

Data Path : Z:\HPCHEM1\BNA M\DATA\BM042816\
 Data File : BM005121.D
 Acq On : 28 Apr 2016 19:39
 Operator : UM/SJ
 Sample : H2639-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BC7N7

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 1 % 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:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM042516.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.822	693	703	709	rBV	26682	60150	6.33%	0.531%
2	6.957	721	726	735	rBB	58881	90925	9.57%	0.802%
3	7.122	748	754	762	rBB	110372	166775	17.56%	1.472%
4	7.322	783	788	795	rBV	128512	199558	21.01%	1.761%
5	7.787	860	867	874	rBV	143033	229676	24.18%	2.027%
6	7.845	874	877	882	rVB	10811	14506	1.53%	0.128%
7	8.357	960	964	969	rVB	6830	9894	1.04%	0.087%
8	8.492	981	987	998	rVB	146814	235151	24.76%	2.075%
9	8.945	1058	1064	1073	rBV	136897	224782	23.67%	1.984%
10	9.098	1086	1090	1093	rBV	7181	9521	1.00%	0.084%
11	9.139	1093	1097	1102	rVB	17816	25247	2.66%	0.223%
12	9.257	1111	1117	1123	rBB2	5562	11998	1.26%	0.106%
13	9.322	1123	1128	1135	rBB2	9093	14862	1.56%	0.131%
14	9.663	1180	1186	1193	rBV	120269	192792	20.30%	1.702%
15	9.904	1221	1227	1233	rBB	18156	29074	3.06%	0.257%
16	10.098	1256	1260	1265	rBB	6665	10053	1.06%	0.089%
17	10.198	1271	1277	1288	rBB	193618	319205	33.61%	2.817%
18	10.575	1335	1341	1348	rBV	222676	382375	40.26%	3.375%
19	10.745	1364	1370	1378	rBB	163281	273482	28.80%	2.414%
20	11.386	1472	1479	1489	rBB	47198	84460	8.89%	0.745%
21	11.686	1525	1530	1539	rBB	25791	43762	4.61%	0.386%
22	12.139	1601	1607	1609	rBV	6842	9583	1.01%	0.085%
23	12.357	1639	1644	1651	rBV	14135	22773	2.40%	0.201%
24	12.698	1695	1702	1706	rBV2	8089	13238	1.39%	0.117%
25	12.933	1737	1742	1749	rVB	12721	18393	1.94%	0.162%
26	13.416	1819	1824	1834	rVB2	6675	12210	1.29%	0.108%
27	13.839	1890	1896	1904	rVB	373719	536672	56.51%	4.737%
28	14.021	1922	1927	1932	rBV	9122	13149	1.38%	0.116%
29	14.121	1938	1944	1953	rVB	429354	621519	65.45%	5.485%
30	14.274	1961	1970	1982	rBV2	7703	24076	2.54%	0.212%
31	14.427	1990	1996	2003	rBV2	367785	564617	59.45%	4.983%
32	14.492	2003	2007	2012	rVB	7948	10970	1.16%	0.097%
33	14.639	2027	2032	2042	rBV	57968	96657	10.18%	0.853%
34	14.851	2058	2068	2075	rBV4	50953	87802	9.25%	0.775%

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35	15.221	2125	2131	2134	rBV	8707	12796	1.35%	0.113%
36	15.351	2148	2153	2157	rBV	17180	25113	2.64%	0.222%
37	15.421	2157	2165	2171	rVV	537534	802535	84.51%	7.083%
38	15.480	2171	2175	2181	rVB	13280	18728	1.97%	0.165%
39	15.545	2181	2186	2196	rBV	198666	282795	29.78%	2.496%
40	16.157	2280	2290	2295	rBV	40616	74294	7.82%	0.656%
41	16.345	2314	2322	2327	rBV2	19211	28266	2.98%	0.249%
42	16.457	2336	2341	2346	rBV	10342	13804	1.45%	0.122%
43	16.704	2375	2383	2390	rBV	78362	137399	14.47%	1.213%
44	16.957	2420	2426	2429	rBV4	7208	12698	1.34%	0.112%
45	17.015	2429	2436	2440	rVV	62770	102930	10.84%	0.908%
46	17.068	2440	2445	2452	rVB	29603	45581	4.80%	0.402%
47	17.174	2458	2463	2468	rBV2	468693	684115	72.04%	6.038%
48	17.215	2468	2470	2474	rVV	20769	26195	2.76%	0.231%
49	17.274	2474	2480	2485	rVV2	625716	949679	100.00%	8.382%
50	17.315	2485	2487	2493	rVV2	54860	66394	6.99%	0.586%
51	17.386	2493	2499	2503	rVV2	9484	20948	2.21%	0.185%
52	17.433	2503	2507	2515	rVV2	30350	49737	5.24%	0.439%
53	17.521	2518	2522	2526	rVV	9402	14368	1.51%	0.127%
54	17.580	2528	2532	2540	rVB	10165	15367	1.62%	0.136%
55	18.068	2607	2615	2622	rBV4	29273	46965	4.95%	0.415%
56	18.139	2623	2627	2632	rVB	9259	10995	1.16%	0.097%
57	18.598	2701	2705	2708	rBV	9853	15295	1.61%	0.135%
58	18.956	2760	2766	2774	rBV3	13689	20523	2.16%	0.181%
59	19.203	2804	2808	2810	rBV	10817	12574	1.32%	0.111%
60	19.233	2810	2813	2827	rVB2	10048	20998	2.21%	0.185%
61	19.351	2827	2833	2844	rBV2	32221	55267	5.82%	0.488%
62	19.486	2849	2856	2865	rVV2	41793	57170	6.02%	0.505%
63	19.574	2865	2871	2882	rVB2	660346	927958	97.71%	8.190%
64	19.803	2907	2910	2921	rVB	9481	12019	1.27%	0.106%
65	20.297	2991	2994	3001	rBV6	7345	17305	1.82%	0.153%
66	21.274	3157	3160	3164	rBV	10574	11764	1.24%	0.104%
67	21.368	3171	3176	3182	rBV	548593	674869	71.06%	5.956%
68	22.421	3351	3355	3360	rVB3	15136	19893	2.09%	0.176%
69	22.933	3438	3442	3448	rVB	20179	30349	3.20%	0.268%
70	23.503	3532	3539	3549	rBV2	355281	669582	70.51%	5.910%
71	23.650	3557	3564	3575	rVB2	283117	543135	57.19%	4.794%

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72	24.162	3647	3651	3658	rVB	26050	48648	5.12%	0.429%
73	24.927	3775	3781	3788	rVB	25570	51987	5.47%	0.459%
74	25.809	3925	3931	3938	rVB2	17525	39354	4.14%	0.347%

