

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF103019\
 Data File : BF117623.D
 Acq On : 30 Oct 2019 14:39
 Operator : JU
 Sample : K5542-07
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 198-90-(0-2)

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-BF101819.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	4.916	478	481	498	rBV	80372	122432	15.72%	1.317%
2	5.357	552	556	575	rBV	427441	633570	81.33%	6.814%
3	6.387	728	731	747	rBV	552449	750739	96.37%	8.074%
4	6.504	747	751	763	rVB	791600	778984	100.00%	8.378%
5	6.728	785	789	798	rBV	230040	215640	27.68%	2.319%
6	6.881	812	815	833	rBV	650628	588320	75.52%	6.328%
7	7.293	882	885	907	rBV	349108	466727	59.91%	5.020%
8	8.010	1003	1007	1023	rBV	223689	277100	35.57%	2.980%
9	9.081	1185	1189	1201	rBV	785309	725519	93.14%	7.803%
10	9.481	1254	1257	1265	rVB	82942	77285	9.92%	0.831%
11	9.757	1301	1304	1308	rBV	278752	277981	35.69%	2.990%
12	9.792	1308	1310	1318	rVB	23364	26120	3.35%	0.281%
13	9.981	1337	1342	1345	rBV2	6125	8591	1.10%	0.092%
14	10.316	1395	1399	1405	rBV2	15186	17289	2.22%	0.186%
15	10.557	1436	1440	1451	rBV	303999	347982	44.67%	3.743%
16	11.110	1530	1534	1537	rBV3	6637	8324	1.07%	0.090%
17	11.151	1537	1541	1545	rVB	6435	8669	1.11%	0.093%
18	11.251	1554	1558	1560	rBV	236193	238446	30.61%	2.565%
19	11.275	1560	1562	1568	rVV	203517	193327	24.82%	2.079%
20	11.328	1568	1571	1577	rVB	46540	50440	6.48%	0.543%
21	11.498	1594	1600	1604	rBV3	13428	21873	2.81%	0.235%
22	11.751	1640	1643	1645	rBV	22636	21141	2.71%	0.227%
23	11.780	1645	1648	1654	rVV2	61938	81793	10.50%	0.880%
24	11.833	1654	1657	1659	rVV	13413	16443	2.11%	0.177%
25	11.863	1659	1662	1665	rVV	38501	46094	5.92%	0.496%
26	11.886	1665	1666	1672	rVB	17019	14798	1.90%	0.159%
