

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF020720\
 Data File : BF118848.D
 Acq On : 7 Feb 2020 14:08
 Operator : CG/JU
 Sample : L1431-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 350

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-BF020320.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	5.434	564	569	572	rBV	4384519	4155261	56.02%	5.593%
2	6.445	736	741	744	rBV	4050538	4104371	55.34%	5.524%
3	6.810	800	803	807	rBV	1547021	1369436	18.46%	1.843%
4	6.969	826	830	834	rBV	4279641	3760821	50.70%	5.062%
5	7.134	854	858	861	rBV	107222	92193	1.24%	0.124%
6	7.375	890	899	902	rBV	2801473	3002590	40.48%	4.041%
7	7.586	930	935	937	rBV	156171	150842	2.03%	0.203%
8	7.610	937	939	943	rVB	134360	112162	1.51%	0.151%
9	7.769	960	966	968	rBV3	73665	111773	1.51%	0.150%
10	7.833	972	977	980	rBV3	184233	222811	3.00%	0.300%
11	7.869	980	983	986	rVB2	81123	84689	1.14%	0.114%
12	8.092	1017	1021	1023	rBV	2043691	1875039	25.28%	2.524%
13	8.116	1023	1025	1028	rVB	659584	541356	7.30%	0.729%
14	8.480	1084	1087	1090	rBV	119952	108276	1.46%	0.146%
15	8.804	1137	1142	1146	rBV	1072700	961243	12.96%	1.294%
16	8.904	1154	1159	1163	rBV	713944	631105	8.51%	0.849%
17	9.169	1199	1204	1208	rBV	5907934	5767230	77.75%	7.762%
18	9.269	1217	1221	1224	rBV	280771	230308	3.11%	0.310%
19	9.363	1232	1237	1243	rVB	135386	169392	2.28%	0.228%
20	9.433	1243	1249	1254	rBV2	208627	295139	3.98%	0.397%
21	9.510	1257	1262	1264	rBV	265079	279195	3.76%	0.376%
22	9.533	1264	1266	1268	rVV	182505	146885	1.98%	0.198%
23	9.563	1268	1271	1278	rVB	260848	241731	3.26%	0.325%
24	9.627	1278	1282	1286	rBV	103753	128119	1.73%	0.172%
25	9.710	1292	1296	1301	rVB	76328	77963	1.05%	0.105%
26	9.851	1313	1320	1322	rBV	2476467	2404157	32.41%	3.236%
27	9.880	1322	1325	1329	rVB	2121726	1781504	24.02%	2.398%
28	10.004	1342	1346	1350	rBV2	73358	78236	1.05%	0.105%
29	10.051	1350	1354	1358	rBV	1232415	1068613	14.41%	1.438%
30	10.392	1408	1412	1418	rBV	1500365	1477664	19.92%	1.989%
31	10.480	1425	1427	1430	rVB	178501	136735	1.84%	0.184%
32	10.551	1437	1439	1444	rVB	209389	186105	2.51%	0.250%
33	10.639	1449	1454	1460	rBV	4375244	4555602	61.42%	6.132%
34	10.757	1470	1474	1479	rVB4	83741	99766	1.35%	0.134%

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 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

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	10.827	1481	1486	1491	rBV	76532	95336	1.29%	0.128%
36	10.945	1498	1506	1508	rBV2	127866	179342	2.42%	0.241%
37	11.186	1543	1547	1549	rBV2	57335	83950	1.13%	0.113%
38	11.227	1551	1554	1561	rVB	416245	374049	5.04%	0.503%
39	11.333	1567	1572	1574	rBV	2787120	2761558	37.23%	3.717%
40	11.357	1574	1576	1580	rVV	4417387	4052841	54.64%	5.455%
41	11.404	1580	1584	1588	rVB	480991	495626	6.68%	0.667%
42	11.563	1607	1611	1620	rBV2	139201	184660	2.49%	0.249%
43	11.786	1646	1649	1653	rBV3	96995	113030	1.52%	0.152%
44	11.833	1653	1657	1659	rBV	241344	211161	2.85%	0.284%
45	11.863	1659	1662	1668	rVV	1129738	1197307	16.14%	1.612%
46	11.910	1668	1670	1673	rVB	110557	93275	1.26%	0.126%
47	11.945	1673	1676	1679	rBV	517492	521569	7.03%	0.702%
48	12.127	1702	1707	1710	rBV2	200411	277020	3.73%	0.373%
49	12.545	1774	1778	1781	rBV	2527804	2235322	30.14%	3.009%
50	12.645	1791	1795	1801	rBV	289070	344346	4.64%	0.463%
51	12.692	1801	1803	1808	rVV	94764	104610	1.41%	0.141%
52	12.774	1810	1817	1819	rVV	1506724	1807749	24.37%	2.433%
53	12.821	1823	1825	1831	rVB3	89481	92426	1.25%	0.124%
54	12.921	1837	1842	1845	rVV	7477304	7417283	100.00%	9.983%
55	12.992	1849	1854	1857	rBV2	56079	86773	1.17%	0.117%
56	13.098	1870	1872	1876	rVB	214162	193715	2.61%	0.261%
57	13.168	1880	1884	1886	rVB	216697	210168	2.83%	0.283%
58	13.204	1886	1890	1892	rBV2	83233	81297	1.10%	0.109%
59	13.715	1973	1977	1980	rBV	66990	88256	1.19%	0.119%
60	13.815	1991	1994	2004	rVB2	554251	692911	9.34%	0.933%
61	13.968	2012	2020	2023	rBV	3118706	3439152	46.37%	4.629%
62	14.133	2045	2048	2054	rVB2	123497	145064	1.96%	0.195%
63	14.439	2096	2100	2105	rBV2	226295	316965	4.27%	0.427%
64	14.745	2149	2152	2155	rVB	242477	235906	3.18%	0.318%
65	14.815	2160	2164	2171	rVB	1227720	1237863	16.69%	1.666%
66	14.998	2191	2195	2198	rBV	147389	222906	3.01%	0.300%
67	15.074	2204	2208	2211	rBV	273359	315490	4.25%	0.425%
68	15.351	2252	2255	2261	rVB	92771	116289	1.57%	0.157%
69	15.415	2261	2266	2270	rBV	1859739	2675635	36.07%	3.601%
70	15.868	2339	2343	2348	rBV	185800	279834	3.77%	0.377%
71	15.933	2350	2354	2362	rVB5	55006	114847	1.55%	0.155%

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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

72	16.292	2409	2415	2422	rBV5	84706	178498	2.41%	0.240%
73	16.362	2423	2427	2433	rVB	102355	171173	2.31%	0.230%
74	16.851	2506	2510	2519	rVV2	47999	109025	1.47%	0.147%
75	16.939	2523	2525	2531	rVB2	59032	103419	1.39%	0.139%
76	17.274	2576	2582	2590	rVB	51714	155628	2.10%	0.209%
77	17.386	2596	2601	2606	rBV7	37334	76015	1.02%	0.102%