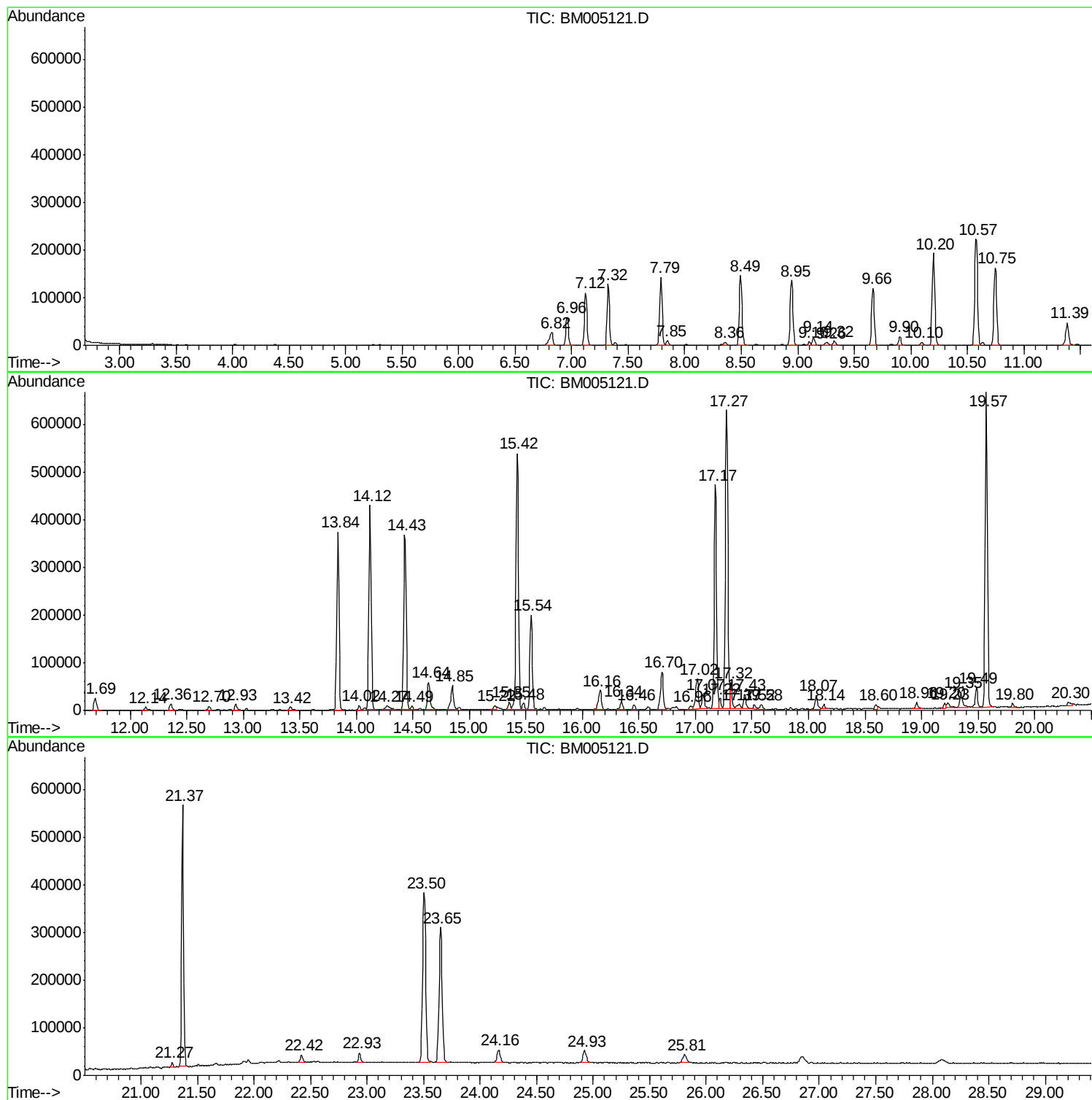
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Instrument :
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM042516.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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Instrument :
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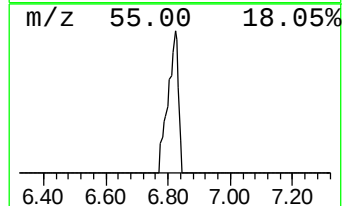
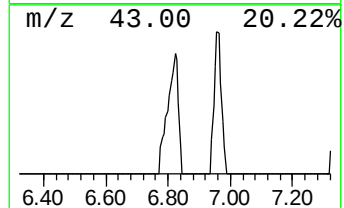
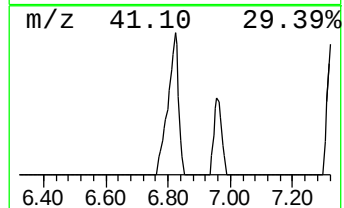
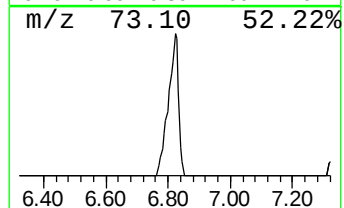
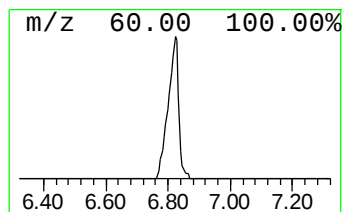
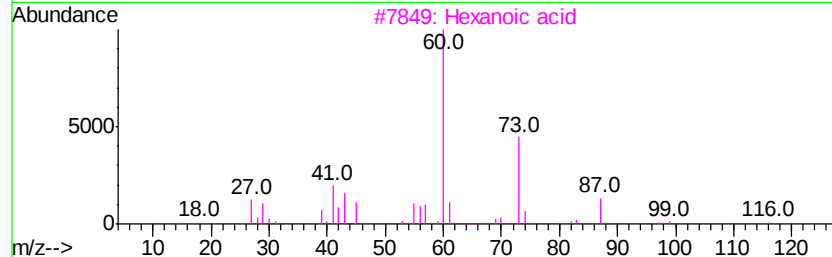
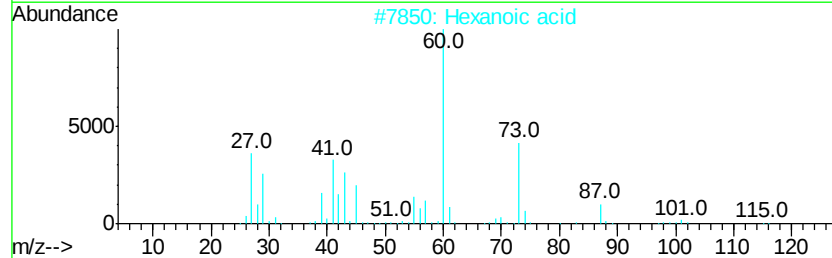
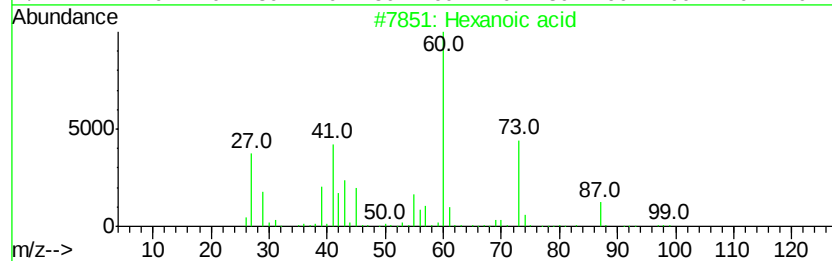
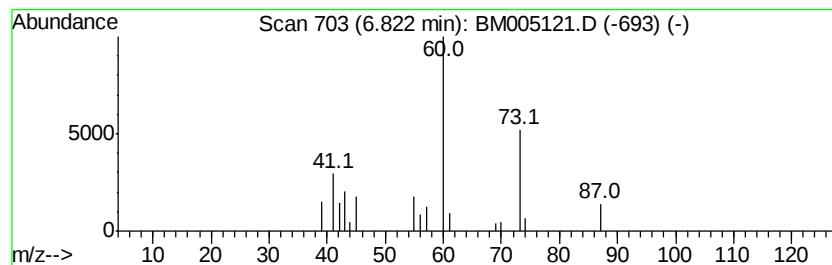
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Hexanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.82	5.24 ng/ul	60150	1,4-Dichlorobenzene-d4	7.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanoic acid	116	C6H12O2	000142-62-1	83
2		Hexanoic acid	116	C6H12O2	000142-62-1	78
3		Hexanoic acid	116	C6H12O2	000142-62-1	78
4		Pentanoic acid	102	C5H10O2	000109-52-4	72
5		Propanedioic acid, propyl-	146	C6H10O4	000616-62-6	56