27	12.045	1690	1693	1700	rBV	16887	18064	2.32%	0.194%
28	12.092	1700	1701	1704	rBV	8204	8093	1.04%	0.087%
29	12.192	1714	1718	1722	rBV3	6687	8679	1.11%	0.093%
30	12.304	1733	1737	1739	rBV	16443	19079	2.45%	0.205%
31	12.327	1739	1741	1745	rVV2	10962	15520	1.99%	0.167%
32	12.398	1749	1753	1757	rVB5	10654	13584	1.74%	0.146%
33	12.463	1760	1764	1769	rBV	320849	293699	37.70%	3.159%
34	12.563	1778	1781	1786	rBV3	11980	15522	1.99%	0.167%

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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
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If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101819.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	12.616	1787	1790	1793	rVV	11778	12169	1.56%	0.131%
36	12.692	1797	1803	1810	rBV	296856	334178	42.90%	3.594%
37	12.745	1810	1812	1818	rVB3	14522	18152	2.33%	0.195%
38	12.827	1822	1826	1831	rVV	648610	544684	69.92%	5.858%
39	12.869	1831	1833	1836	rVV	10699	10310	1.32%	0.111%
40	12.910	1836	1840	1845	rVV6	16850	28731	3.69%	0.309%
41	13.004	1852	1856	1857	rBV3	16869	22802	2.93%	0.245%
42	13.022	1857	1859	1862	rVB2	30384	31308	4.02%	0.337%
43	13.092	1868	1871	1873	rVB	19074	17727	2.28%	0.191%
44	13.127	1873	1877	1885	rBV4	29956	41521	5.33%	0.447%
45	13.222	1890	1893	1895	rBV	22884	21737	2.79%	0.234%
46	13.251	1895	1898	1904	rVV4	16518	21513	2.76%	0.231%
47	13.298	1904	1906	1909	rVB	18024	12980	1.67%	0.140%
48	13.398	1920	1923	1925	rBV3	8660	8148	1.05%	0.088%
49	13.510	1938	1942	1944	rBV4	10361	12320	1.58%	0.133%
50	13.551	1945	1949	1954	rVB	23083	29874	3.83%	0.321%
51	13.645	1963	1965	1967	rBV2	19171	20705	2.66%	0.223%
52	13.663	1967	1968	1971	rVB2	19813	14747	1.89%	0.159%
