

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040616\
 Data File : BF086116.D
 Acq On : 6 Apr 2016 23:53
 Operator : UM/SJ
 Sample : H2096-02 5X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-B8COMP

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	3.550	125	128	136	rVB	5656	13143	1.21%	0.082%
2	4.681	225	227	230	rVB2	16195	16281	1.50%	0.101%
3	4.899	245	246	251	rVB	10519	14148	1.30%	0.088%
4	5.333	282	284	299	rVB	335162	425335	39.17%	2.651%
5	6.396	374	377	385	rBV	254822	431811	39.77%	2.692%
6	6.510	385	387	390	rVV	505800	488246	44.97%	3.044%
7	6.579	390	393	395	rVB	13363	19827	1.83%	0.124%
8	6.739	405	407	410	rBV	592097	587231	54.08%	3.661%
9	6.899	419	421	423	rBV	424491	380892	35.08%	2.374%
10	7.310	455	457	463	rBV	277778	279148	25.71%	1.740%
11	8.030	517	520	522	rBV	848979	740148	68.17%	4.614%
12	8.453	555	557	562	rVB	12380	16692	1.54%	0.104%
13	8.625	570	572	574	rVB	28786	23222	2.14%	0.145%
14	8.750	579	583	584	rBV	11352	14923	1.37%	0.093%
15	8.933	595	599	602	rBV4	4406	11998	1.10%	0.075%
16	9.059	605	610	612	rVV	8691	13771	1.27%	0.086%
17	9.105	612	614	616	rVV	589987	470815	43.36%	2.935%
18	9.185	619	621	626	rVV	28591	36631	3.37%	0.228%
19	9.448	642	644	645	rBV	9643	15507	1.43%	0.097%
20	9.505	647	649	653	rVB	58146	60576	5.58%	0.378%
21	9.642	659	661	663	rBV	27608	23747	2.19%	0.148%
22	9.722	666	668	669	rBV	34334	38422	3.54%	0.240%
23	9.779	671	673	676	rVB	714068	621484	57.24%	3.874%
24	9.985	689	691	694	rBV3	6907	15694	1.45%	0.098%
25	10.042	694	696	700	rVB	10758	21855	2.01%	0.136%
26	10.191	707	709	710	rBV2	6880	11920	1.10%	0.074%
27	10.213	710	711	713	rVV	57368	42894	3.95%	0.267%
28	10.431	726	730	733	rBV	20771	35604	3.28%	0.222%
29	10.568	740	742	748	rVB2	197013	223024	20.54%	1.390%
30	10.693	751	753	757	rBV	50556	94710	8.72%	0.590%
31	10.762	757	759	762	rBV4	5979	15961	1.47%	0.099%
32	10.899	767	771	774	rBV4	14589	34389	3.17%	0.214%
33	11.025	778	782	784	rBV3	13172	29145	2.68%	0.182%
34	11.139	789	792	793	rVV	41239	60001	5.53%	0.374%

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

35	11.162	793	794	797	rVB	55828	51762	4.77%	0.323%
36	11.265	801	803	805	rBV	656437	620532	57.15%	3.868%
37	11.345	806	810	811	rVB3	15209	33197	3.06%	0.207%
38	11.448	816	819	820	rBV2	10526	15199	1.40%	0.095%
39	11.516	822	825	826	rBV3	15452	25452	2.34%	0.159%
40	11.562	826	829	830	rBV	75129	65625	6.04%	0.409%
41	11.596	830	832	837	rBV2	23698	38706	3.56%	0.241%
42	11.676	837	839	842	rBV	16588	24470	2.25%	0.153%
43	11.722	842	843	844	rBV	14180	14149	1.30%	0.088%
44	11.791	845	849	851	rBV3	35299	74596	6.87%	0.465%
45	11.871	854	856	858	rBV	31967	37132	3.42%	0.231%
46	11.939	860	862	863	rBV	12994	15547	1.43%	0.097%
47	11.962	863	864	866	rBV	92218	90264	8.31%	0.563%
48	12.088	868	875	877	rBV6	25513	83085	7.65%	0.518%
49	12.191	882	884	887	rVB2	32660	37647	3.47%	0.235%
50	12.236	887	888	892	rBV4	28825	59644	5.49%	0.372%
51	12.316	892	895	896	rBV	59841	73756	6.79%	0.460%
52	12.351	896	898	901	rVB	133155	141912	13.07%	0.885%
53	12.408	901	903	904	rBV	21653	37031	3.41%	0.231%
54	12.476	907	909	911	rBV	96594	93744	8.63%	0.584%
55	12.579	915	918	921	rBV4	36137	102618	9.45%	0.640%
56	12.625	921	922	924	rBV	29191	49044	4.52%	0.306%
57	12.671	924	926	927	rBV	96463	79005	7.28%	0.492%
58	12.728	927	931	932	rVB2	146147	246590	22.71%	1.537%
59	12.762	932	934	936	rVB2	22475	26696	2.46%	0.166%
60	12.842	939	941	946	rVB	400541	407434	37.52%	2.540%
61	13.082	959	962	964	rBV	230029	263213	24.24%	1.641%
62	13.334	982	984	986	rBV2	46799	72415	6.67%	0.451%
63	13.379	986	988	990	rVV	77250	80264	7.39%	0.500%
64	13.425	990	992	993	rBV	265868	265312	24.43%	1.654%
65	13.699	1014	1016	1019	rBV	63239	124491	11.47%	0.776%
66	13.757	1019	1021	1025	rVB	246285	363822	33.51%	2.268%
67	13.905	1031	1034	1035	rBV	506024	691139	63.65%	4.308%
68	14.065	1046	1048	1052	rVB	418066	441905	40.70%	2.755%
69	14.374	1073	1075	1076	rBV	377415	394042	36.29%	2.456%
70	14.671	1098	1101	1103	rBV3	291290	472782	43.54%	2.947%
71	14.980	1125	1128	1130	rVB	282538	417818	38.48%	2.605%

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72	15.311	1154	1157	1164	rVB2	547175	1085798	100.00%	6.768%
73	15.711	1189	1192	1195	rBV	291210	448712	41.33%	2.797%
74	16.168	1229	1232	1241	rVB	233533	526083	48.45%	3.279%
75	17.403	1338	1340	1342	rVB2	109529	129690	11.94%	0.808%
76	17.471	1342	1346	1347	rBV4	82901	242239	22.31%	1.510%
77	17.574	1354	1355	1358	rBV3	69112	179335	16.52%	1.118%
78	17.620	1358	1359	1360	rVV	120643	93605	8.62%	0.583%
79	17.677	1360	1364	1365	rVV3	102525	216686	19.96%	1.351%
80	17.711	1365	1367	1368	rVV2	92677	112355	10.35%	0.700%
81	17.746	1368	1370	1372	rVV3	92050	165165	15.21%	1.030%
82	17.791	1372	1374	1375	rVV2	113042	159747	14.71%	0.996%
83	17.860	1378	1380	1381	rVV2	108144	133927	12.33%	0.835%
84	17.906	1382	1384	1385	rVV	126034	136381	12.56%	0.850%
85	17.951	1385	1388	1390	rVV3	133341	312008	28.74%	1.945%
86	17.997	1390	1392	1395	rVV4	70597	174655	16.09%	1.089%
87	18.043	1395	1396	1402	rVB2	111260	187472	17.27%	1.169%
88	18.351	1421	1423	1426	rBV3	54618	88547	8.16%	0.552%
89	18.454	1430	1432	1433	rVB2	81231	85531	7.88%	0.533%
90	18.511	1436	1437	1439	rBV2	42380	44652	4.11%	0.278%
91	19.094	1486	1488	1491	rBV3	41161	58257	5.37%	0.363%