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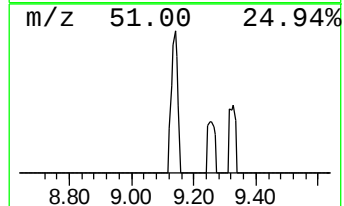
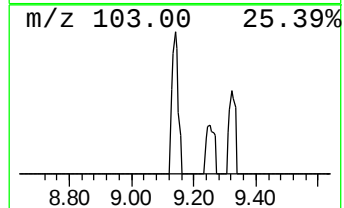
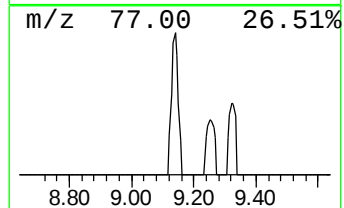
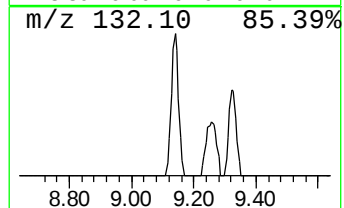
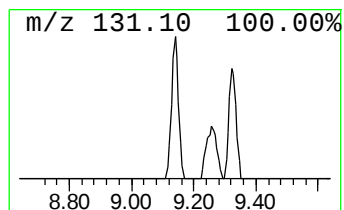
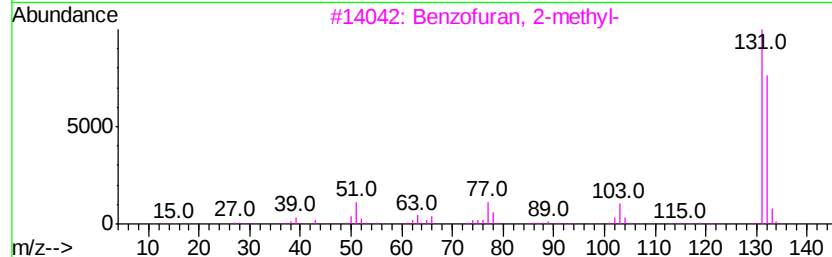
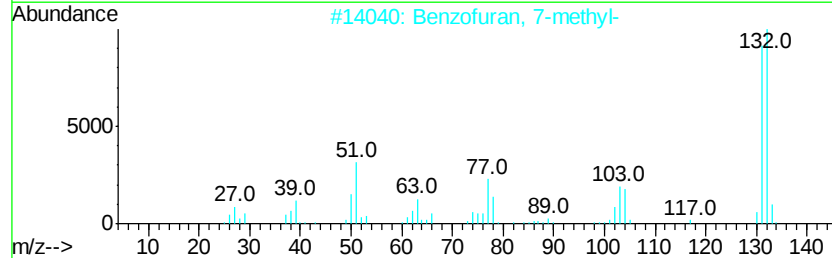
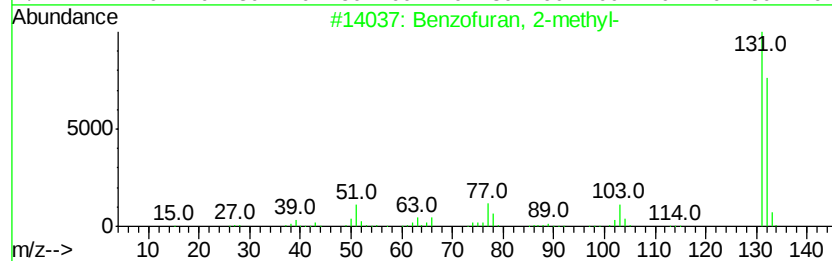
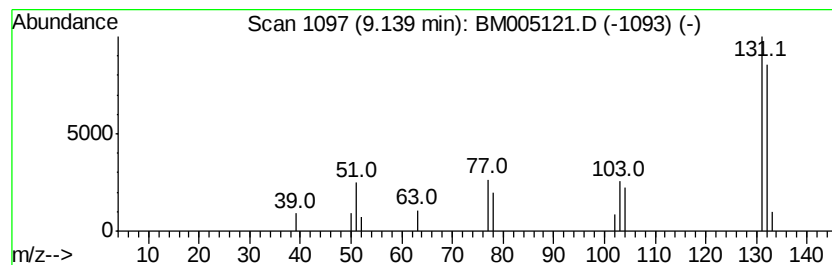
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzofuran, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.14	2.20 ng/ul	25247	1,4-Dichlorobenzene-d4	7.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
2		Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	91
3		Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
4		Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	90
5		3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	90