53	13.698	1971	1974	1976	rBV	20411	19666	2.52%	0.212%
54	13.733	1976	1980	1982	rBV3	41799	60877	7.81%	0.655%
55	13.857	1998	2001	2003	rBV	40364	35691	4.58%	0.384%
56	13.892	2003	2007	2010	rVV2	311710	427632	54.90%	4.599%
57	13.921	2010	2012	2019	rVB	148590	144543	18.56%	1.555%
58	13.974	2019	2021	2023	rBV3	8777	7791	1.00%	0.084%
59	14.004	2023	2026	2028	rVV	18386	17909	2.30%	0.193%
60	14.245	2065	2067	2069	rBV3	8642	9020	1.16%	0.097%
61	14.269	2069	2071	2077	rBV3	39087	69214	8.89%	0.744%
62	14.345	2082	2084	2087	rBV3	18381	25185	3.23%	0.271%
63	14.639	2132	2134	2136	rVB2	22730	20962	2.69%	0.225%
64	14.710	2143	2146	2149	rVB	39775	42401	5.44%	0.456%
65	14.774	2155	2157	2160	rVB4	16971	15940	2.05%	0.171%
66	14.921	2179	2182	2185	rBV	119629	163248	20.96%	1.756%
67	14.951	2185	2187	2191	rVB3	83309	92605	11.89%	0.996%
68	15.027	2197	2200	2210	rVB2	32981	55777	7.16%	0.600%
69	15.216	2228	2232	2235	rBV	65267	76883	9.87%	0.827%
70	15.274	2239	2242	2245	rBV	92993	105758	13.58%	1.137%
71	15.333	2248	2252	2256	rVB	134315	158585	20.36%	1.706%

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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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101819.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

72	15.715	2313	2317	2320	rVB3	24778	31924	4.10%	0.343%
73	16.180	2395	2396	2401	rVB5	17651	17385	2.23%	0.187%
74	16.762	2491	2495	2500	rVB2	25507	46408	5.96%	0.499%
75	17.186	2562	2567	2574	rBV2	22151	40810	5.24%	0.439%

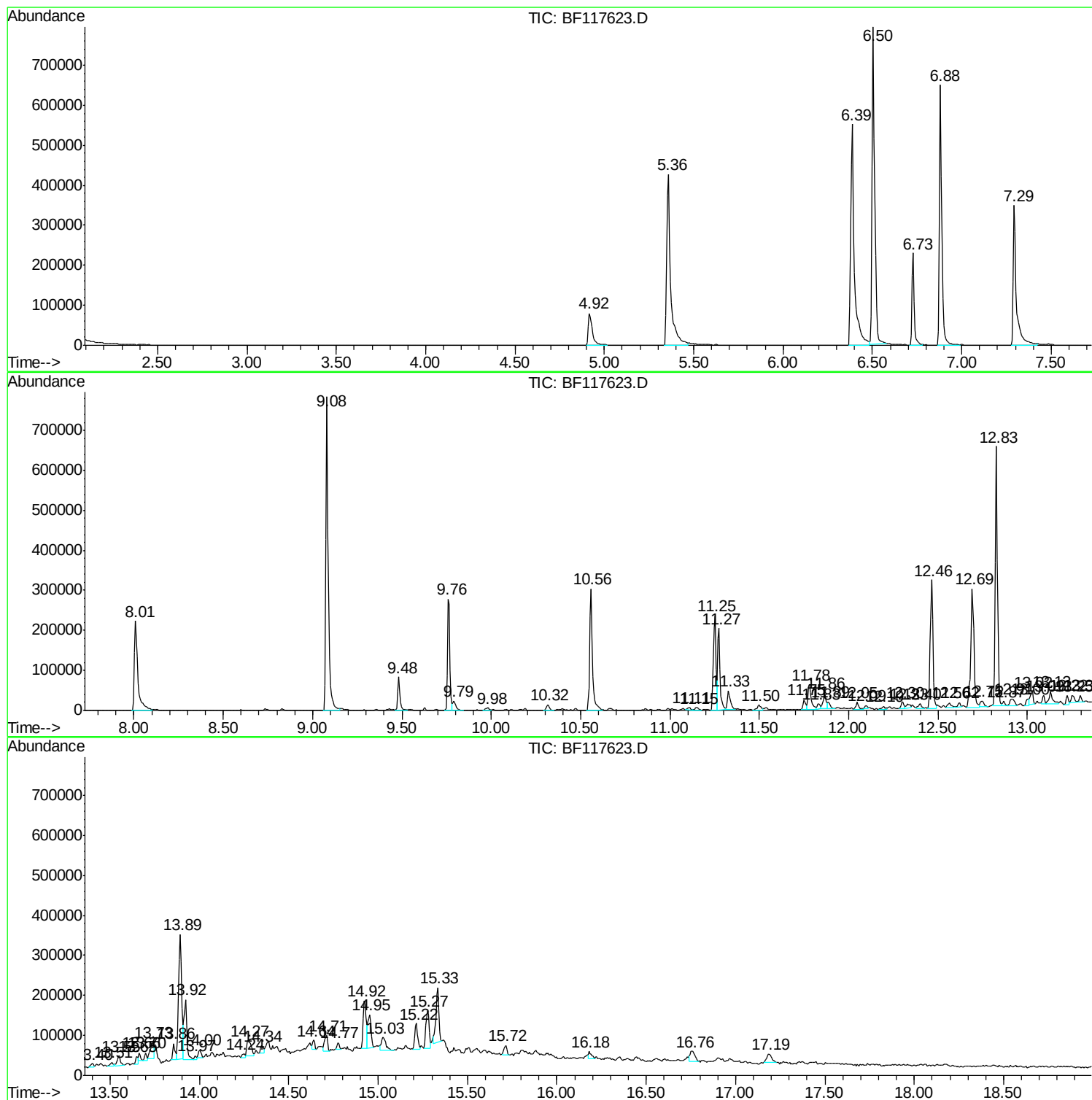
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Instrument :
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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ALS Vial : 10 Sample Multiplier: 1