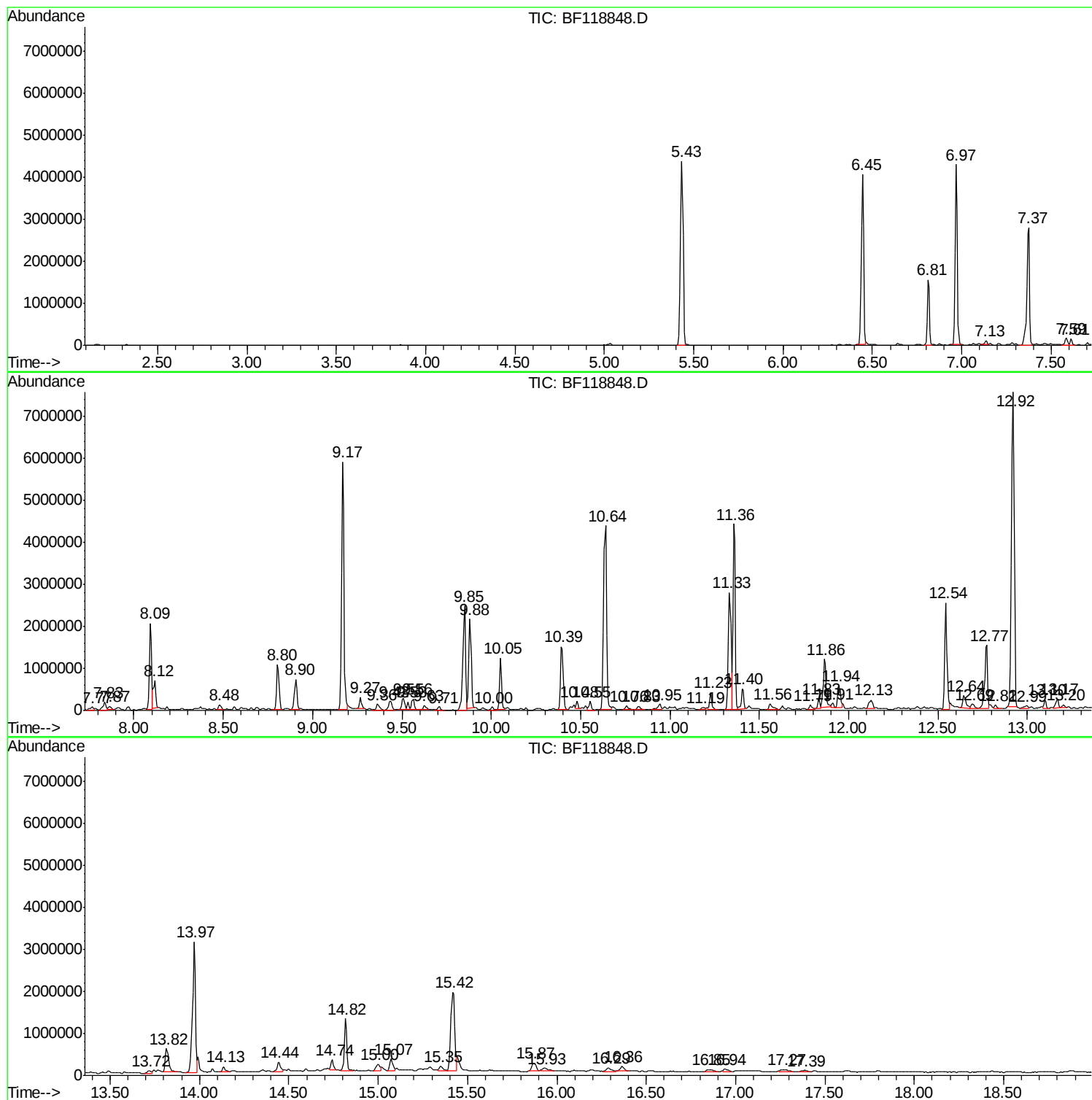
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Sample : L1431-04
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
350

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF020320.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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Data File : BF118848.D
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Sample : L1431-04
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
350

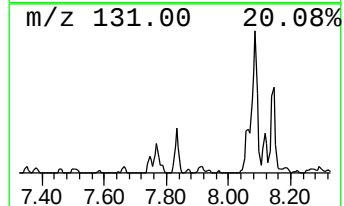
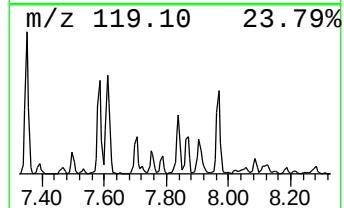
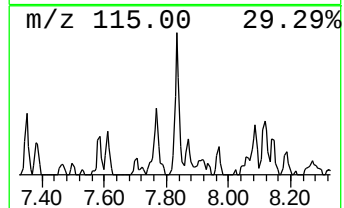
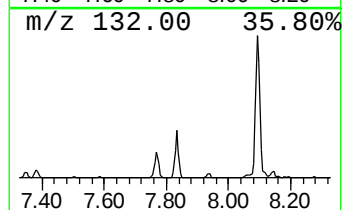
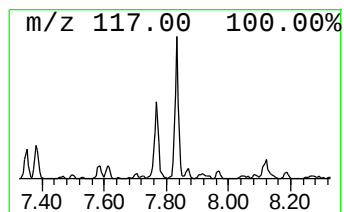
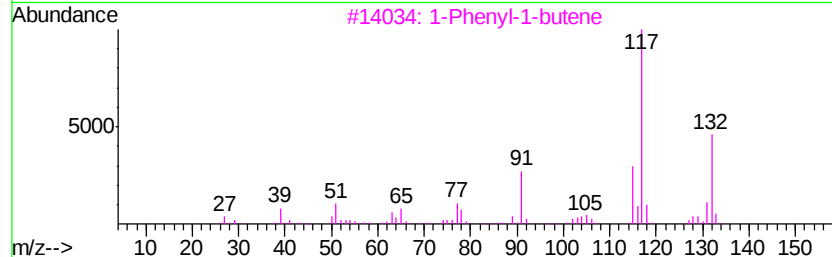
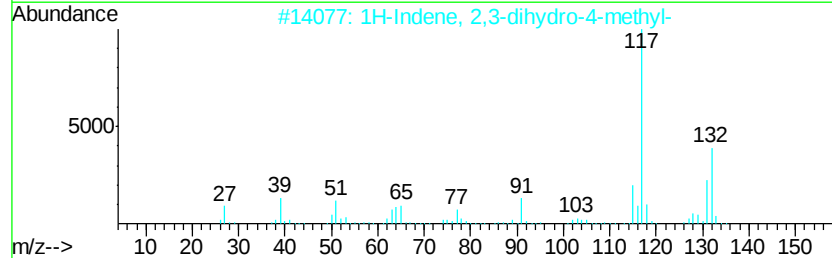
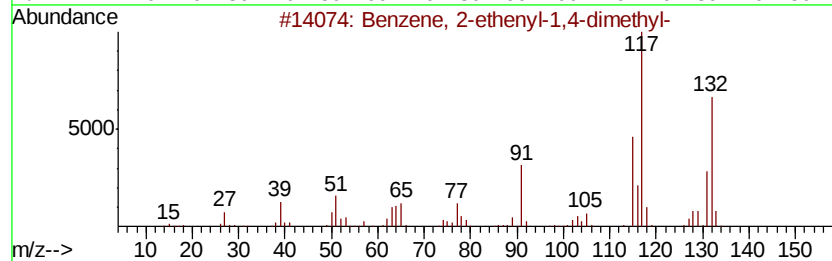
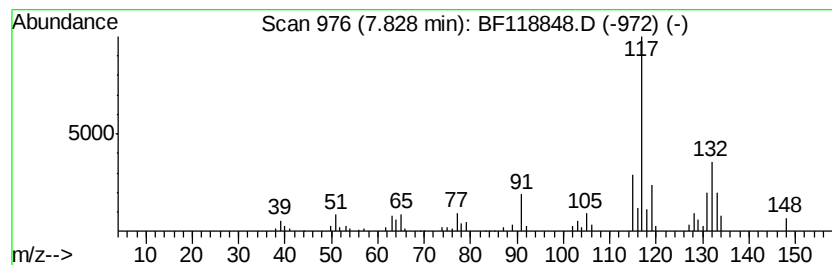
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF020320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.83	2.38 ng	222811	Naphthalene-d8	8.09

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
2			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	64
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	64
4			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	64
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	60