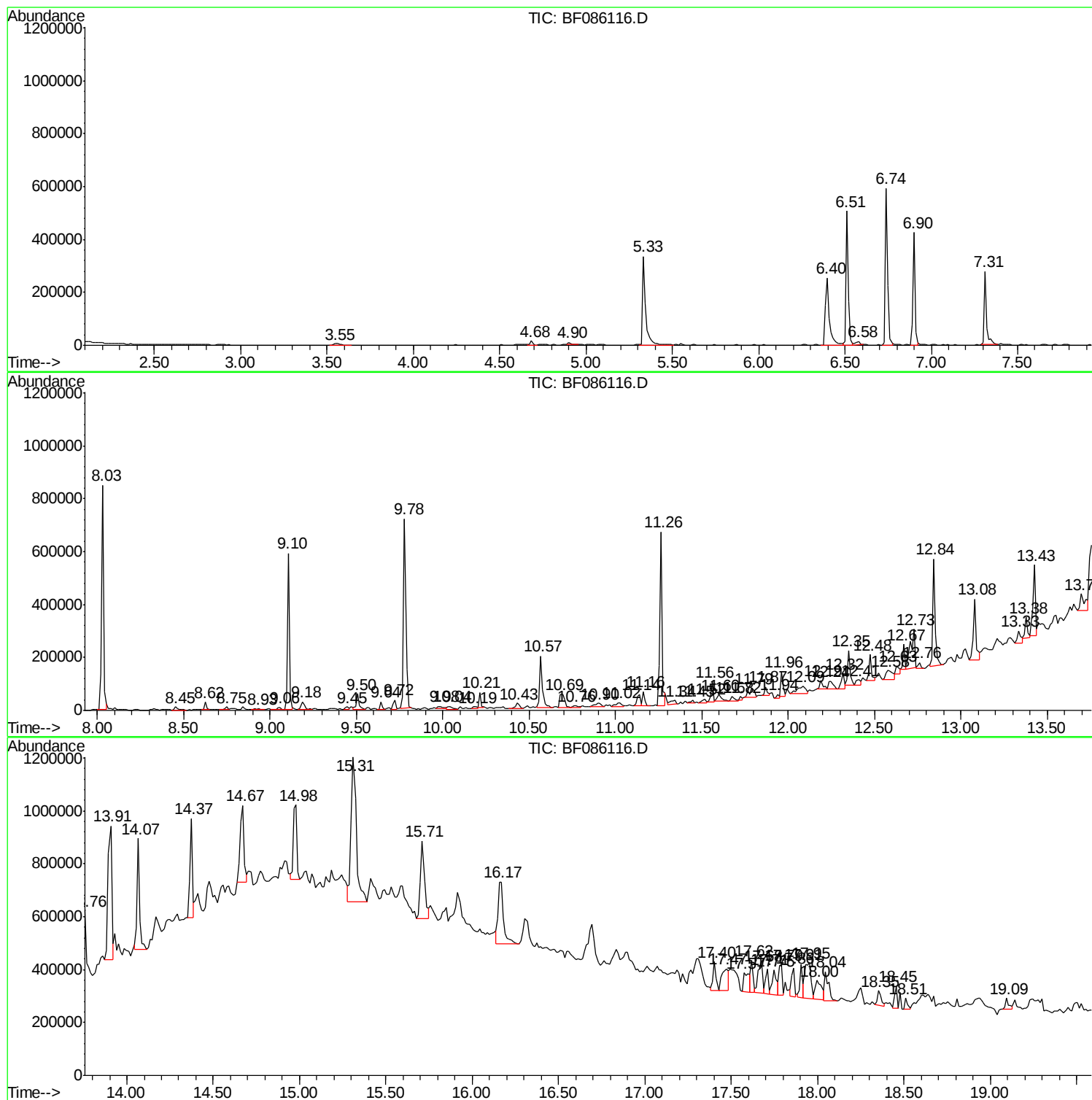
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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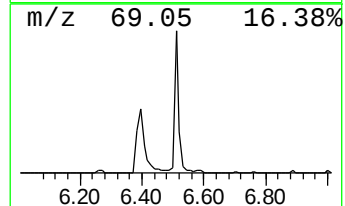
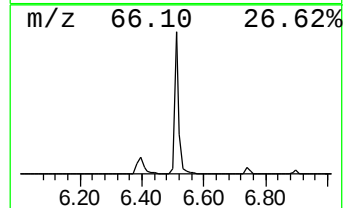
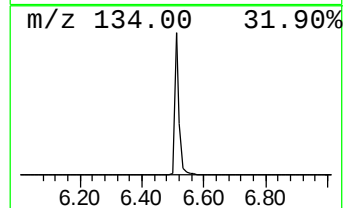
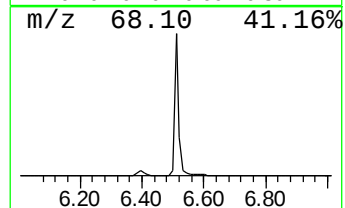
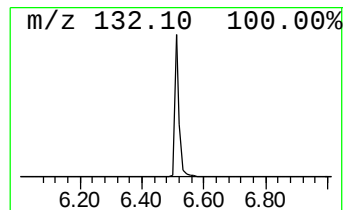
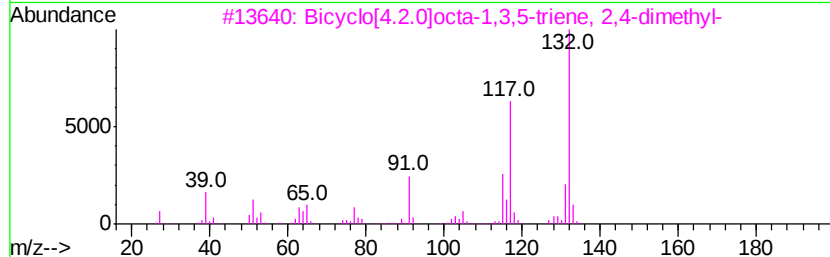
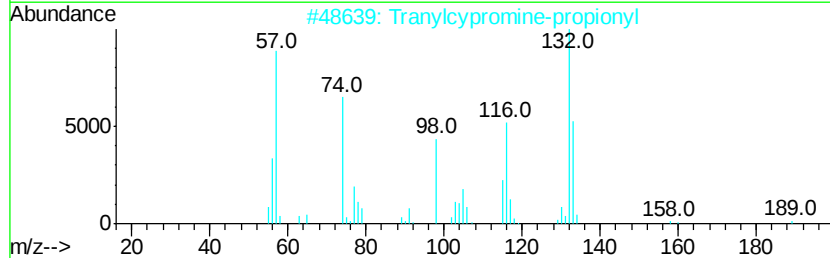
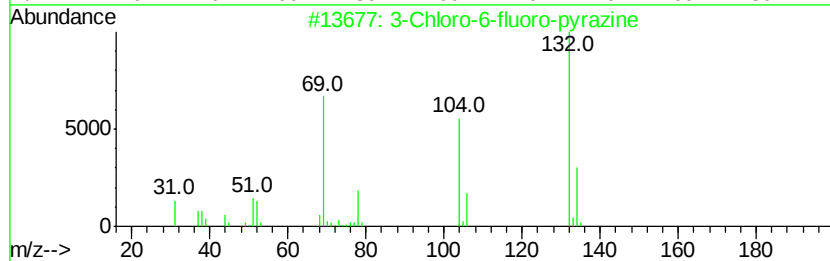
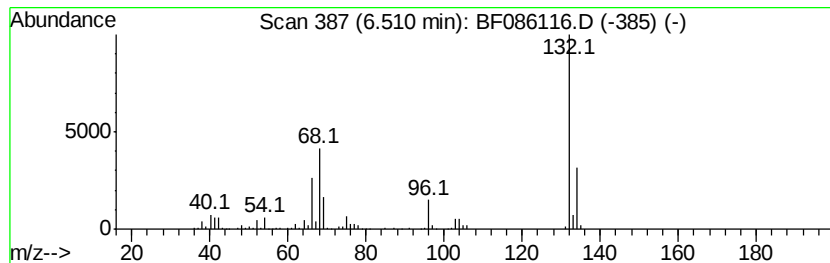
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.51 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	16.63 ng	488246	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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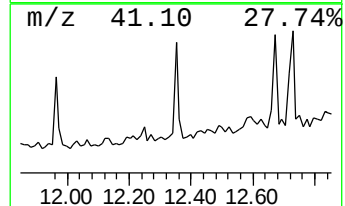
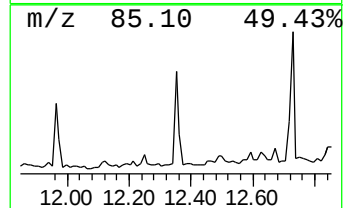
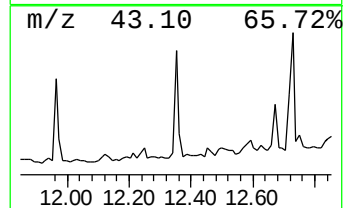
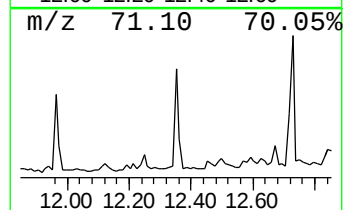
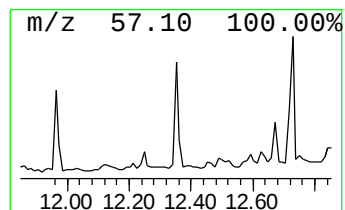
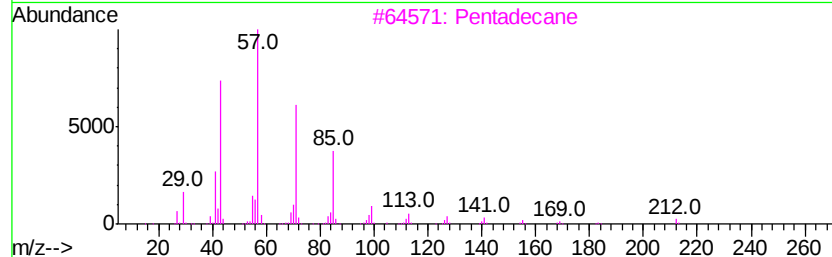
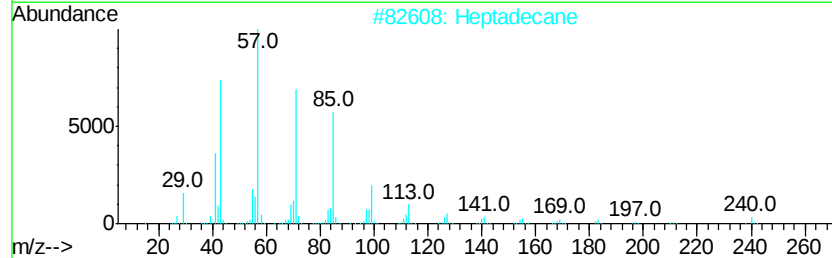
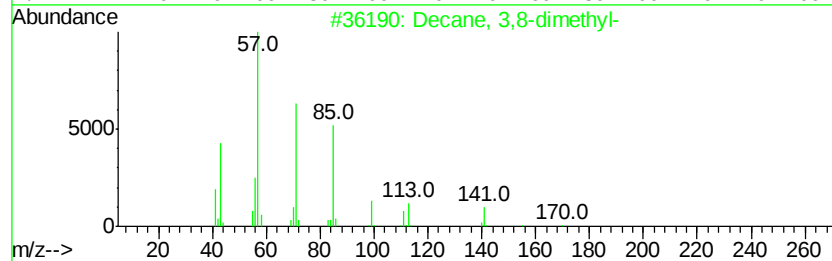
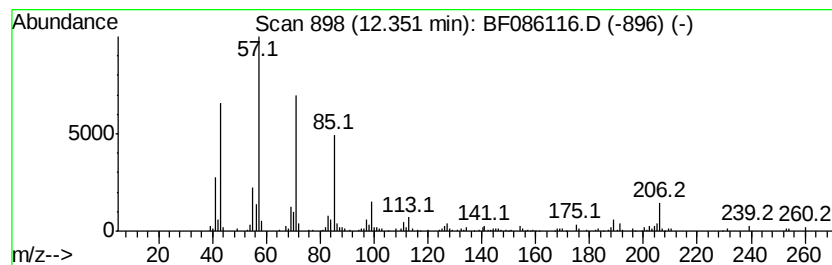
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Decane, 3,8-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.35	4.57 ng	141912	Phenanthrene-d10	11.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	74
2		Heptadecane	240	C17H36	000629-78-7	74
3		Pentadecane	212	C15H32	000629-62-9	72
4		Eicosane, 10-methyl-	296	C21H44	054833-23-7	72
5		Eicosane	282	C20H42	000112-95-8	72