Instrument :
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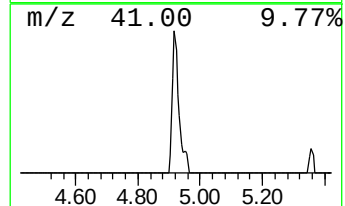
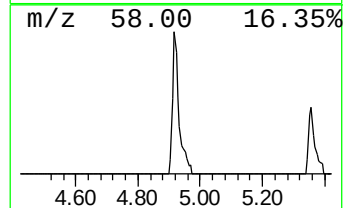
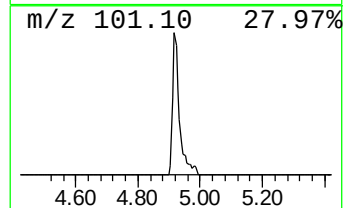
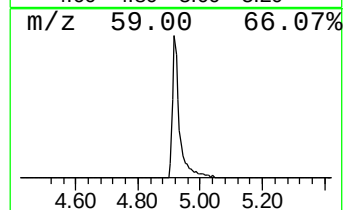
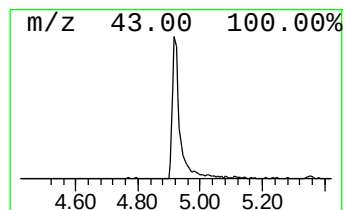
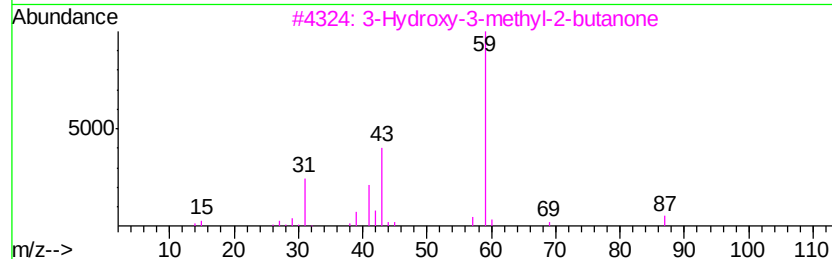
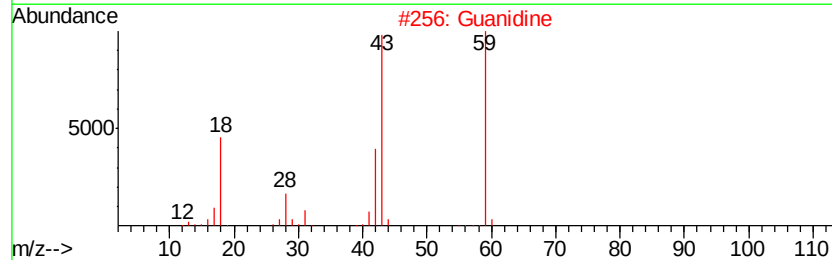
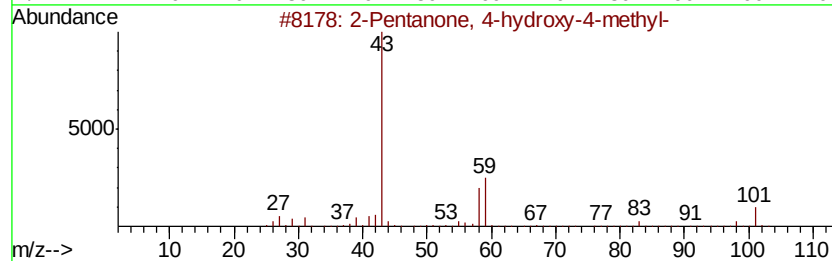
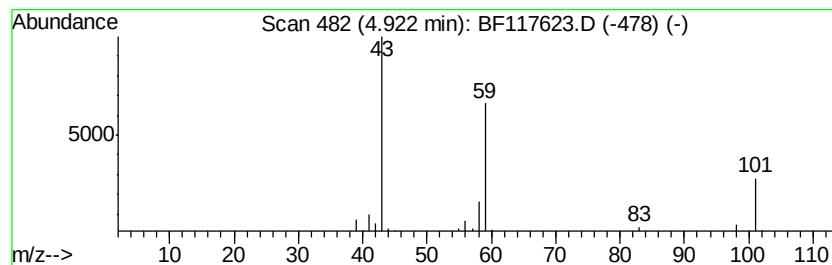
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.92	11.36 ng	122432	1,4-Dichlorobenzene-d4	6.73

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Guanidine	59	CH5N3	000113-00-8	43
3			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	25
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	25
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17



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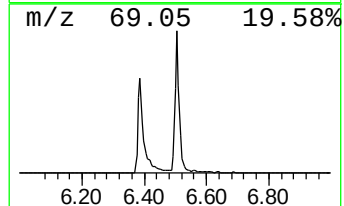
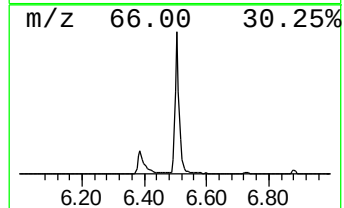
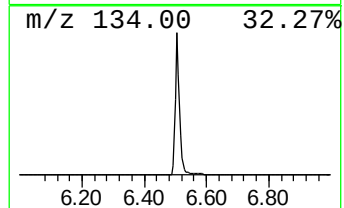
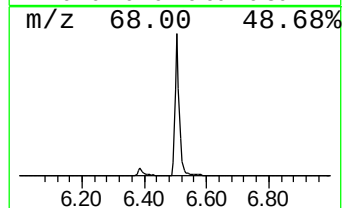
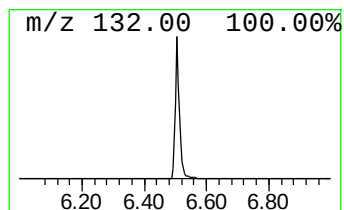
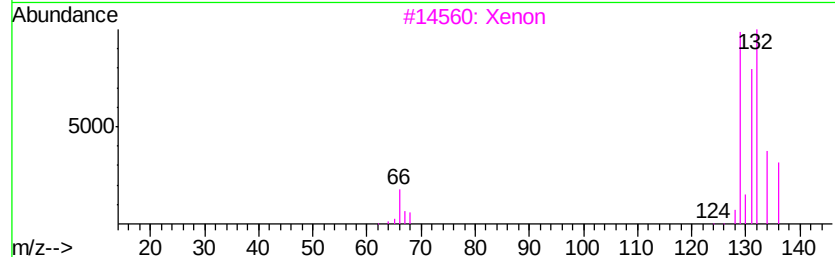
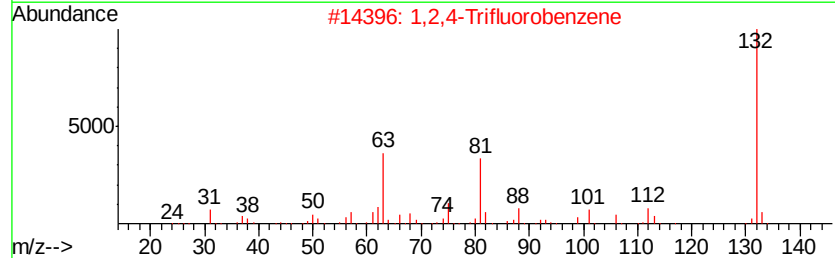
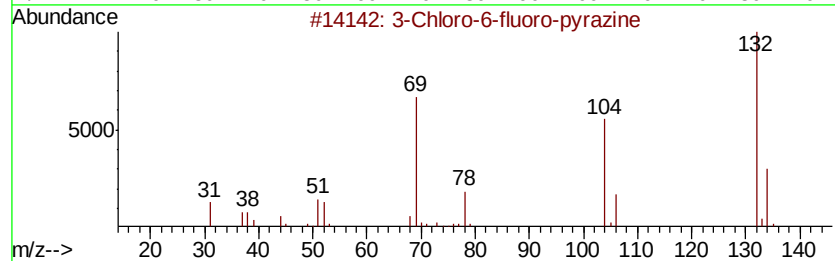
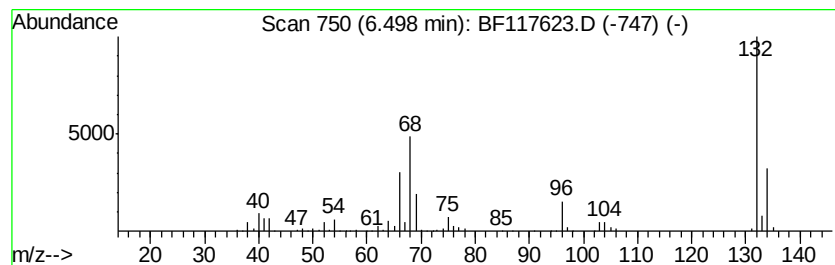
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.50 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.50	72.25 ng	778984	1,4-Dichlorobenzene-d4	6.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9
3		Xenon	132	Xe	007440-63-3	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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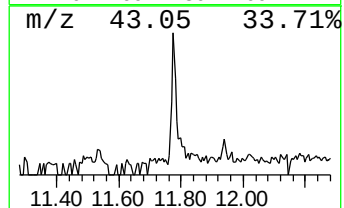
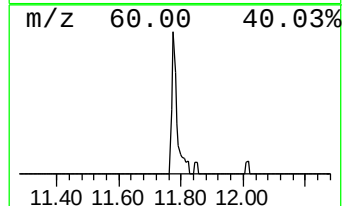
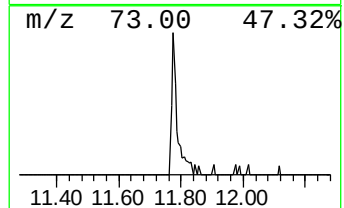
