

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040416\
 Data File : BF086052.D
 Acq On : 4 Apr 2016 22:10
 Operator : UM/SJ
 Sample : H2054-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-B11COMP

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:\HPCHEM1\BNA_F\METHODS\8270-BF040216.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	2.612	39	46	50	rVB	30887	63361	3.27%	0.208%
2	3.595	129	132	141	rVB	91920	176293	9.11%	0.579%
3	4.281	188	192	200	rBV2	29473	60270	3.11%	0.198%
4	4.841	238	241	244	rVB2	31214	44077	2.28%	0.145%
5	4.921	247	248	254	rVB	73314	118115	6.10%	0.388%
6	5.115	263	265	267	rBV	28642	29567	1.53%	0.097%
7	5.195	270	272	274	rBV	38482	58710	3.03%	0.193%
8	5.310	279	282	285	rBV	136133	236544	12.22%	0.777%
9	5.367	285	287	296	rVV	1073291	1694715	87.54%	5.566%
10	5.493	296	298	301	rVV	24083	43611	2.25%	0.143%
11	5.573	303	305	310	rVV	180116	206643	10.67%	0.679%
12	5.641	310	311	314	rVV	42348	53374	2.76%	0.175%
13	5.744	317	320	323	rVB2	32810	47649	2.46%	0.156%
14	5.893	331	333	336	rBV2	54984	76411	3.95%	0.251%
15	5.996	339	342	345	rVB2	67152	89141	4.60%	0.293%
16	6.053	345	347	348	rBV	37725	36696	1.90%	0.121%
17	6.190	356	359	360	rBV2	32429	56868	2.94%	0.187%
18	6.213	360	361	363	rVV2	56937	59372	3.07%	0.195%
19	6.281	365	367	372	rVV	96792	137101	7.08%	0.450%
20	6.350	372	373	376	rVV2	30093	49755	2.57%	0.163%
21	6.418	376	379	384	rVV	1310600	1814323	93.72%	5.959%
22	6.533	387	389	392	rVV	1821613	1759437	90.88%	5.779%
23	6.578	392	393	397	rVB2	141132	235610	12.17%	0.774%
24	6.716	401	405	407	rBV2	33028	66745	3.45%	0.219%
25	6.761	407	409	412	rBV	789008	734652	37.95%	2.413%
26	6.819	412	414	415	rBV	75159	53732	2.78%	0.176%
27	6.921	420	423	425	rBV	1200119	1473292	76.10%	4.839%
28	7.036	430	433	435	rVB2	49779	84052	4.34%	0.276%
29	7.081	435	437	441	rBV2	62255	122522	6.33%	0.402%
30	7.150	441	443	447	rBV3	70774	115824	5.98%	0.380%
31	7.230	447	450	451	rBV2	28436	33746	1.74%	0.111%
32	7.299	453	456	457	rBV3	46591	54550	2.82%	0.179%
33	7.333	457	459	461	rVV	1123694	1142665	59.02%	3.753%
34	7.367	461	462	463	rVB	162271	111237	5.75%	0.365%

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35	7.413	463	466	468	rBV4	47763	51497	2.66%	0.169%
36	7.481	470	472	474	rVB3	29130	41921	2.17%	0.138%
37	7.561	474	479	481	rVB3	45244	124450	6.43%	0.409%
38	7.664	481	488	490	rBV4	35413	89015	4.60%	0.292%
39	7.801	496	500	503	rBV4	58154	130136	6.72%	0.427%
40	7.870	503	506	509	rVB4	42942	77316	3.99%	0.254%
41	7.996	515	517	519	rBV2	28127	57905	2.99%	0.190%
42	8.041	519	521	526	rVV2	684409	1321032	68.24%	4.339%
43	8.110	526	527	531	rVB	94133	108299	5.59%	0.356%
44	8.179	531	533	535	rBV2	23998	38516	1.99%	0.127%
45	8.339	544	547	551	rVV5	67619	104372	5.39%	0.343%
46	8.430	553	555	556	rVV	74777	66651	3.44%	0.219%
47	8.476	556	559	561	rVV	179552	206062	10.64%	0.677%
48	8.647	572	574	576	rBV	170352	207917	10.74%	0.683%
49	8.762	582	584	586	rVB	458802	355675	18.37%	1.168%
50	8.853	590	592	594	rBV	210915	289635	14.96%	0.951%
51	8.956	599	601	603	rBV3	30626	42556	2.20%	0.140%
52	9.002	603	605	607	rVB2	49125	66479	3.43%	0.218%
53	9.047	607	609	610	rBV2	23726	35528	1.84%	0.117%
54	9.070	610	611	613	rVB	139062	106461	5.50%	0.350%
55	9.127	613	616	618	rBV	1824607	1935999	100.00%	6.359%
56	9.207	621	623	626	rBV2	352484	346140	17.88%	1.137%
57	9.322	631	633	637	rVB	68397	88049	4.55%	0.289%
58	9.390	637	639	641	rBV	125480	168139	8.68%	0.552%
59	9.459	643	645	646	rBV	225138	191645	9.90%	0.629%
60	9.482	646	647	649	rVV	134212	135284	6.99%	0.444%
61	9.527	649	651	652	rVB2	296620	375218	19.38%	1.232%
62	9.573	654	655	660	rVB3	107896	165183	8.53%	0.543%
63	9.665	660	663	665	rVB2	52499	68223	3.52%	0.224%
64	9.699	665	666	668	rBV	21704	29566	1.53%	0.097%
65	9.733	668	669	671	rVB	327656	256043	13.23%	0.841%
66	9.802	671	675	676	rBV	791887	899663	46.47%	2.955%
67	9.836	676	678	682	rVB3	64802	110980	5.73%	0.365%
68	9.905	682	684	686	rVB	80264	96464	4.98%	0.317%
69	9.962	686	689	691	rBV4	88727	168140	8.68%	0.552%
70	10.007	691	693	695	rBV2	102816	145731	7.53%	0.479%
71	10.042	695	696	701	rVB3	100176	170671	8.82%	0.561%

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72	10.122	701	703	704 rBV	78484	122875	6.35%	0.404%
73	10.202	708	710	711 rBV2	138434	151741	7.84%	0.498%
74	10.236	711	713	714 rVB	277837	277447	14.33%	0.911%
75	10.293	714	718	720 rBV4	30562	79291	4.10%	0.260%
76	10.339	720	722	723 rBV	157323	145734	7.53%	0.479%
77	10.453	729	732	735 rBV	148152	211366	10.92%	0.694%
78	10.499	735	736	738 rVB	86824	99233	5.13%	0.326%
79	10.533	738	739	741 rBV2	37018	61793	3.19%	0.203%
80	10.590	741	744	746 rVV	718460	758805	39.19%	2.492%
81	10.625	746	747	750 rVB2	44364	59171	3.06%	0.194%
82	10.705	752	754	758 rVV	364869	576254	29.77%	1.893%
83	10.899	769	771	772 rBV2	51287	66307	3.42%	0.218%
84	11.025	780	782	786 rBV4	76148	141682	7.32%	0.465%
85	11.150	791	793	794 rVV	260584	233926	12.08%	0.768%
86	11.185	794	796	798 rVB2	135476	166222	8.59%	0.546%
87	11.288	802	805	808 rBV	524068	1057927	54.65%	3.475%
88	11.573	829	830	832 rBV	146805	174601	9.02%	0.573%
89	11.813	850	851	853 rVB	263471	202788	10.47%	0.666%
90	11.893	856	858	859 rBV	209982	198167	10.24%	0.651%
91	11.985	864	866	868 rBV	240788	243190	12.56%	0.799%
92	12.339	895	897	898 rBV	188868	262643	13.57%	0.863%
93	12.373	898	900	901 rVV	289354	284862	14.71%	0.936%
94	12.499	909	911	913 rBV	382292	385445	19.91%	1.266%
95	12.728	928	931	934 rBV2	294455	670808	34.65%	2.203%
96	12.865	941	943	947 rVB	1084979	1195990	61.78%	3.928%
97	13.105	962	964	966 rBV2	240064	295898	15.28%	0.972%
98	13.448	992	994	996 rBV2	273968	403822	20.86%	1.326%
99	15.962	1212	1214	1222 rVB	404317	1153971	59.61%	3.790%
100	16.362	1247	1249	1254 rVB	418952	751760	38.83%	2.469%