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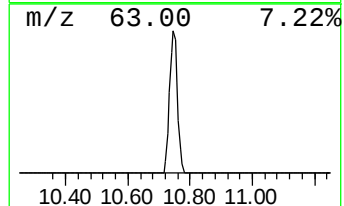
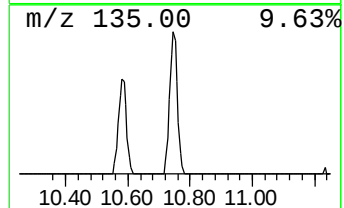
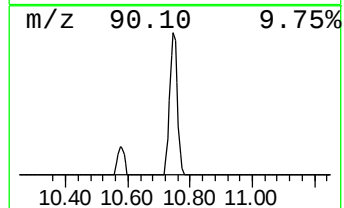
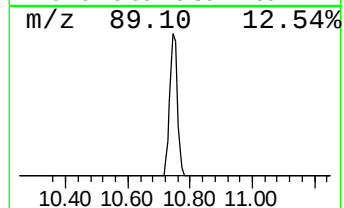
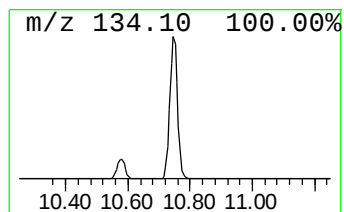
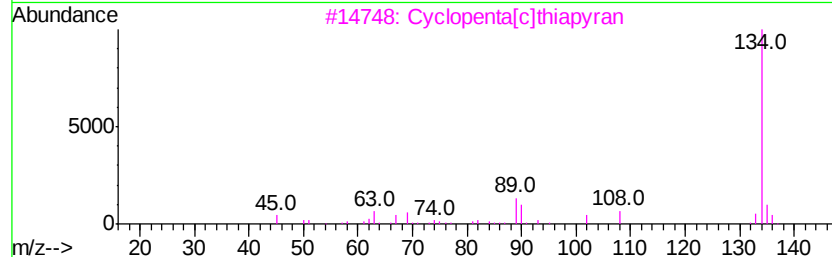
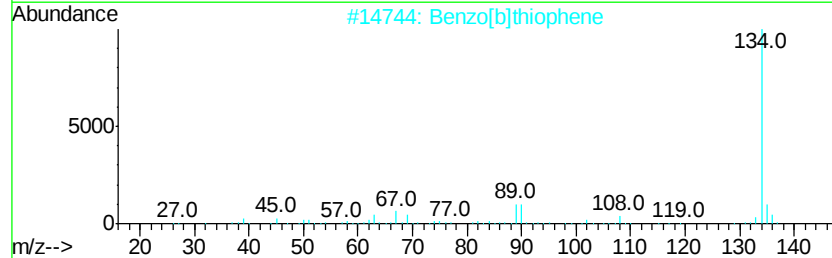
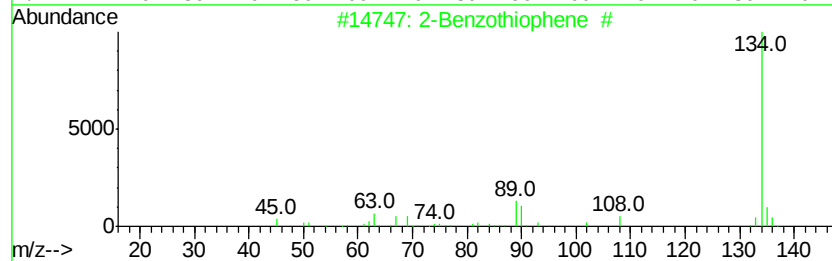
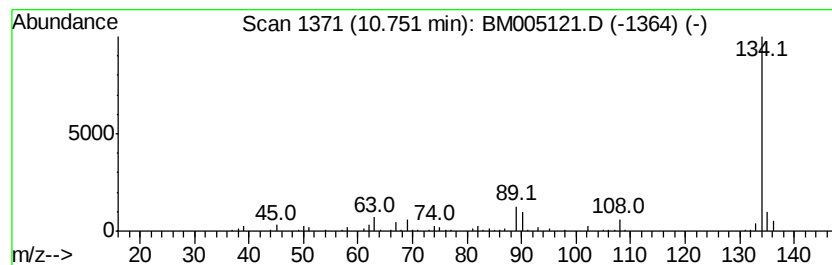
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Peak Number 3 2-Benzothiophene # Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.75	14.30 ng/ul	273482	Napthalene-d8	10.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Benzothiophene #	134	C8H6S	000270-82-6	97
2		Benzo[b]thiophene	134	C8H6S	000095-15-8	95
3		Cyclopenta[c]thiopyran	134	C8H6S	000270-63-3	94
4		Benzo[b]thiophene	134	C8H6S	000095-15-8	94
5		Benzo[b]thiophene	134	C8H6S	000095-15-8	90



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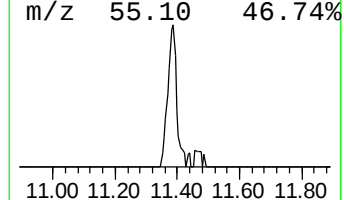
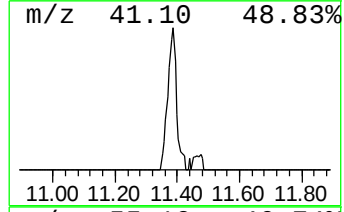
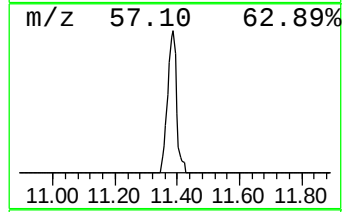
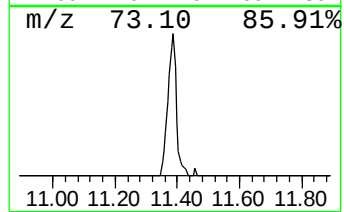
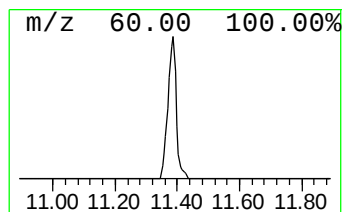
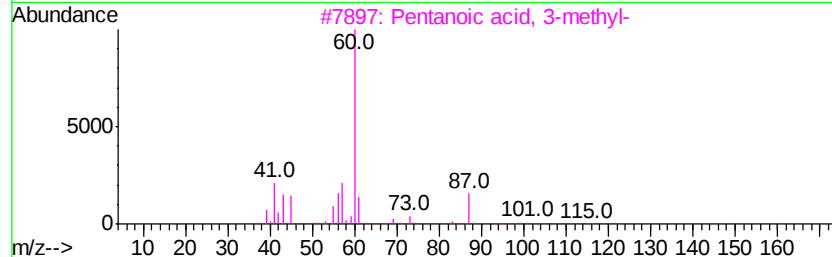
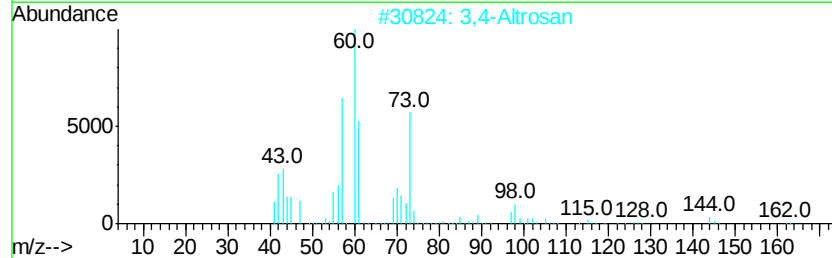
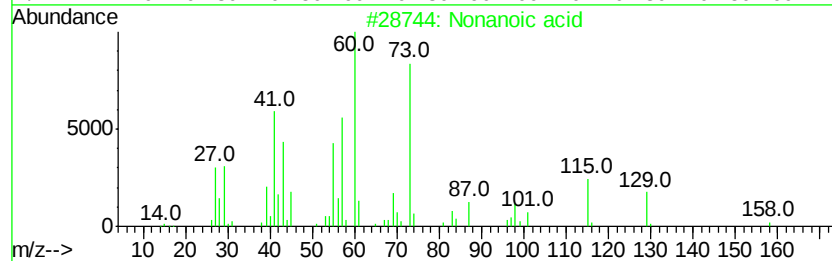
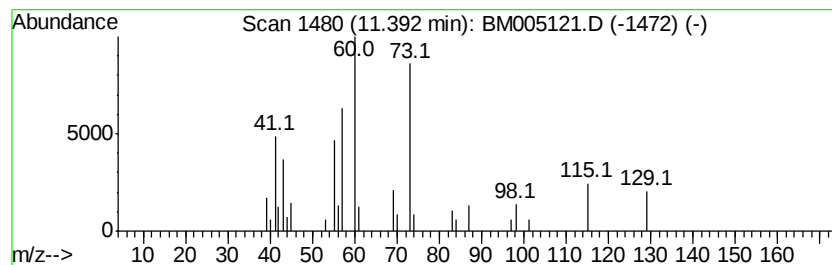
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TIC Integration Parameters: LSCINT.P