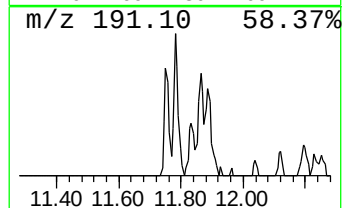
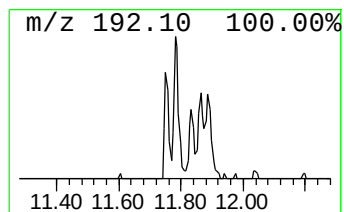
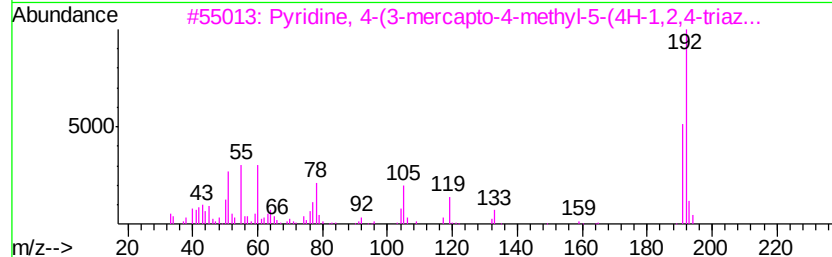
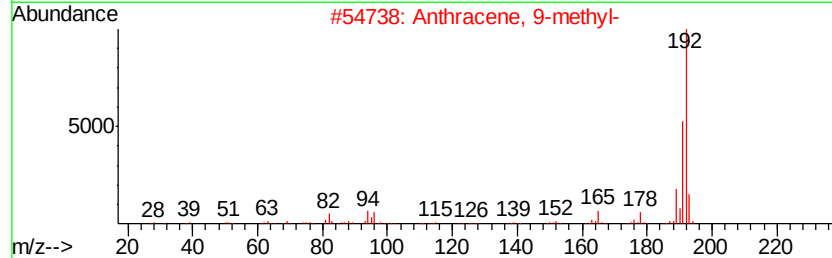
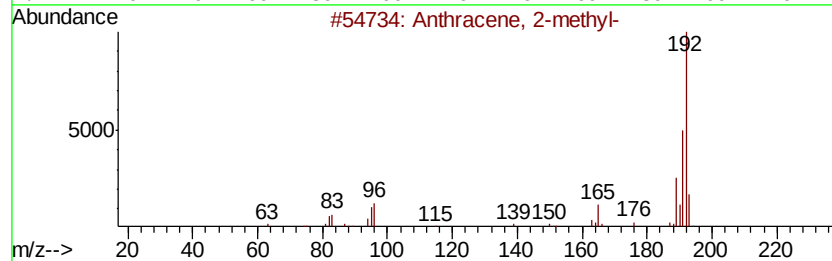
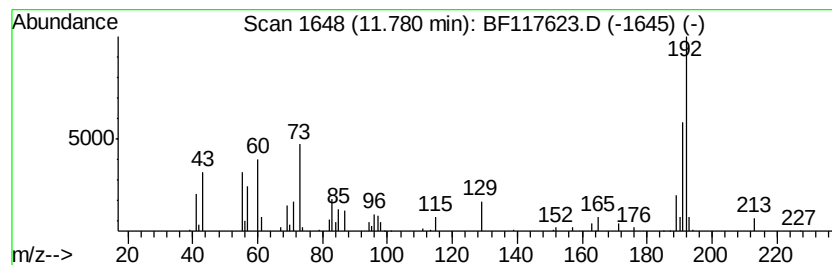
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Anthracene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.78	6.86 ng	81793	Phenanthrene-d10	11.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	62
2		Anthracene, 9-methyl-	192	C15H12	000779-02-2	60
3		Pyridine, 4-(3-mercapto-4-methyl...	192	C8H8N4S	003652-32-2	58
4		Anthracene, 1-methyl-	192	C15H12	000610-48-0	55
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	53



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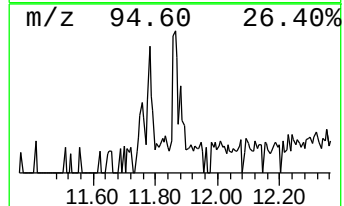
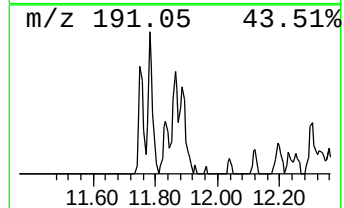
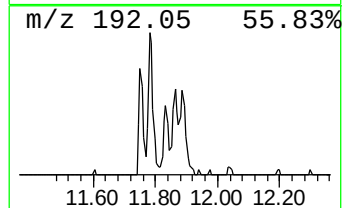
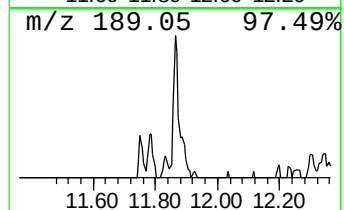
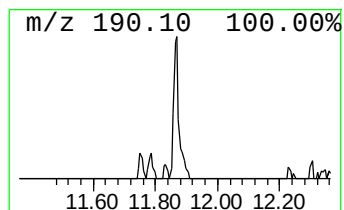
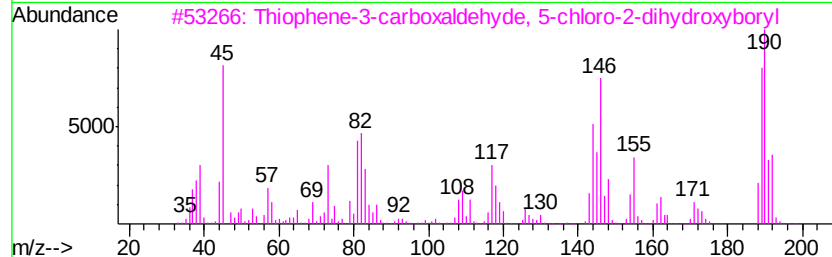
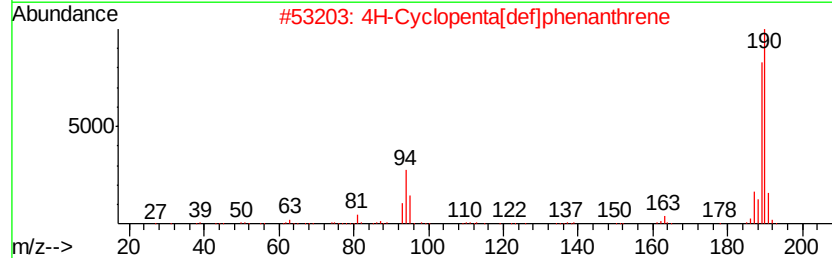
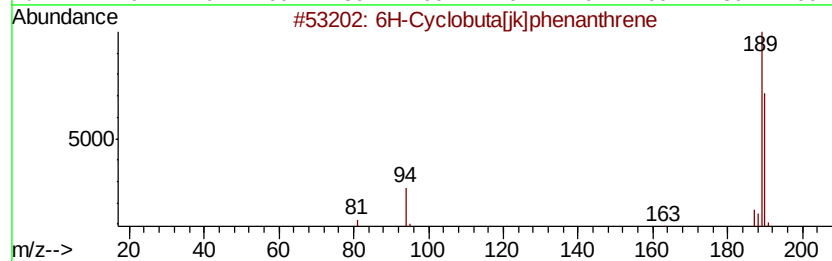
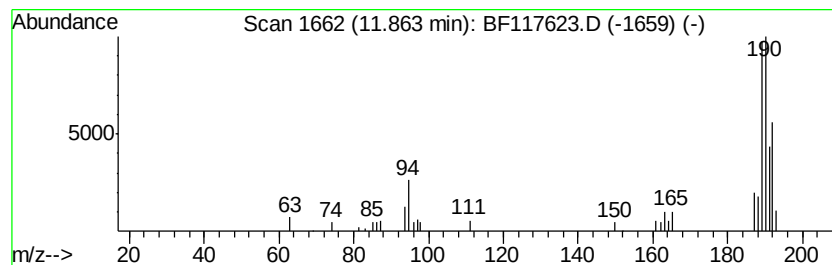
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 6H-Cyclobuta[jk]phenanthrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.86	3.87 ng	46094	Phenanthrene-d10	11.25

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	58
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
3		Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	50
4		1H-Pyrazole-4-carboxaldehyde, 1-...	190	C10H7FN2O	1000338-05-6	38
5		Methyl diselenide	190	C2H6Se2	007101-31-7	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103019\
Data File : BF117623.D
Acq On : 30 Oct 2019 14:39
Operator : JU