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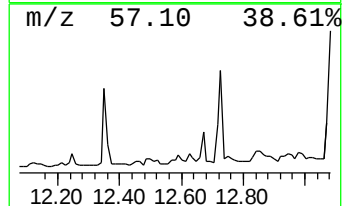
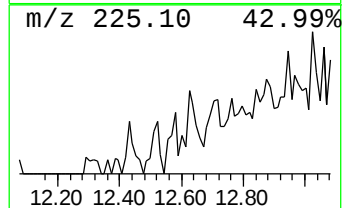
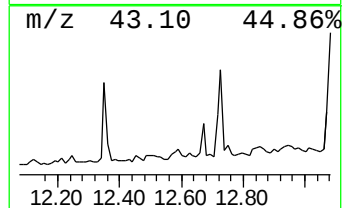
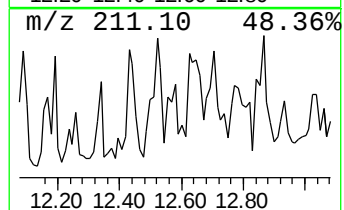
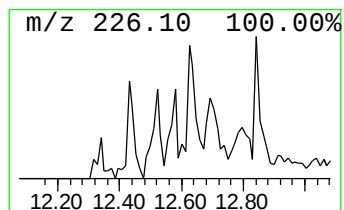
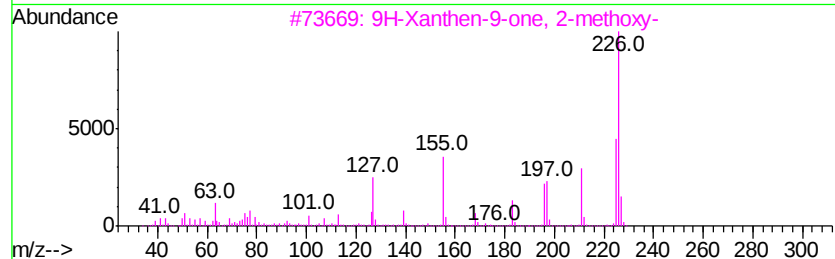
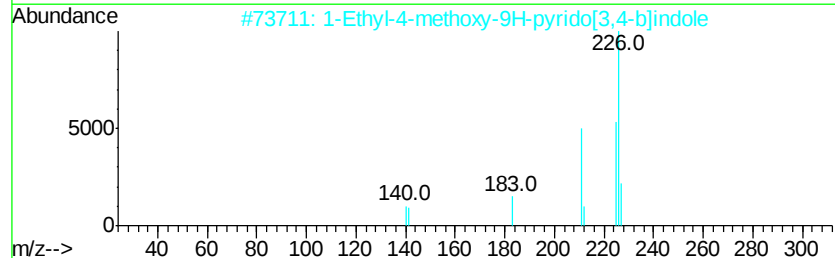
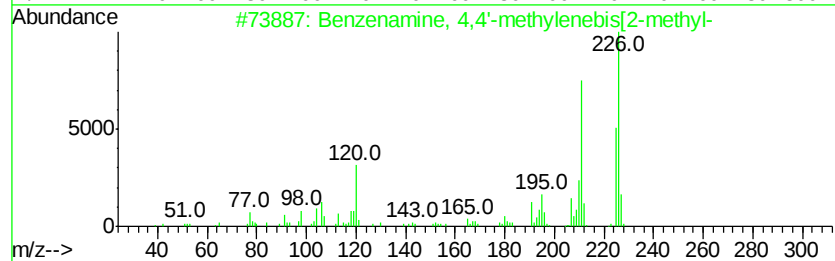
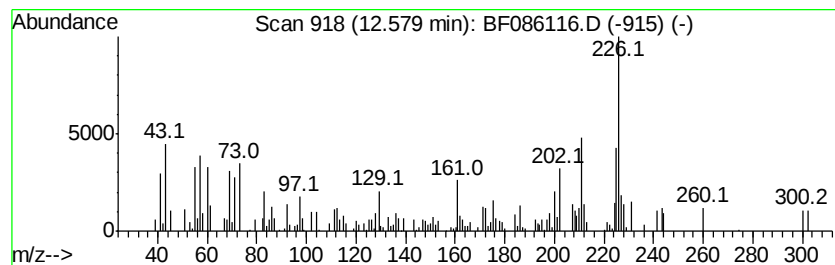
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzenamine, 4,4'-methylene... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.58	3.31 ng	102618	Phenanthrene-d10	11.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenamine, 4,4'-methylenebis[2...	226	C15H18N2	000838-88-0	89
2		1-Ethyl-4-methoxy-9H-pyrido[3,4-...	226	C14H14N2O	026585-14-8	58
3		9H-Xanthen-9-one, 2-methoxy-	226	C14H10O3	001214-20-6	50
4		N,N-Dimethylindoaniline	226	C14H14N2O	002150-58-5	49
5		Acridin-9-amine, 1,2,3,4-tetrahy...	226	C15H18N2	297758-19-1	47