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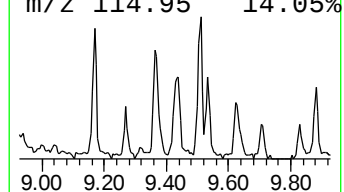
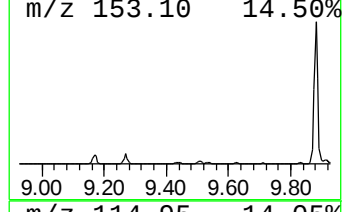
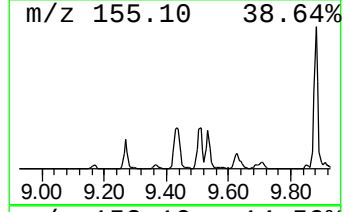
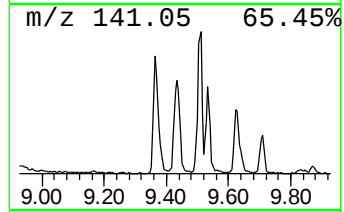
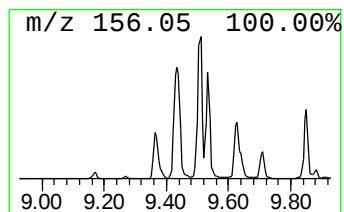
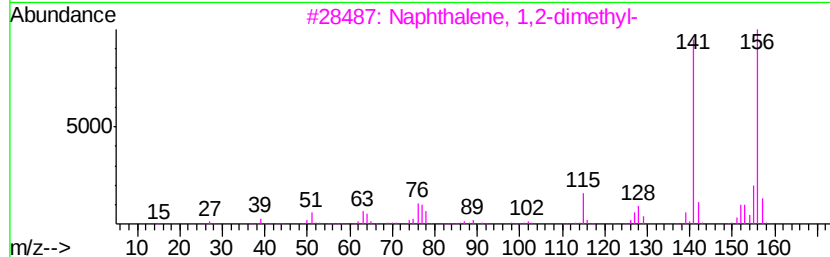
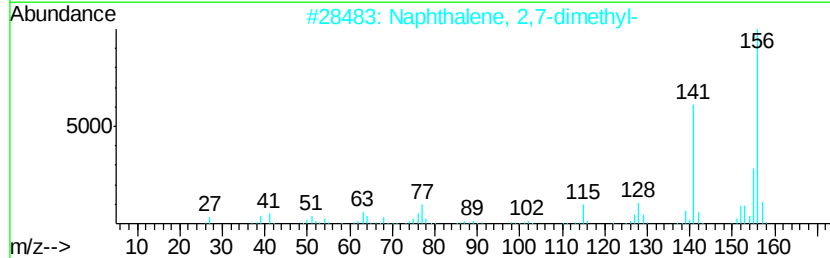
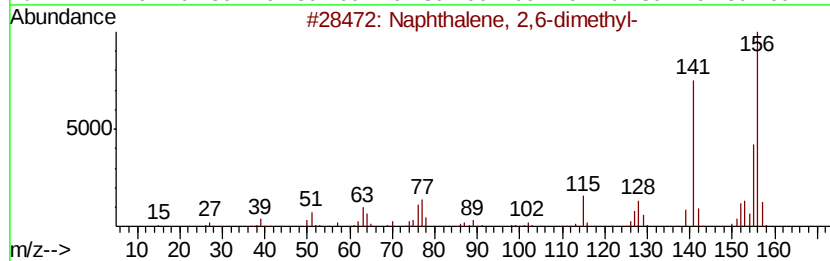
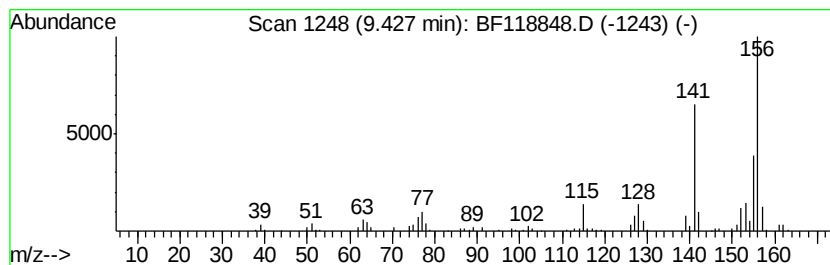
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF020320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Naphthalene, 2,6-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
9.43	2.46 ng	295139	Acenaphthene-d10	9.85		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
2		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3		Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	96
4		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
5		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	96



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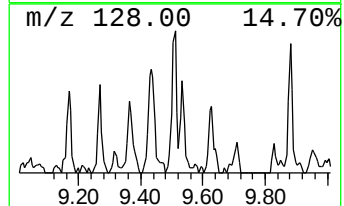
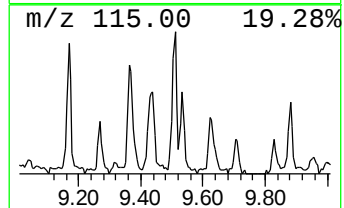
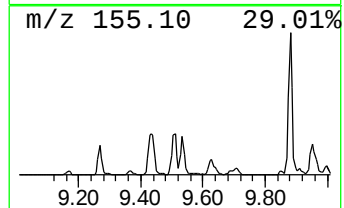
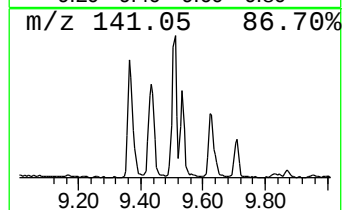
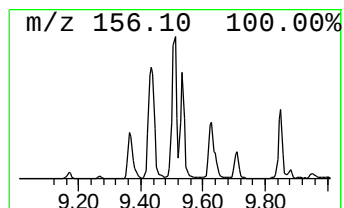
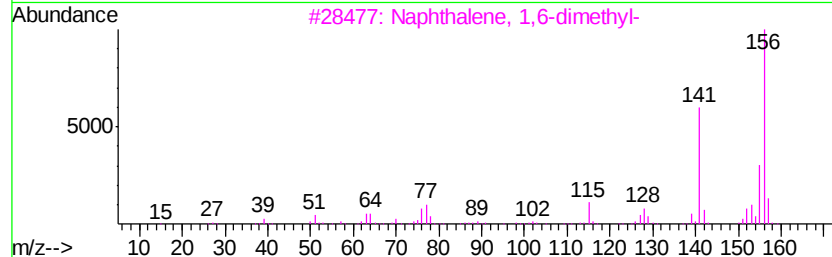
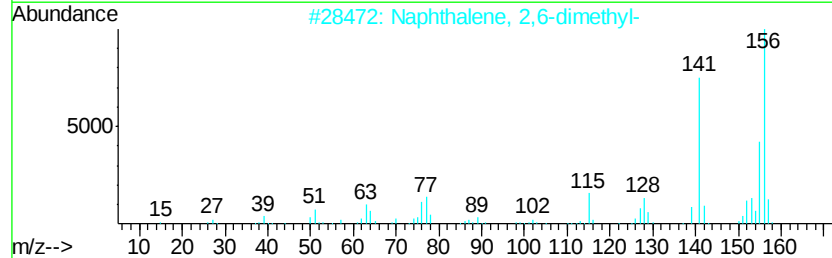
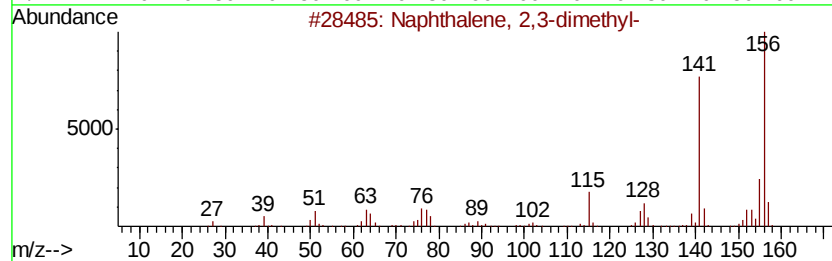
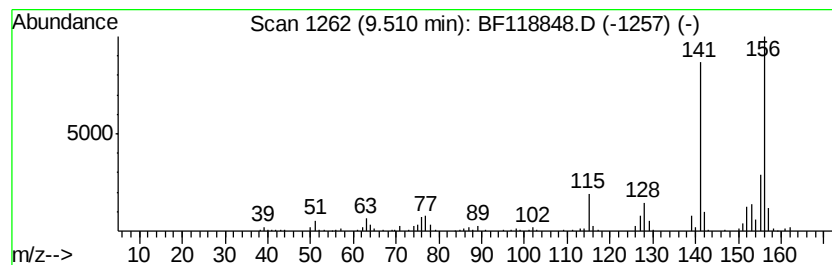
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Naphthalene, 2,3-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.51	2.32 ng	279195	Acenaphthene-d10	9.85
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97
2		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 97
3		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 97
4		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 96
5		Naphthalene, 1,3-dimethyl-	156 C12H12	000575-41-7 96