Sample : K5542-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

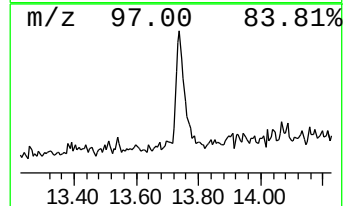
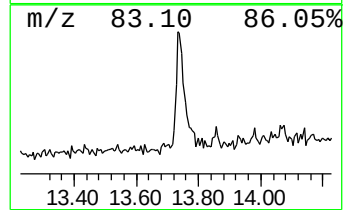
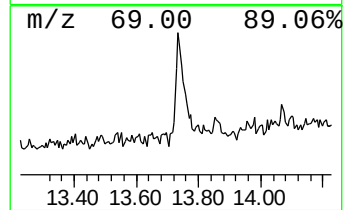
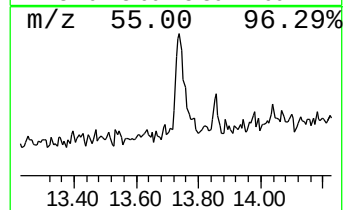
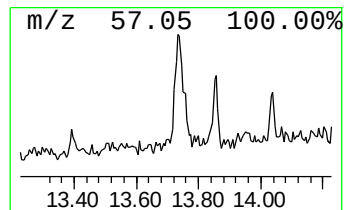
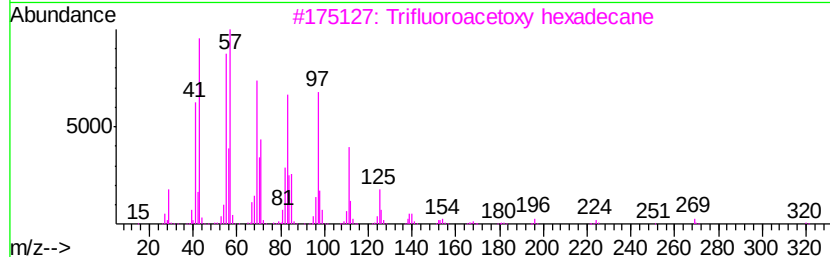
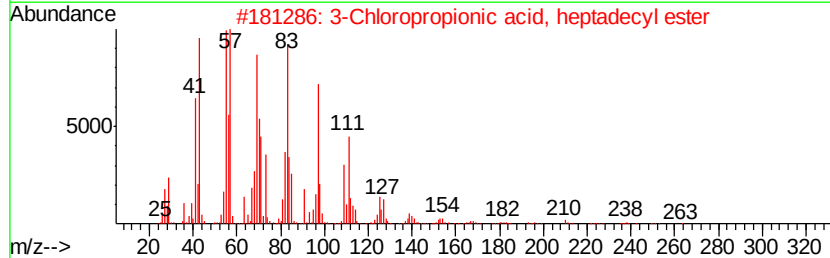
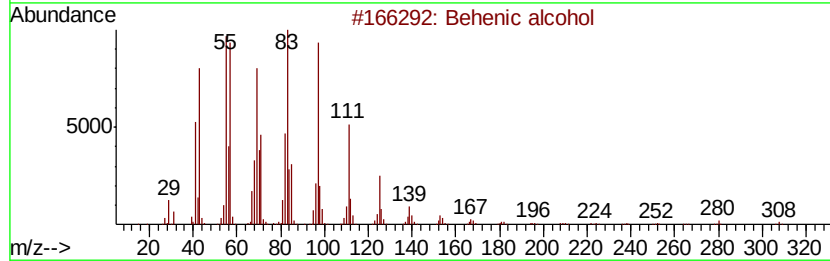
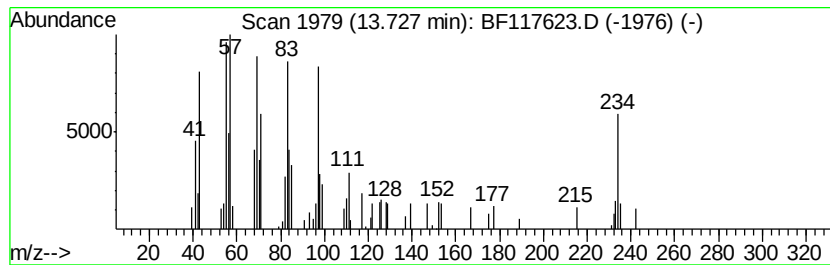
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Behenic alcohol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.73	2.85 ng	60877	Chrysene-d12	13.89		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Behenic alcohol	326	C22H46O	000661-19-8	93
2		3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	91
3		Trifluoroacetoxv hexadecane	338	C18H33F3O2	006222-03-3	87
4		Isoheptadecanol	256	C17H36O	057289-07-3	83
5		Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103019\
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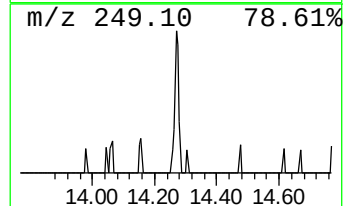
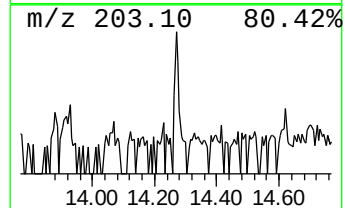
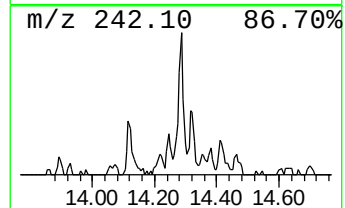
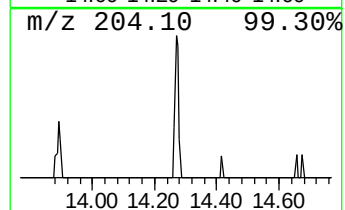
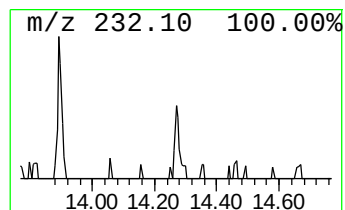
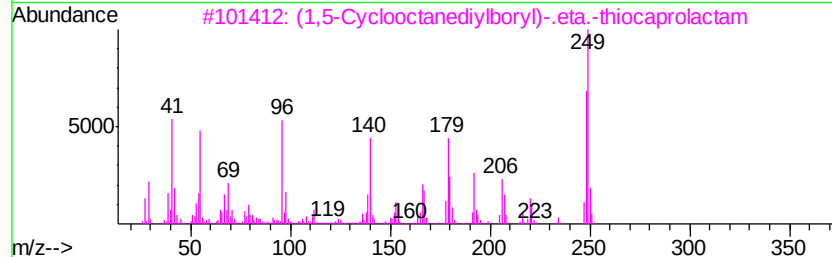
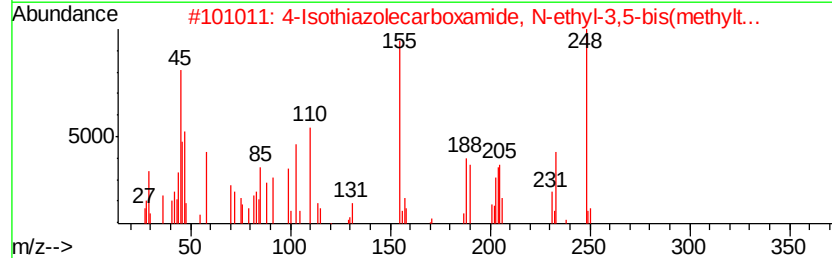
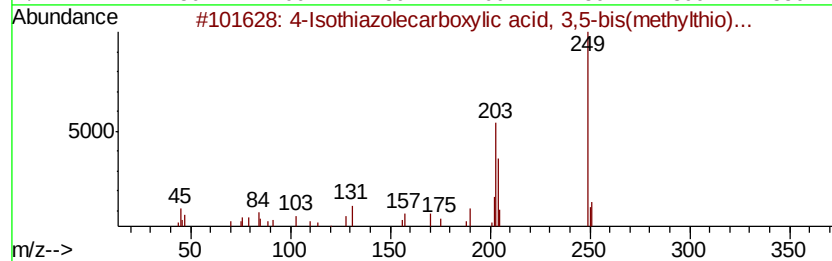
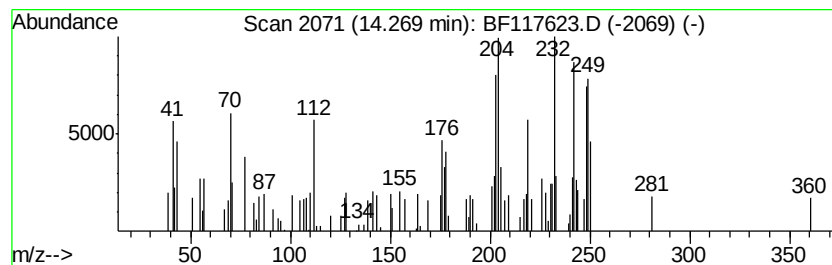
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown14.27 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	3.24 ng	69214	Chrysene-d12	13.89

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4		4-Isothiazolecarboxylic acid, 3,...	249	C8H11N02S3	002273-03-2	22
2	4		4-Isothiazolecarboxamide, N-ethy...	248	C8H12N20S3	037572-35-3	22
