

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF101718\
 Data File : BF109939.D
 Acq On : 17 Oct 2018 23:56
 Operator : JU/SJ
 Sample : J5539-05
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-1-6-8FT

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101518.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.275	27	32	39	rVB	459998	622437	3.94%	0.167%
2	4.351	376	385	389	rBV2	703689	987795	6.25%	0.264%
3	4.669	434	439	448	rVB	753119	907534	5.74%	0.243%
4	4.769	448	456	461	rBV	391535	517011	3.27%	0.138%
5	5.051	500	504	511	rVB	767071	817124	5.17%	0.219%
6	5.128	511	517	520	rBV3	742594	1336995	8.46%	0.358%
7	5.157	520	522	527	rVB	1465913	1508180	9.54%	0.404%
8	5.210	527	531	535	rBV	2281289	2453996	15.53%	0.657%
9	5.269	537	541	544	rVB2	458760	618382	3.91%	0.165%
10	5.410	558	565	568	rBV2	1745678	2665849	16.87%	0.713%
11	5.481	573	577	581	rBV2	782929	812028	5.14%	0.217%
12	5.593	591	596	599	rBV2	4691303	5253435	33.24%	1.406%
13	5.681	608	611	614	rBV	1131286	1230525	7.79%	0.329%
14	5.757	621	624	629	rBV3	1217599	2007765	12.71%	0.537%
15	5.804	629	632	635	rVB	2380546	1969436	12.46%	0.527%
16	5.846	635	639	642	rBV	1559993	1656300	10.48%	0.443%
17	5.910	646	650	654	rBV	924353	906339	5.74%	0.243%
18	6.010	658	667	670	rBV3	2819131	4075963	25.79%	1.091%
19	6.045	670	673	677	rBV3	1261713	1870076	11.83%	0.500%
20	6.128	684	687	688	rBV	893207	753764	4.77%	0.202%
21	6.146	688	690	693	rVB2	3501832	3071529	19.44%	0.822%
22	6.181	693	696	699	rBV2	1137311	1334849	8.45%	0.357%
23	6.228	699	704	711	rBV3	5072300	10817447	68.45%	2.895%
24	6.287	711	714	716	rVB	3798488	3395480	21.49%	0.909%
25	6.322	716	720	722	rBV4	1058058	1496272	9.47%	0.400%
26	6.351	722	725	727	rVB	949230	863820	5.47%	0.231%
27	6.369	727	728	732	rVB3	740226	706529	4.47%	0.189%
28	6.422	732	737	739	rBV3	3040308	4725778	29.91%	1.265%
29	6.475	743	746	749	rVB	3314148	3307372	20.93%	0.885%
30	6.522	749	754	762	rBV4	6579834	10655944	67.43%	2.852%
31	6.598	762	767	770	rBV2	4691494	5719901	36.20%	1.531%
32	6.634	770	773	775	rVV	2294276	2030643	12.85%	0.543%
33	6.657	775	777	779	rVV2	3013992	2507862	15.87%	0.671%
34	6.681	779	781	784	rVB	1708530	1446892	9.16%	0.387%

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	6.822	801	805	807	rBV2	3149848	4008926	25.37%	1.073%
36	6.840	807	808	811	rVB2	3115375	2270822	14.37%	0.608%
37	6.869	811	813	815	rBV	1341778	1317969	8.34%	0.353%
38	6.898	815	818	822	rVV4	1666165	2985993	18.90%	0.799%
39	6.951	822	827	828	rVV5	3147197	4982148	31.53%	1.333%
40	6.998	830	835	837	rVB2	8624997	11728418	74.22%	3.139%
41	7.040	837	842	845	rBV3	2337434	3036736	19.22%	0.813%
42	7.093	845	851	854	rBV3	2875931	5411030	34.24%	1.448%
43	7.134	854	858	861	rBV3	6605627	9054463	57.30%	2.423%
44	7.163	861	863	866	rVB	3202480	2794449	17.68%	0.748%
45	7.210	869	871	873	rBV	2611796	2408812	15.24%	0.645%
46	7.234	873	875	876	rBV2	1682445	1456748	9.22%	0.390%
47	7.257	876	879	882	rVB3	5582325	6642923	42.04%	1.778%
48	7.304	882	887	889	rBV3	2002866	2407043	15.23%	0.644%
49	7.328	889	891	894	rBV3	1926641	1889528	11.96%	0.506%
50	7.392	894	902	904	rBV3	6062713	9240467	58.48%	2.473%
51	7.463	910	914	917	rBV5	2260016	3118884	19.74%	0.835%
52	7.522	918	924	927	rVV3	9276551	15802340	100.00%	4.229%
53	7.563	928	931	934	rVV3	5262571	6842158	43.30%	1.831%
54	7.592	934	936	938	rVB	2004084	1474795	9.33%	0.395%
55	7.628	938	942	946	rVB4	6474769	11376396	71.99%	3.045%
56	7.692	946	953	956	rBV4	4838662	11787906	74.60%	3.155%
57	7.763	962	965	968	rVB3	4957690	5378019	34.03%	1.439%
58	7.804	968	972	977	rVB4	6572426	8994279	56.92%	2.407%
59	7.869	980	983	985	rBV2	3366431	3889502	24.61%	1.041%
60	7.892	985	987	989	rVV	4055708	3735123	23.64%	1.000%
61	7.922	989	992	994	rVV3	6011890	6420479	40.63%	1.718%
62	7.945	994	996	1000	rVB2	3674623	3804653	24.08%	1.018%
63	7.981	1000	1002	1004	rBV2	1211569	1056073	6.68%	0.283%
64	8.016	1004	1008	1010	rBV2	5694371	6477465	40.99%	1.734%
65	8.092	1017	1021	1023	rBV3	3494724	4681028	29.62%	1.253%
66	8.116	1023	1025	1027	rVB	2885630	2200312	13.92%	0.589%
67	8.145	1027	1030	1033	rBV2	4255446	5655198	35.79%	1.513%
68	8.181	1033	1036	1040	rBV5	1890860	2494680	15.79%	0.668%
69	8.234	1043	1045	1047	rBV2	2411685	2301650	14.57%	0.616%
70	8.269	1047	1051	1053	rBV4	3894045	5043510	31.92%	1.350%
71	8.345	1058	1064	1066	rVV2	6509159	12615787	79.83%	3.376%

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72	8.369	1066	1068	1070	rVB2	3041553	2611243	16.52%	0.699%
73	8.434	1076	1079	1081	rBV2	2271684	2184738	13.83%	0.585%
74	8.457	1081	1083	1085	rBV3	2507565	2836010	17.95%	0.759%
75	8.481	1085	1087	1092	rVB3	4451958	5822365	36.84%	1.558%
76	8.528	1093	1095	1097	rVV	1858138	1450228	9.18%	0.388%
77	8.575	1098	1103	1107	rVB5	5116120	8797045	55.67%	2.354%
78	8.669	1116	1119	1121	rVB	3691662	3738682	23.66%	1.001%
79	8.716	1121	1127	1131	rBV2	6477684	12896647	81.61%	3.451%
80	8.792	1137	1140	1143	rBV2	4306032	5398115	34.16%	1.445%
81	8.881	1152	1155	1158	rBV3	2604085	3444620	21.80%	0.922%
82	9.039	1180	1182	1184	rVB2	2544623	2043994	12.93%	0.547%
83	9.075	1184	1188	1189	rBV2	2449586	3124065	19.77%	0.836%
84	9.463	1249	1254	1258	rBV5	3507228	6523792	41.28%	1.746%
85	10.392	1408	1412	1414	rBV	2686759	4609040	29.17%	1.233%
86	11.780	1645	1648	1654	rVB3	2668096	3344428	21.16%	0.895%
87	12.163	1711	1713	1716	rVB2	1984956	1955424	12.37%	0.523%
88	12.604	1784	1788	1790	rVB	3303870	3764396	23.82%	1.007%
89	12.680	1799	1801	1804	rBV2	1310788	1479155	9.36%	0.396%
90	12.816	1822	1824	1827	rBV3	1115359	931411	5.89%	0.249%
91	12.998	1851	1855	1858	rVB2	2149002	2692259	17.04%	0.721%
92	13.033	1858	1861	1864	rVB	2165460	2257511	14.29%	0.604%
93	13.080	1867	1869	1872	rVB3	1200970	1244046	7.87%	0.333%
94	13.116	1872	1875	1877	rBV2	2706680	2534167	16.04%	0.678%
95	13.980	2020	2022	2026	rVB2	628662	638785	4.04%	0.171%
96	14.163	2050	2053	2059	rVB	1639564	1489674	9.43%	0.399%
97	14.927	2179	2183	2190	rVB	518856	558528	3.53%	0.149%
98	15.286	2240	2244	2249	rVB	606397	658774	4.17%	0.176%
99	15.686	2306	2312	2320	rVB2	1494713	2114676	13.38%	0.566%
100	16.151	2386	2391	2397	rBV	462201	677434	4.29%	0.181%

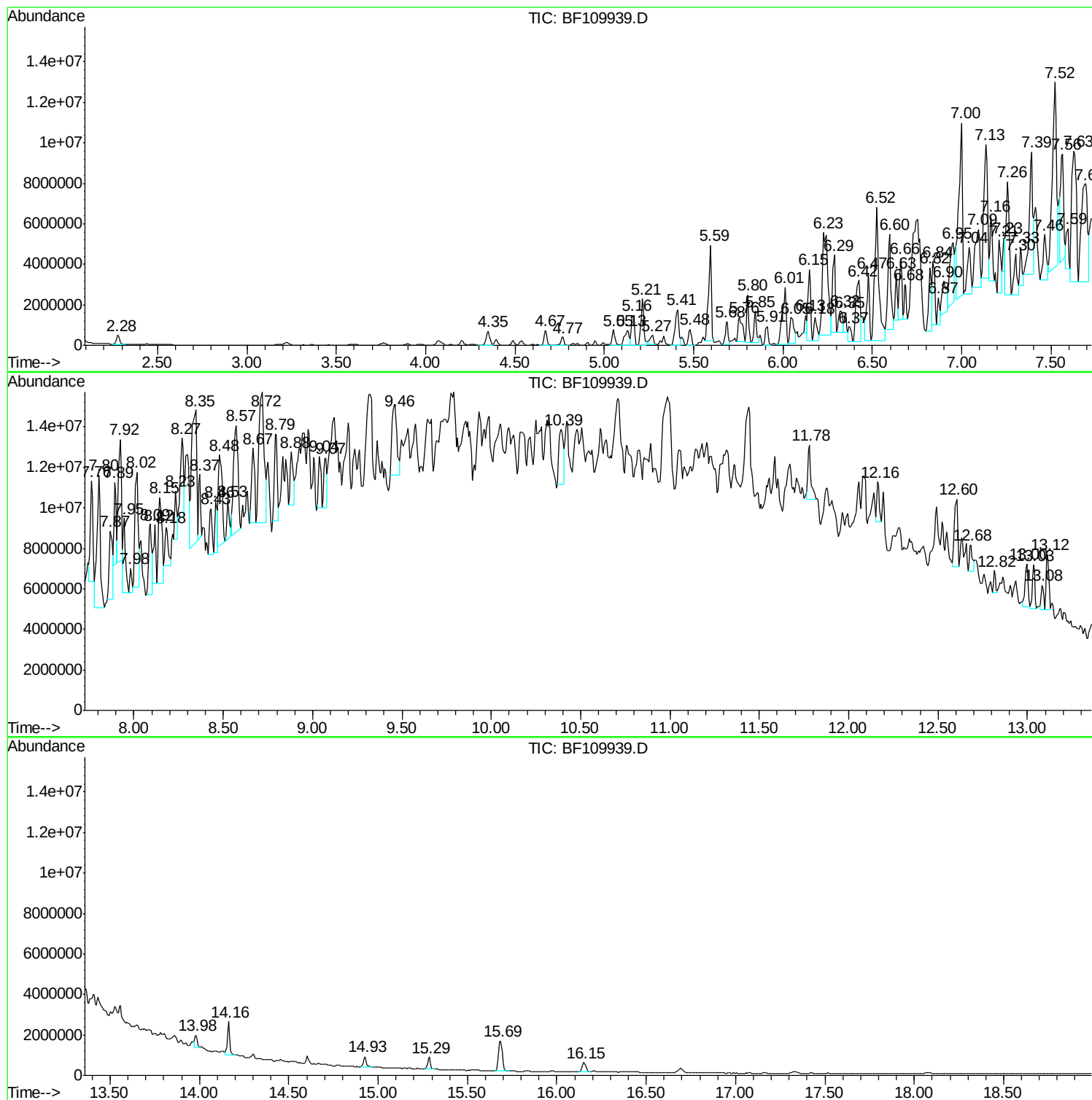
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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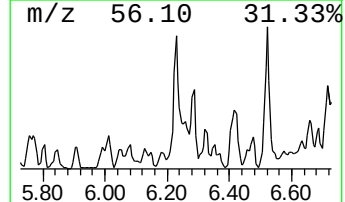
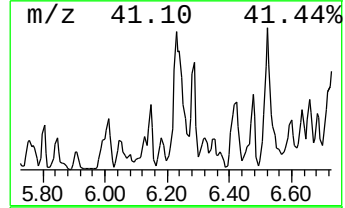
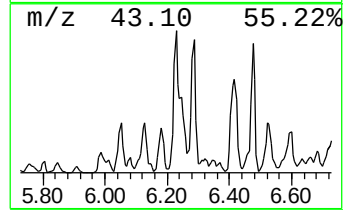
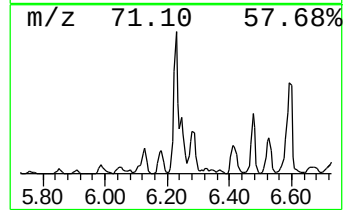
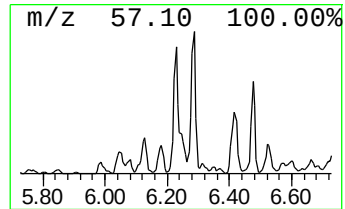
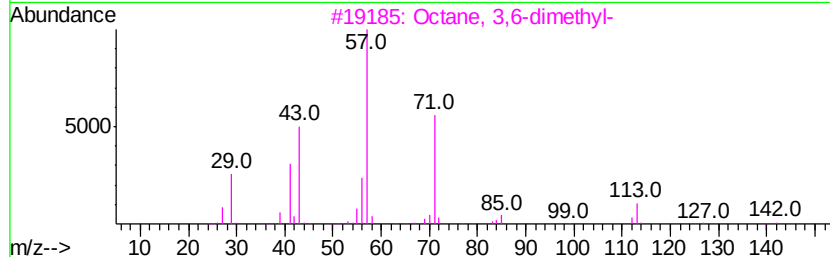
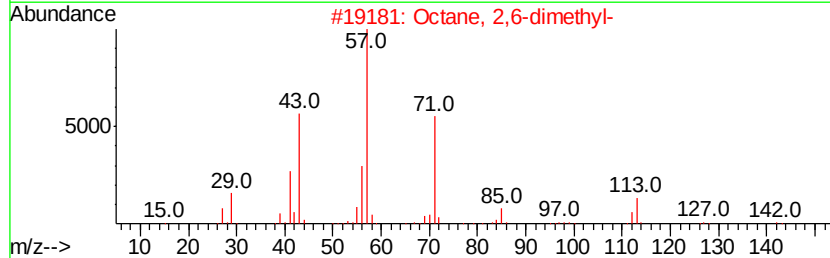
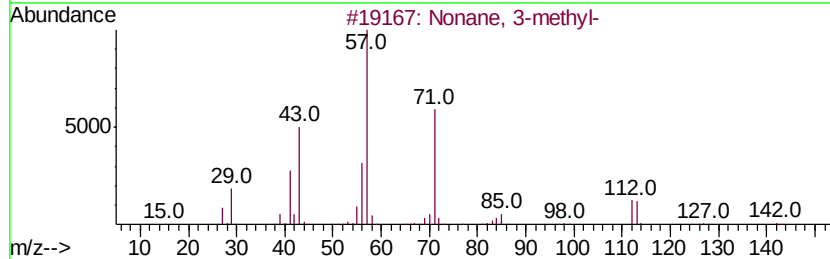
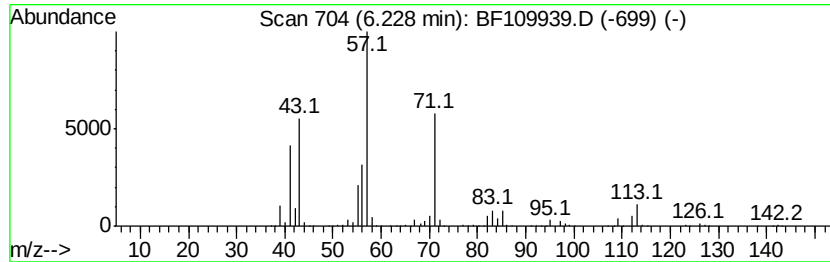
Instrument :
BNA_F
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SB-1-6-8FT