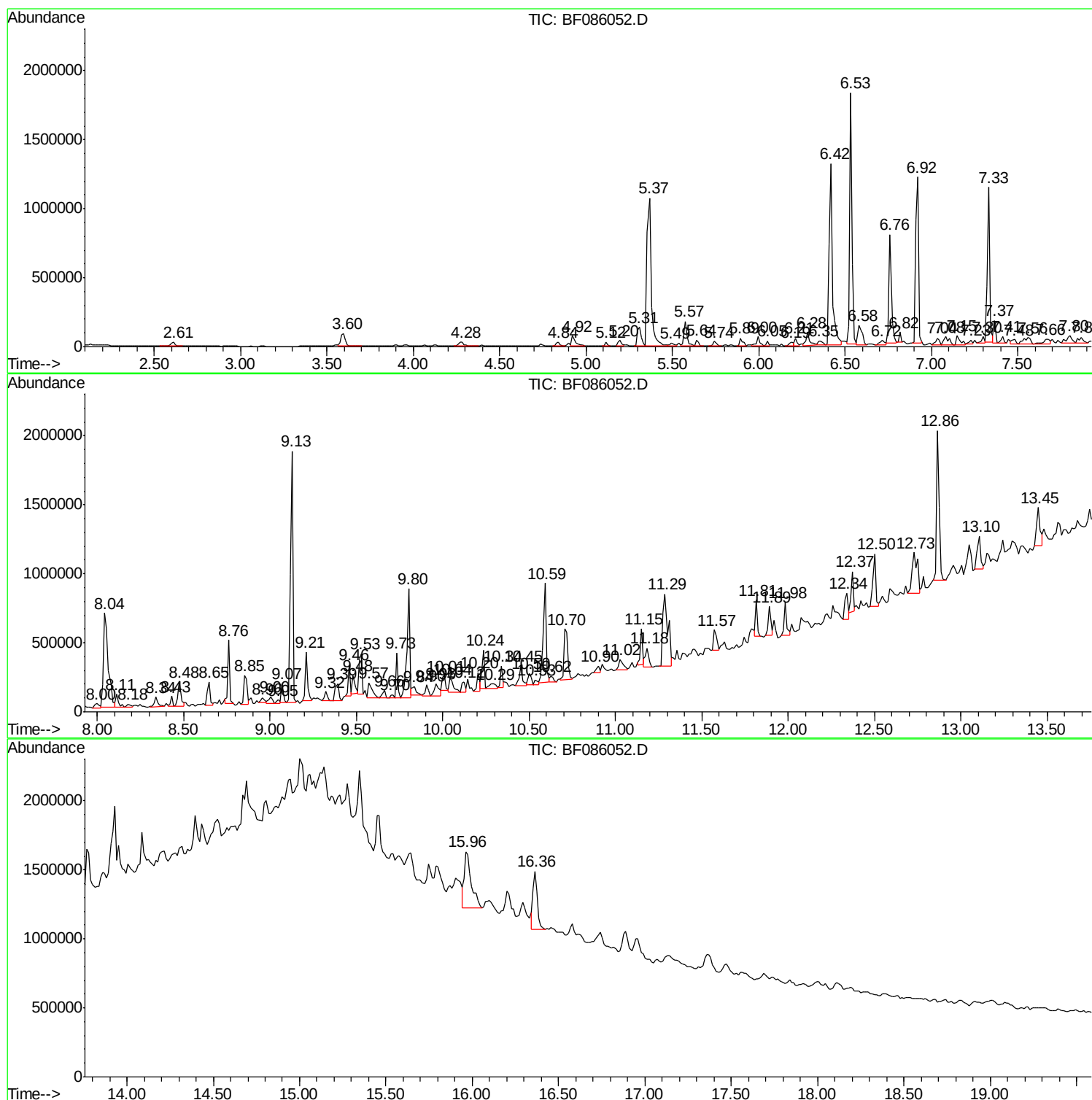
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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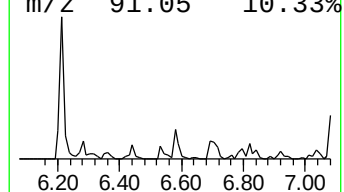
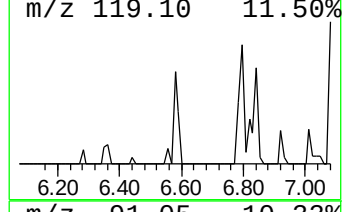
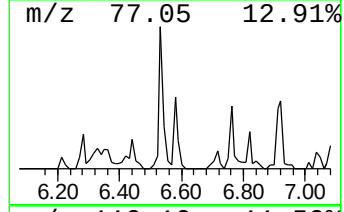
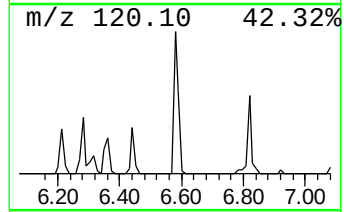
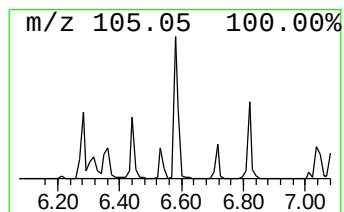
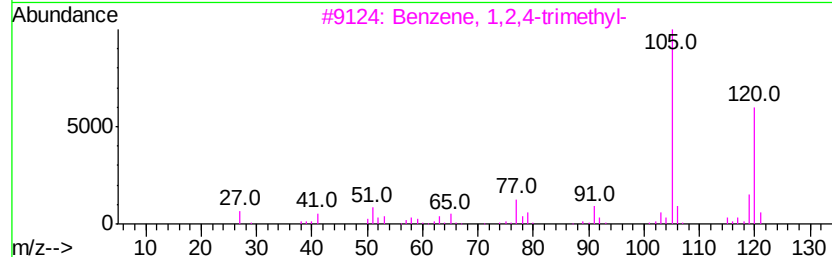
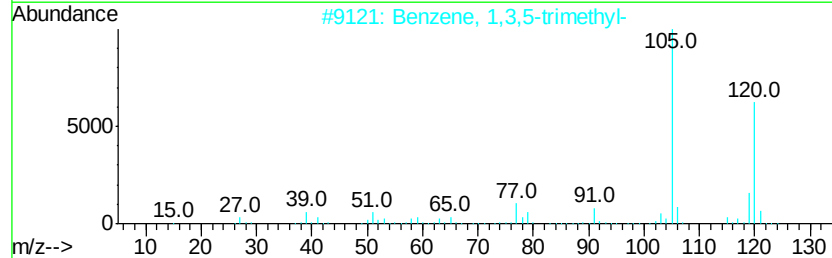
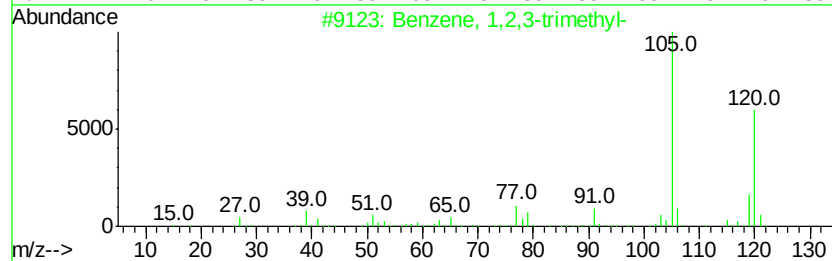
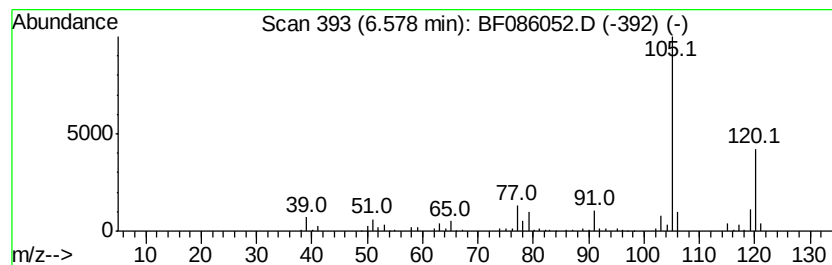
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	6.41 ng	235610	1,4-Dichlorobenzene-d4	6.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	95
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90