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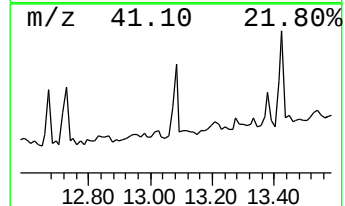
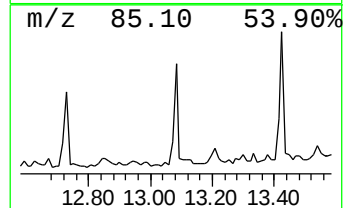
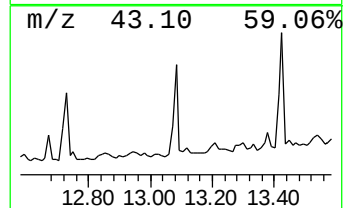
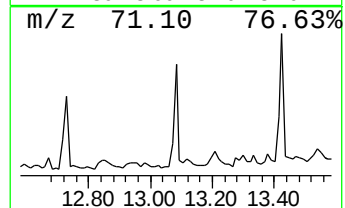
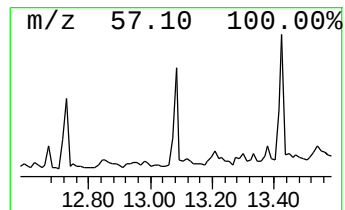
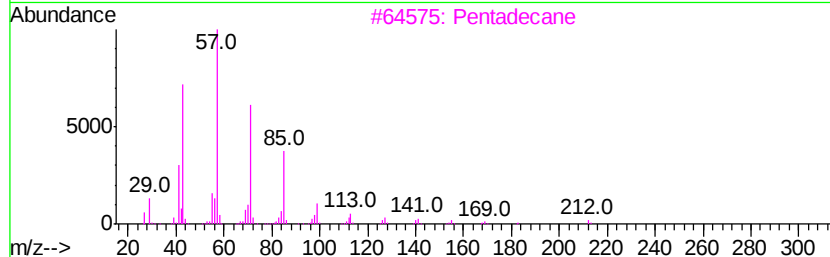
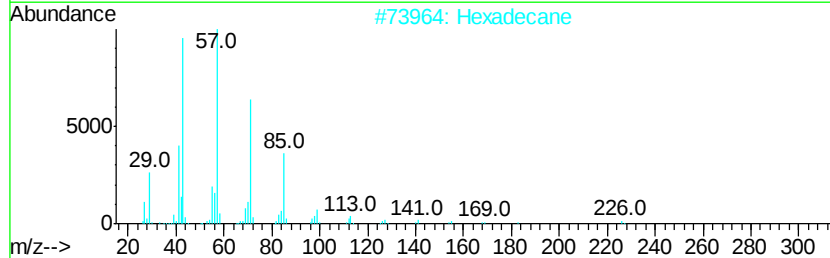
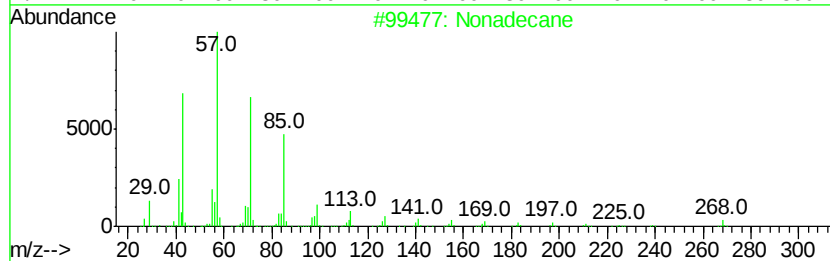
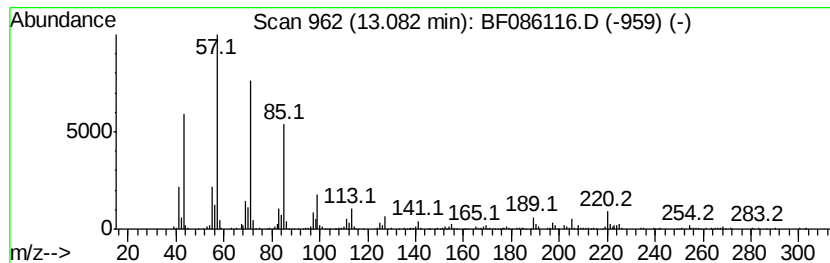
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Nonadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.08	7.62 ng	263213	Chrysene-d12	13.91

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	93
2	Hexadecane		226	C16H34	000544-76-3	93
3	Pentadecane		212	C15H32	000629-62-9	91
4	Tridecane, 6-propyl-		226	C16H34	055045-10-8	87
5	Heneicosane		296	C21H44	000629-94-7	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
Data File : BF086116.D
Acq On : 6 Apr 2016 23:53
Operator : UM/SJ
Sample : H2096-02 5X
Misc :
ALS Vial : 27 Sample Multiplier: 1