Peak Number 4 Nonanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.39	4.42 ng/ul	84460	Napthalene-d8	10.57

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonanoic acid		158	C9H18O2	000112-05-0	91
2	3,4-Altrosan		162	C6H10O5	1000129-76-4	47
3	Pentanoic acid, 3-methyl-		116	C6H12O2	000105-43-1	43
4	Octanoic Acid		144	C8H16O2	000124-07-2	42
5	Hexanoic acid		116	C6H12O2	000142-62-1	40



Data Path : Z:\HPCHEM1\BNA M\DATA\BM042816\
Data File : BM005121.D
Acq On : 28 Apr 2016 19:39
Operator : UM/SJ
Sample : H2639-04
Misc :
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ClientSampleId :
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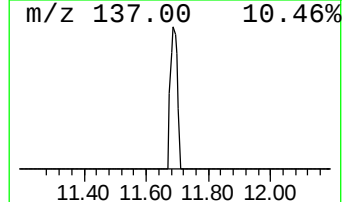
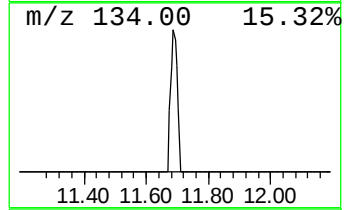
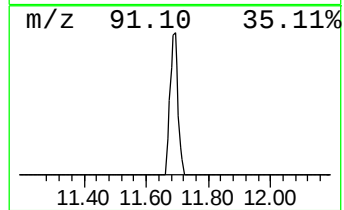
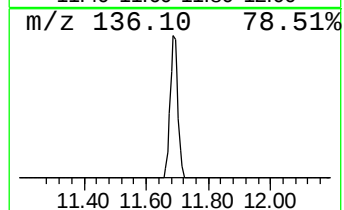
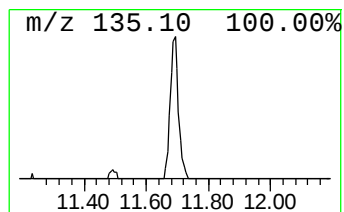
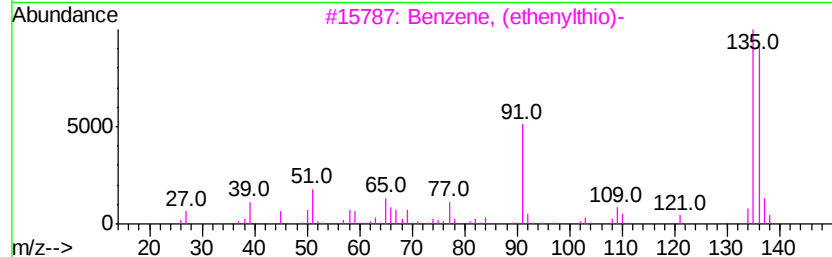
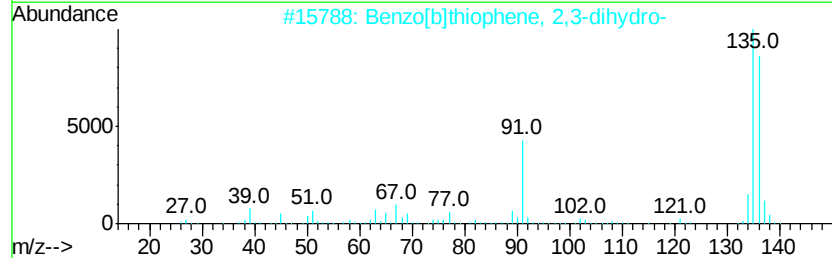
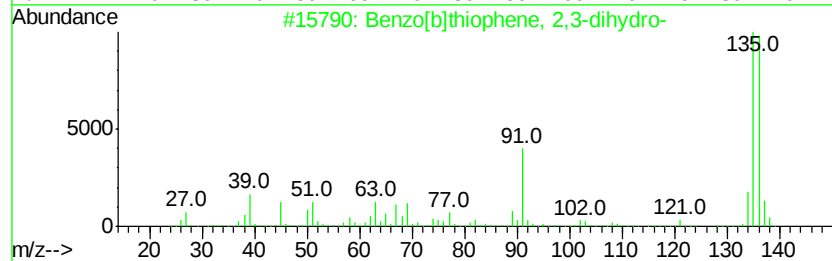
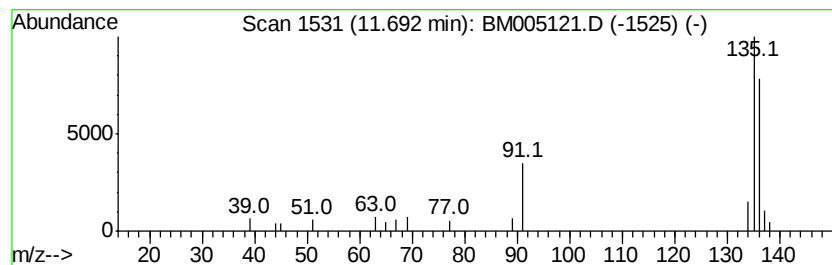
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzo[b]thiophene, 2,3-dihy... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.69	2.29 ng/ul	43762	Napthalene-d8	10.57

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	91
2		Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	86
3		Benzene, (ethenylthio)-	136	C8H8S	001822-73-7	78
4		Benzene, (ethenylthio)-	136	C8H8S	001822-73-7	78
5		Benzo[c]thiophene, 1,3-dihydro-	136	C8H8S	002471-92-3	72