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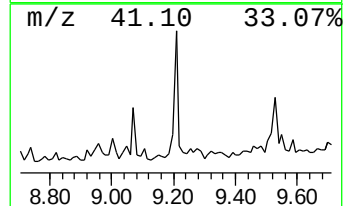
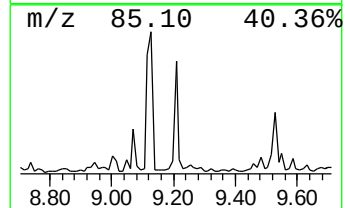
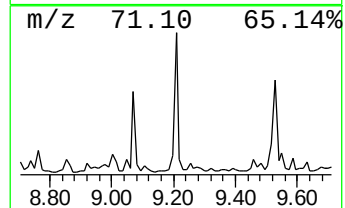
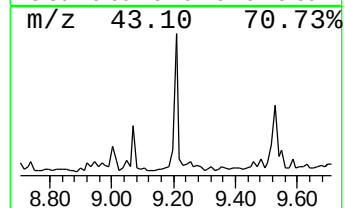
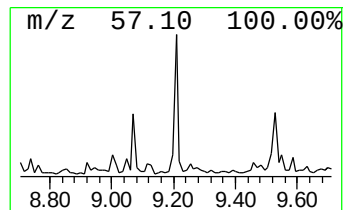
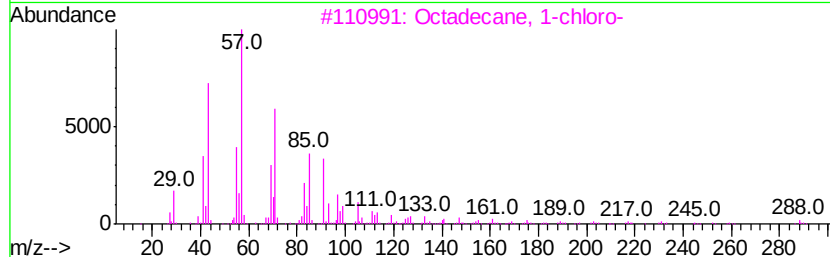
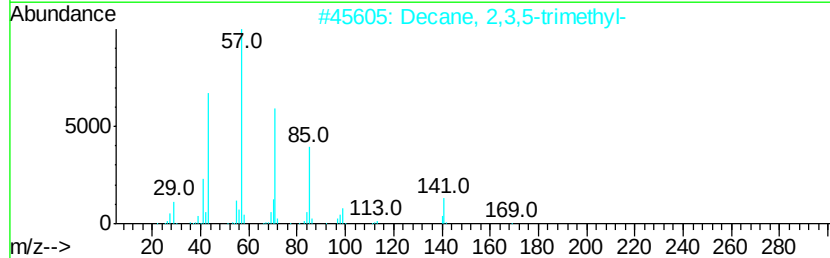
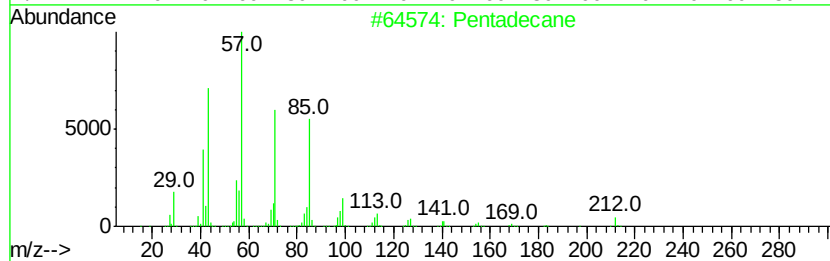
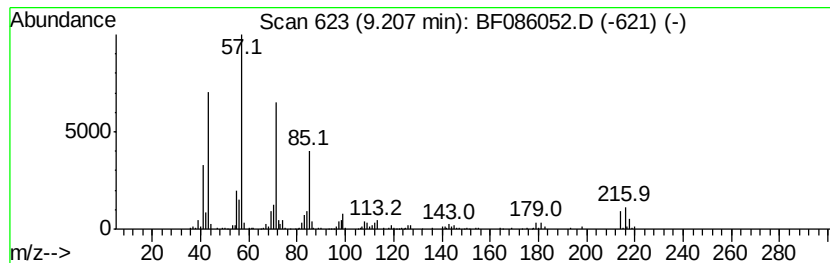
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Pentadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.21	7.69 ng	346140	Acenaphthene-d10	9.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane	212	C15H32	000629-62-9	64
2		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	64
3		Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	59
4		Decane, 2,3,6-trimethyl-	184	C13H28	062238-12-4	59
5		Tridecane	184	C13H28	000629-50-5	59



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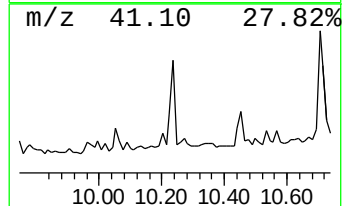
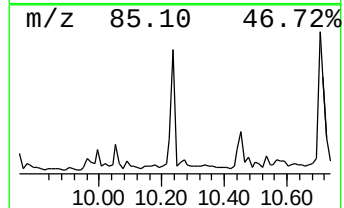
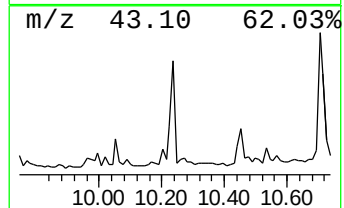
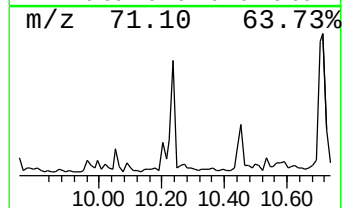
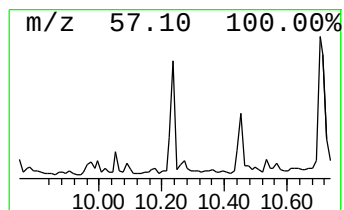
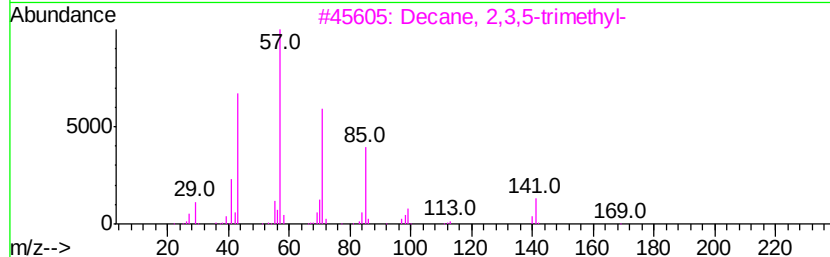
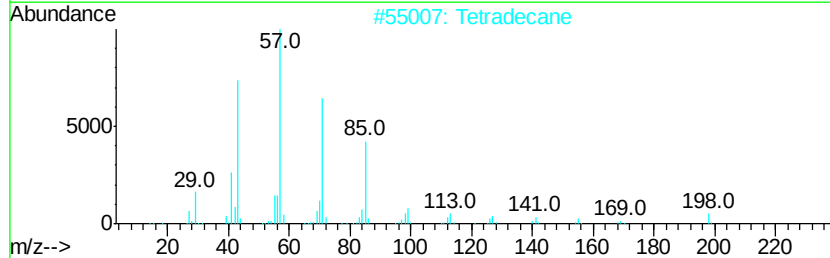
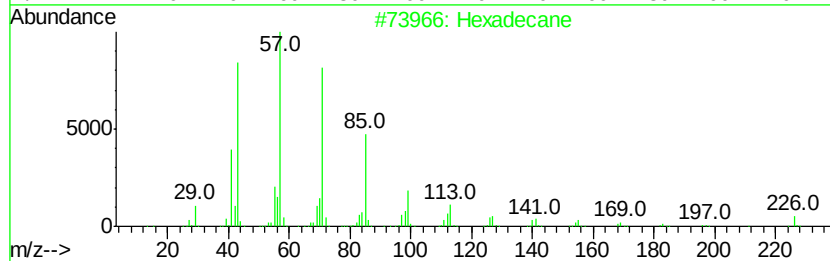
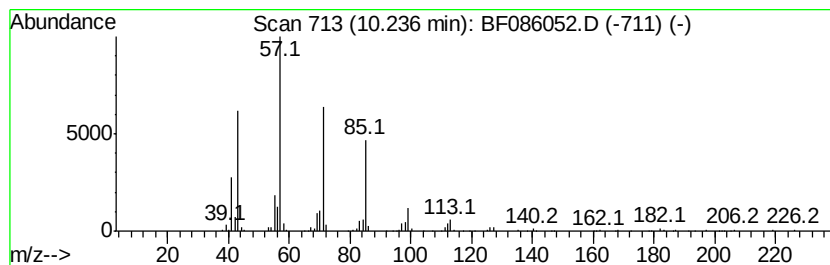
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.24	6.17 ng	277447	Acenaphthene-d10	9.80	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	94
2	Tetradecane	198	C14H30	000629-59-4	90
3	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	90
4	Pentadecane	212	C15H32	000629-62-9	86
5	Heptadecane	240	C17H36	000629-78-7	83



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ALS Vial : 18 Sample Multiplier: 1