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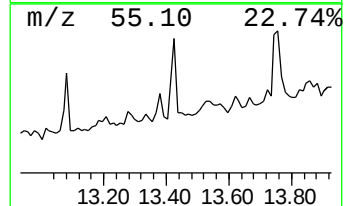
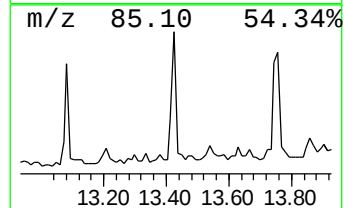
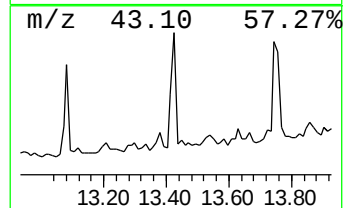
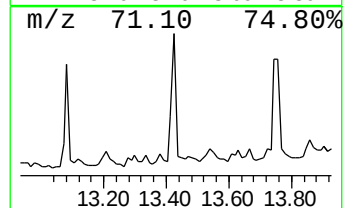
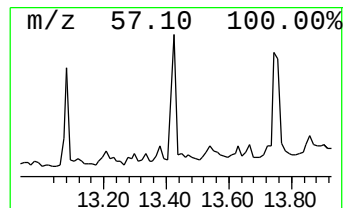
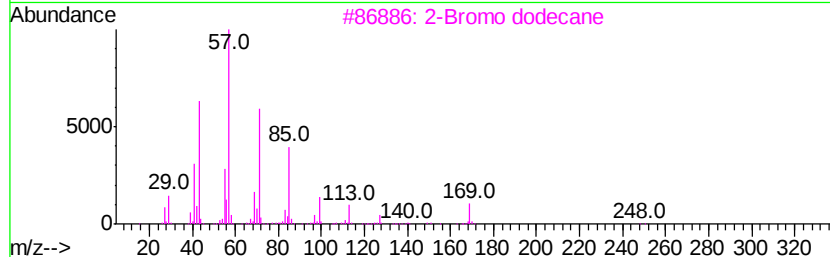
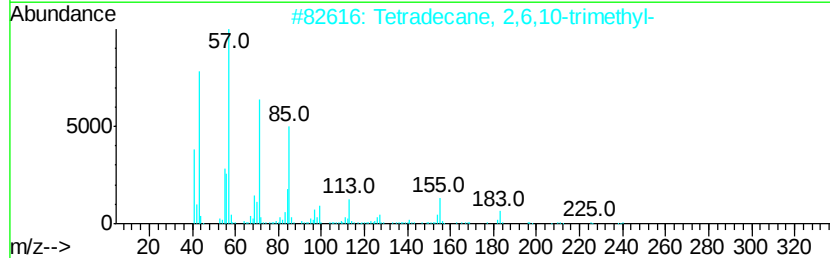
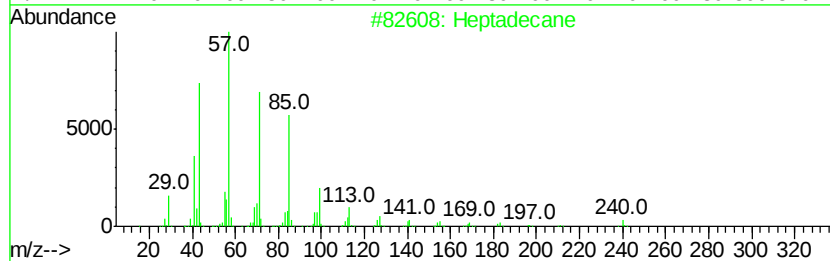
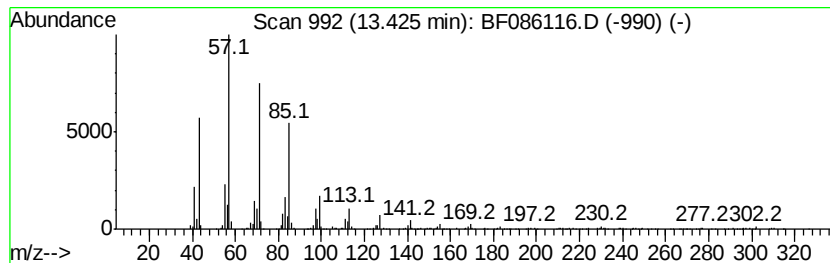
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Heptadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.43	7.68 ng	265312	Chrysene-d12	13.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	94
2		Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	93
3		2-Bromo dodecane	248	C12H25Br	013187-99-0	93
4		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
5		Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040616\
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Misc :
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Instrument :
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ClientSampleId :
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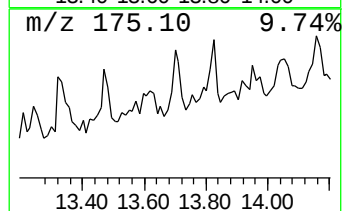
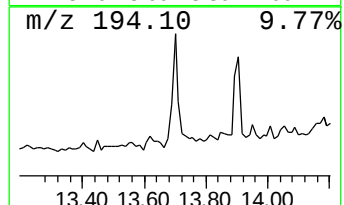
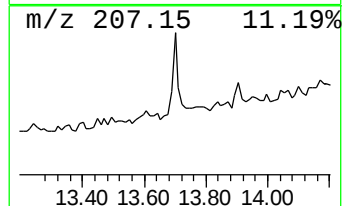
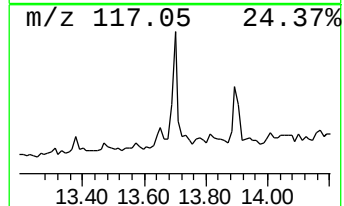
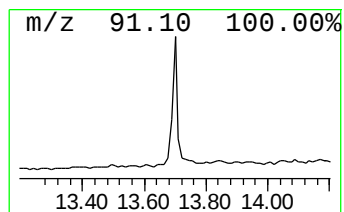
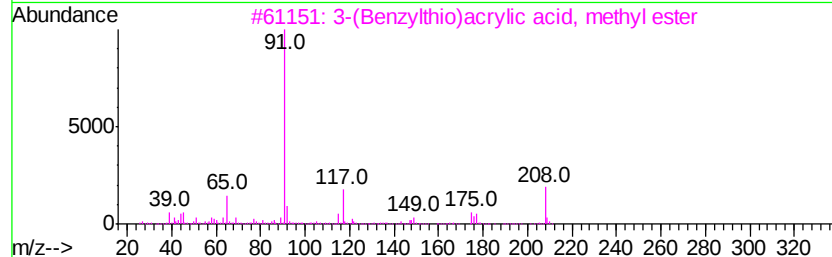
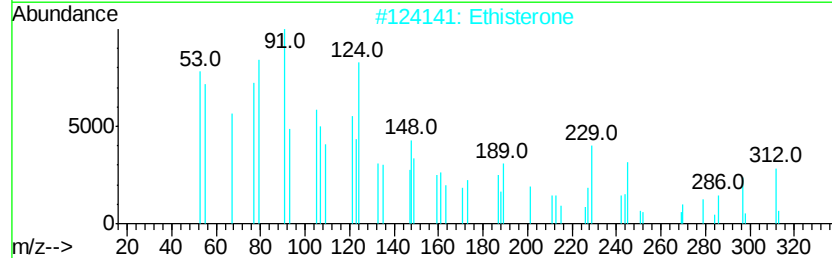
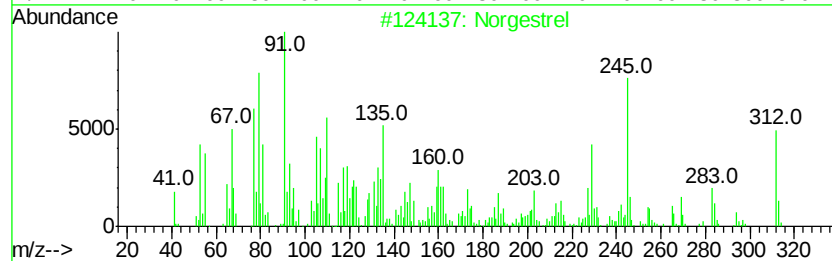
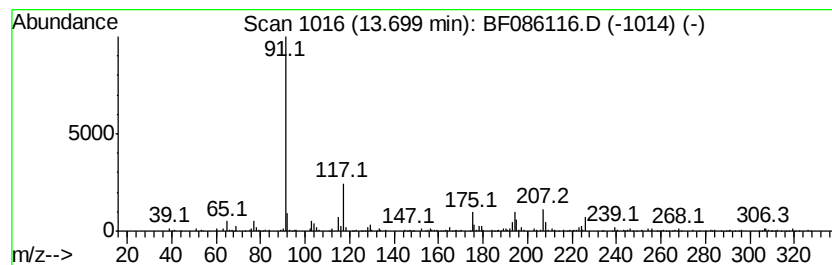
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Norgestrel Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.70	3.60 ng	124491	Chrysene-d12	13.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nordestrel	312	C21H28O2	006533-00-2	90
2		Ethisterone	312	C21H28O2	000434-03-7	59
3		3-(Benzylthio)acrylic acid, meth...	208	C11H12O2S	1000192-96-8	53
4		5(2H)-Oxazolone, 4-(phenylmethyl)-	175	C10H9N02	049656-77-1	50
5		Benzylidene-L-ornithine	220	C12H16N2O2	025693-00-9	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
Data File : BF086116.D
Acq On : 6 Apr 2016 23:53
Operator : UM/SJ
Sample : H2096-02 5X
Misc :
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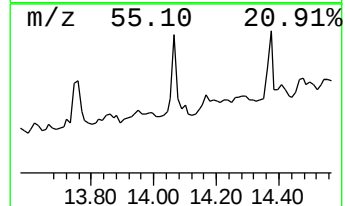
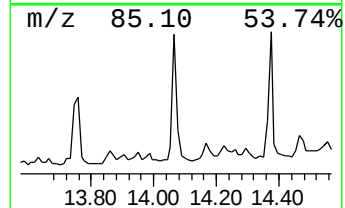
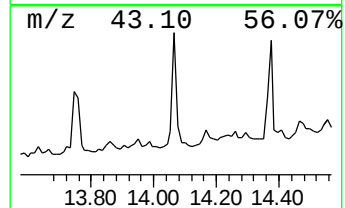
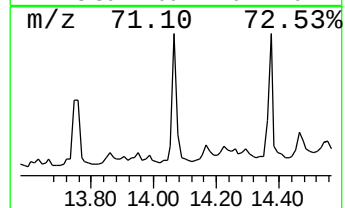
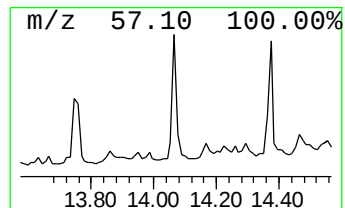
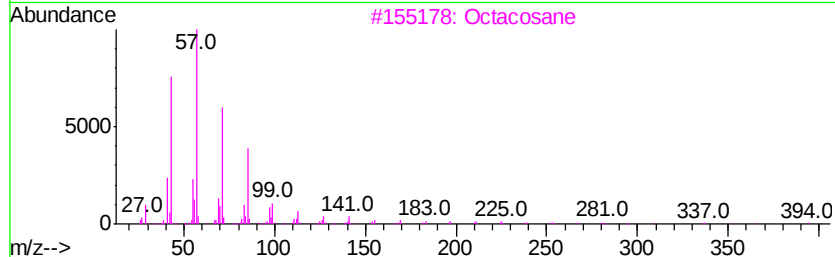
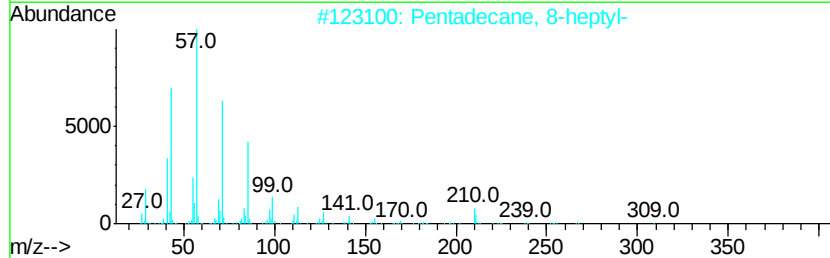
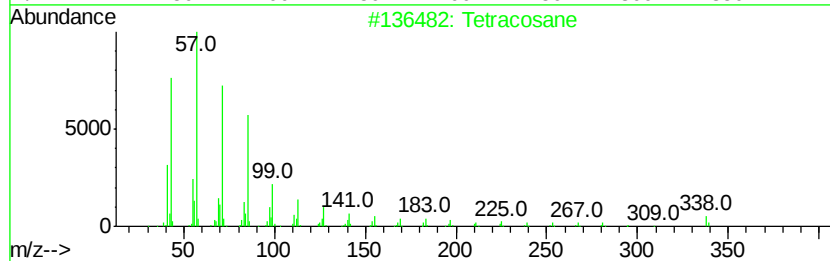
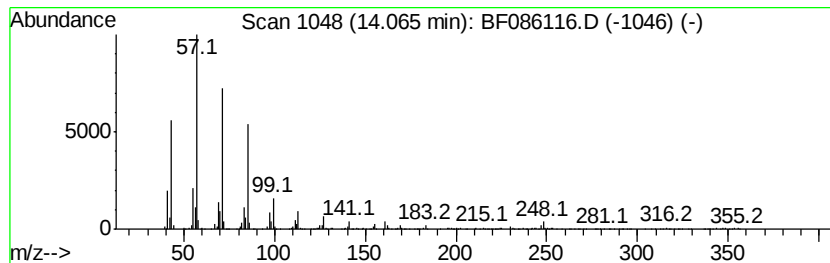
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetracosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.07	12.79 ng	441905	Chrysene-d12	13.91