Data Path : Z:\HPCHEM1\BNA M\DATA\BM042816\
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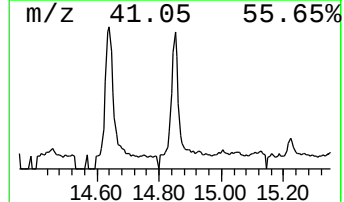
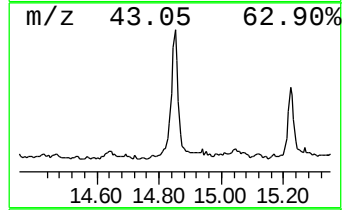
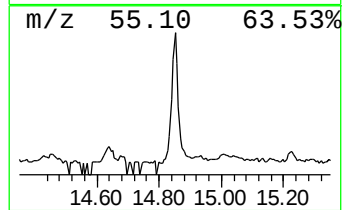
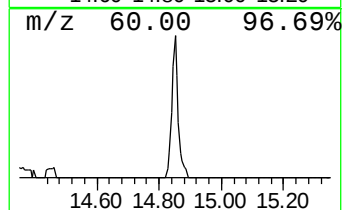
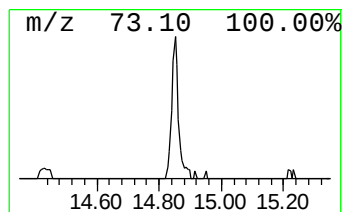
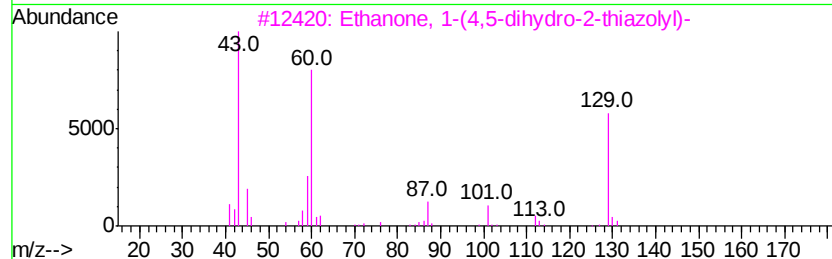
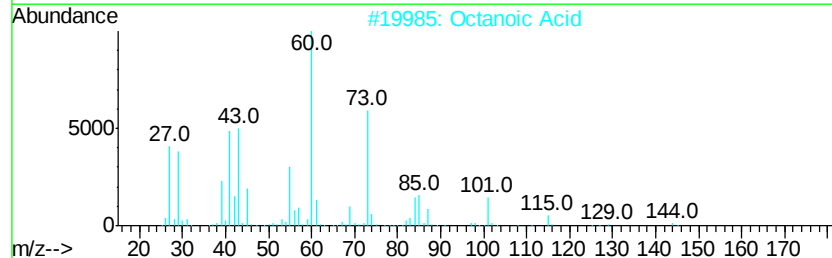
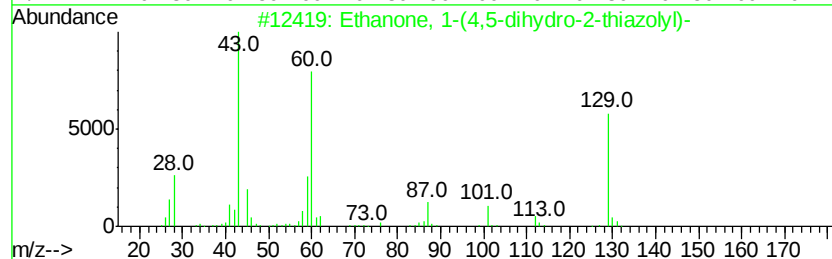
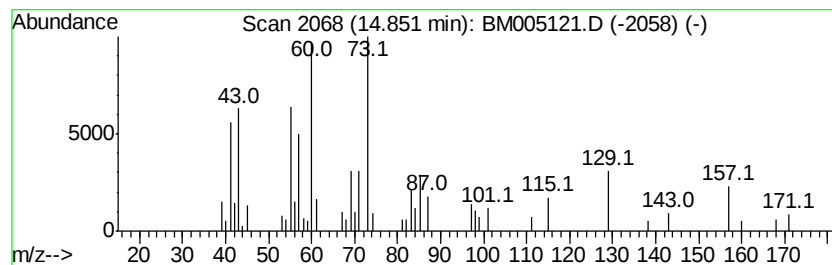
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	3.11 ng/ul	87802	Acenaphthene-d10	14.43

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Ethanone, 1-(4,5-dihydro-2-thiaz...	129	C5H7NOS	029926-41-8	43
2		Octanoic Acid	144	C8H16O2	000124-07-2	43
3		Ethanone, 1-(4,5-dihydro-2-thiaz...	129	C5H7NOS	029926-41-8	43
4		Nonanoic acid	158	C9H18O2	000112-05-0	42
5		D-Glucose, 4-O-.alpha.-D-glucopy...	342	C12H22O11	000069-79-4	40