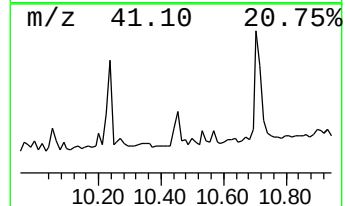
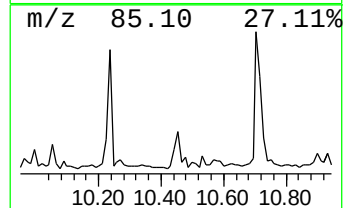
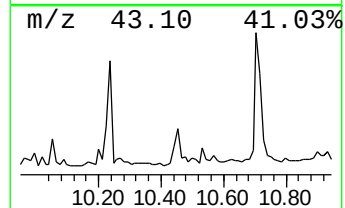
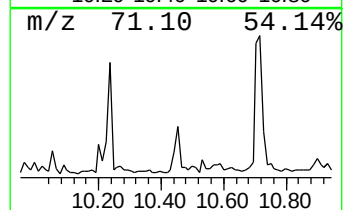
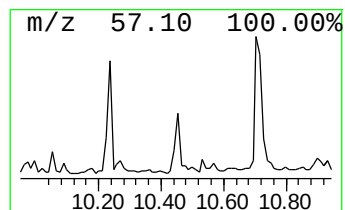
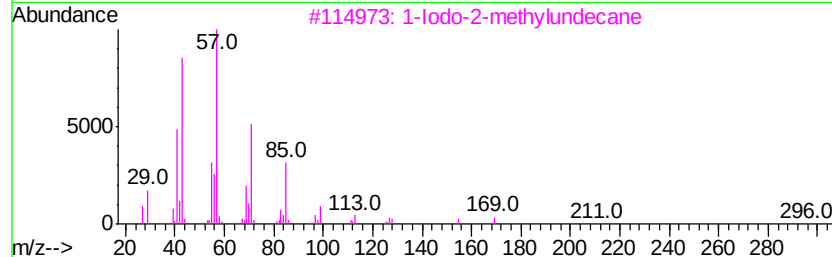
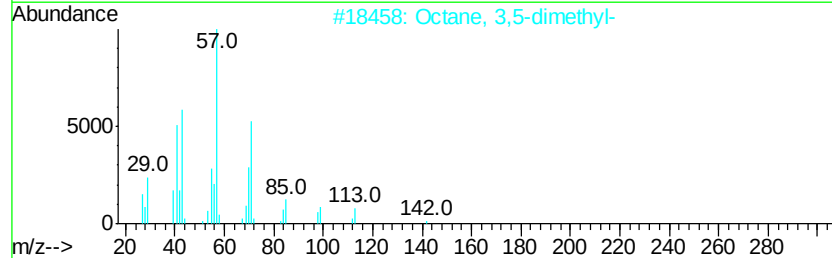
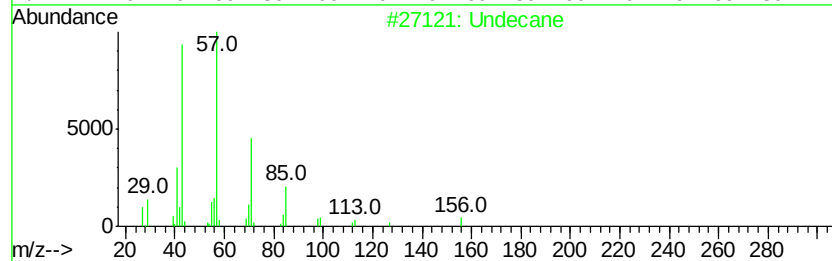
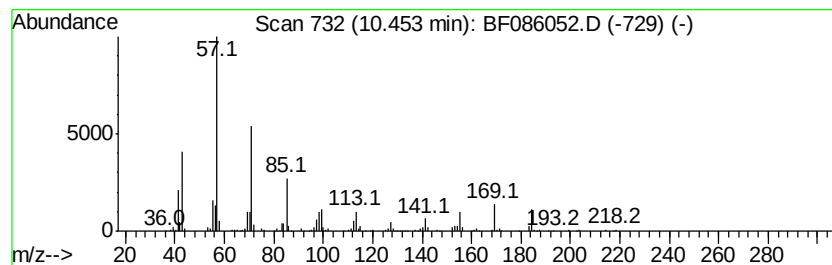
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Undecane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.45	4.70 ng	211366	Acenaphthene-d10	9.80
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Undecane		156 C11H24	001120-21-4 68
2	Octane, 3,5-dimethyl-		142 C10H22	015869-93-9 58
3	1-Iodo-2-methylundecane		296 C12H25I	073105-67-6 53
4	Dodecane, 3-methyl-		184 C13H28	017312-57-1 53
5	Dodecane, 1-iodo-		296 C12H25I	004292-19-7 52



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040416\
Data File : BF086052.D
Acq On : 4 Apr 2016 22:10
Operator : UM/SJ
Sample : H2054-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

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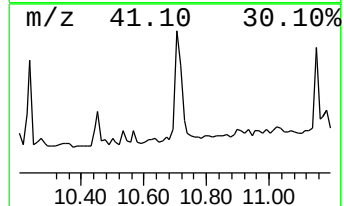
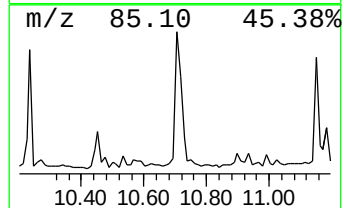
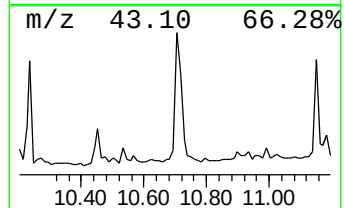
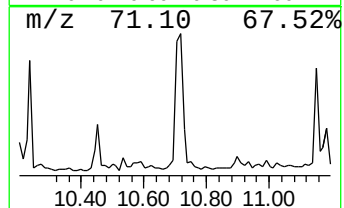
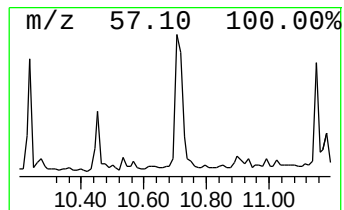
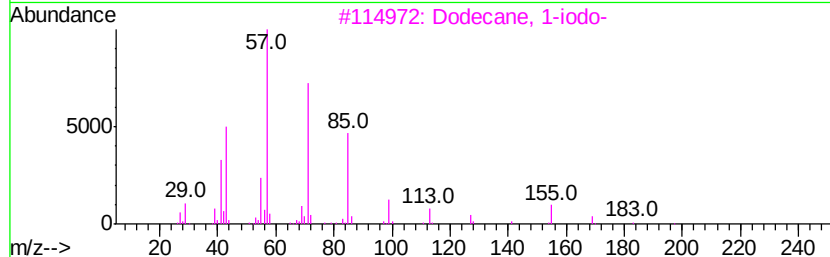
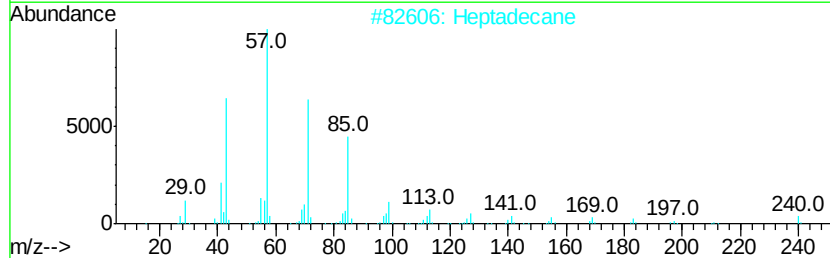
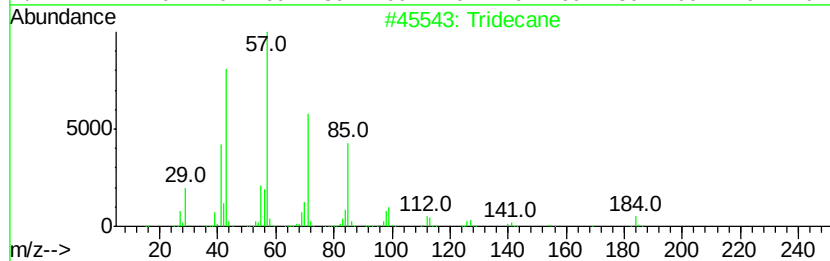
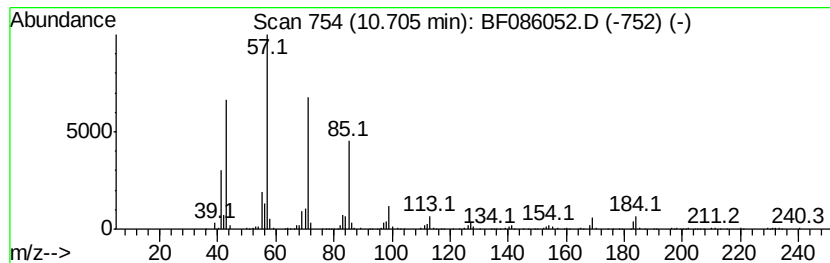
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tridecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.70	10.89 ng	576254	Phenanthrene-d10	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane	184	C13H28	000629-50-5	87
2		Heptadecane	240	C17H36	000629-78-7	87
3		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	83
4		Pentadecane	212	C15H32	000629-62-9	80
5		Hexadecane	226	C16H34	000544-76-3	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040416\
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Sample : H2054-02
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ALS Vial : 18 Sample Multiplier: 1