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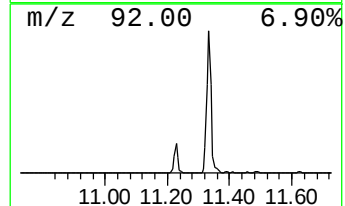
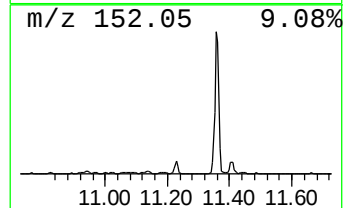
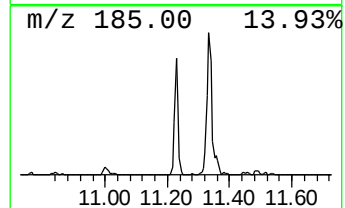
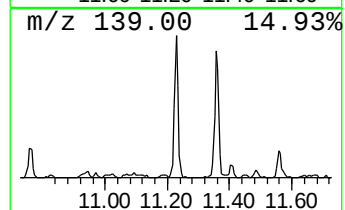
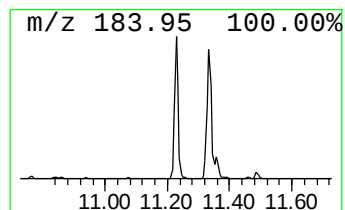
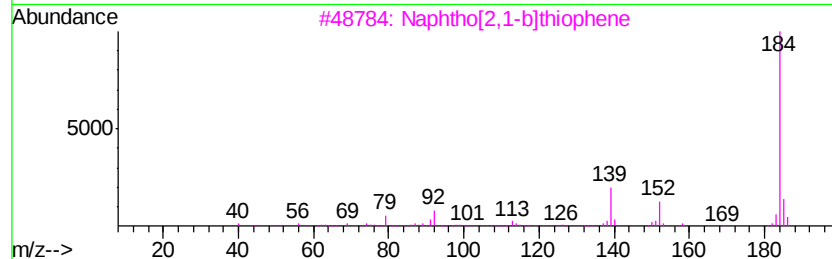
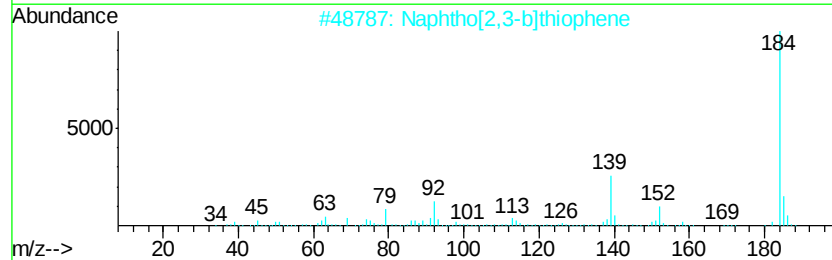
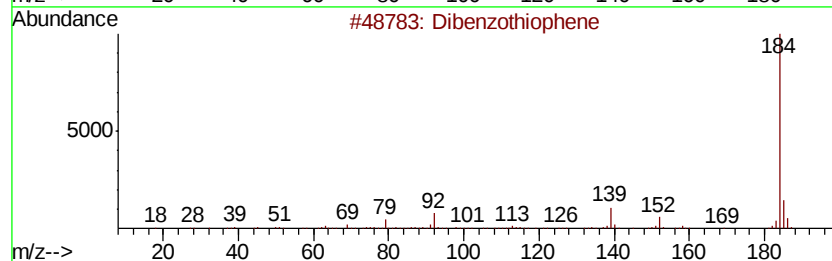
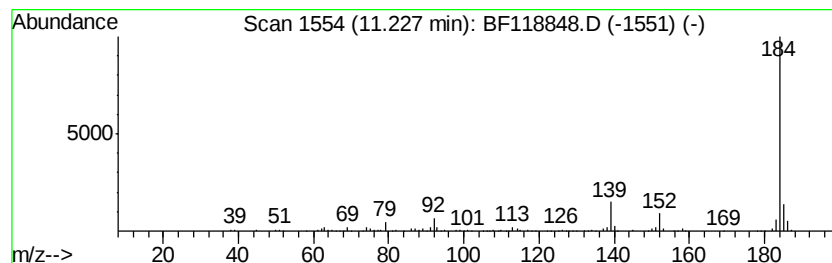
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Dibenzothiophene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.23	2.71 ng	374049	Phenanthrene-d10		11.33
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzothiophene	184	C12H8S	000132-65-0	97
2	Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	95
3	Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	95
4	Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	95
5	1,1'-Biphenyl, 2-methoxy-	184	C13H12O	000086-26-0	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF020720\
Data File : BF118848.D
Acq On : 7 Feb 2020 14:08
Operator : CG/JU
Sample : L1431-04
Misc :
ALS Vial : 5 Sample Multiplier: 1

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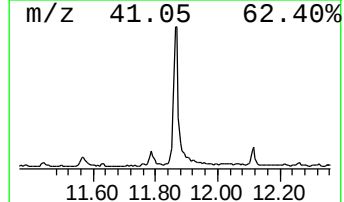
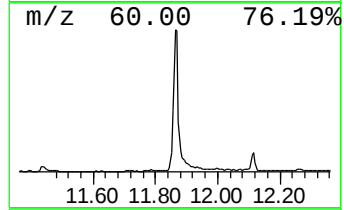
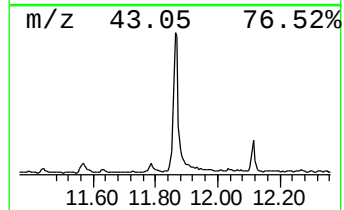
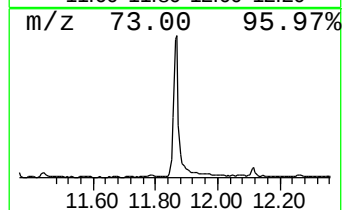
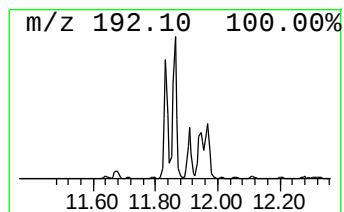
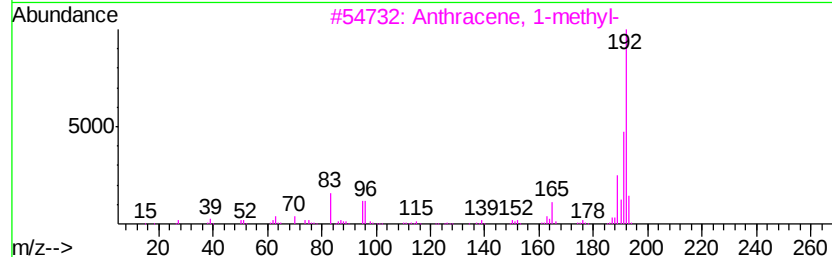
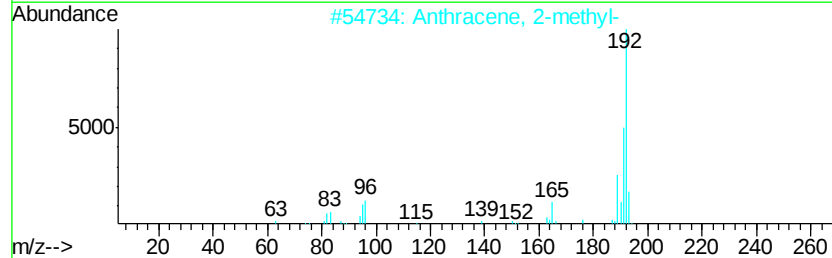
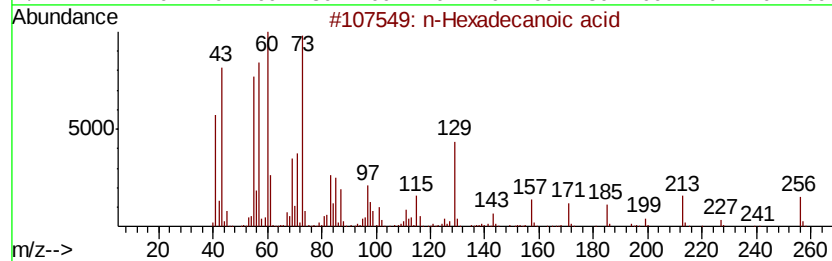
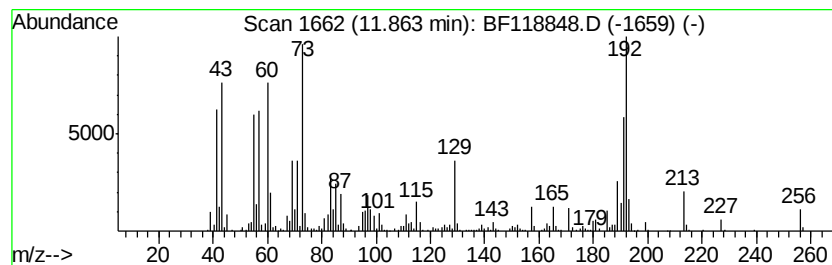
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.86	8.67 ng	1197310	Phenanthrene-d10	11.33

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
3		Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	92
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	92