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Misc :
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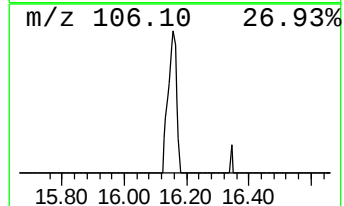
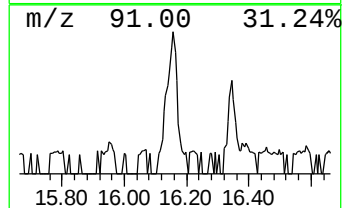
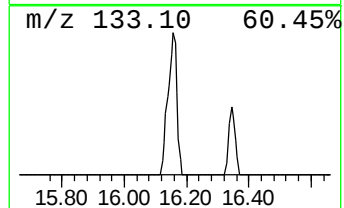
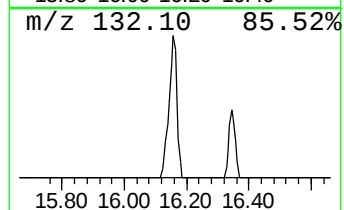
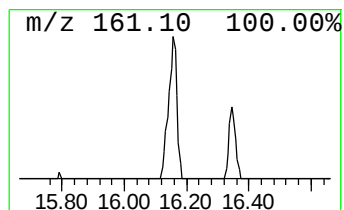
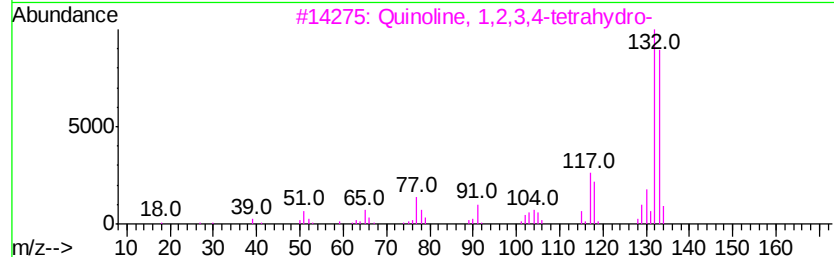
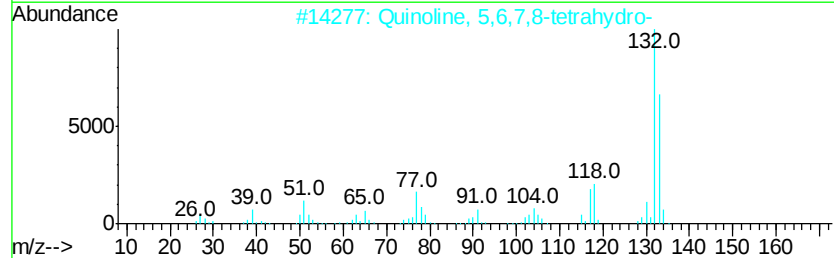
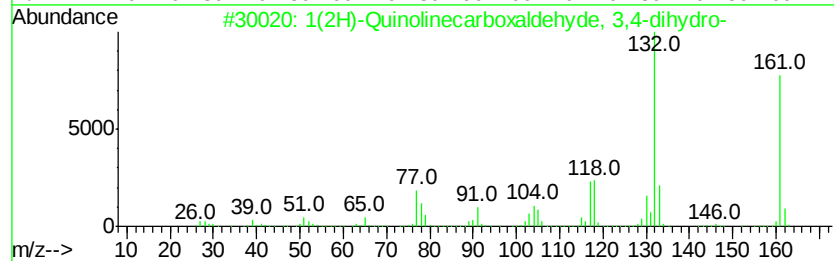
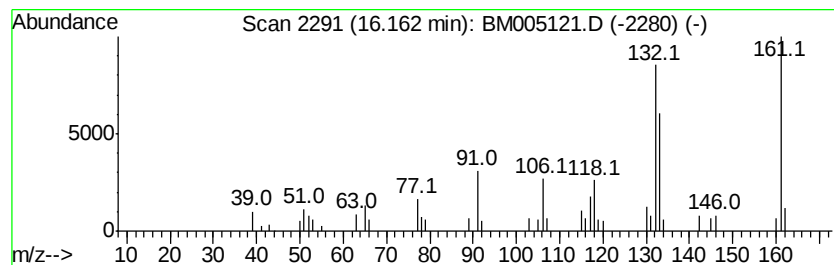
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 1(2H)-Quinolinecarboxaldehy... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.16	2.17 ng/ul	74294	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1(2H)-Quinolinecarboxaldehyde, 3...	161	C10H11NO	002739-16-4	68
2		Quinoline, 5,6,7,8-tetrahydro-	133	C9H11N	010500-57-9	55
3		Quinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000635-46-1	50
4		Cyclopropanamine, 2-phenyl-, tra...	133	C9H11N	000155-09-9	49
5		1,4-Dihydrobenzo[e][1,2,4]triaz...	161	C8H7N3O	057711-25-8	49



Data Path : Z:\HPCHEM1\BNA M\DATA\BM042816\
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Sample : H2639-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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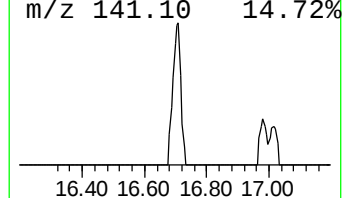
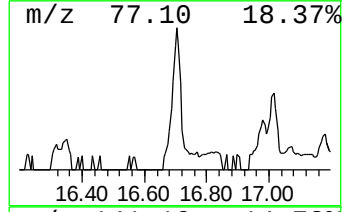
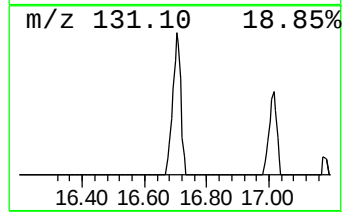
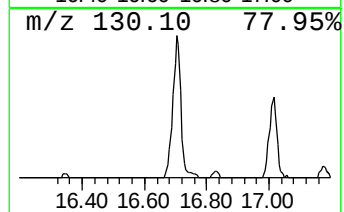
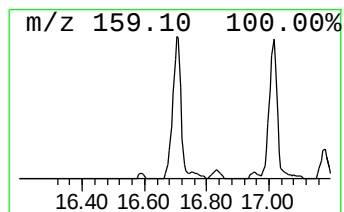
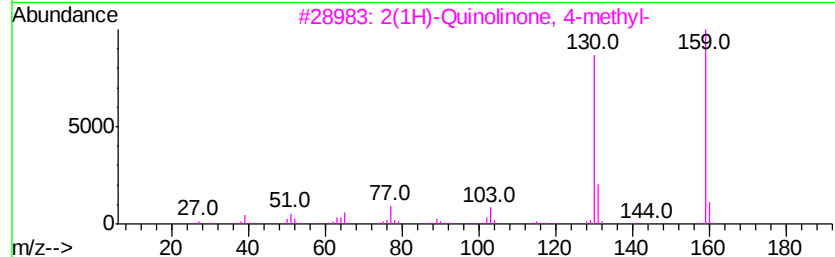
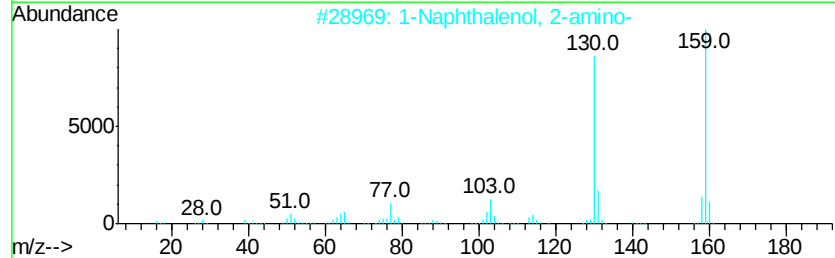
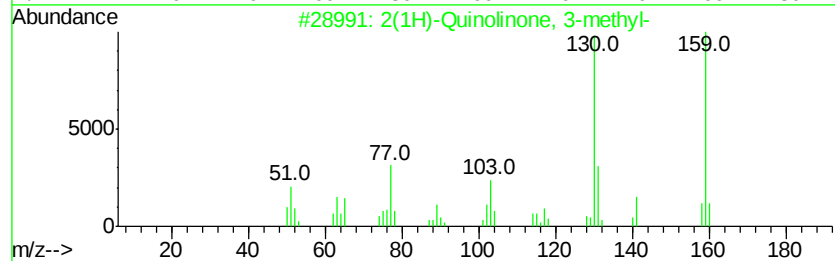
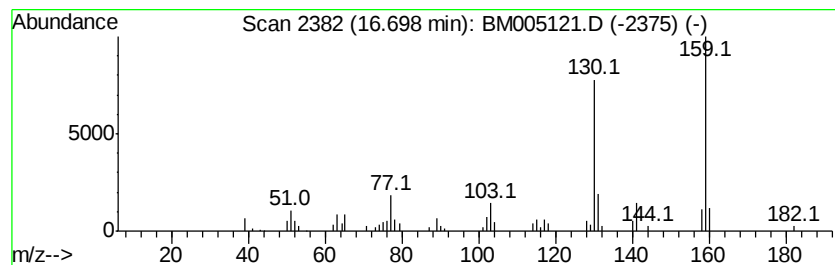
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 2(1H)-Quinolinone, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.70	4.02 ng/ul	137399	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2(1H)-Quinolinone, 3-methyl-	159	C10H9NO	002721-59-7	97
2		1-Naphthalenol, 2-amino-	159	C10H9NO	000606-41-7	93
3		2(1H)-Quinolinone, 4-methyl-	159	C10H9NO	000607-66-9	91
4		5-Amino-1-naphthol	159	C10H9NO	000083-55-6	90
5		2(1H)-Quinolinone, 4-methyl-	159	C10H9NO	000607-66-9	87