3	(1,5-Cyclooctanediylborvl)-.eta...			249	C14H24BNS	1000161-66-4	18
4	2,1,3-Benzothiadiazole, 6-triflu...			249	C7H2F3N302S	1000276-24-3	18
5	Ethyl 5,6-dimethoxyindole-2-carb...			249	C13H15N04	016382-18-6	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103019\
 Data File : BF117623.D
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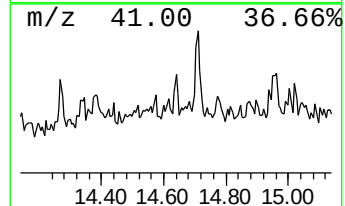
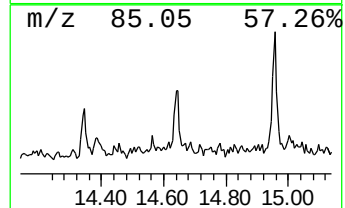
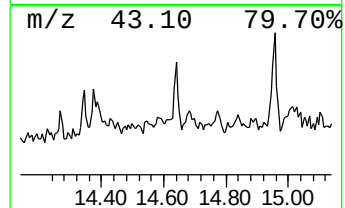
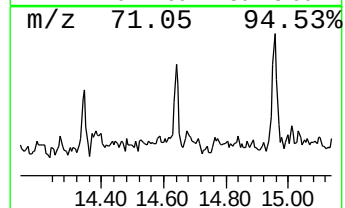
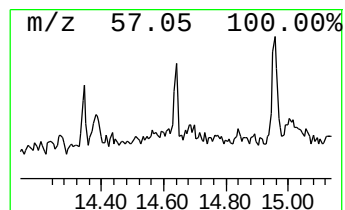
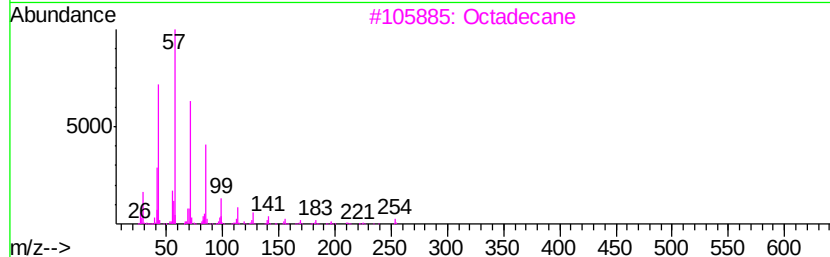
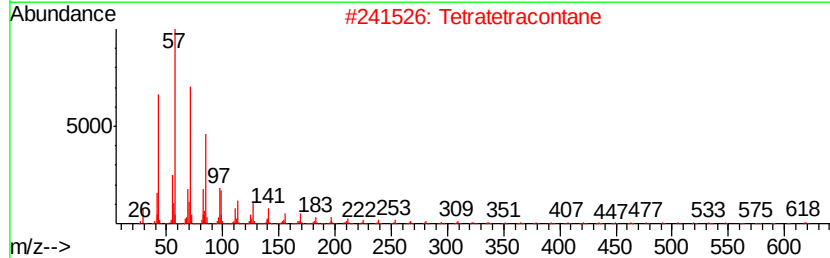
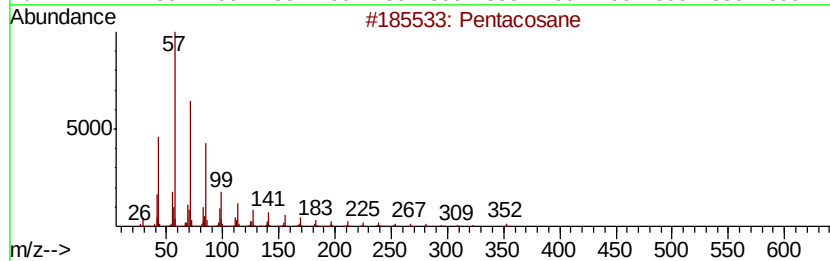
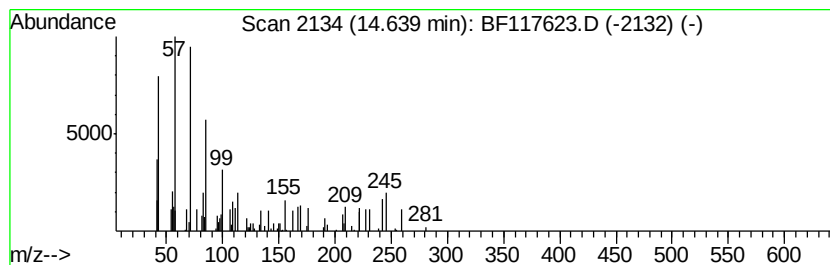
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 8 Pentacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.64	2.64 ng	20962	Perylene-d12	15.33

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentacosane		352	C25H52	000629-99-2	58
2	Tetratetracontane		619	C44H90	007098-22-8	53
3	Octadecane		254	C18H38	000593-45-3	53
4	Tetracosane		338	C24H50	000646-31-1	53
5	Octacosane		394	C28H58	000630-02-4	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103019\
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Misc :
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Instrument :
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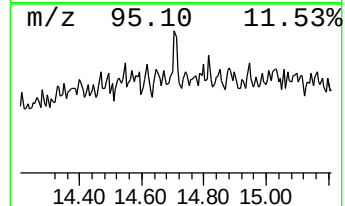
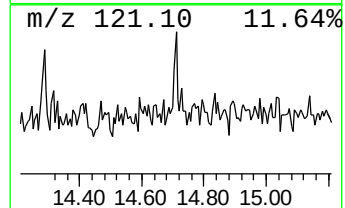
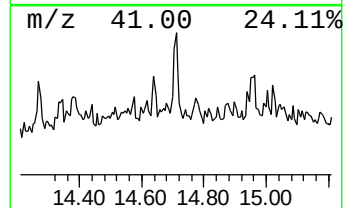
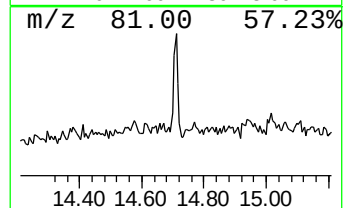
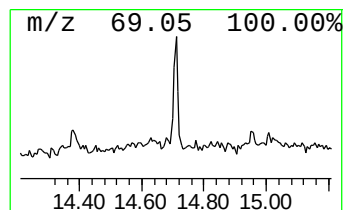
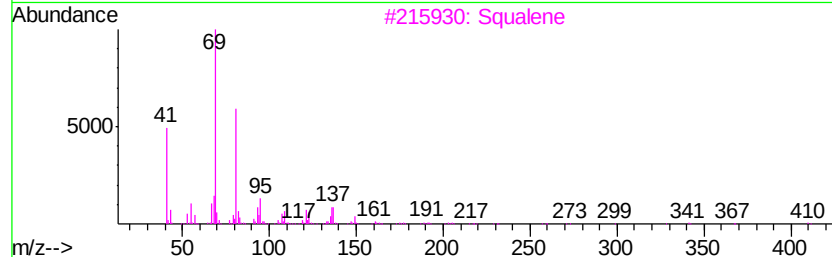
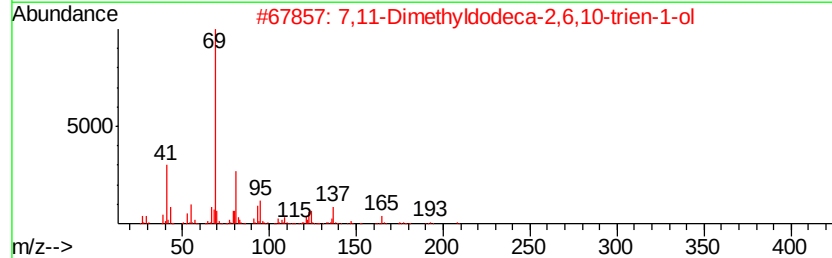
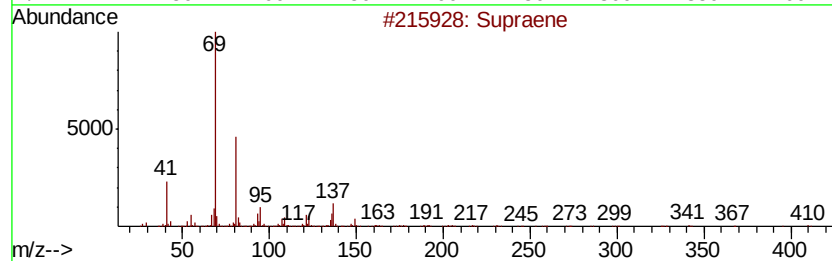
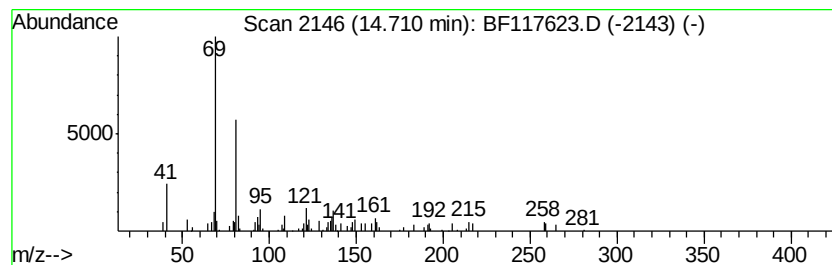