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF020720\
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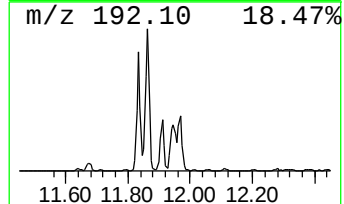
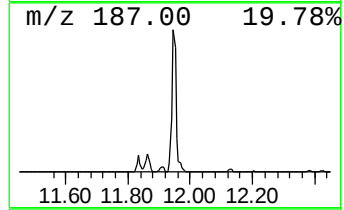
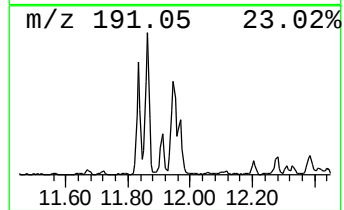
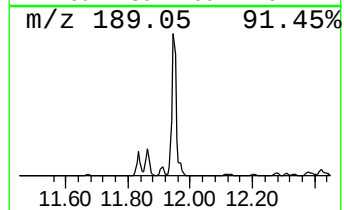
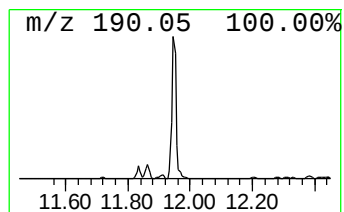
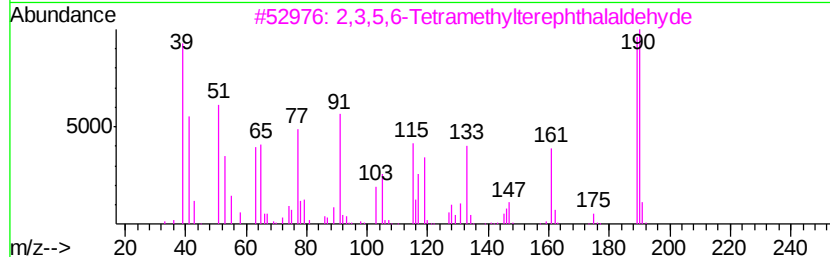
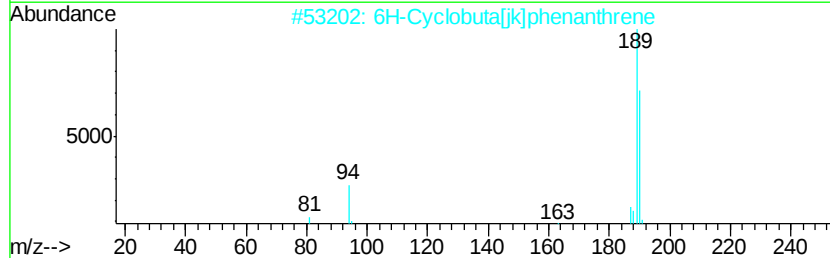
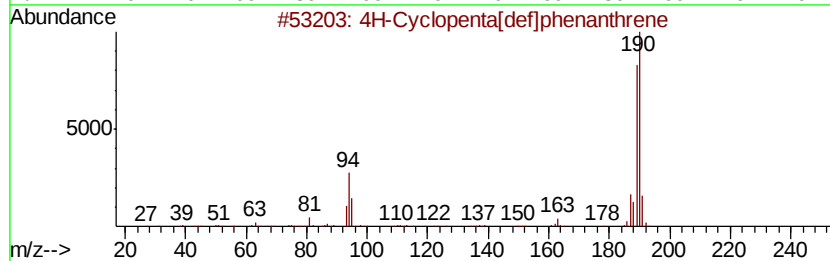
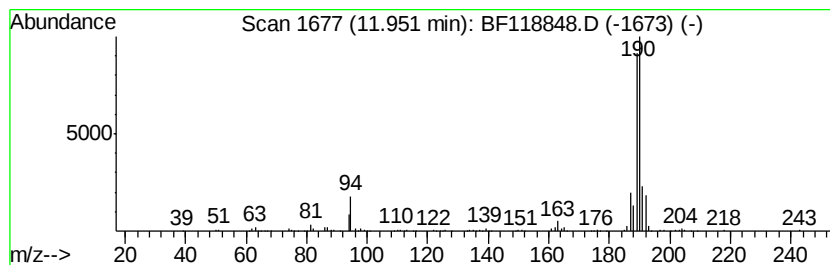
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.95	3.78 ng	521569	Phenanthrene-d10	11.33		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2		6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	72
3		2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	50
4		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45
5		Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	40