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Nonane, 3-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD		R.T.	
6.23	18.45 ng	10817400	1,4-Dichlorobenzene-d4		6.98	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 3-methyl-	142	C10H22	005911-04-6	90
2		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	83
3		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	78
4		Tridecane, 7-methyl-	198	C14H30	026730-14-3	64
5		Sulfurous acid, butyl octyl ester	250	C12H26O3S	1000309-17-5	64



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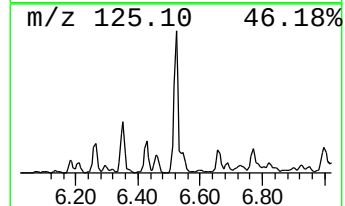
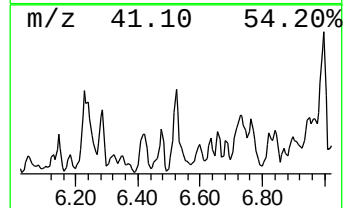
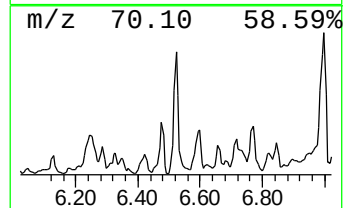
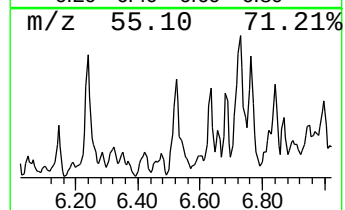
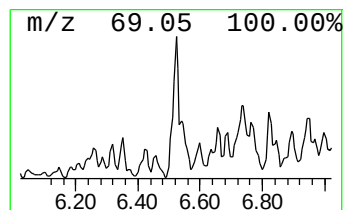
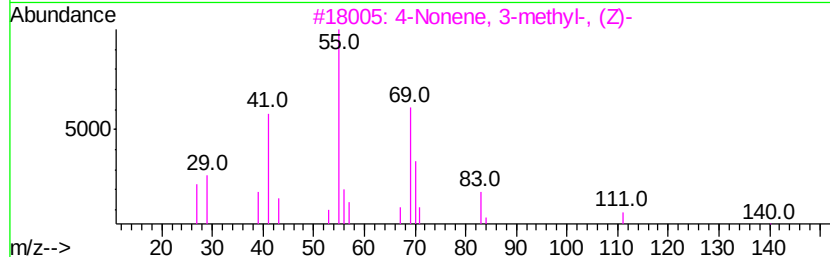
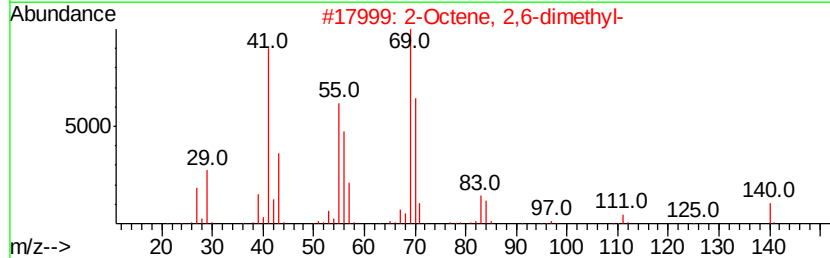
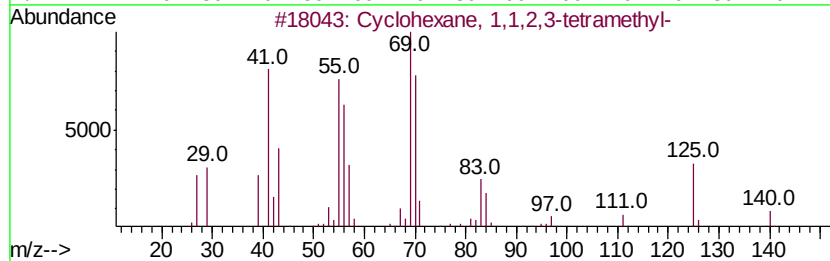
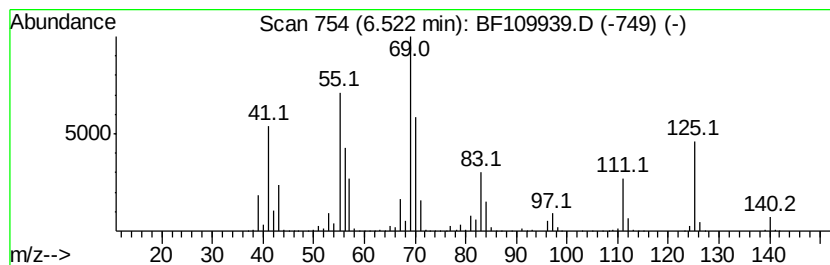
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclohexane, 1,1,2,3-tetram... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.52	18.17 ng	10655900	1,4-Dichlorobenzene-d4	6.98

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1,2,3-tetramethyl-	140	C10H20	006783-92-2	74
2			2-Octene, 2,6-dimethyl-	140	C10H20	004057-42-5	58
3			4-Nonene, 3-methyl-, (Z)-	140	C10H20	063830-69-3	58
4			1-Hexene, 3,3-dimethyl-	112	C8H16	003404-77-1	46
5			4-Octene, 2,6-dimethyl-, [S-(E)]-	140	C10H20	062960-76-3	43



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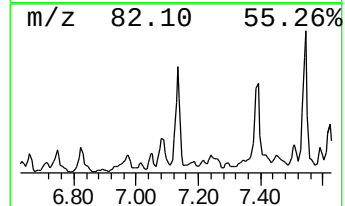
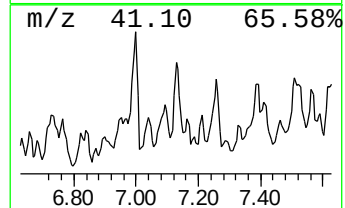
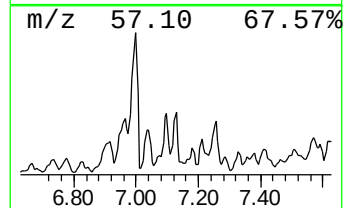
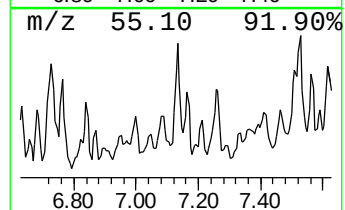
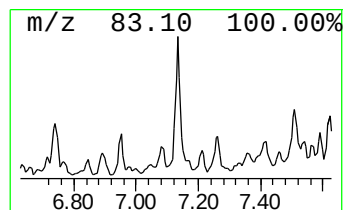
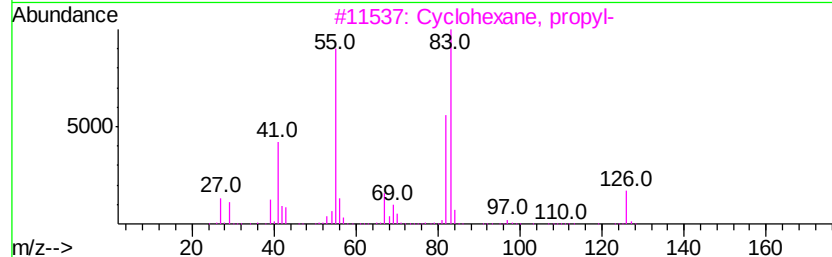
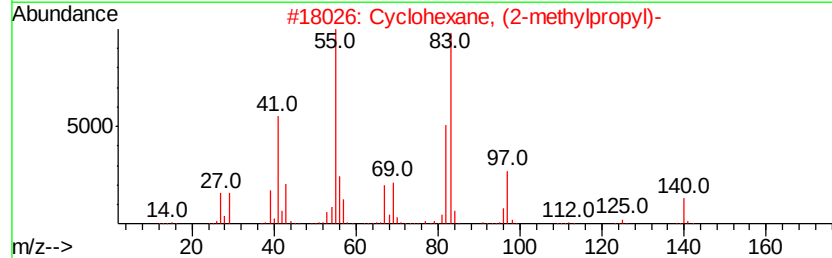
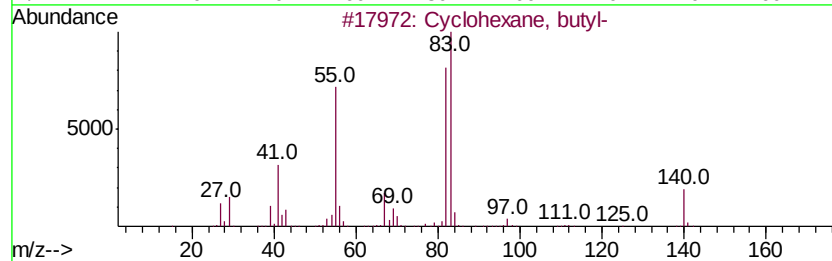
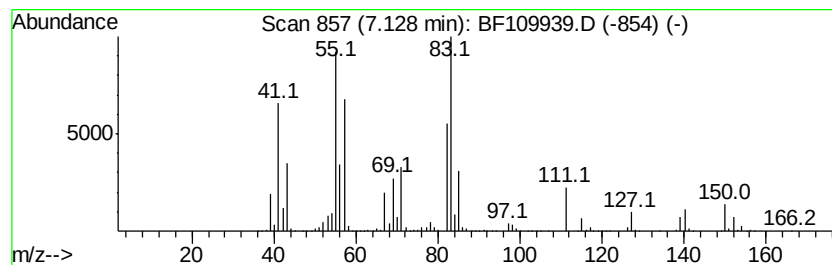
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Cyclohexane, butyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.13	15.44 ng	9054460	1,4-Dichlorobenzene-d4	6.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, butyl-	140	C10H20	001678-93-9	76
2		Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	58
3		Cyclohexane, propyl-	126	C9H18	001678-92-8	58
4		Cyclohexane, octyl-	196	C14H28	001795-15-9	53
5		Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	53