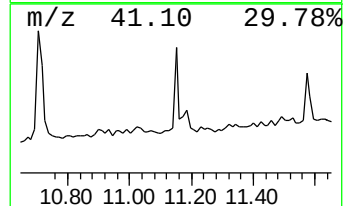
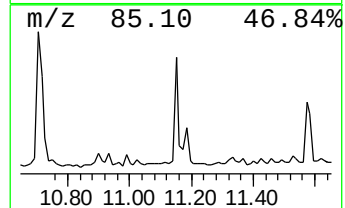
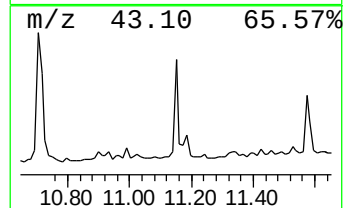
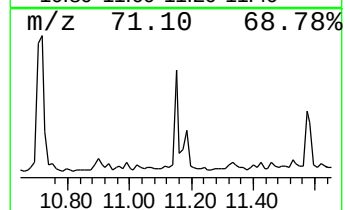
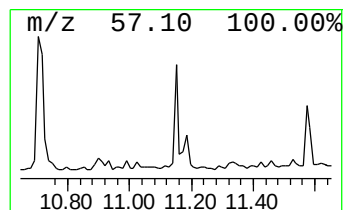
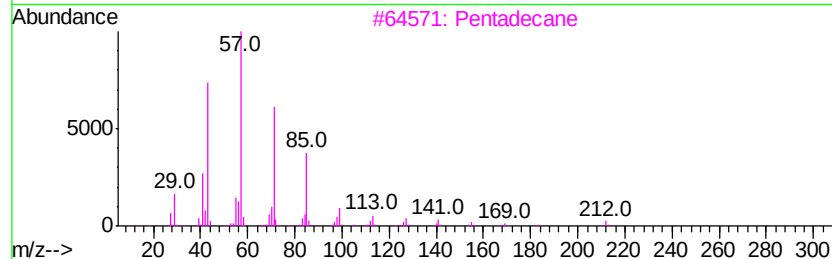
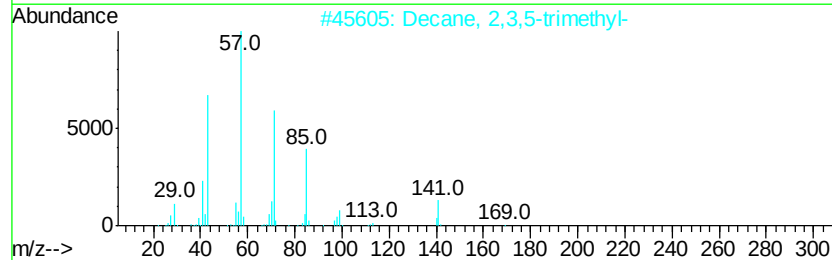
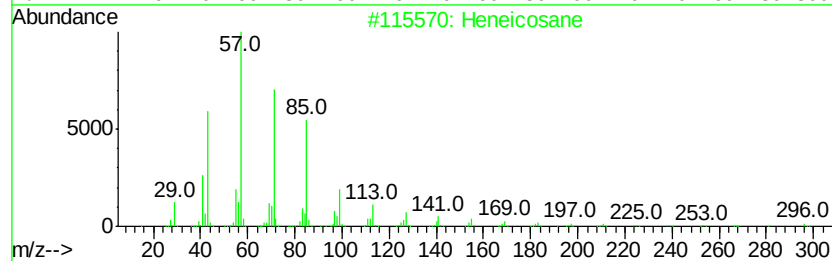
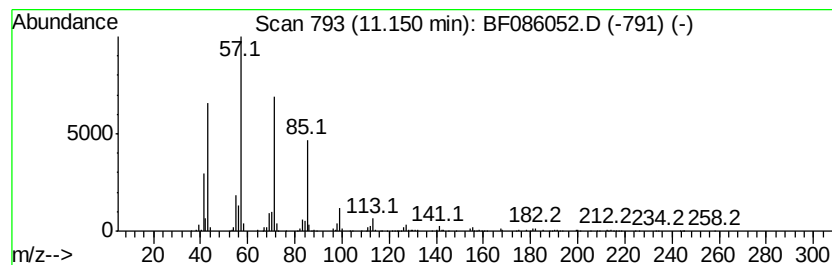
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heneicosane Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.15	4.42 ng	233926	Phenanthrene-d10		11.29
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	90
2	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	90
3	Pentadecane	212	C15H32	000629-62-9	90
4	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	80
5	Eicosane	282	C20H42	000112-95-8	80



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040416\
Data File : BF086052.D
Acq On : 4 Apr 2016 22:10
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Misc :
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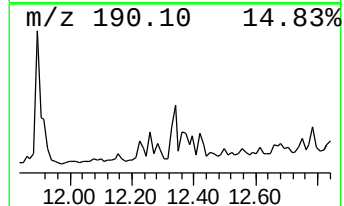
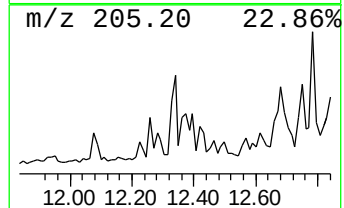
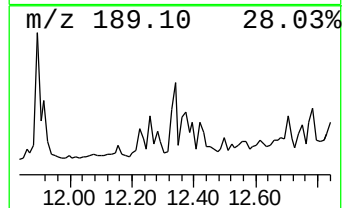
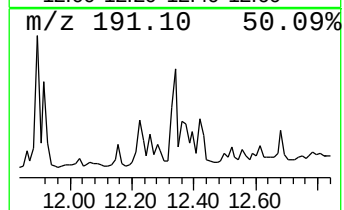
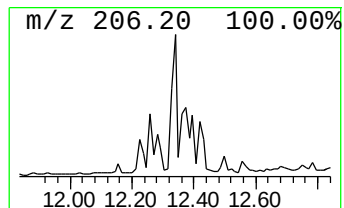
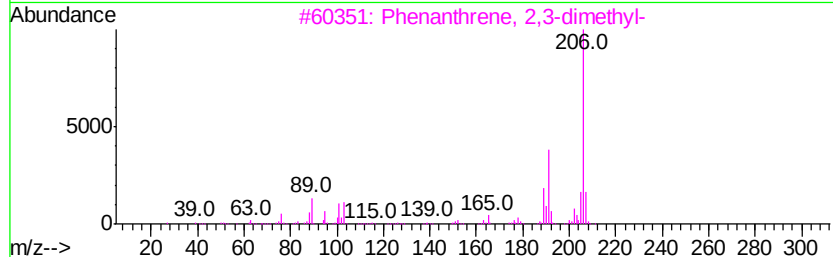
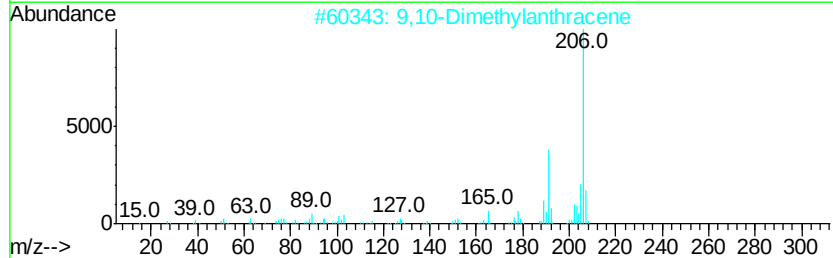
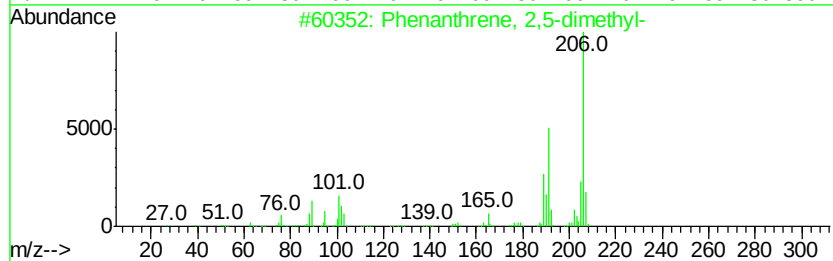
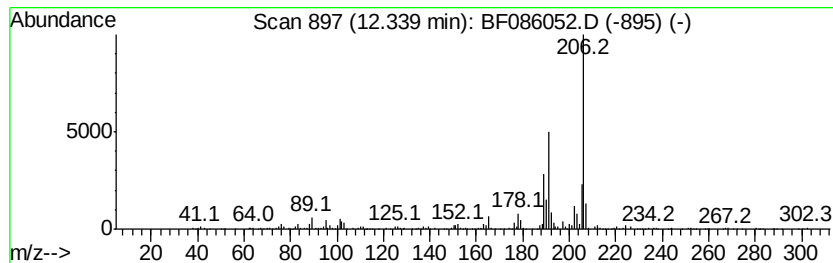
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Phenanthrene, 2,5-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.34	4.97 ng	262643	Phenanthrene-d10	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2		9,10-Dimethylantracene	206	C16H14	000781-43-1	89
3		Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	87
4		di-p-Tolylacetylene	206	C16H14	002789-88-0	83
5		1,3-Diphenyl-3-methylcyclopropene	206	C16H14	086544-79-8	74