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	91
2		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	87
3		Octacosane	394	C28H58	000630-02-4	87
4		Heneicosane	296	C21H44	000629-94-7	87
5		Nonadecane	268	C19H40	000629-92-5	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
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Acq On : 6 Apr 2016 23:53
Operator : UM/SJ
Sample : H2096-02 5X
Misc :
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Instrument :
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ClientSampleId :
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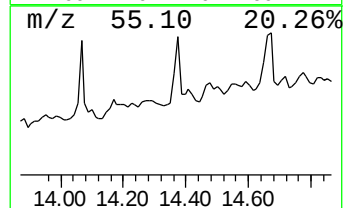
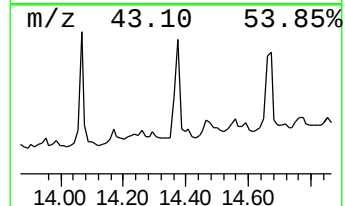
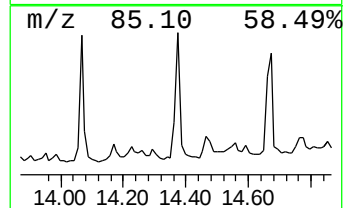
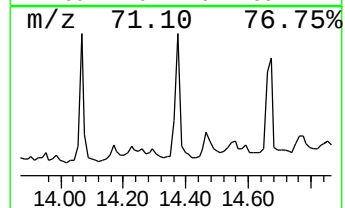
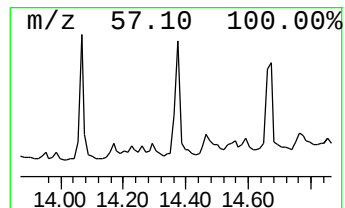
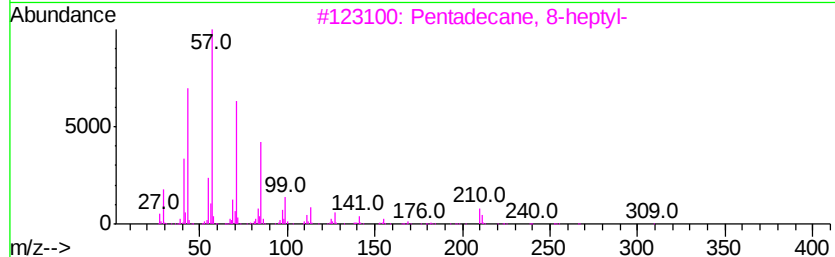
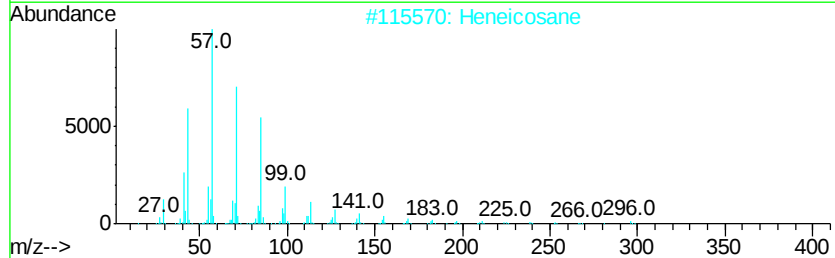
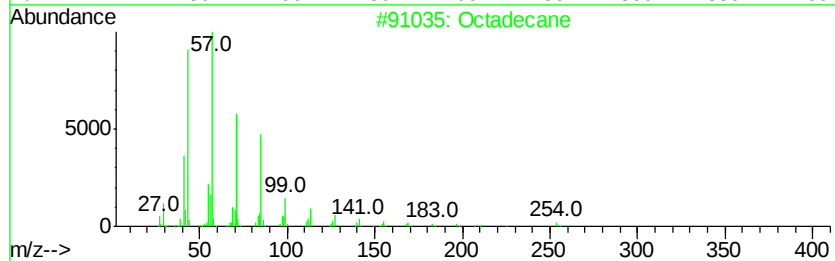
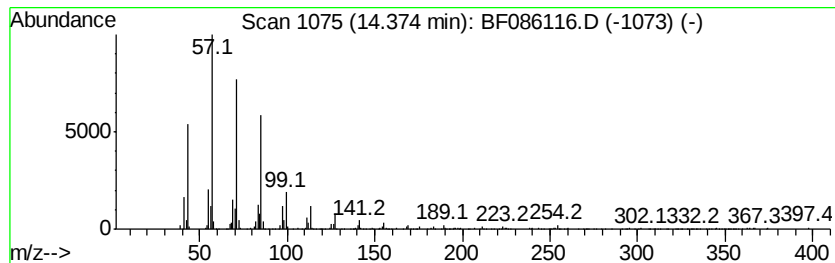
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Octadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.37	11.40 ng	394042	Chrysene-d12	13.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	94
2		Heneicosane	296	C21H44	000629-94-7	91
3		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
4		Tetradecane	198	C14H30	000629-59-4	90
5		Eicosane	282	C20H42	000112-95-8	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
Data File : BF086116.D
Acq On : 6 Apr 2016 23:53
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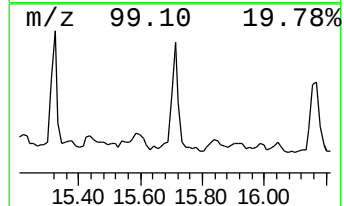
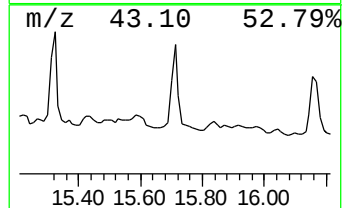
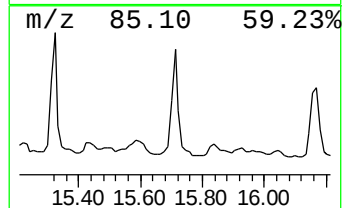
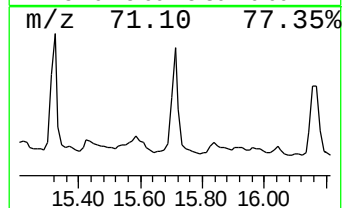
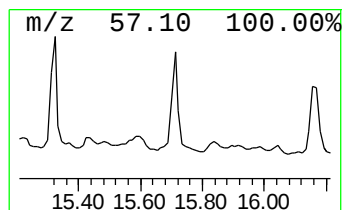
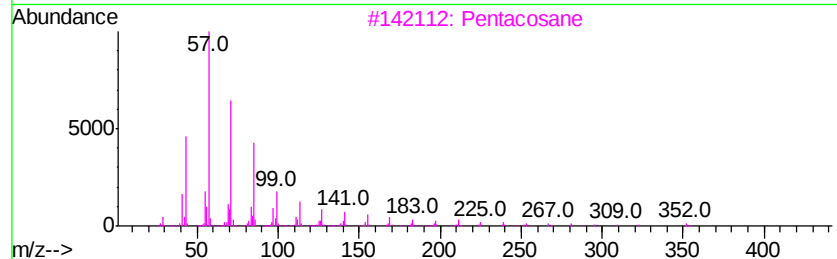
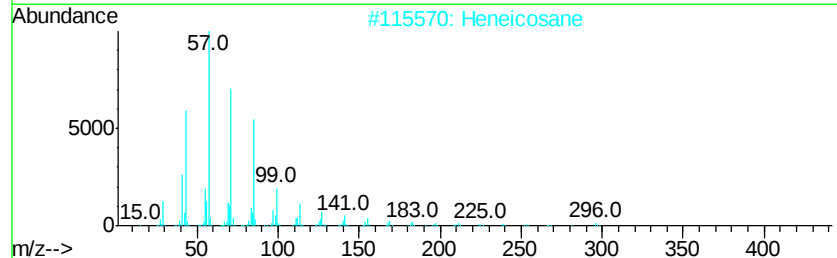
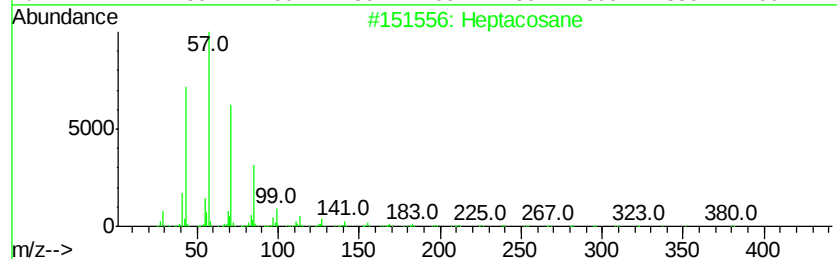
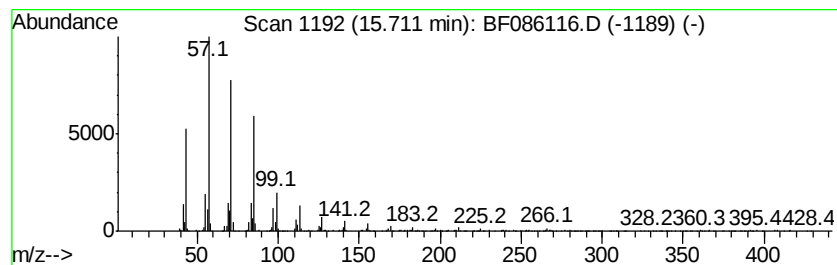
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heptacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.71	8.27 ng	448712	Perylene-d12	15.31