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Operator : JU/SJ
Sample : J5539-05
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ClientSampleId :
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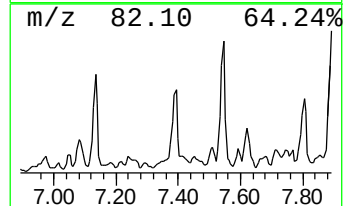
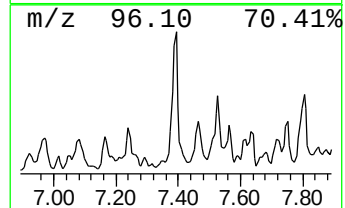
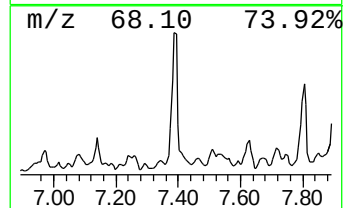
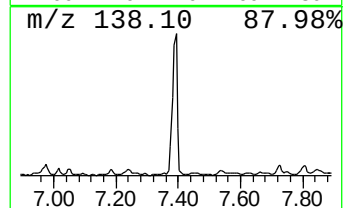
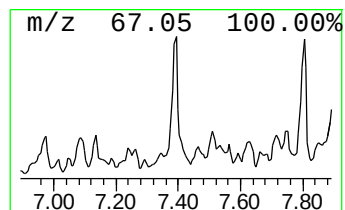
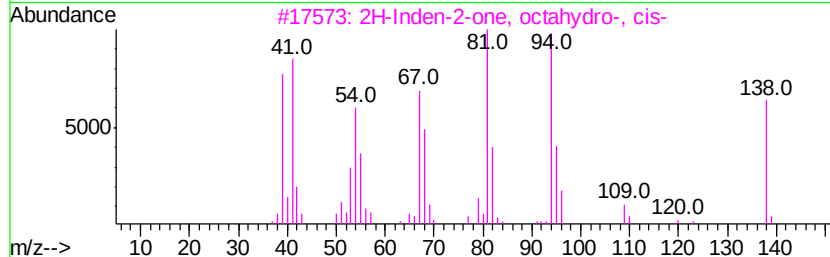
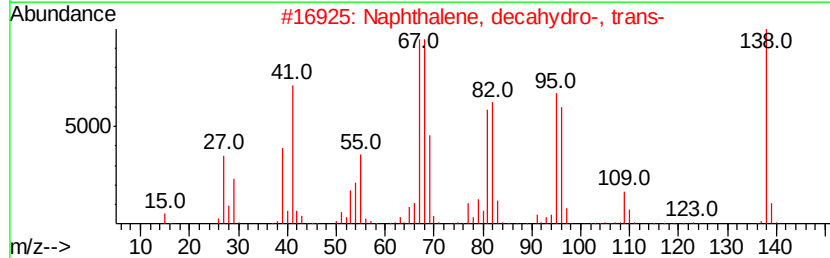
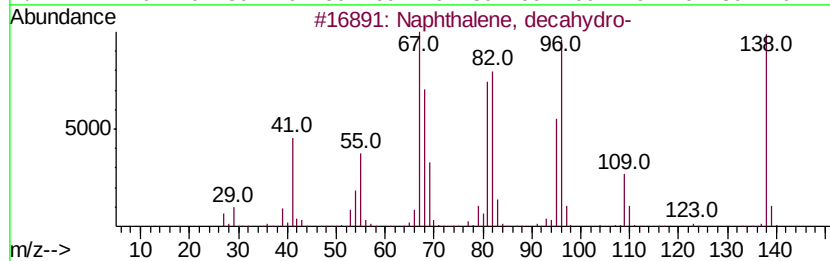
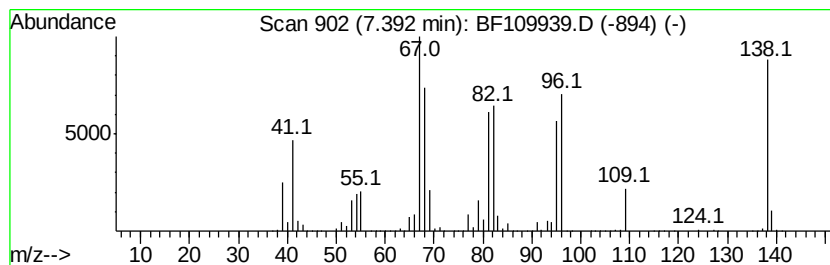
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Naphthalene, decahydro- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.39	15.76 ng	9240470	1,4-Dichlorobenzene-d4	6.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, decahydro-	138	C10H18	000091-17-8	94
2		Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	91
3		2H-Inden-2-one, octahydro-, cis-	138	C9H14O	005689-04-3	86
4		Spiro[4.5]decane	138	C10H18	000176-63-6	76
5		Naphthalene, decahydro-, cis-	138	C10H18	000493-01-6	76



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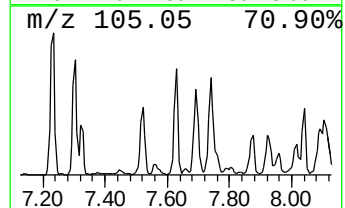
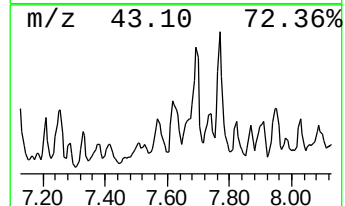
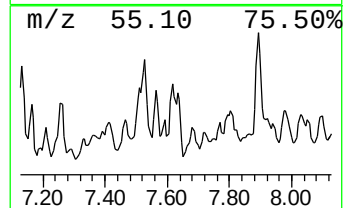
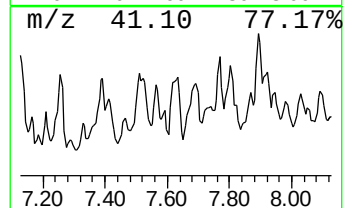
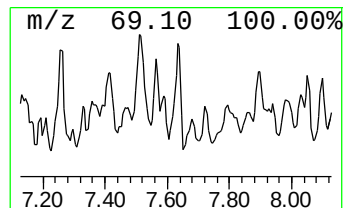
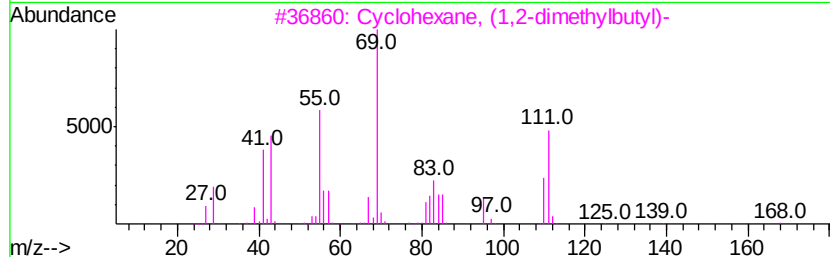
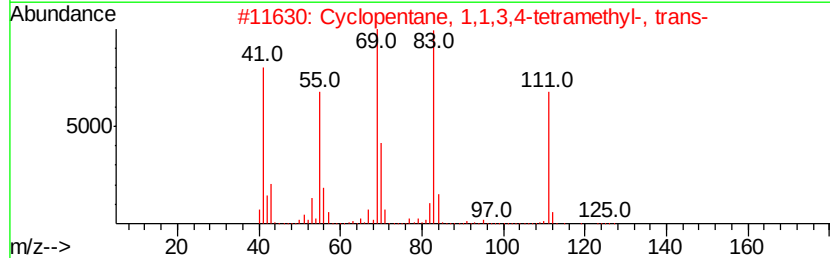
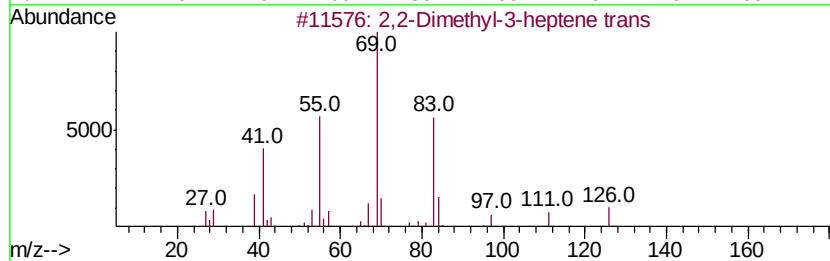
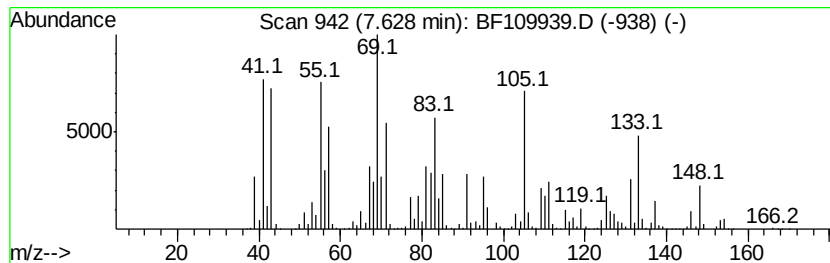
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown7.63 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.63	19.40 ng	11376400	1,4-Dichlorobenzene-d4	6.98