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040416\
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Sample : H2054-02
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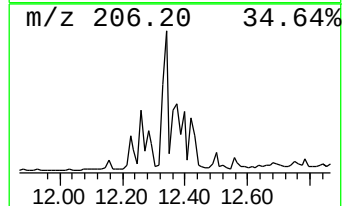
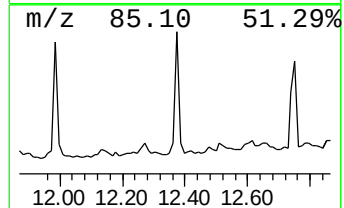
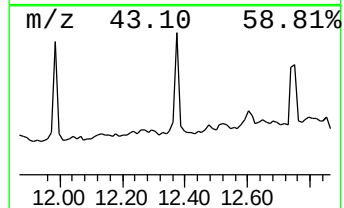
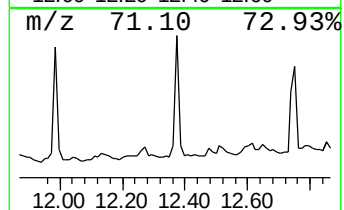
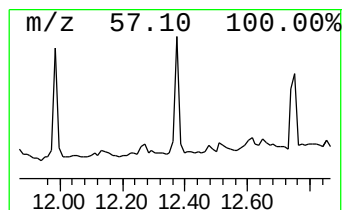
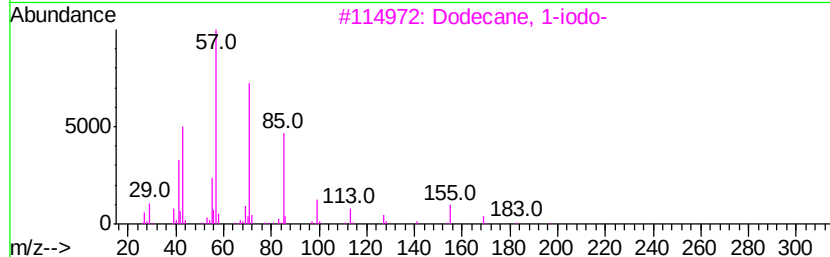
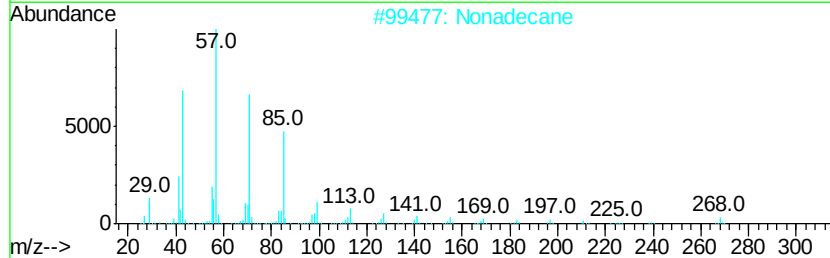
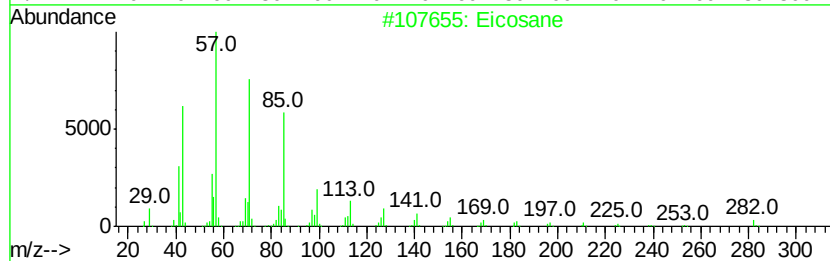
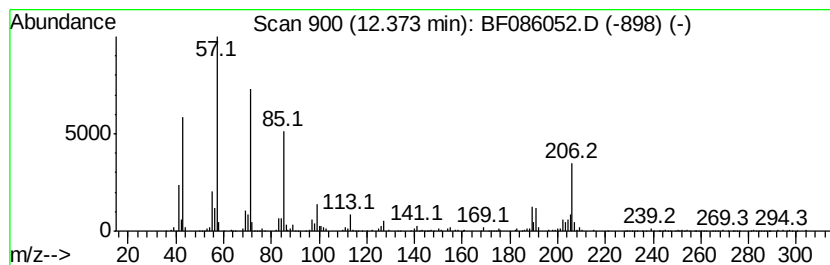
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Eicosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.37	5.39 ng	284862	Phenanthrene-d10		11.29
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	83
2	Nonadecane	268	C19H40	000629-92-5	52
3	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	52
4	Heneicosane	296	C21H44	000629-94-7	52
5	1-Iodoundecane	282	C11H23I	004282-44-4	52