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane	380	C27H56	000593-49-7	93
2			Heneicosane	296	C21H44	000629-94-7	91
3			Pentacosane	352	C25H52	000629-99-2	91
4			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	86
5			Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
Data File : BF086116.D
Acq On : 6 Apr 2016 23:53
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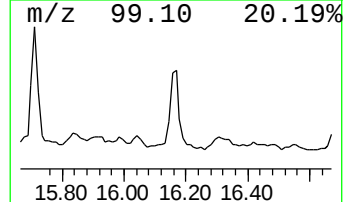
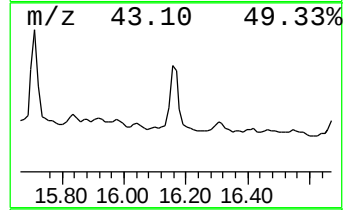
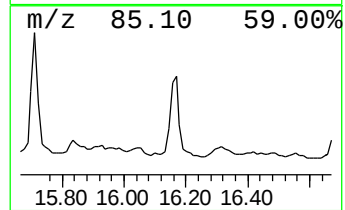
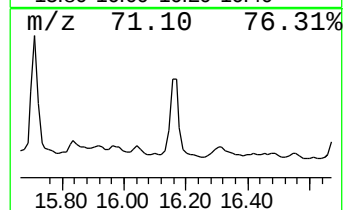
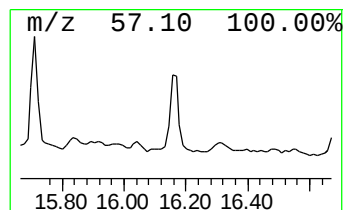
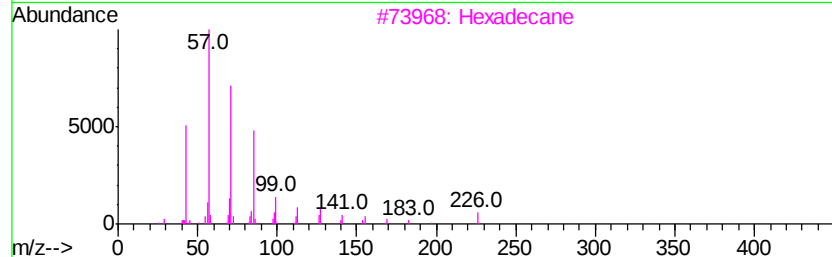
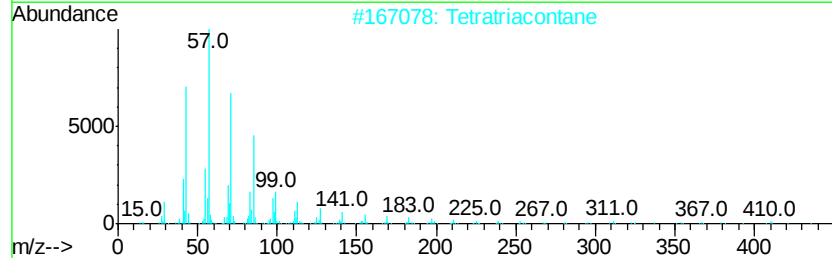
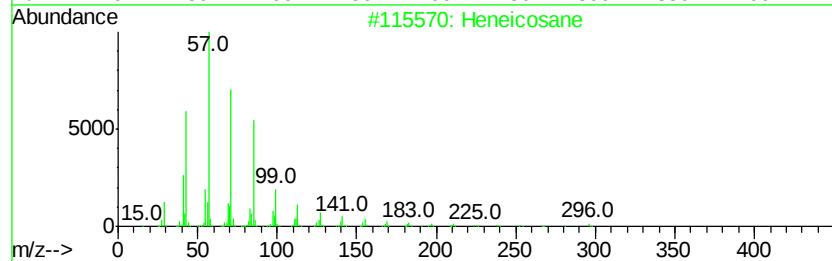
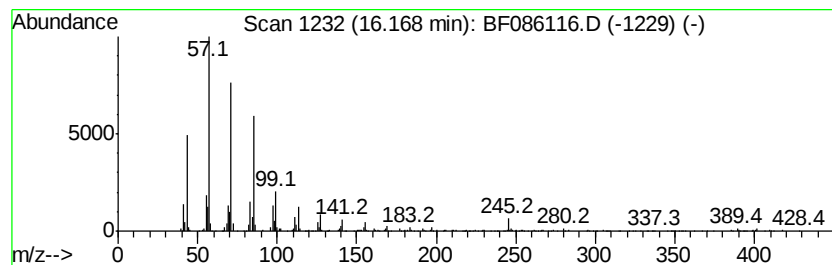
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Heneicosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.17	9.69 ng	526083	Perylene-d12	15.31

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	90
2			Tetratriacontane	479	C34H70	014167-59-0	90
3			Hexadecane	226	C16H34	000544-76-3	90
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	89
5			Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
Data File : BF086116.D
Acq On : 6 Apr 2016 23:53
Operator : UM/SJ
Sample : H2096-02 5X
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TEC-B8COMP