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2-Dimethyl-3-heptene trans			126	C9H18	019550-75-5	30
2	Cyclopentane, 1,1,3,4-tetramethyl-			126	C9H18	020309-77-7	27
3	Cyclohexane, (1,2-dimethylbutyl)-			168	C12H24	061142-37-8	27
4	Cyclohexane, 1,1,2-trimethyl-			126	C9H18	007094-26-0	25
5	Cyclohexane, 1,1,3,5-tetramethyl-			140	C10H20	050876-31-8	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101718\
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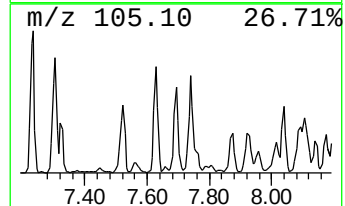
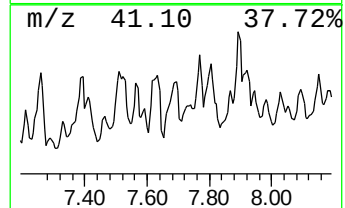
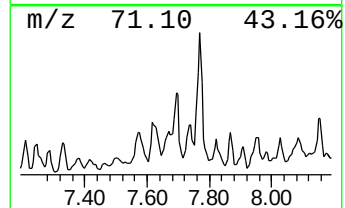
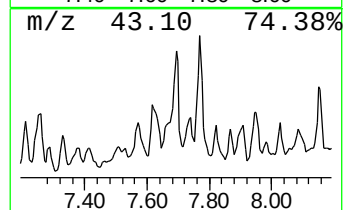
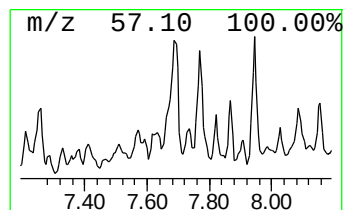
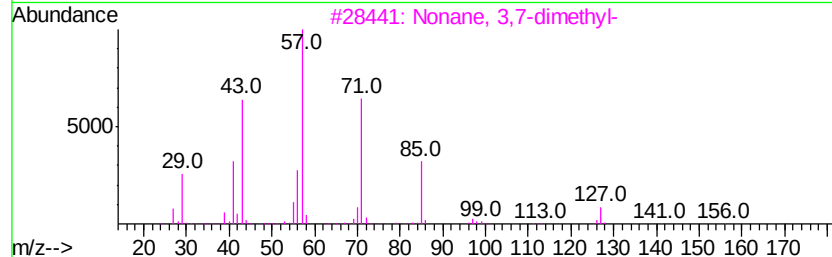
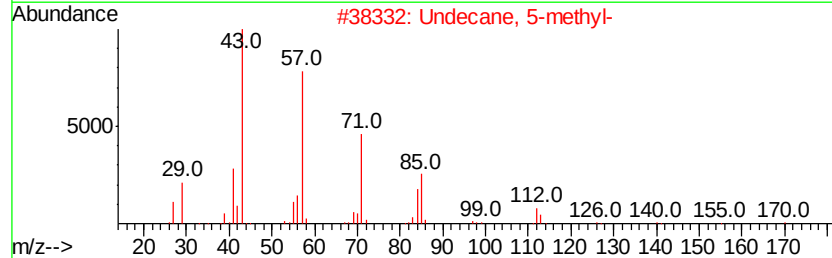
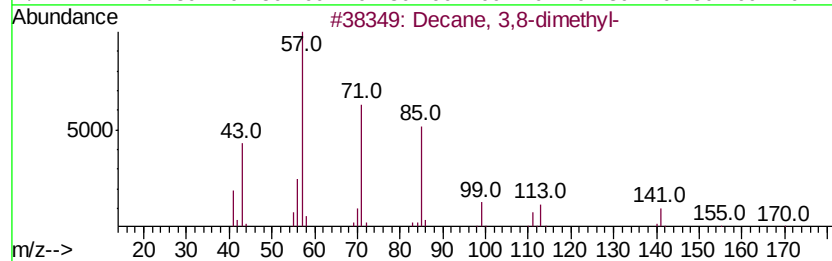
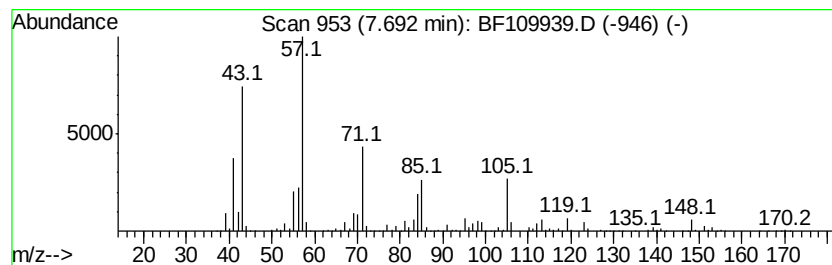
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Decane, 3,8-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.69	46.74 ng	11787900	Napthalene-d8	8.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	62
2		Undecane, 5-methyl-	170	C12H26	001632-70-8	62
3		Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	58
4		2,6-Dimethyldecane	170	C12H26	013150-81-7	53
5		Dodecane	170	C12H26	000112-40-3	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101718\
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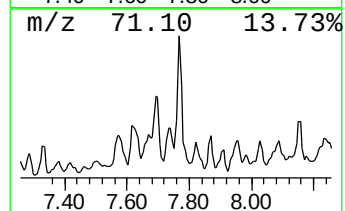
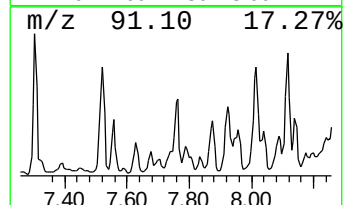
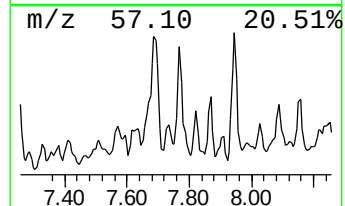
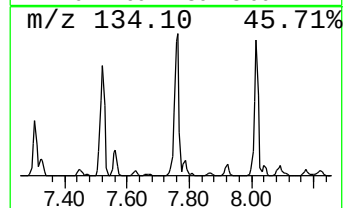
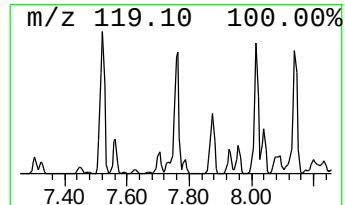
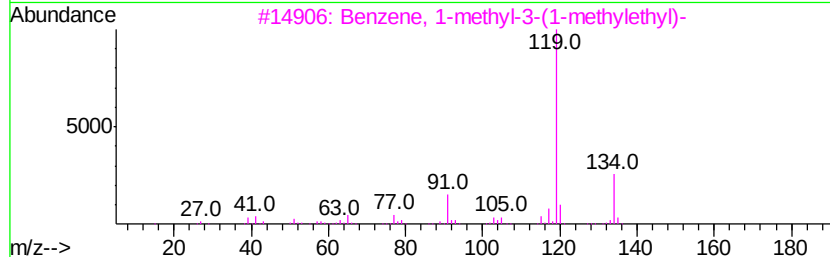
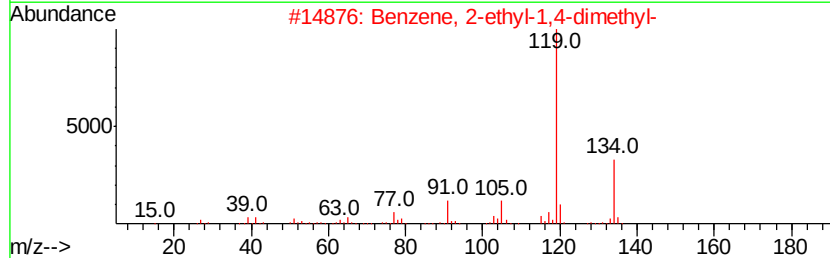
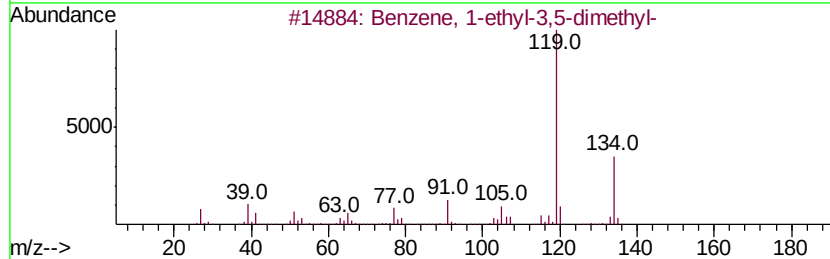
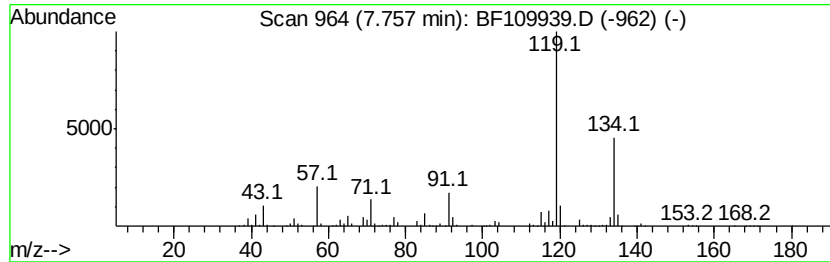
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.76	21.33 ng	5378020	Napthalene-d8	8.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	70
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	70
3		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	70
4		p-Cymene	134	C10H14	000099-87-6	70
5		o-Cymene	134	C10H14	000527-84-4	70



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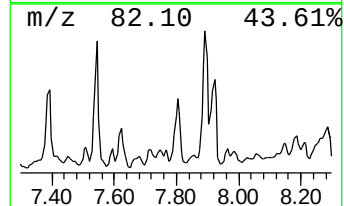
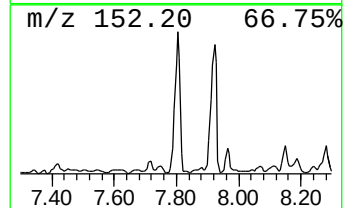
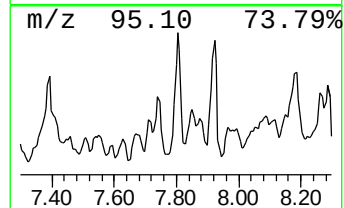
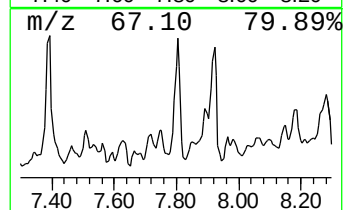
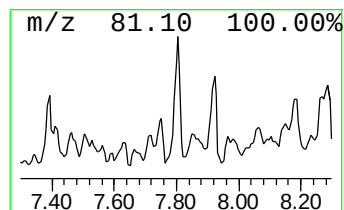
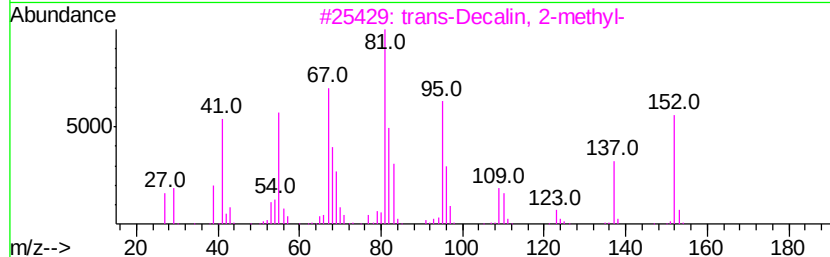
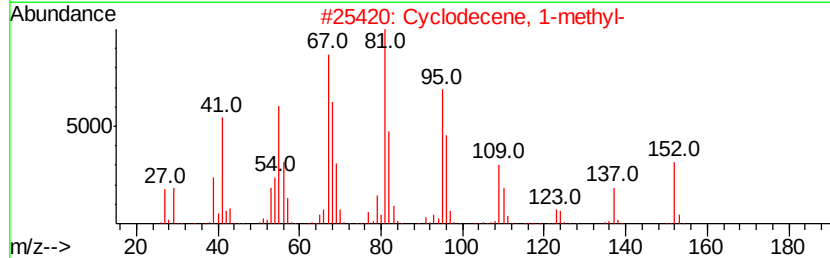
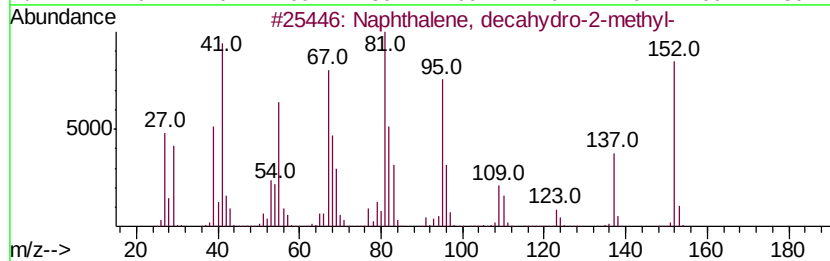
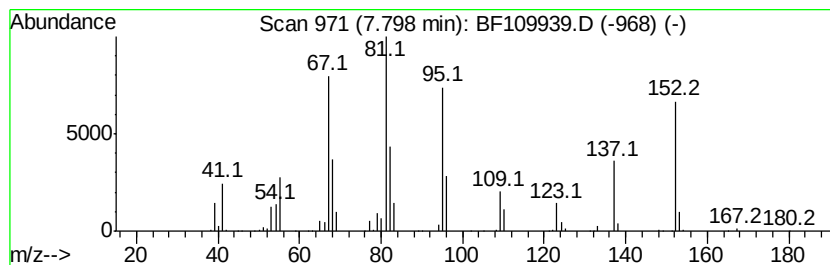
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Naphthalene, decahydro-2-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.80	35.67 ng	8994280	Naphthalene-d8	8.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	91
2		Cyclodecene, 1-methyl-	152	C11H20	066633-38-3	87
3		trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	87
4		trans-4a-Methyl-decahydronaphtha...	152	C11H20	002547-27-5	76
5		Cyclohexanone, 2-methyl-5-(1-met...	152	C10H16O	005948-04-9	70



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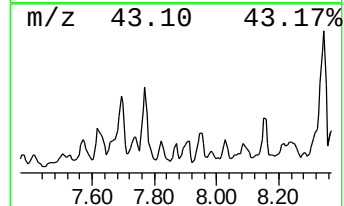
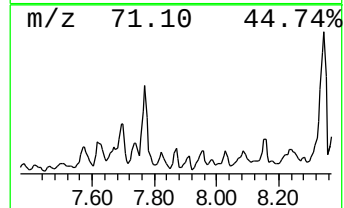
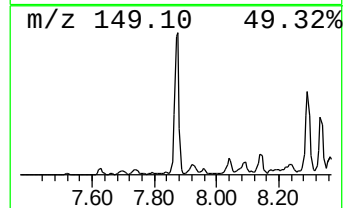
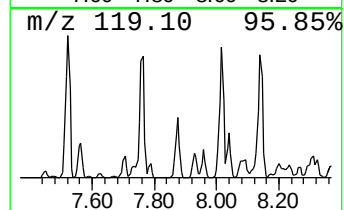
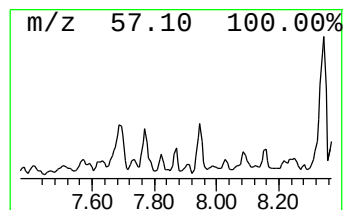
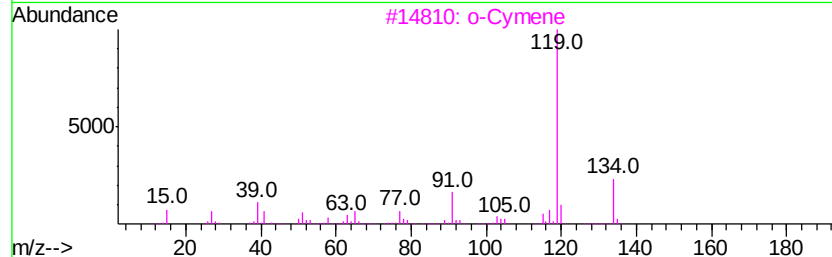
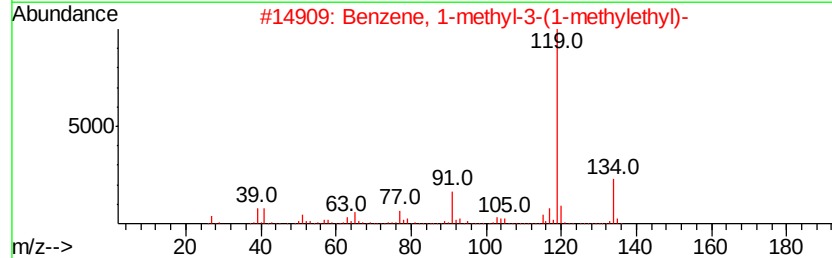
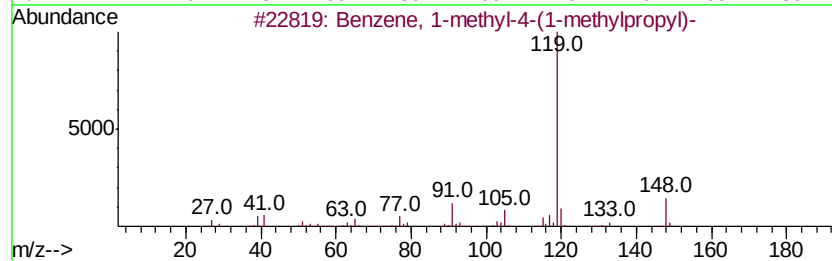
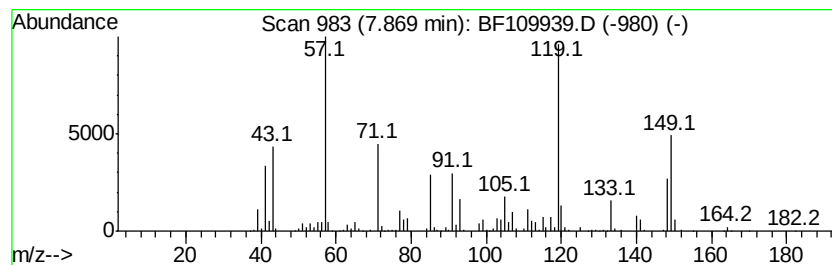
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown7.87 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.87	15.42 ng	3889500	Napthalene-d8	8.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	46
2		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	38
3		o-Cymene	134	C10H14	000527-84-4	38
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	38
5		Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	38