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040416\
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Sample : H2054-02
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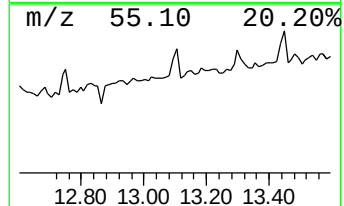
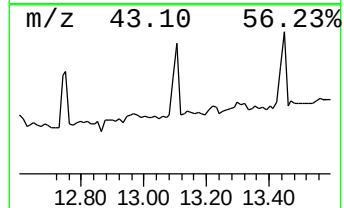
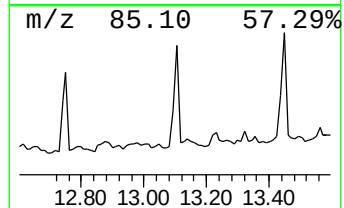
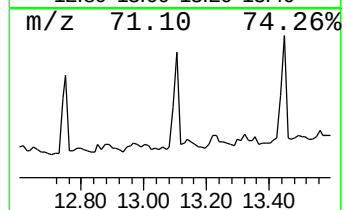
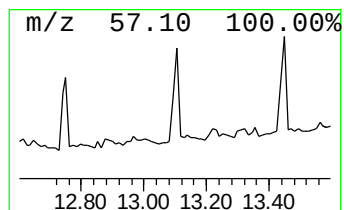
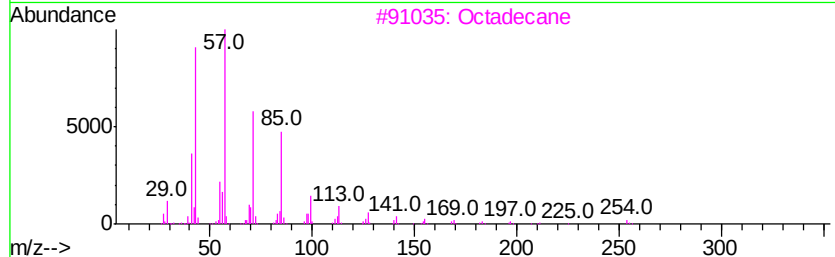
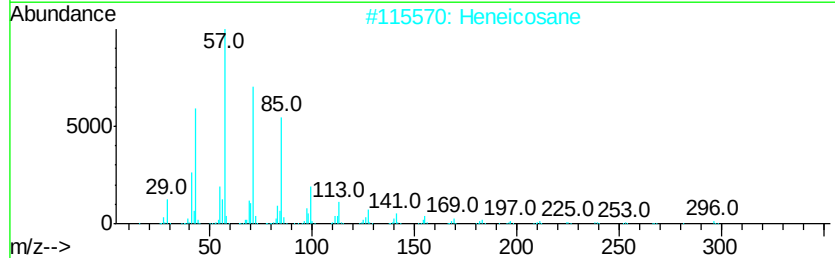
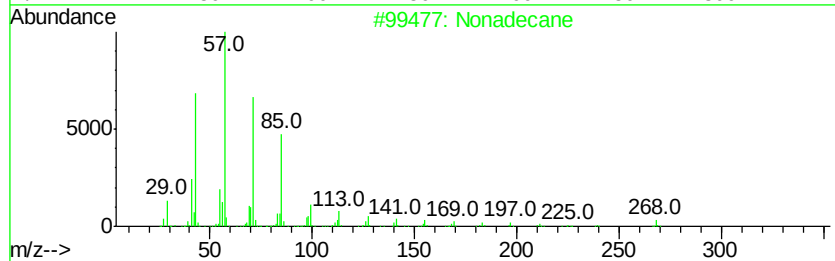
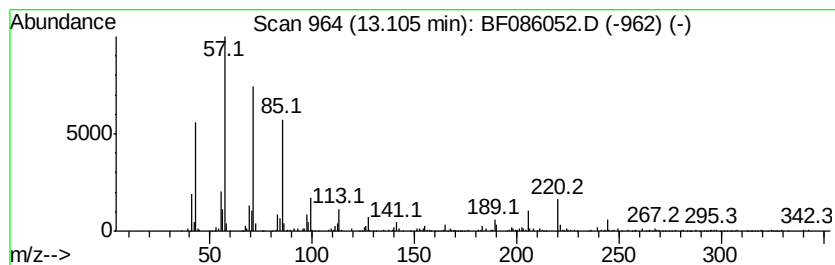
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Nonadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.10	14.65 ng	295898	Chrysene-d12	13.93	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	95
2	Heneicosane	296	C21H44	000629-94-7	87
3	Octadecane	254	C18H38	000593-45-3	83
4	Eicosane	282	C20H42	000112-95-8	83
5	Heptadecane	240	C17H36	000629-78-7	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040416\
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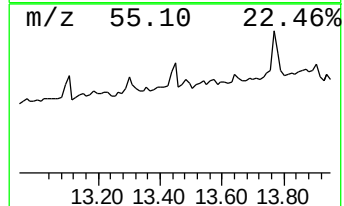
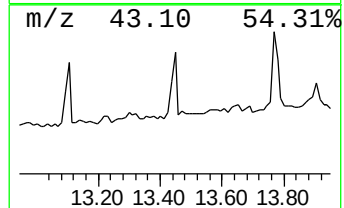
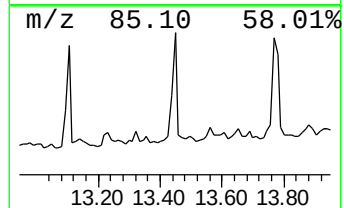
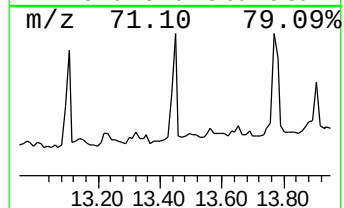
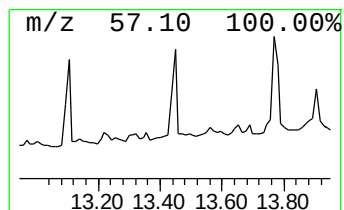
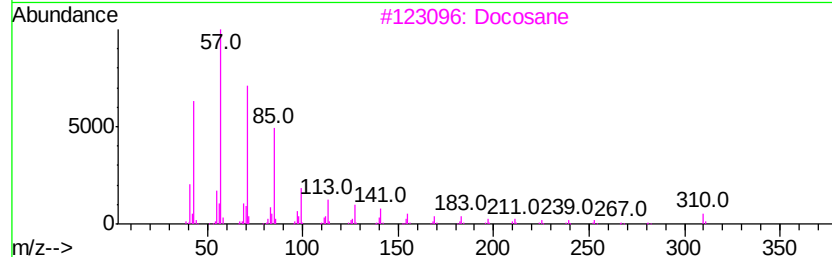
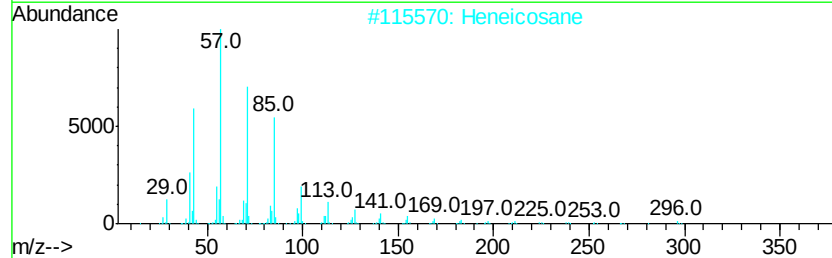
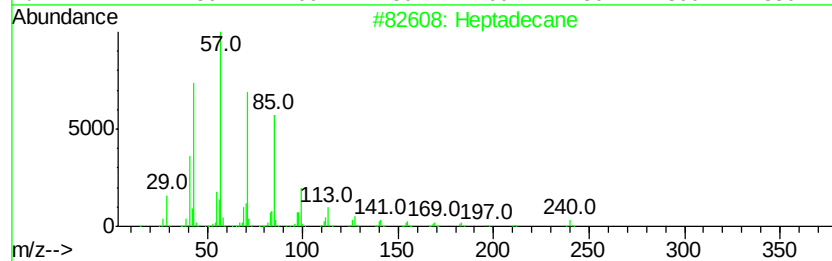
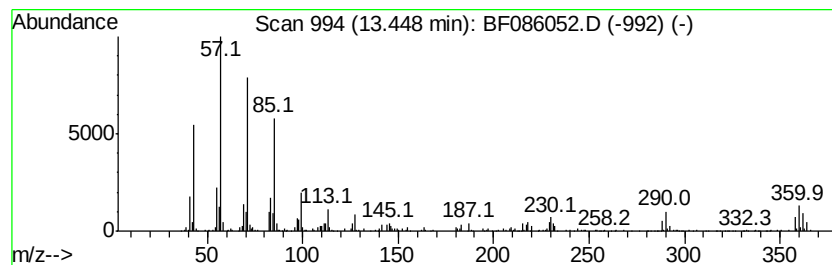
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Heptadecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.45	20.00 ng	403822	Chrysene-d12	13.93	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	70
2	Heneicosane	296	C21H44	000629-94-7	62
3	Docosane	310	C22H46	000629-97-0	62
4	Hexadecane	226	C16H34	000544-76-3	58
5	Pentacosane	352	C25H52	000629-99-2	58



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040416\
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ALS Vial : 18 Sample Multiplier: 1

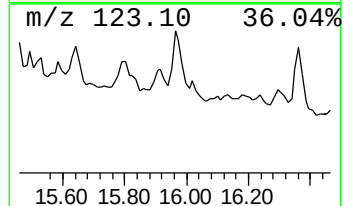
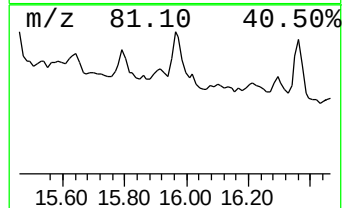
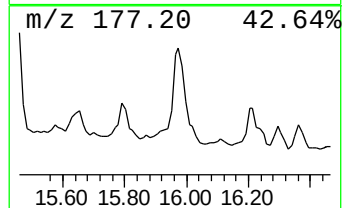
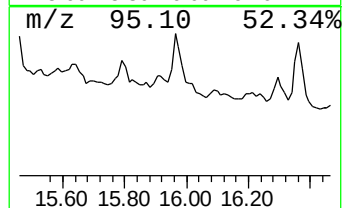
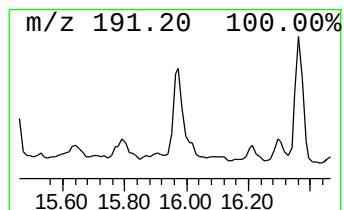
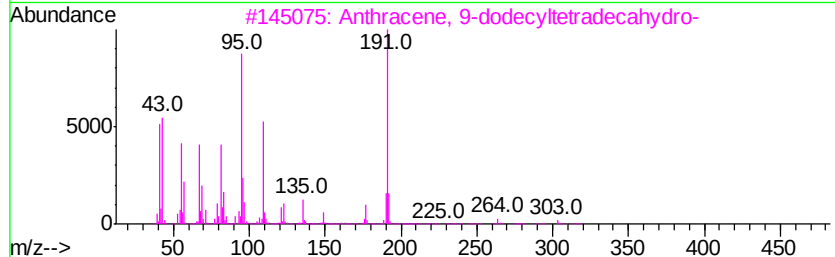
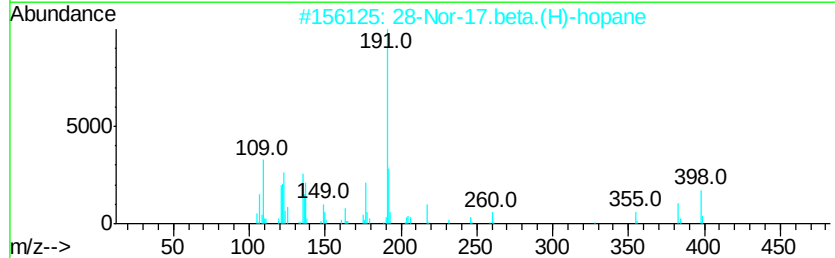
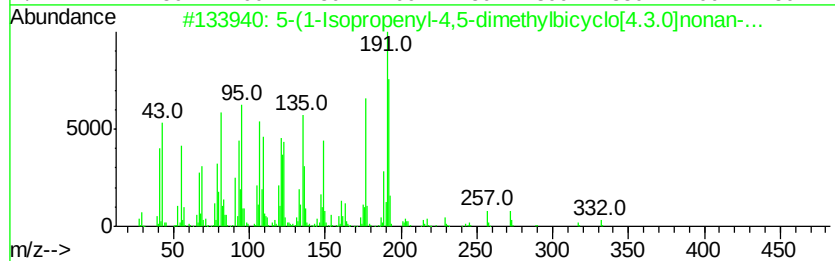
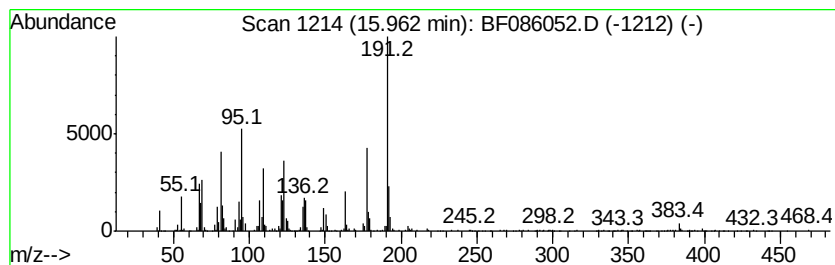
Instrument :
BNA_F
ClientSampleId :
TEC-B11COMP

