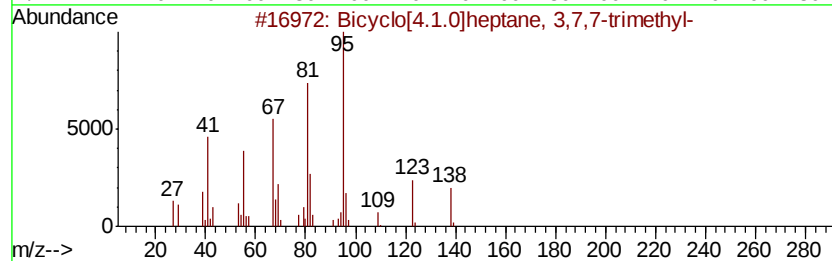
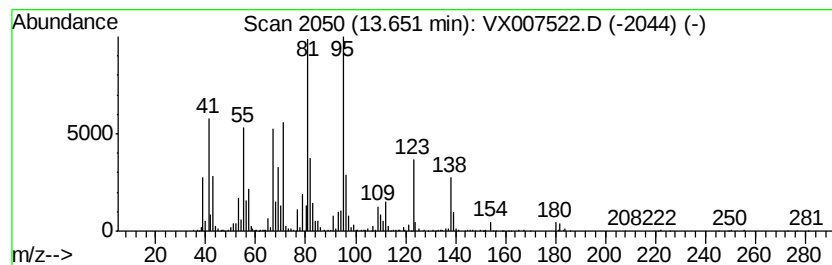
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Peak Number 1 Bicyclo[4.1.0]heptane, 3,7,... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.65	9.66 ug/l	418736	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Bicyclo[4.1.0]heptane, 3,7,7-tri...	138 C10H18	000554-59-6 94
2		Bicyclo[4.1.0]heptane, 3,7,7-tri...	138 C10H18	018968-23-5 94
3		Cyclohexane, 1-methyl-4-(1-methyl...	138 C10H18	001879-07-8 90
4		Cyclohexene, 4-methyl-1-(1-methyl...	138 C10H18	000500-00-5 68
5		L-Menthyl chloroformate	218 C11H19ClO2	014602-86-9 68



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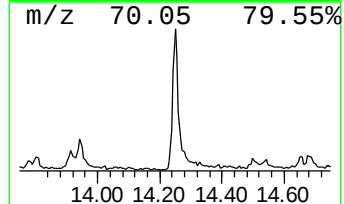
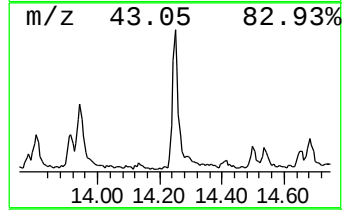
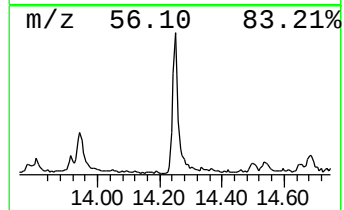
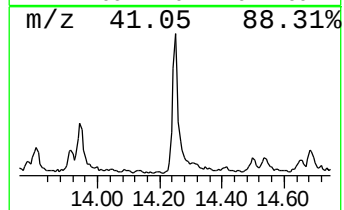
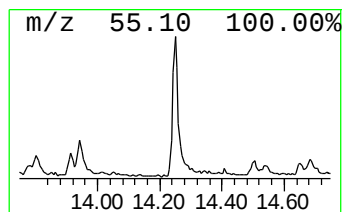
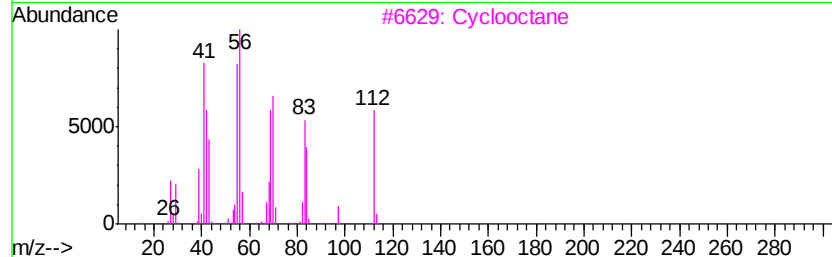
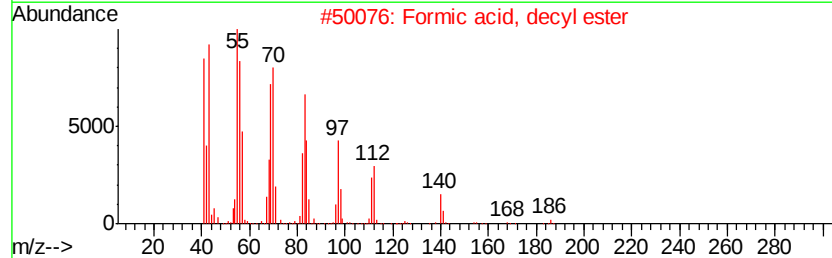
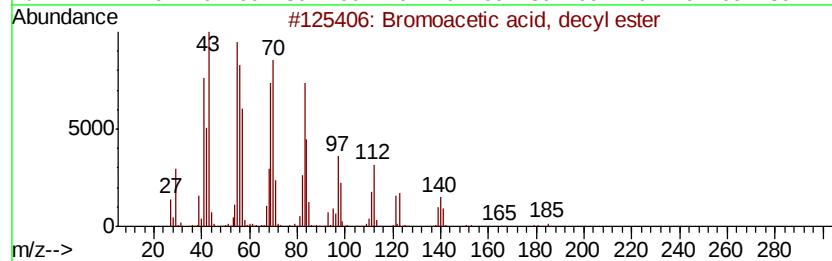
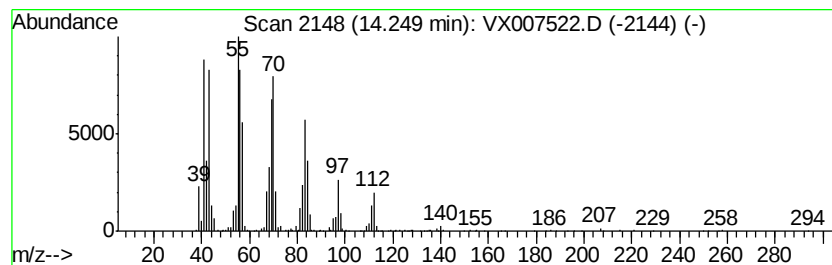
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Peak Number 2 Bromoacetic acid, decyl ester Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.25	9.04 ug/l	391789	1,4-Dichlorobenzene-d4		12.08	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Bromoacetic acid, decyl ester	278	C12H23BrO2	005436-93-1	87
2		Formic acid, decyl ester	186	C11H22O2	005451-52-5	86
3		Cvclooctane	112	C8H16	000292-64-8	81
4		1-Decanol	158	C10H22O	000112-30-1	80
5		2-Dodecene, (Z)-	168	C12H24	007206-26-0	72



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Bicyclo[4.1.0]hep...	13.65	9.7	ug/l	418736	4	12.08	2166320	50.0
Bromoacetic acid,...	14.25	9.0	ug/l	391789	4	12.08	2166320	50.0