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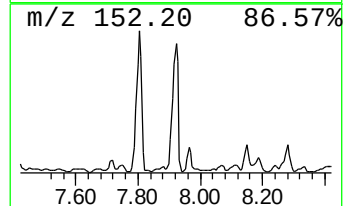
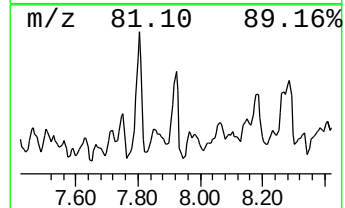
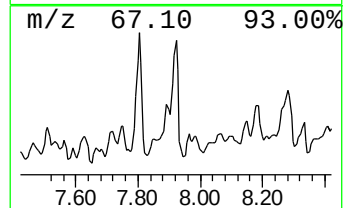
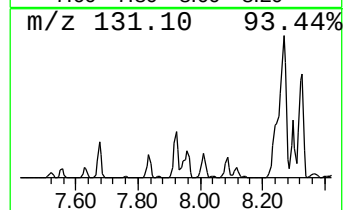
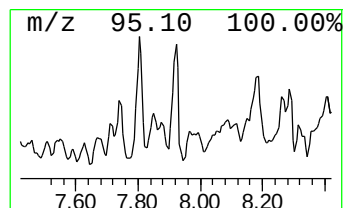
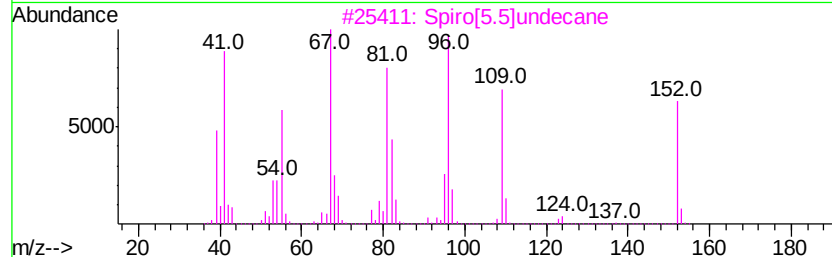
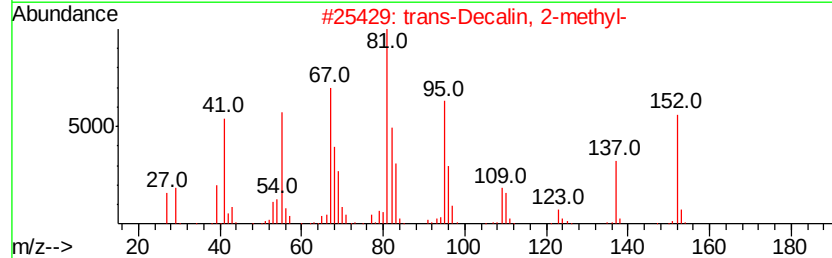
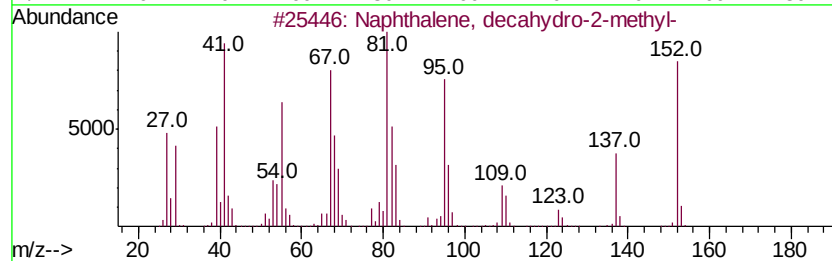
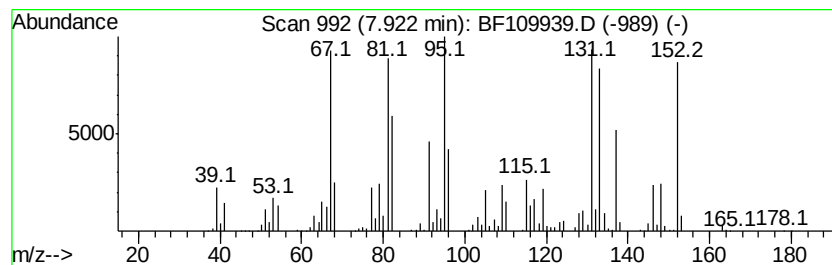
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown7.92 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.92	25.46 ng	6420480	Naphthalene-d8	8.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	49
2		trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	45
3		Spiro[5.5]undecane	152	C11H20	000180-43-8	38
4		Decalin, anti-1-methyl-, cis-	152	C11H20	1000158-89-0	38
5		trans-4a-Methyl-decahydronaphtha...	152	C11H20	002547-27-5	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101718\
Data File : BF109939.D
Acq On : 17 Oct 2018 23:56
Operator : JU/SJ
Sample : J5539-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

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ClientSampleId :
SB-1-6-8FT

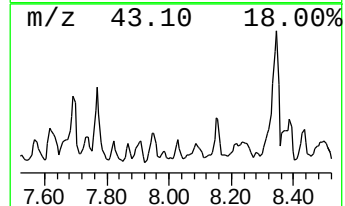
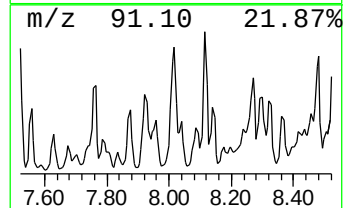
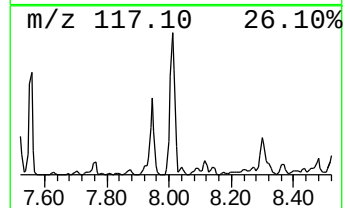
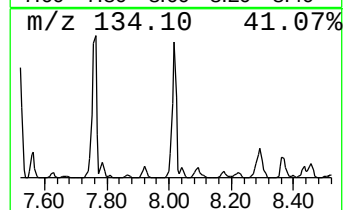
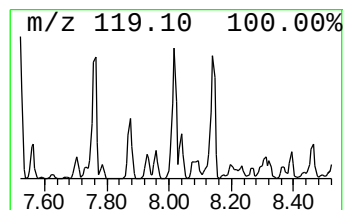
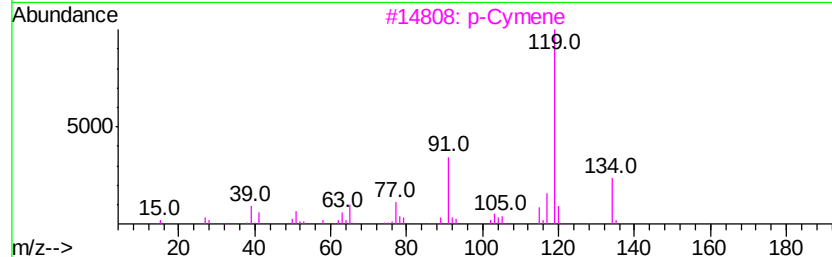
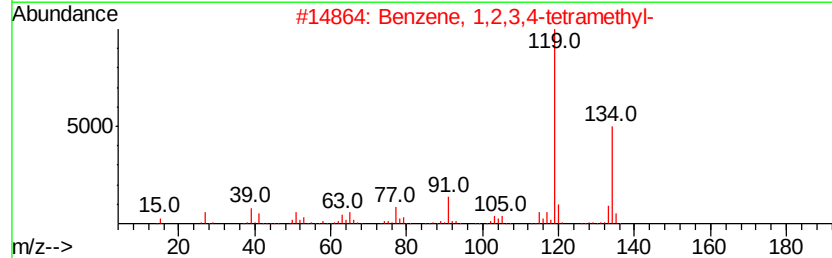
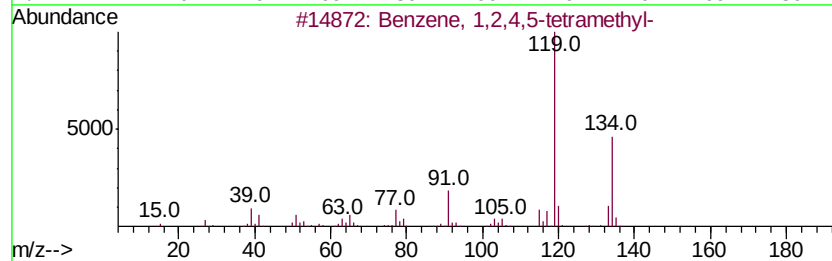
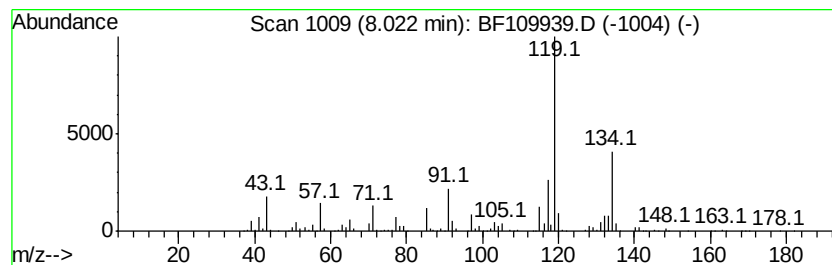
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.02	25.69 ng	6477470	Napthalene-d8	8.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	64
3		p-Cymene	134	C10H14	000099-87-6	60
4		o-Cymene	134	C10H14	000527-84-4	60
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	60