Data Path : Z:\HPCHEM1\BNA M\DATA\BM042816\
Data File : BM005121.D
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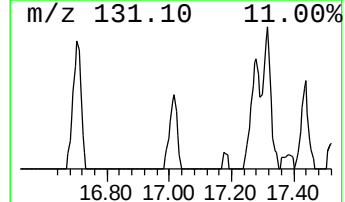
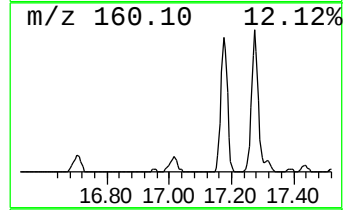
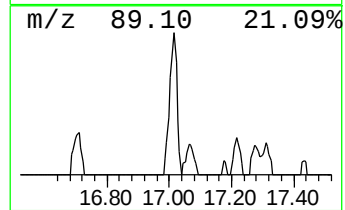
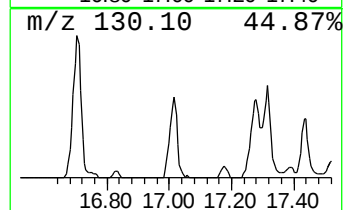
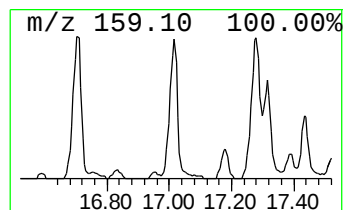
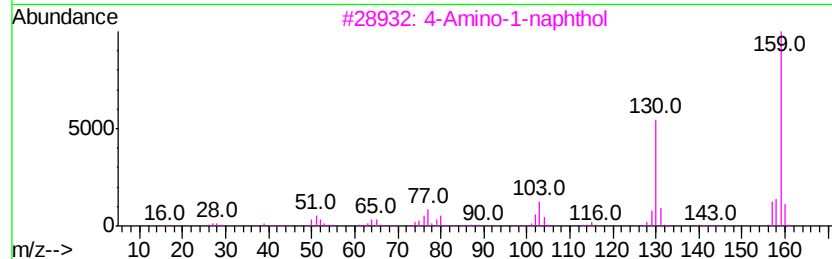
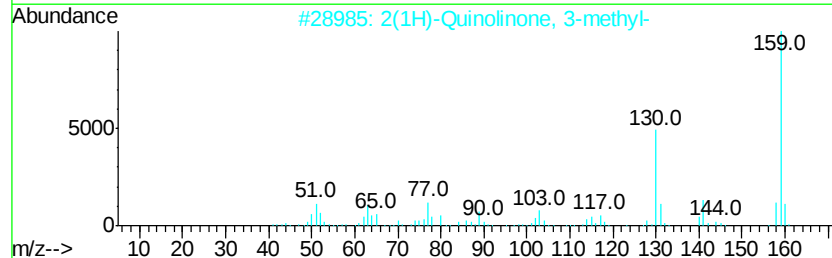
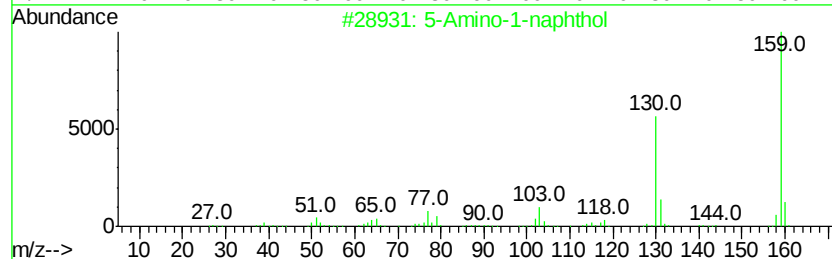
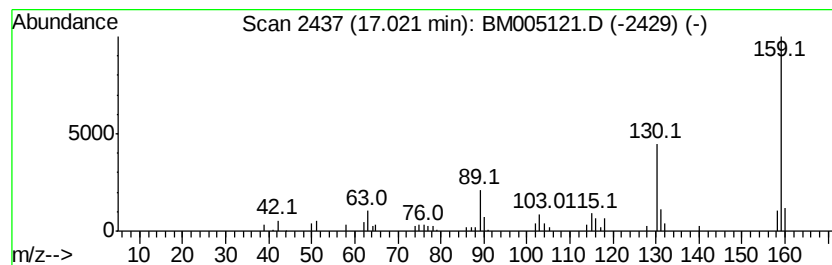
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 5-Amino-1-naphthol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.02	3.01 ng/ul	102930	Phenanthrene-d10	17.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Amino-1-naphthol		159	C10H9NO	000083-55-6	87
2	2(1H)-Quinolinone, 3-methyl-		159	C10H9NO	002721-59-7	83
3	4-Amino-1-naphthol		159	C10H9NO	002834-90-4	81
4	2(1H)-Quinolinone, 1-methyl-		159	C10H9NO	000606-43-9	80
5	2(1H)-Quinolinone, 1-methyl-		159	C10H9NO	000606-43-9	80



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM042816\
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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Hexanoic acid	6.82	5.2	ng/ul	60150	1	7.79	229676	20.0
Benzofuran, 2-met...	9.14	2.2	ng/ul	25247	1	7.79	229676	20.0
2-Benzothiophene #	10.75	14.3	ng/ul	273482	2	10.57	382375	20.0
Nonanoic acid	11.39	4.4	ng/ul	84460	2	10.57	382375	20.0
Benzo[b]thiophene...	11.69	2.3	ng/ul	43762	2	10.57	382375	20.0
unknown-01	14.85	3.1	ng/ul	87802	3	14.43	564617	20.0
1(2H)-Quinolineca...	16.16	2.2	ng/ul	74294	4	17.17	684115	20.0
2(1H)-Quinolinone...	16.70	4.0	ng/ul	137399	4	17.17	684115	20.0
5-Amino-1-naphthol	17.02	3.0	ng/ul	102930	4	17.17	684115	20.0