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 5-(1-Isopropenyl-4,5-dimeth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
15.96	20.00 ng	1153970	Perylene-d12	15.35		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-(1-Isopropenyl-4,5-dimethylbic...	332	C22H36O2	1000195-69-8	50	
2	28-Nor-17.beta.(H)-hopane	398	C29H50	036728-72-0	42	
3	Anthracene, 9-dodecyltetradecahv...	360	C26H48	055401-75-7	38	
4	1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	32	
5	2-(2-((2-Chloro-6-fluorobenzyl)s...	334	C17H16ClFN2S	1000298-18-2	27	



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040416\
Data File : BF086052.D
Acq On : 4 Apr 2016 22:10
Operator : UM/SJ
Sample : H2054-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TEC-B11COMP

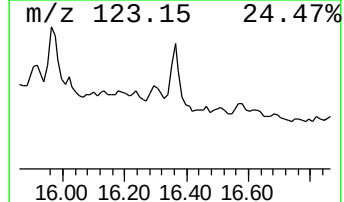
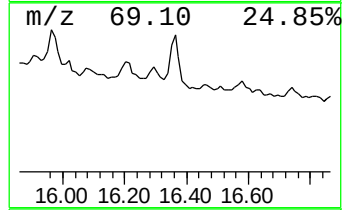
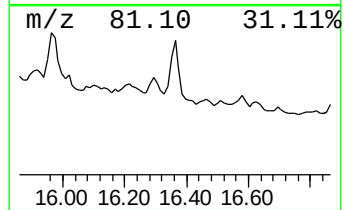
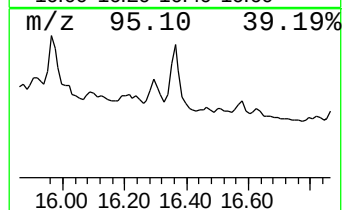
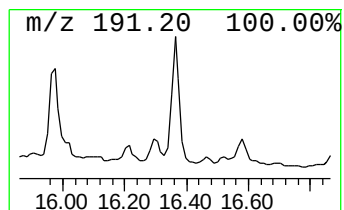
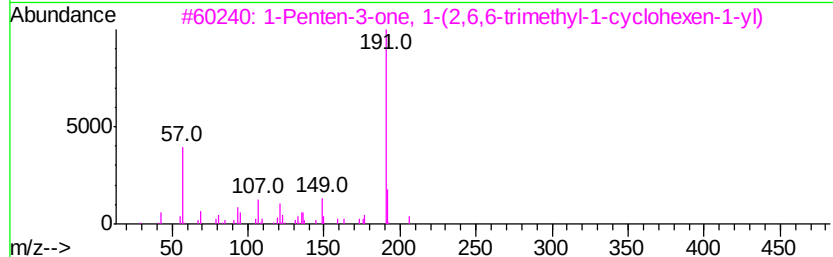
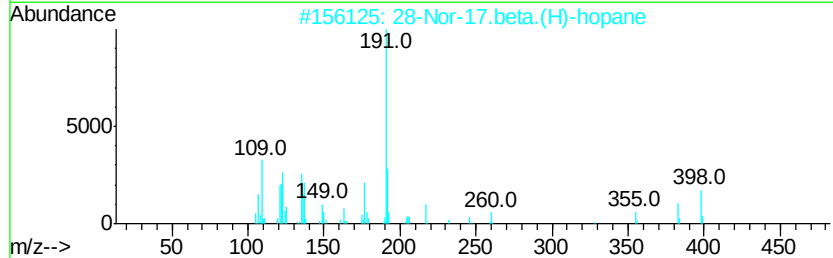
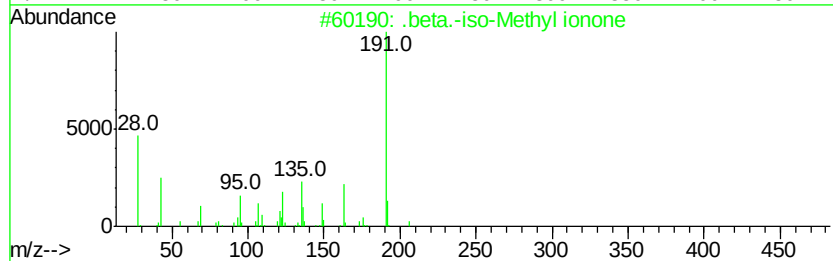
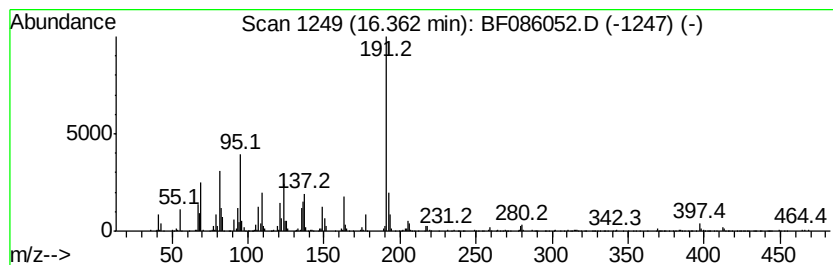
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 .beta.-iso-Methyl ionone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.36	13.03 ng	751760	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	64
2		28-Nor-17.beta.(H)-hopane	398	C29H50	036728-72-0	50
3		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	43
4		1-Cyclohexene, 1,3,3-trimethyl-2...	206	C14H22O	1000197-08-4	42
5		Anthracene, 9-dodecyltetradecahy...	360	C26H48	055401-75-7	42



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040416\
Data File : BF086052.D
Acq On : 4 Apr 2016 22:10
Operator : UM/SJ
Sample : H2054-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TEC-B11COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
Benzene, 1,2,3-tr...	6.58	6.4	ng	235610	1	6.76	734652	20.0
Pentadecane	9.21	7.7	ng	346140	3	9.80	899663	20.0
Hexadecane	10.24	6.2	ng	277447	3	9.80	899663	20.0
Undecane	10.45	4.7	ng	211366	3	9.80	899663	20.0
Tridecane	10.70	10.9	ng	576254	4	11.29	1057930	20.0
Heneicosane	11.15	4.4	ng	233926	4	11.29	1057930	20.0
Phenanthrene, 2,5...	12.34	5.0	ng	262643	4	11.29	1057930	20.0
Eicosane	12.37	5.4	ng	284862	4	11.29	1057930	20.0
Nonadecane	13.10	14.7	ng	295898	5	13.93	403822	20.0
Heptadecane	13.45	20.0	ng	403822	5	13.93	403822	20.0
5-(1-Isopropenyl-...	15.96	20.0	ng	1153970	6	15.35	1153970	20.0
.beta.-iso-Methyl...	16.36	13.0	ng	751760	6	15.35	1153970	20.0