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101718\
Data File : BF109939.D
Acq On : 17 Oct 2018 23:56
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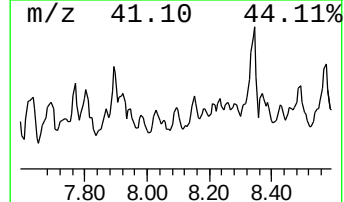
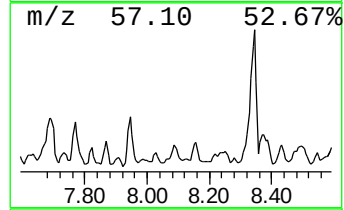
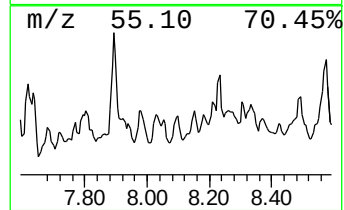
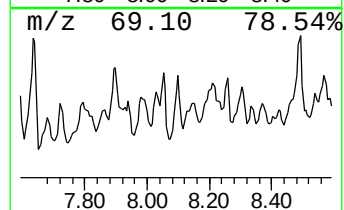
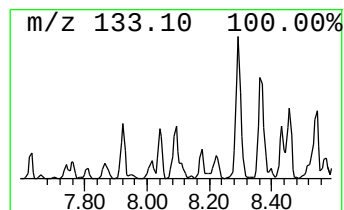
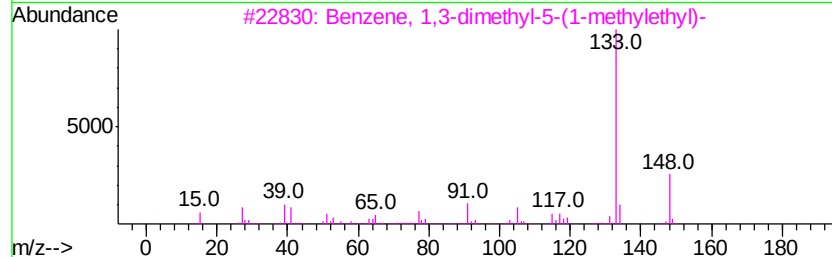
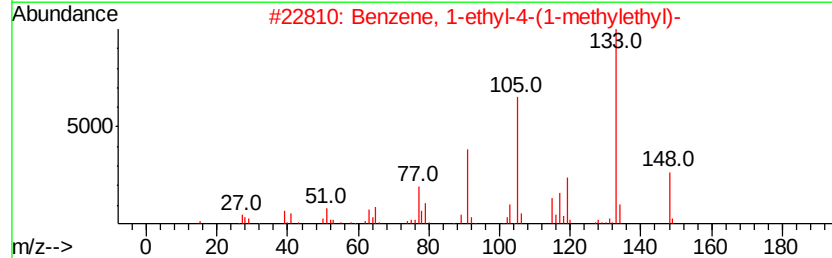
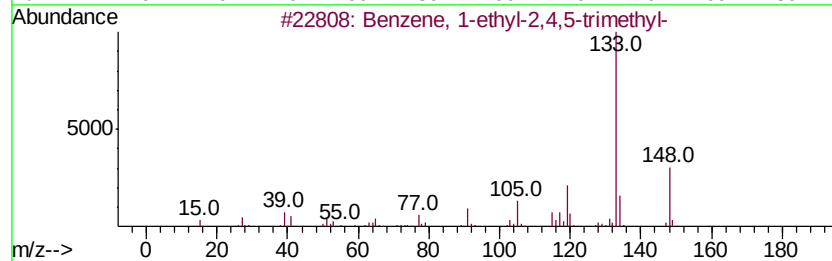
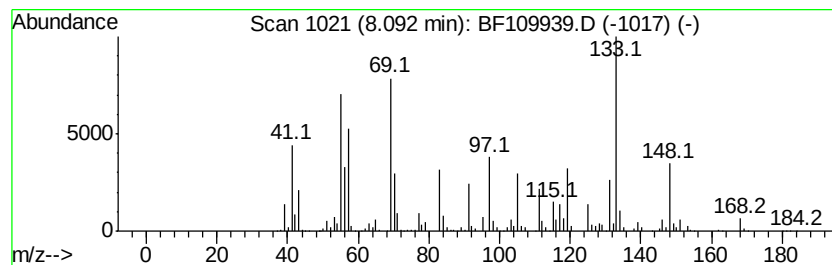
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown8.09 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.09	18.56 ng	4681030	Napthalene-d8	8.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4,5-trimethyl-	148	C11H16	017851-27-3	46
2		Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	46
3		Benzene, 1,3-dimethyl-5-(1-methv...	148	C11H16	004706-90-5	41
4		Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	38
5		Ethanone, 1-(3,4-dimethylphenyl)-	148	C10H12O	003637-01-2	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101718\
Data File : BF109939.D
Acq On : 17 Oct 2018 23:56
Operator : JU/SJ
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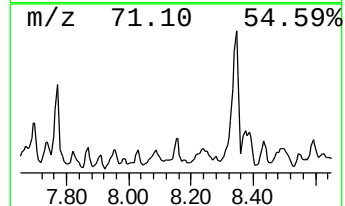
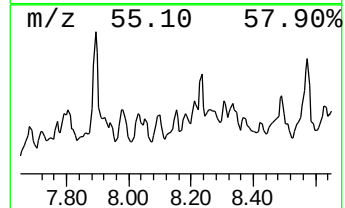
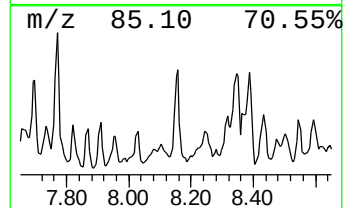
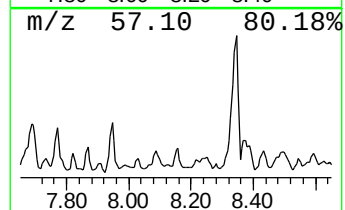
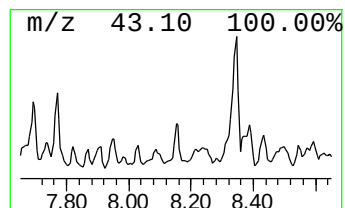
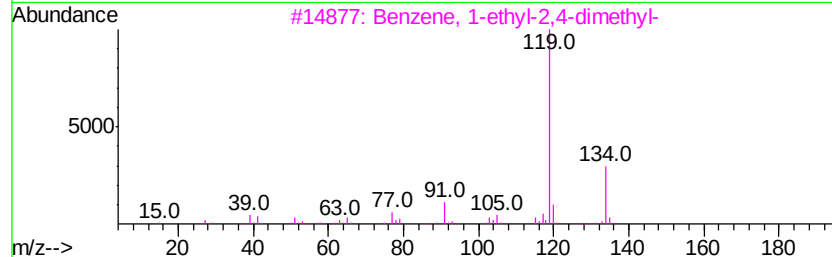
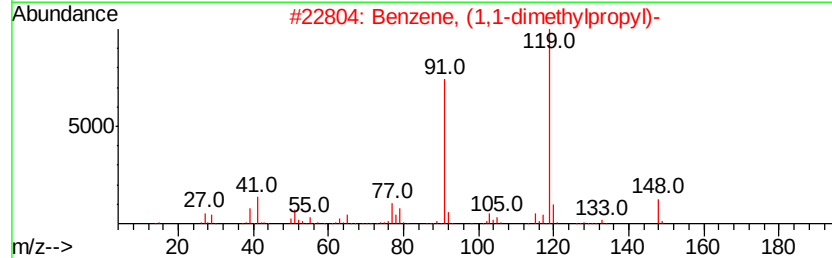
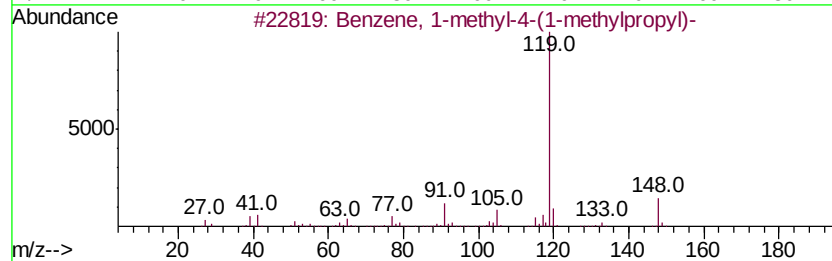
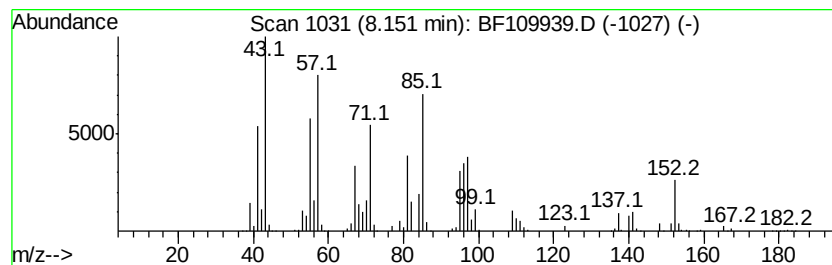
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzene, 1-methyl-4-(1-meth... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.15	22.43 ng	5655200	Napthalene-d8	8.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	64
2		Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	52
3		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	38
4		1,3-Benzenediol, 0-(2-chlorobenz...	366	C21H15ClO4	1000345-54-5	38
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101718\
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Acq On : 17 Oct 2018 23:56
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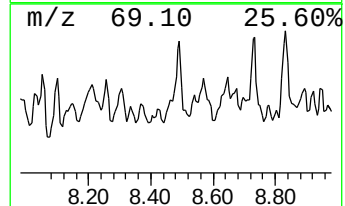
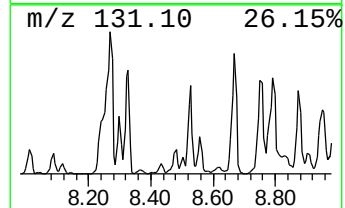
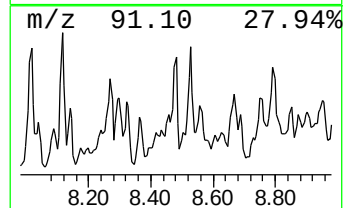
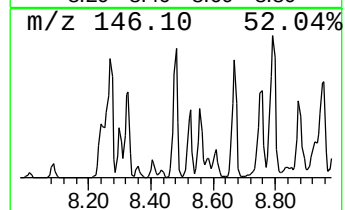
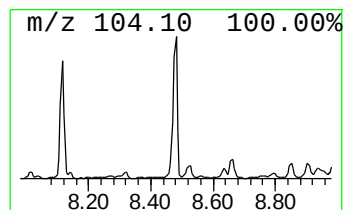
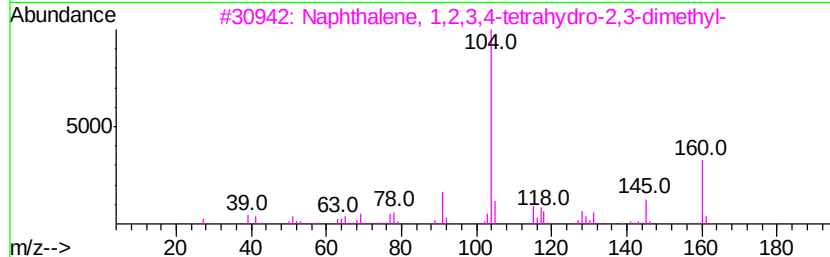
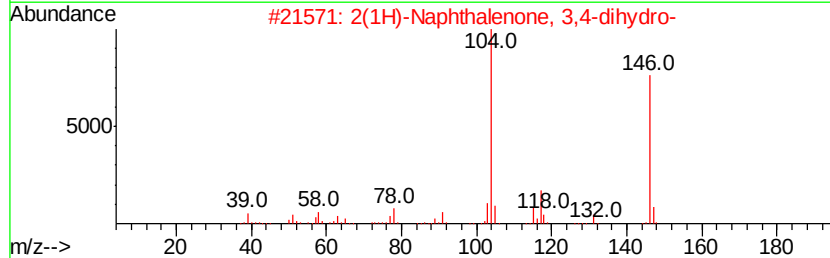
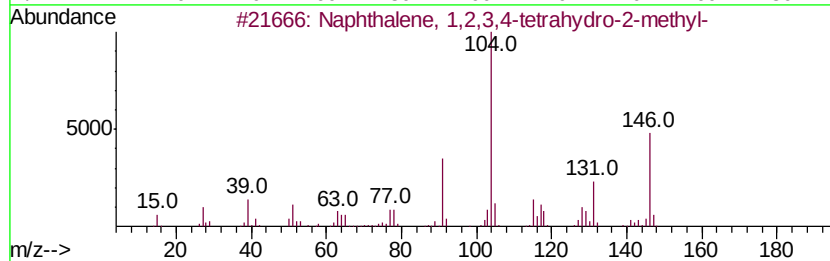
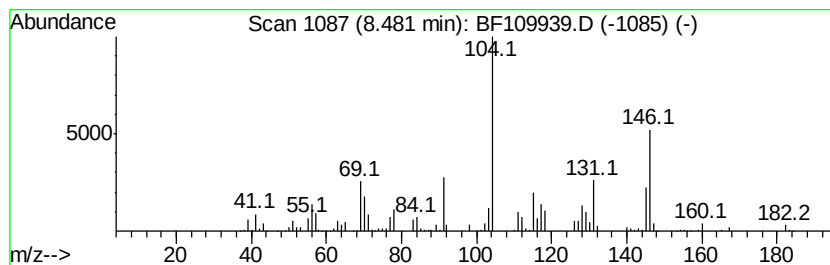
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Naphthalene, 1,2,3,4-tetra... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.48	23.09 ng	5822370	Naphthalene-d8	8.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	87
2		2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	47
3		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-92-1	47
4		Naphthalene, 2-ethyl-1,2,3,4-tet...	160	C12H16	032367-54-7	38
5		2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101718\
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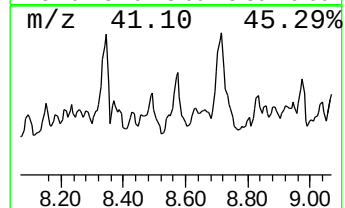
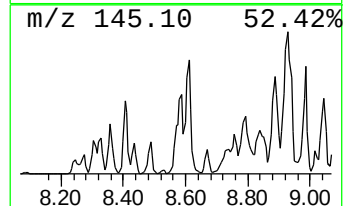
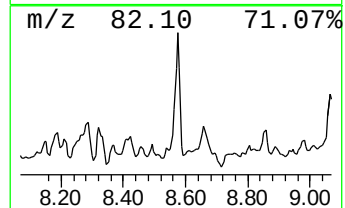
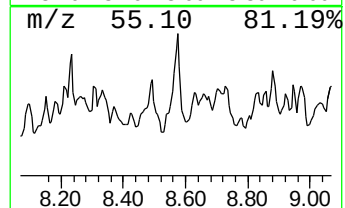
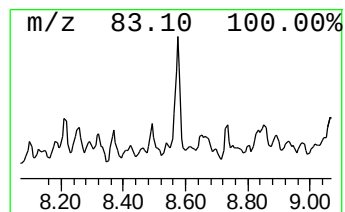
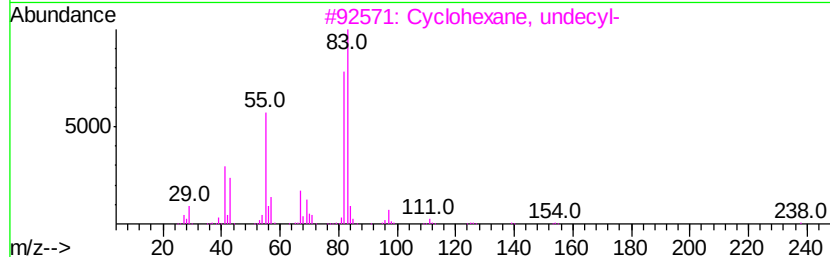
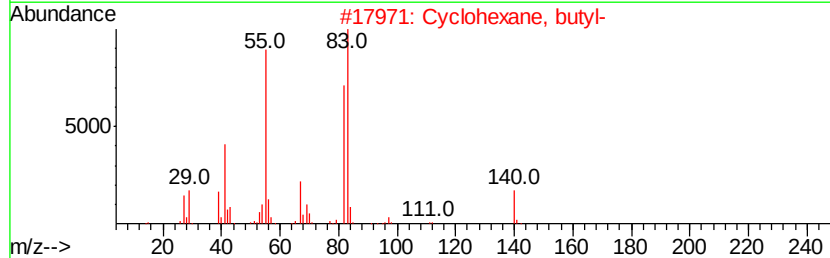
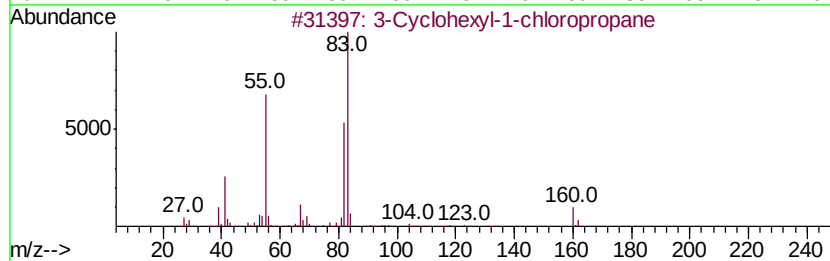
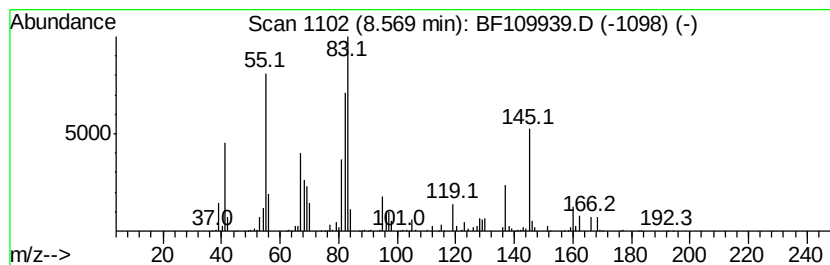
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ClientSampled :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 3-Cyclohexyl-1-chloropropane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
8.57	34.88 ng	8797050	Napthalene-d8	8.28		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Cyclohexyl-1-chloropropane	160	C9H17Cl	001124-62-5	64
2		Cyclohexane, butyl-	140	C10H20	001678-93-9	50
3		Cyclohexane, undecyl-	238	C17H34	054105-66-7	47
4		Cyclohexane, hexyl-	168	C12H24	004292-75-5	47
5		Cyclohexane, propyl-	126	C9H18	001678-92-8	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101718\
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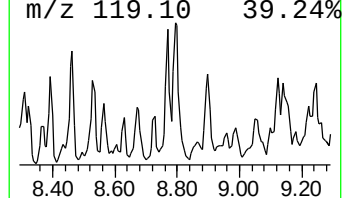
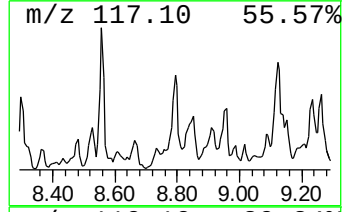
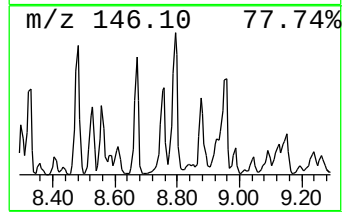
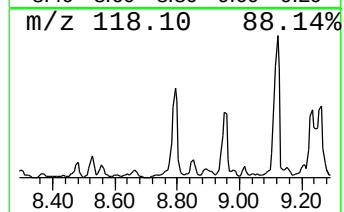
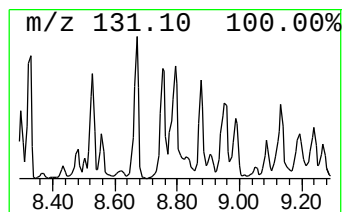
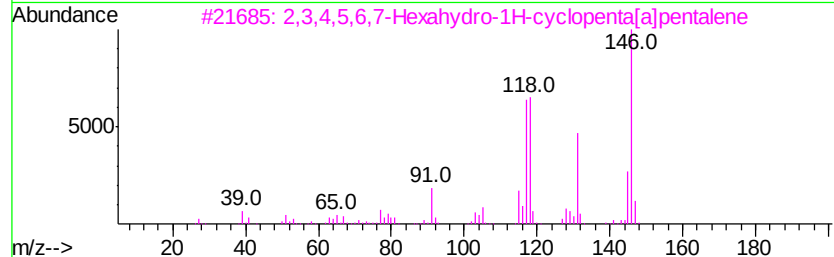
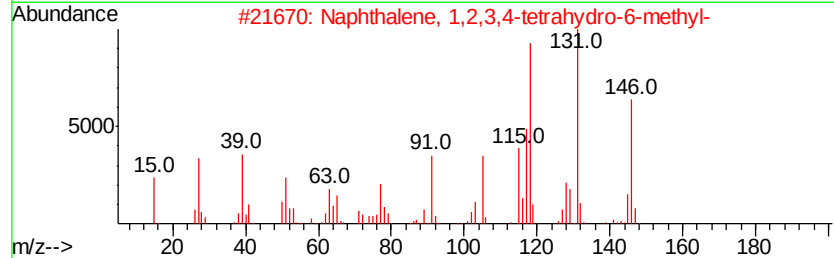
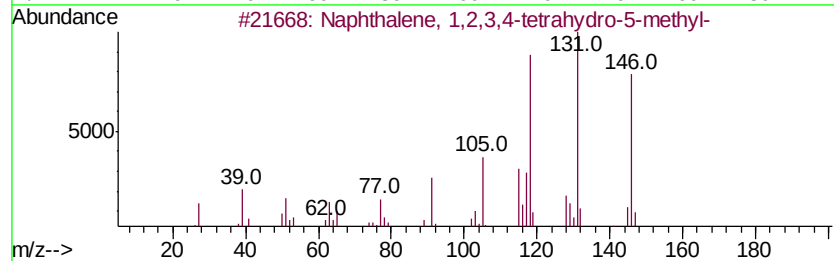
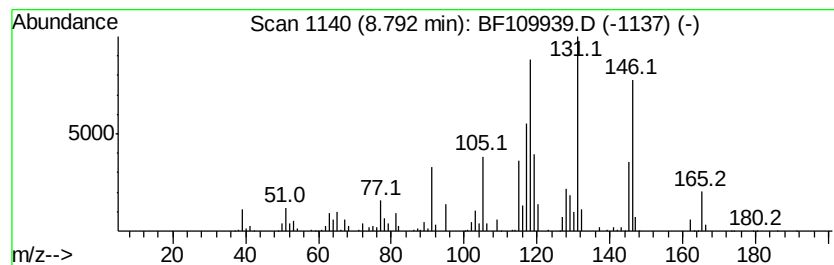
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Naphthalene, 1,2,3,4-tetra... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.		
8.79	21.41 ng	5398120	Naphthalene-d8	8.28		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	95
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	95
3		2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1000189-31-0	89
4		1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	62
5		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101718\
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ALS Vial : 17 Sample Multiplier: 1