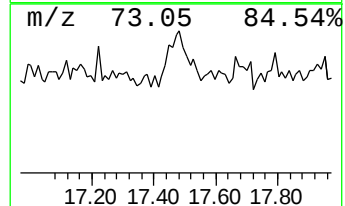
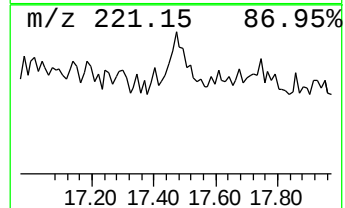
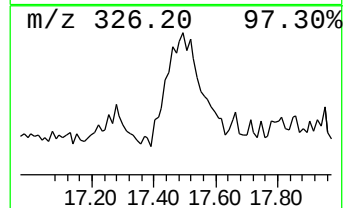
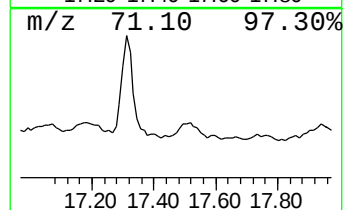
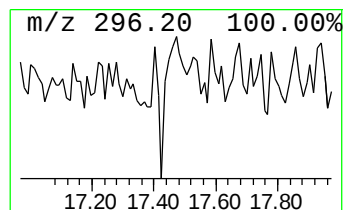
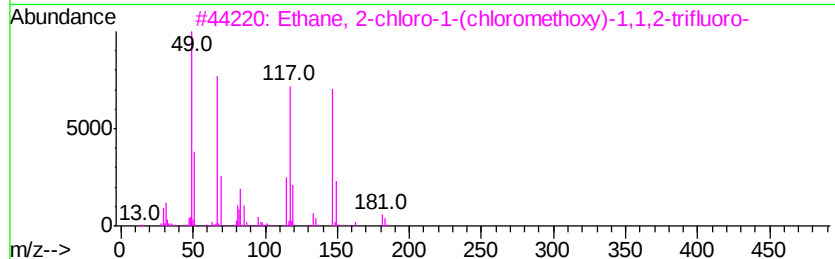
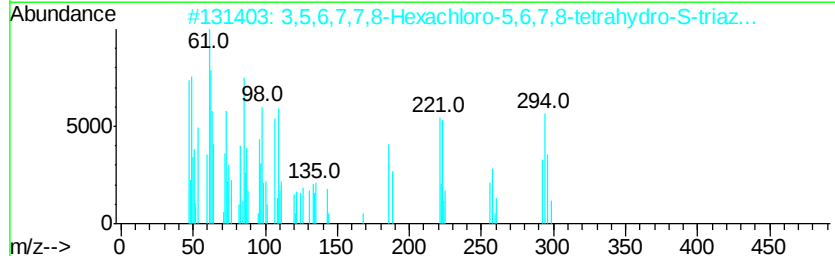
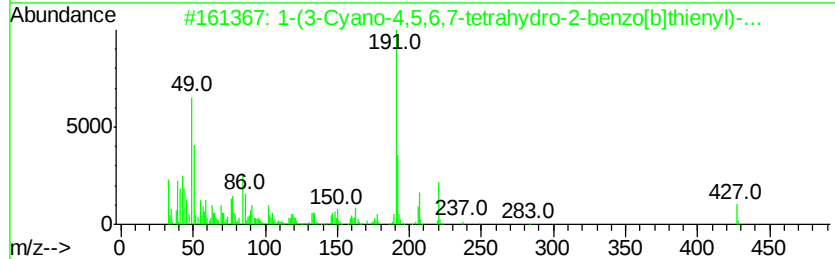
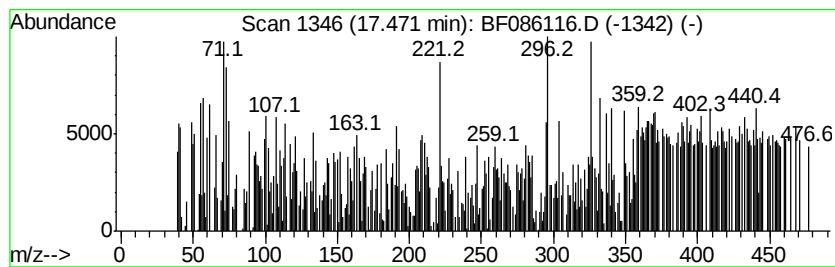
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown17.47 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.47	4.46 ng	242239	Perylene-d12	15.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-(3-Cyano-4,5,6,7-tetrahydro-2-...	427	C21H21N3O3S2	1000225-91-0	20
2		3,5,6,7,7,8-Hexachloro-5,6,7,8-t...	327	C6H3Cl6N3	022841-85-6	18
3		Ethane, 2-chloro-1-(chloromethox...	182	C3H3Cl2F3O	000428-92-2	14
4		Pyridine-2,5-dicarboxylic acid, ...	311	C13H21N04Si2	1000153-38-6	10
5		Acetamide, N-(3,4,5-tribromophen...	403	C8H5Br3ClNO	1000276-95-3	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
Data File : BF086116.D
Acq On : 6 Apr 2016 23:53
Operator : UM/SJ
Sample : H2096-02 5X
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TEC-B8COMP

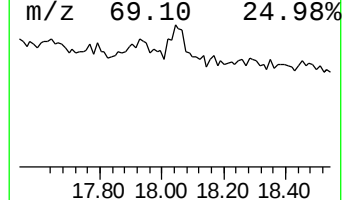
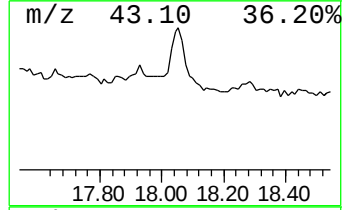
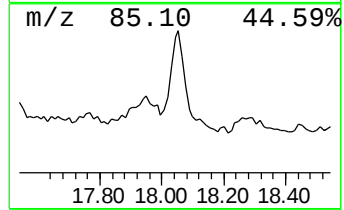
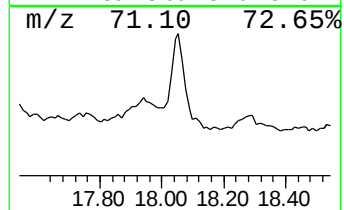
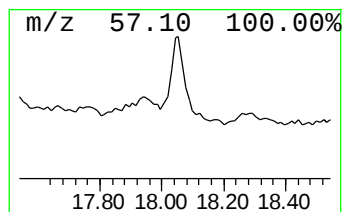
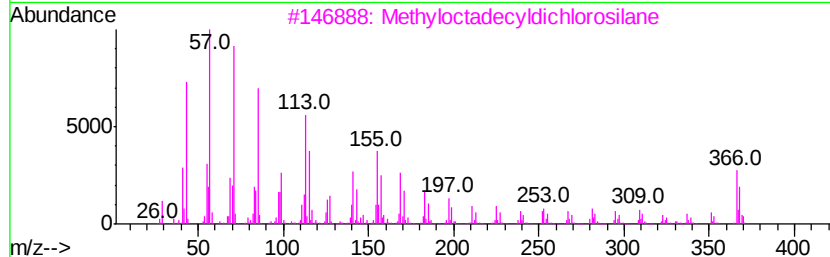
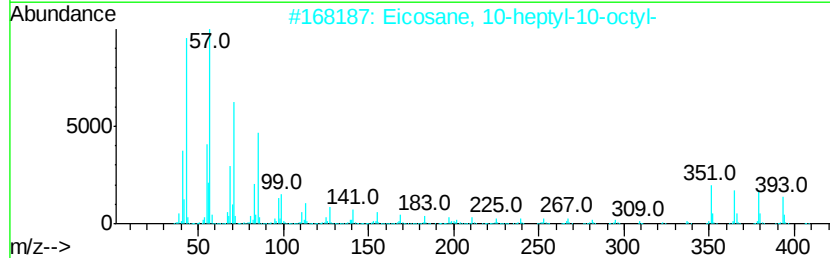
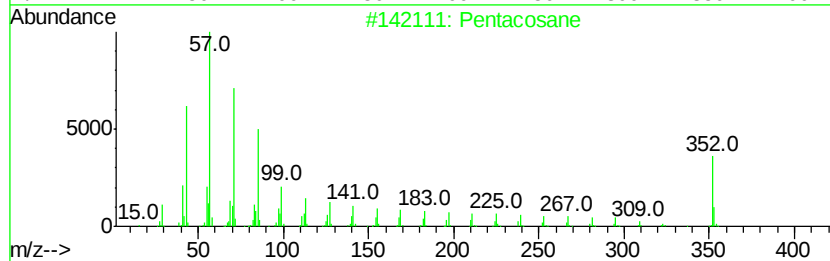
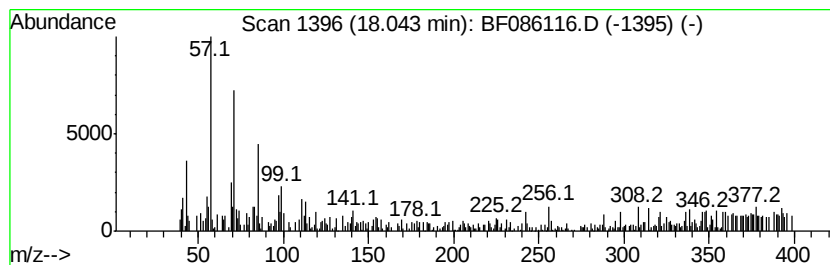
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 unknown18.04 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.04	3.45 ng	187472	Perylene-d12	15.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentacosane	352	C25H52	000629-99-2	30
2		Eicosane, 10-heptyl-10-octyl-	493	C35H72	055470-98-9	27
3		Methyloctadecyldichlorosilane	366	C19H40Cl2Si	005157-75-5	22
4		Hexacosane	366	C26H54	000630-01-3	22
5		Hexacosane, 9-octyl-	479	C34H70	055429-83-9	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
Data File : BF086116.D
Acq On : 6 Apr 2016 23:53
Operator : UM/SJ
Sample : H2096-02 5X
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TEC-B8COMP

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
unknown6.51	6.51	16.6	ng	488246	1	6.74	587231	20.0
Decane, 3,8-dimet...	12.35	4.6	ng	141912	4	11.26	620532	20.0
Benzenamine, 4,4'...	12.58	3.3	ng	102618	4	11.26	620532	20.0
Nonadecane	13.08	7.6	ng	263213	5	13.91	691139	20.0
Heptadecane	13.43	7.7	ng	265312	5	13.91	691139	20.0
Norgestrel	13.70	3.6	ng	124491	5	13.91	691139	20.0
Tetracosane	14.07	12.8	ng	441905	5	13.91	691139	20.0
Octadecane	14.37	11.4	ng	394042	5	13.91	691139	20.0
Heptacosane	15.71	8.3	ng	448712	6	15.31	1085800	20.0
Heneicosane	16.17	9.7	ng	526083	6	15.31	1085800	20.0
unknown17.47	17.47	4.5	ng	242239	6	15.31	1085800	20.0
unknown18.04	18.04	3.5	ng	187472	6	15.31	1085800	20.0