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF020720\
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Sample : L1431-04
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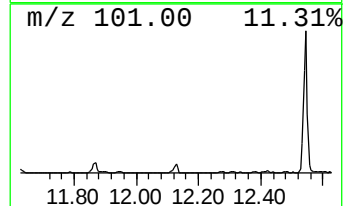
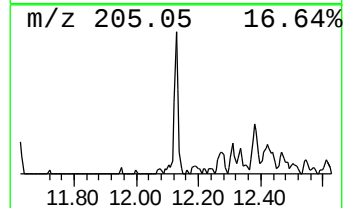
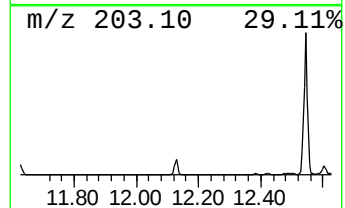
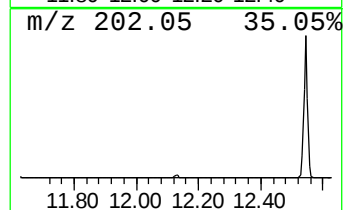
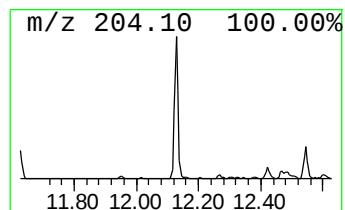
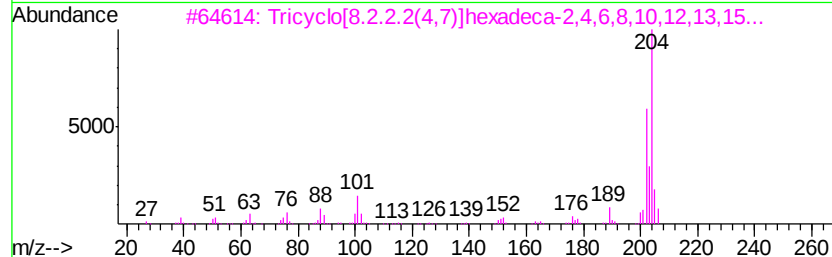
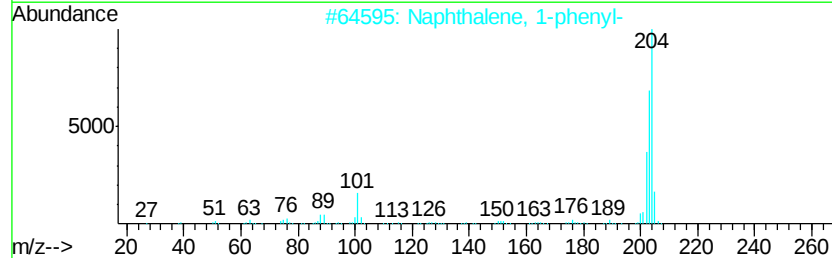
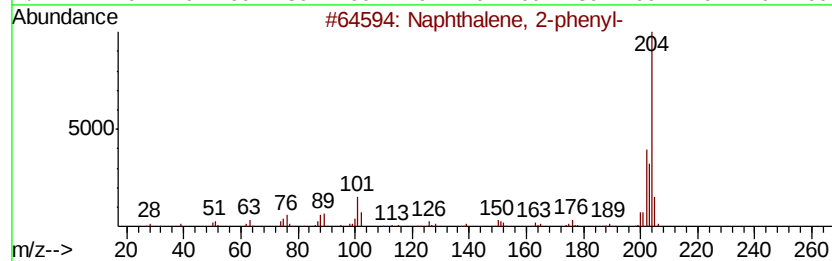
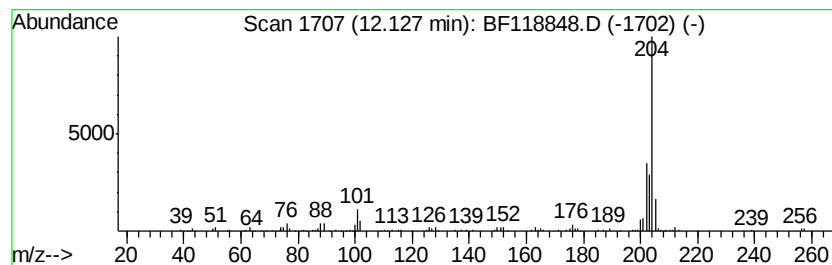
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Naphthalene, 2-phenyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.13	2.01 ng	277020	Phenanthrene-d10	11.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	96
2		Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	93
3		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	91
4		5,16[1',2'1:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	90
5		1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF020720\
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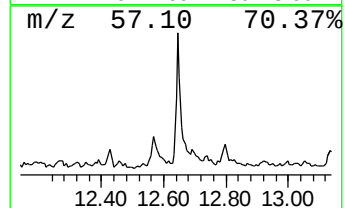
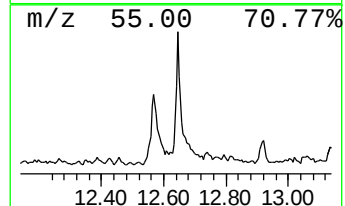
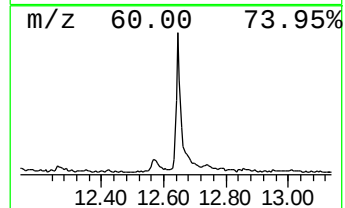
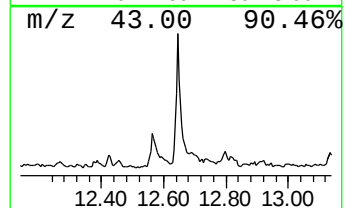
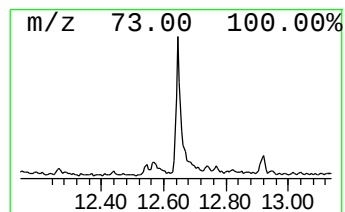
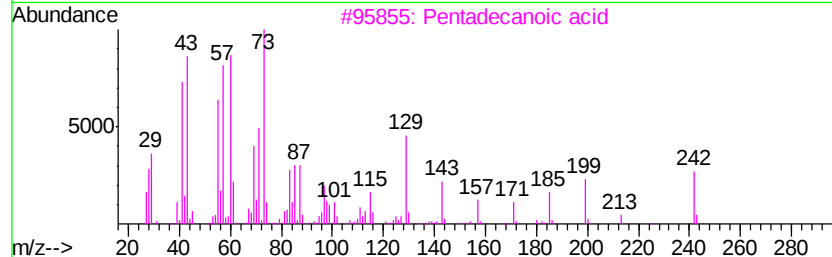
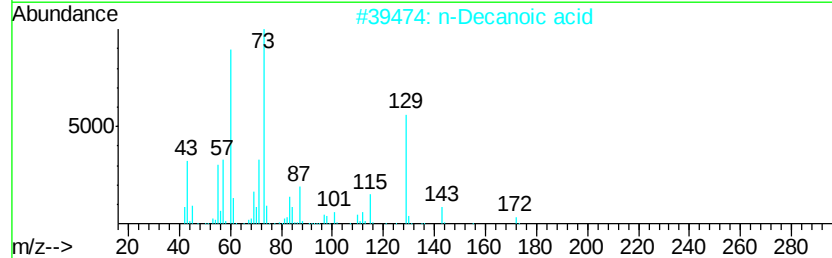
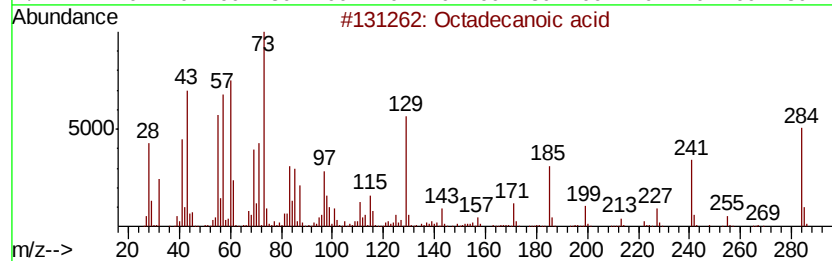
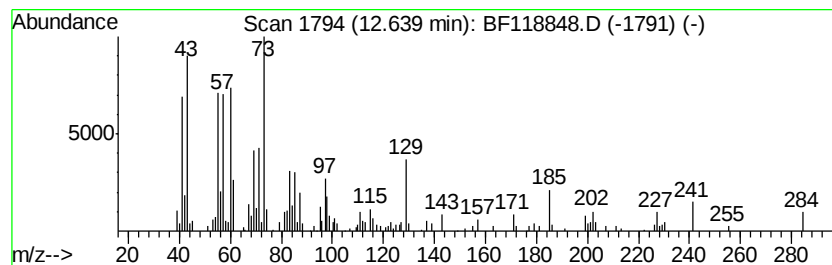
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.64	2.49 ng	344346	Phenanthrene-d10		11.33	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		n-Decanoic acid	172	C10H20O2	000334-48-5	90
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	83
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	64
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF020720\
Data File : BF118848.D
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Instrument :
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ClientSampleId :
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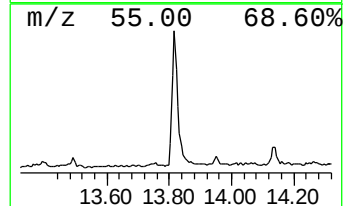
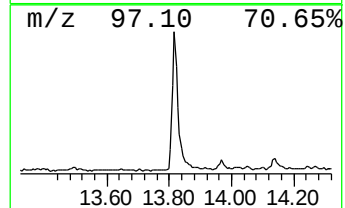
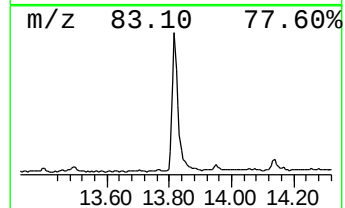
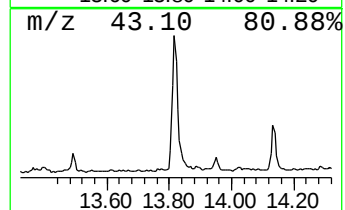
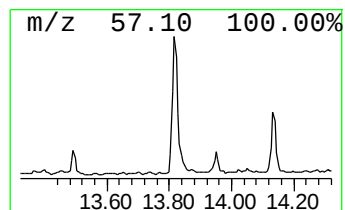
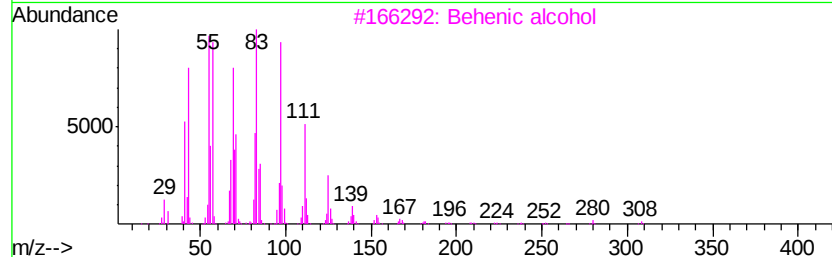
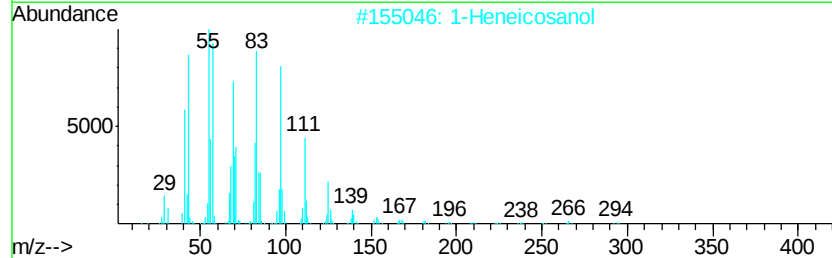
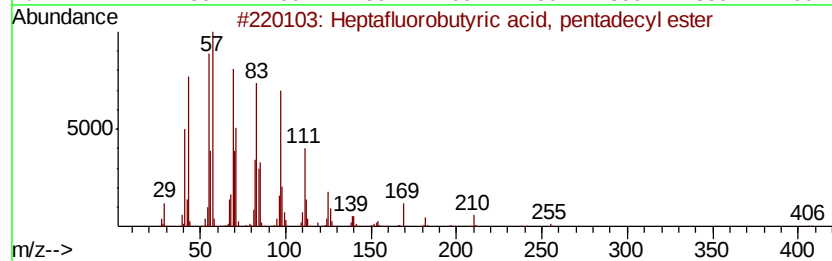
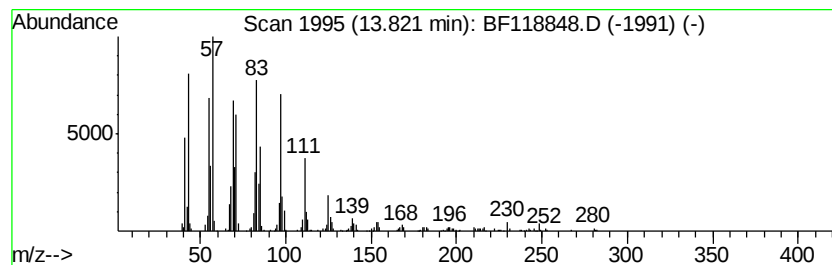
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptafluorobutyric acid, pe... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.82	4.03 ng	692911	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
2		1-Heneicosanol	312	C21H44O	015594-90-8	93
3		Behenic alcohol	326	C22H46O	000661-19-8	93
4		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
5		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF020720\
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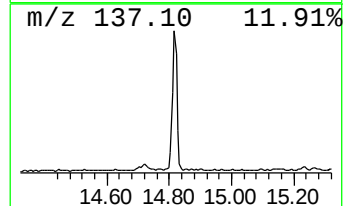
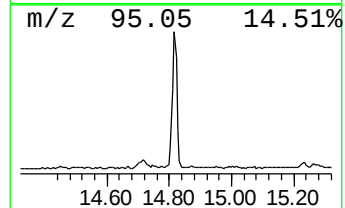
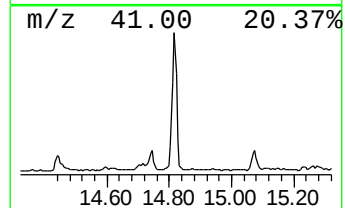
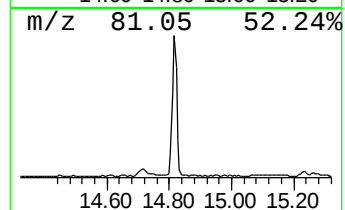
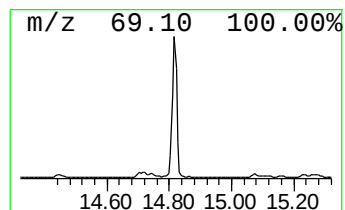
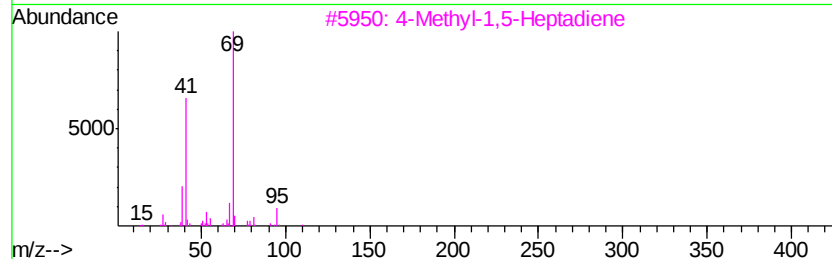
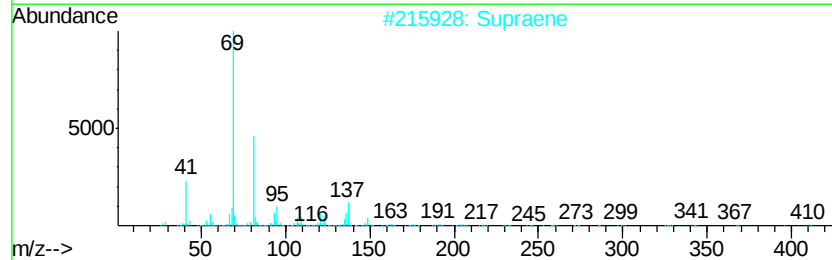
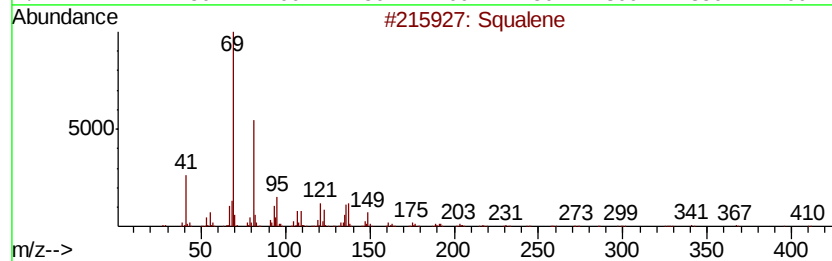
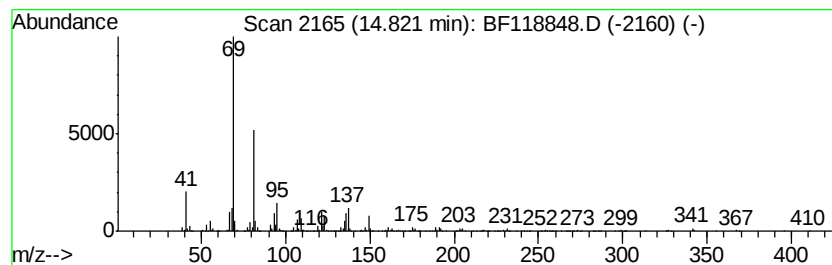
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Squalene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	9.25 ng	1237860	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	96
2		Supraene	410	C30H50	007683-64-9	91
3		4-Methyl-1,5-Heptadiene	110	C8H14	000998-94-7	52
4		2,6-Octadiene, 4-methyl-	124	C9H16	074498-94-5	50
5		1,5-Heptadiene, 2,6-dimethyl-	124	C9H16	006709-39-3	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF020720\
Data File : BF118848.D
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Sample : L1431-04
Misc :
ALS Vial : 5 Sample Multiplier: 1