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SB-1-6-8FT

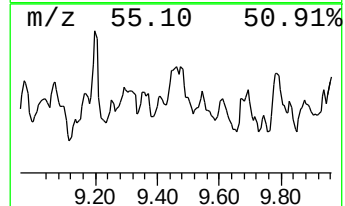
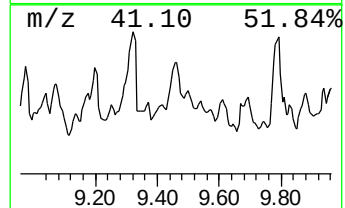
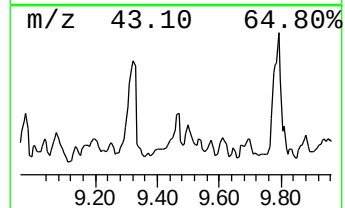
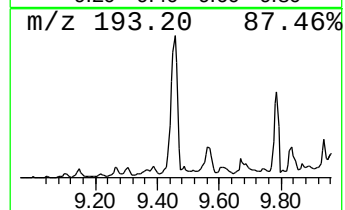
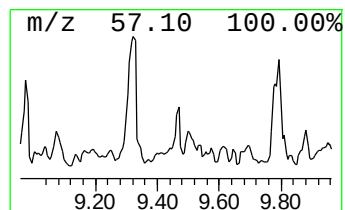
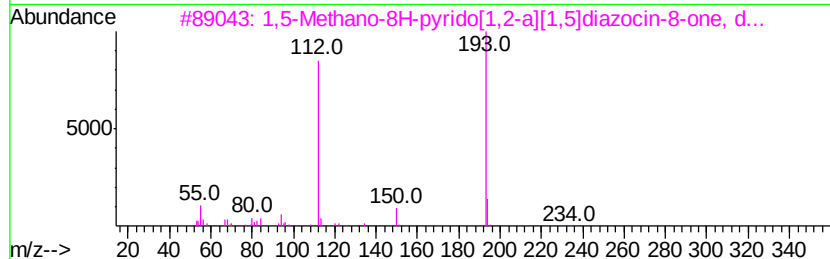
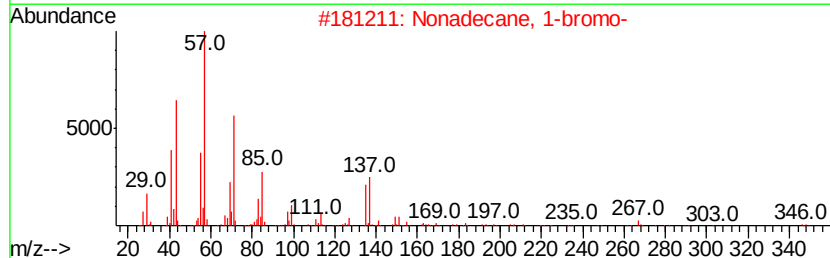
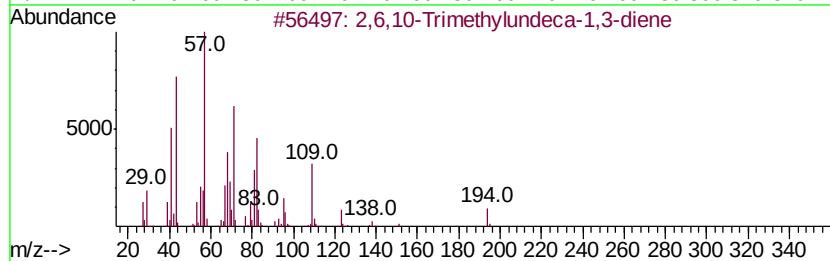
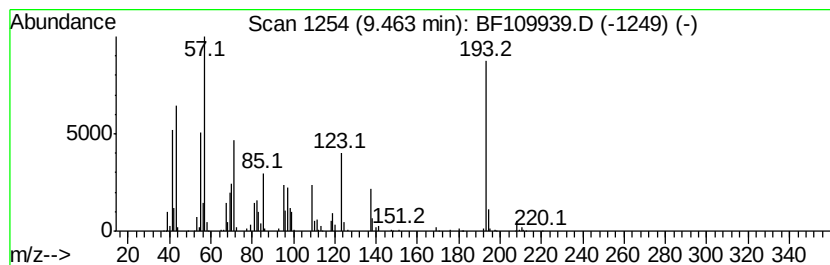
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown9.46 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.46	28.31 ng	6523790	Acenaphthene-d10	10.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,10-Trimethylundeca-1,3-diene			194	C14H26	1000222-09-1	18
2	Nonadecane, 1-bromo-			346	C19H39Br	004434-66-6	14
3	1,5-Methano-8H-pyrido[1,2-a][1,5...			234	C14H22N2O	000550-43-6	14
4	3,4-Undecadiene-2,10-dione, 6,6-...			208	C13H20O2	052588-78-0	14
5	Cyclohexanol, 2-(2-ethylhexyl)-			212	C14H28O	077189-05-0	11



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101718\
Data File : BF109939.D
Acq On : 17 Oct 2018 23:56
Operator : JU/SJ
Sample : J5539-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

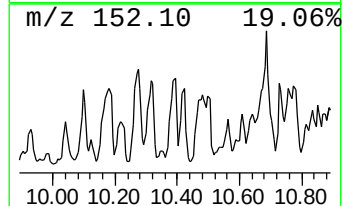
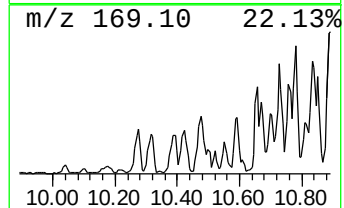
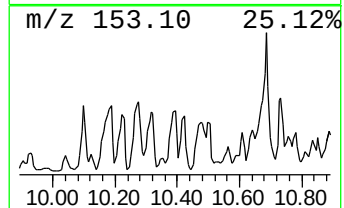
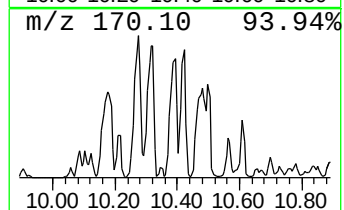
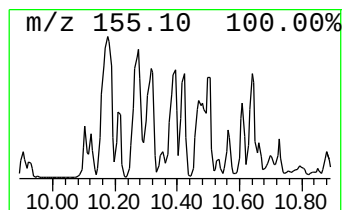
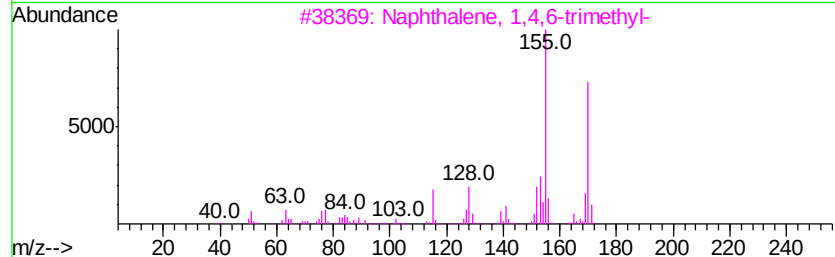
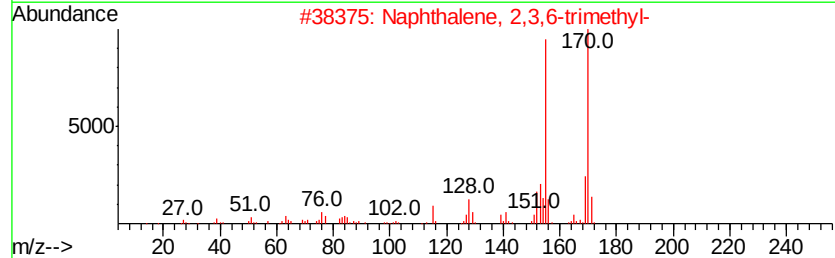
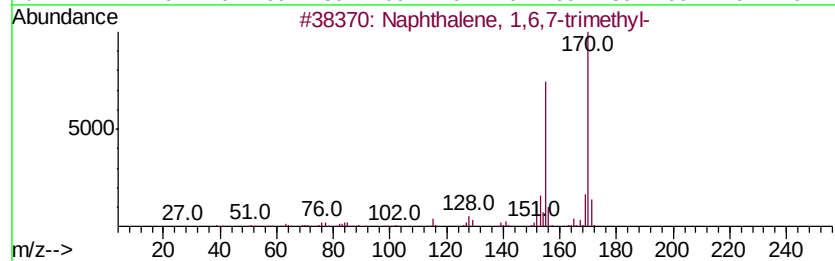
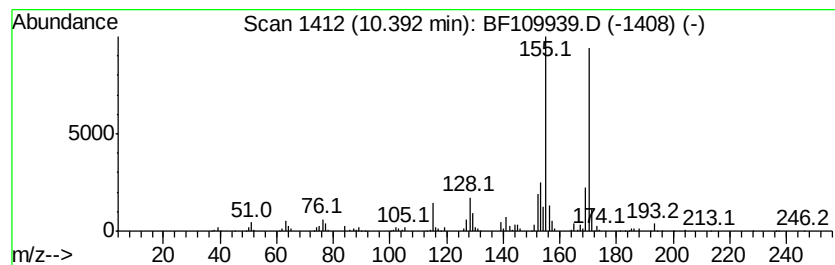
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ClientSampleId :
SB-1-6-8FT