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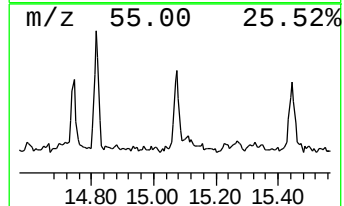
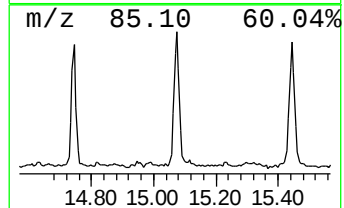
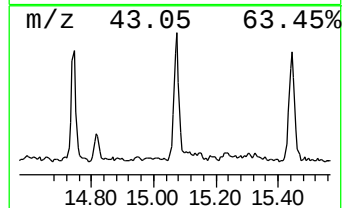
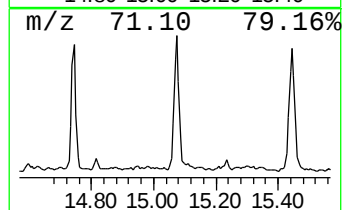
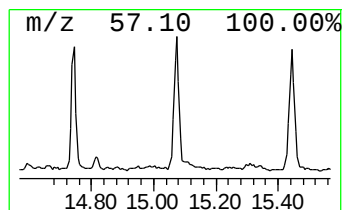
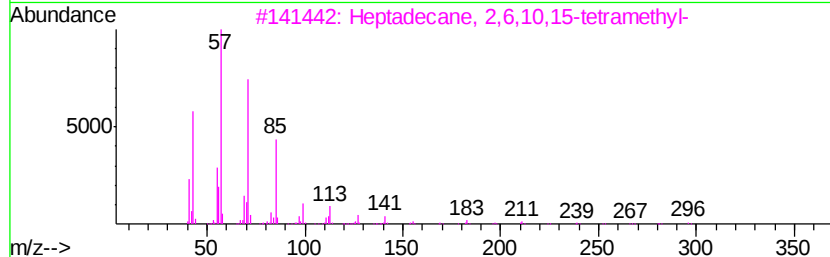
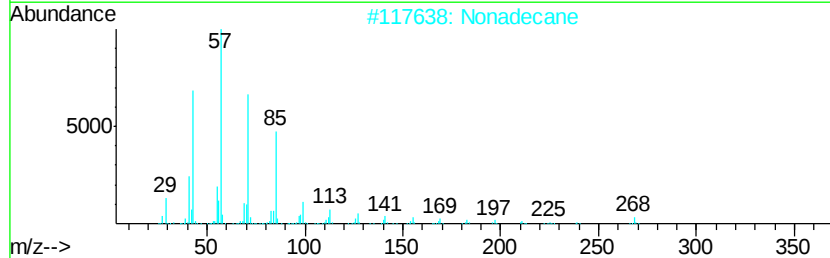
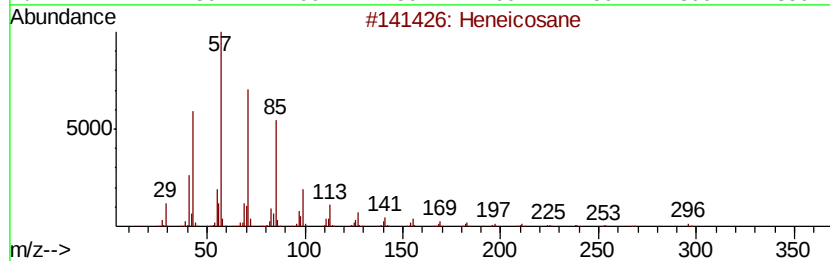
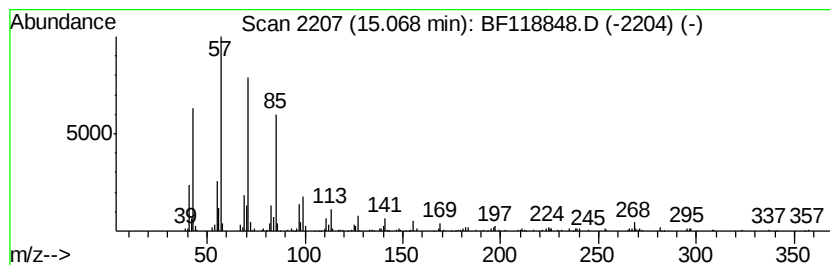
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	2.36 ng	315490	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	99
2		Nonadecane	268	C19H40	000629-92-5	96
3		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
4		2-methyloctacosane	408	C29H60	1000376-72-8	90
5		Hexacosane	366	C26H54	000630-01-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF020720\
Data File : BF118848.D
Acq On : 7 Feb 2020 14:08
Operator : CG/JU
Sample : L1431-04
Misc :
ALS Vial : 5 Sample Multiplier: 1

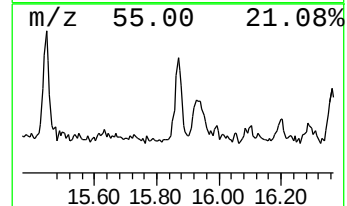
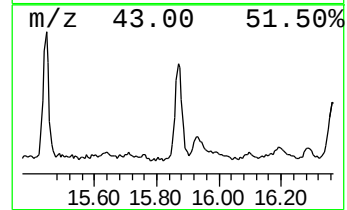
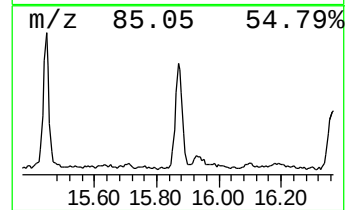
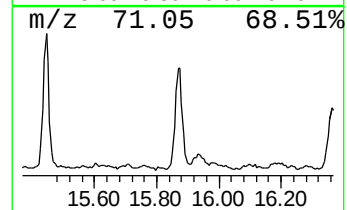
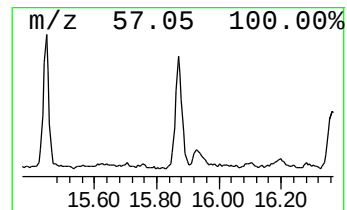
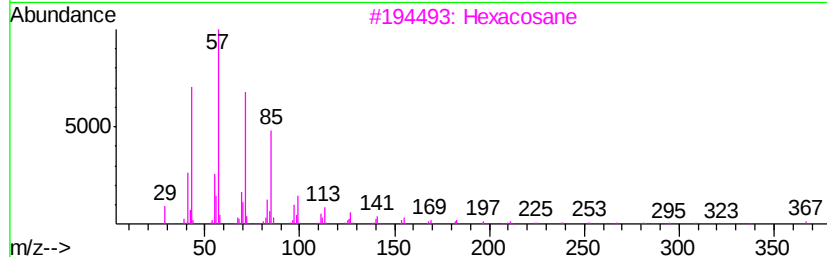
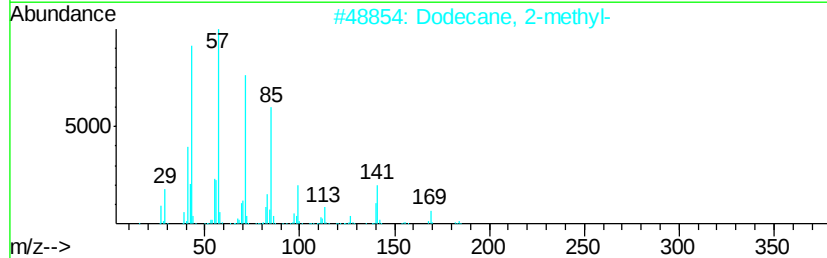
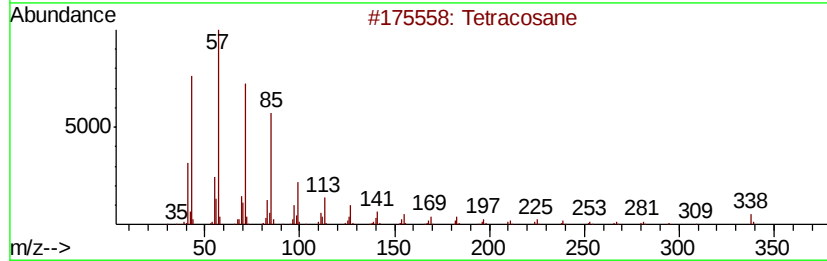
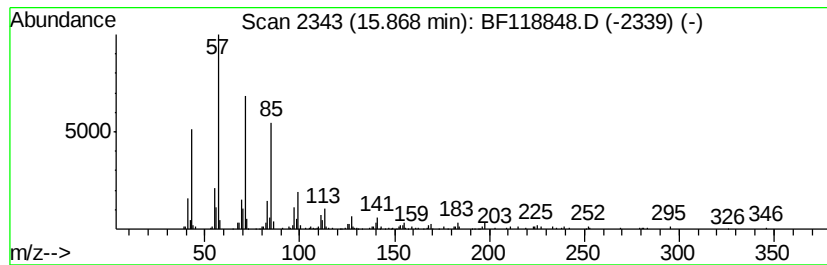
Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF020320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tetracosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.87	2.09 ng	279834	Perylene-d12	15.42
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tetracosane	338 C24H50	000646-31-1 91
2		Dodecane, 2-methyl-	184 C13H28	001560-97-0 90
3		Hexacosane	366 C26H54	000630-01-3 90
4		Heneicosane	296 C21H44	000629-94-7 90
5		Octadecane	254 C18H38	000593-45-3 87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF020720\
 Data File : BF118848.D
 Acq On : 7 Feb 2020 14:08
 Operator : CG/JU
 Sample : L1431-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 350

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF020320.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
Benzene, 2-etheny...	7.83	2.4	ng	222811	2	8.09	1875040	20.0
Naphthalene, 2,6-...	9.43	2.5	ng	295139	3	9.85	2404160	20.0
Naphthalene, 2,3-...	9.51	2.3	ng	279195	3	9.85	2404160	20.0
Dibenzothiophene	11.23	2.7	ng	374049	4	11.33	2761560	20.0
n-Hexadecanoic acid	11.86	8.7	ng	1197310	4	11.33	2761560	20.0
4H-Cyclopenta[def...	11.95	3.8	ng	521569	4	11.33	2761560	20.0
Naphthalene, 2-ph...	12.13	2.0	ng	277020	4	11.33	2761560	20.0
Octadecanoic acid	12.64	2.5	ng	344346	4	11.33	2761560	20.0
Heptafluorobutyri...	13.82	4.0	ng	692911	5	13.97	3439150	20.0
Squalene	14.82	9.3	ng	1237860	6	15.42	2675640	20.0
Heneicosane	15.07	2.4	ng	315490	6	15.42	2675640	20.0
Tetracosane	15.87	2.1	ng	279834	6	15.42	2675640	20.0