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Naphthalene, 1,6,7-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.39	20.00 ng	4609040	Acenaphthene-d10	10.06	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
2	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	96
3	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	94
4	Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	91
5	4,6,8-Trimethylazulene	170	C13H14	000941-81-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101718\
Data File : BF109939.D
Acq On : 17 Oct 2018 23:56
Operator : JU/SJ
Sample : J5539-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

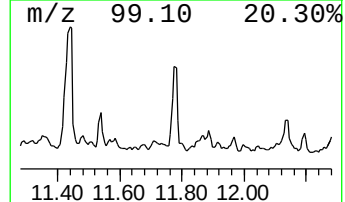
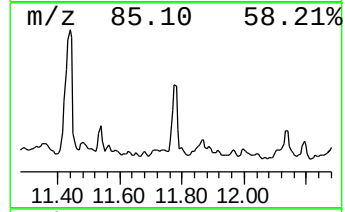
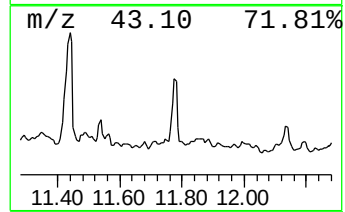
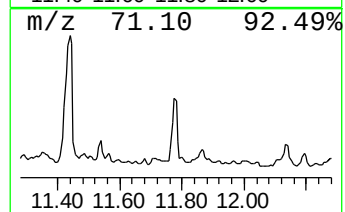
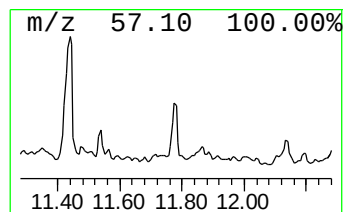
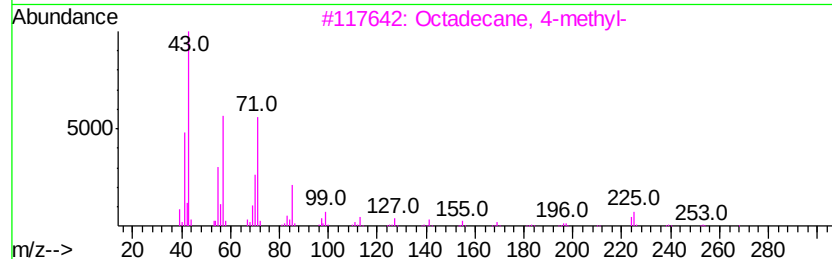
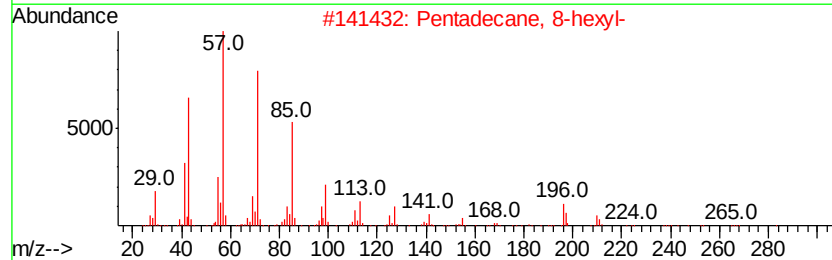
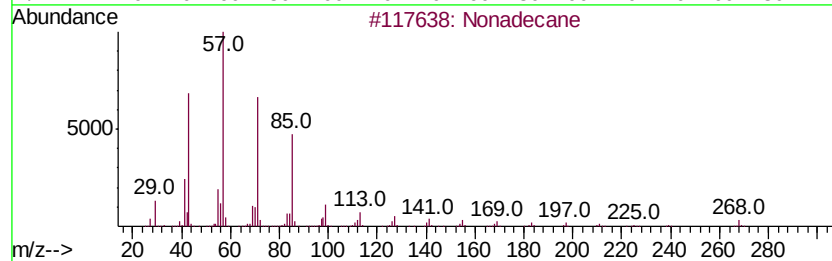
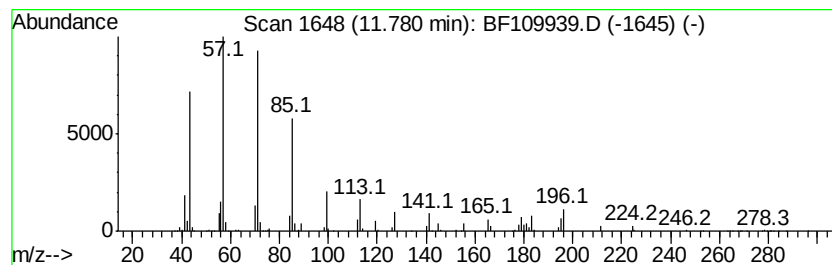
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Nonadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.78	20.00 ng	3344430	Phenanthrene-d10		11.56	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	90
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	90
3		Octadecane, 4-methyl-	268	C19H40	010544-95-3	87
4		Tridecane, 7-hexyl-	268	C19H40	007225-66-3	87
5		Octadecane, 2-methyl-	268	C19H40	001560-88-9	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101718\
Data File : BF109939.D
Acq On : 17 Oct 2018 23:56
Operator : JU/SJ
Sample : J5539-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-1-6-8FT

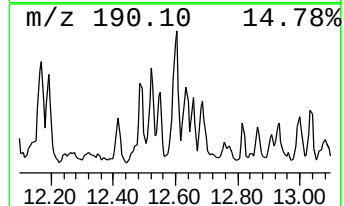
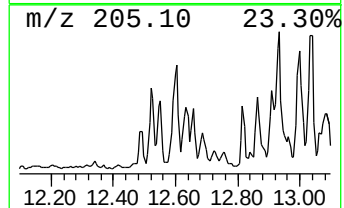
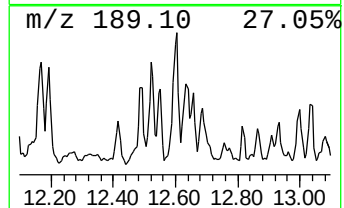
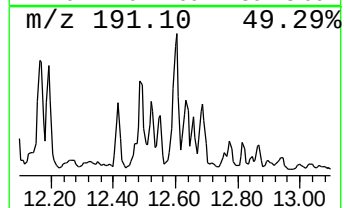
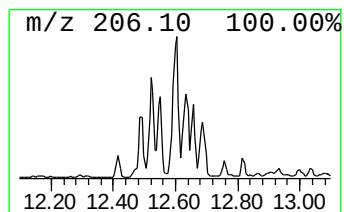
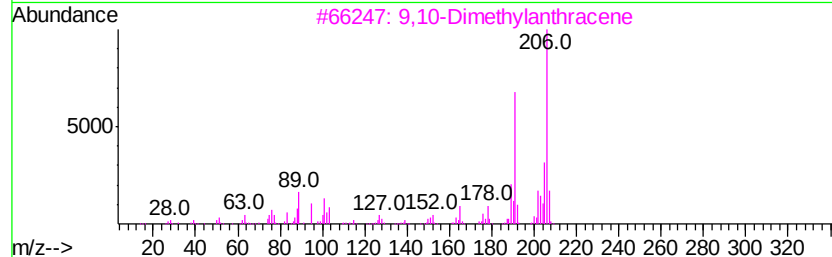
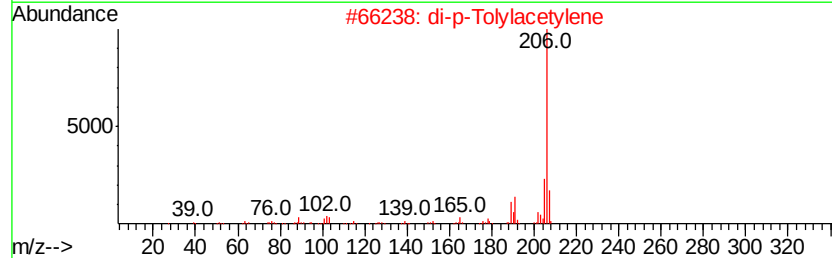
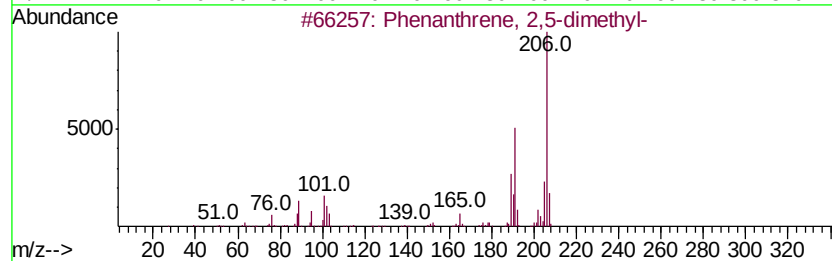
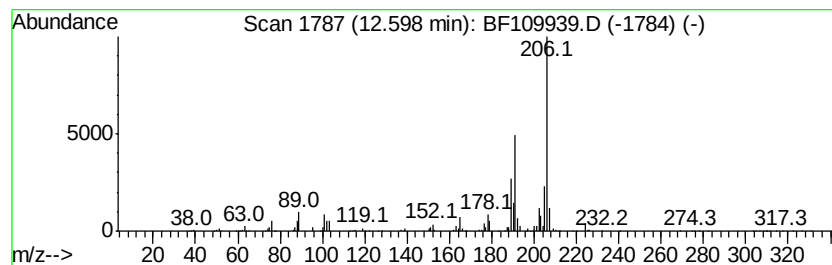
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Phenanthrene, 2,5-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.60	22.51 ng	3764400	Phenanthrene-d10	11.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	94
2		di-p-Tolylacetylene	206	C16H14	002789-88-0	93
3		9,10-Dimethylantracene	206	C16H14	000781-43-1	91
4		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	90
5		Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	89



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF101718\
 Data File : BF109939.D
 Acq On : 17 Oct 2018 23:56
 Operator : JU/SJ
 Sample : J5539-05
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-1-6-8FT

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Nonane, 3-methyl-	6.23	18.4	ng	10817400	1	6.98	11728400	20.0
Cyclohexane, 1,1,...	6.52	18.2	ng	10655900	1	6.98	11728400	20.0
Cyclohexane, butyl-	7.13	15.4	ng	9054460	1	6.98	11728400	20.0
Naphthalene, deca...	7.39	15.8	ng	9240470	1	6.98	11728400	20.0
unknown7.63	7.63	19.4	ng	11376400	1	6.98	11728400	20.0
Decane, 3,8-dimet...	7.69	46.7	ng	11787900	2	8.28	5043510	20.0
Benzene, 1-ethyl-...	7.76	21.3	ng	5378020	2	8.28	5043510	20.0
Naphthalene, deca...	7.80	35.7	ng	8994280	2	8.28	5043510	20.0
unknown7.87	7.87	15.4	ng	3889500	2	8.28	5043510	20.0
unknown7.92	7.92	25.5	ng	6420480	2	8.28	5043510	20.0
Benzene, 1,2,4,5-...	8.02	25.7	ng	6477470	2	8.28	5043510	20.0
unknown8.09	8.09	18.6	ng	4681030	2	8.28	5043510	20.0
Benzene, 1-methyl...	8.15	22.4	ng	5655200	2	8.28	5043510	20.0
Naphthalene, 1,2,...	8.48	23.1	ng	5822370	2	8.28	5043510	20.0
3-Cyclohexyl-1-ch...	8.57	34.9	ng	8797050	2	8.28	5043510	20.0
Naphthalene, 1,2,...	8.79	21.4	ng	5398120	2	8.28	5043510	20.0
unknown9.46	9.46	28.3	ng	6523790	3	10.06	4609040	20.0
Naphthalene, 1,6,...	10.39	20.0	ng	4609040	3	10.06	4609040	20.0
Nonadecane	11.78	20.0	ng	3344430	4	11.56	3344430	20.0
Phenanthrene, 2,5...	12.60	22.5	ng	3764400	4	11.56	3344430	20.0