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Supraene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.71	5.35 ng	42401	Perylene-d12	15.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Supraene	410	C30H50	007683-64-9	68
2		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	62
3		Squalene	410	C30H50	000111-02-4	59
4		(.+/-)-Lavandulol, pentafluorop...	300	C13H17F5O2	1000352-63-1	59
5		.beta.-D-Mannofuranoside, farnesyl-	384	C21H36O6	1000155-15-5	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103019\
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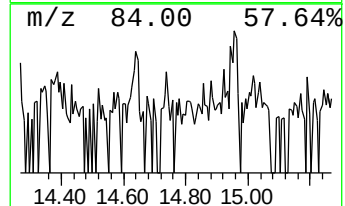
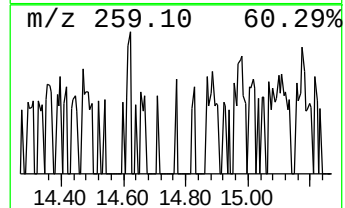
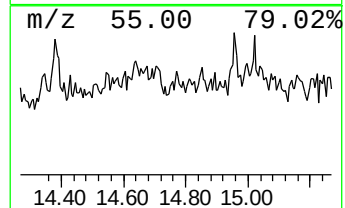
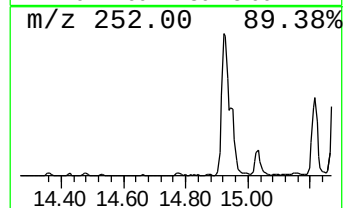
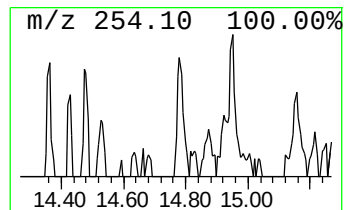
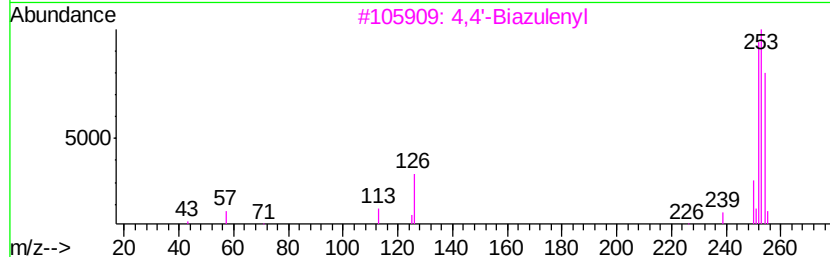
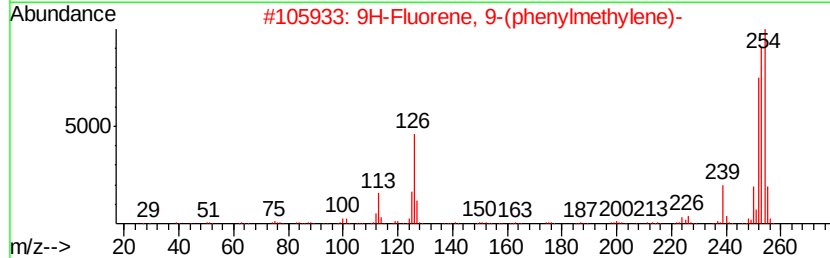
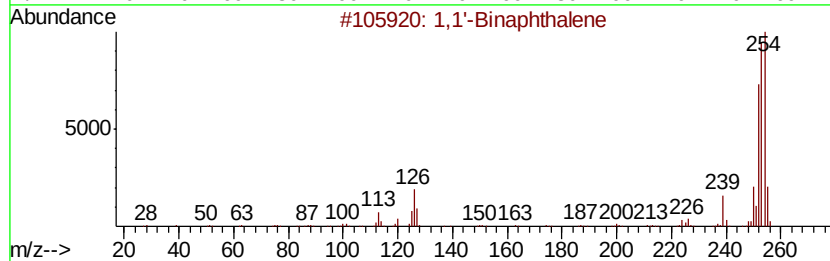
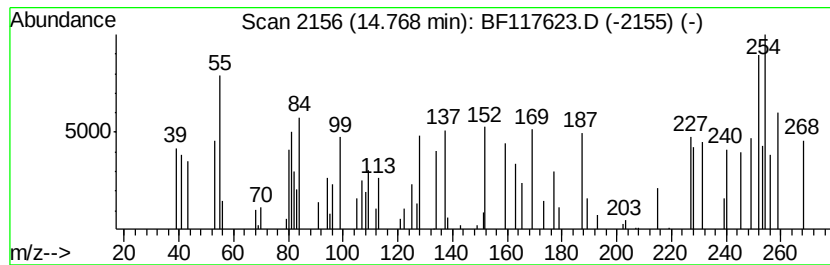
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1,1'-Binaphthalene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.77	2.01 ng	15940	Perylene-d12	15.33	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Binaphthalene	254	C20H14	000604-53-5	53
2	9H-Fluorene, 9-(phenylmethylene)-	254	C20H14	001836-87-9	52
3	4,4'-Biazulenvl	254	C20H14	094053-54-0	52
4	4,6'-Biazulenvl	254	C20H14	094154-49-1	52
5	1,2'-Binaphthalene	254	C20H14	004325-74-0	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103019\
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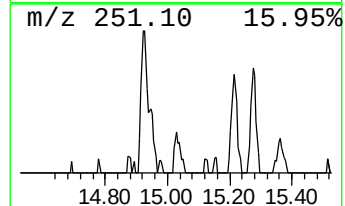
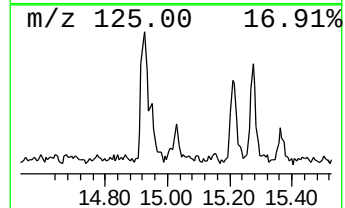
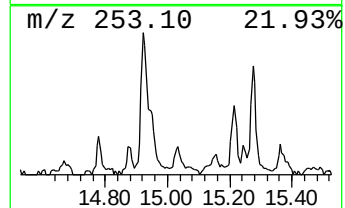
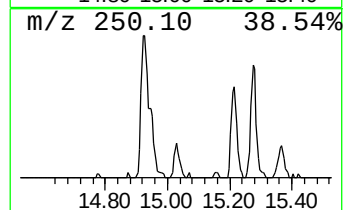
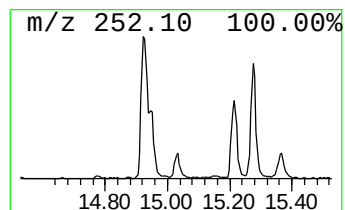
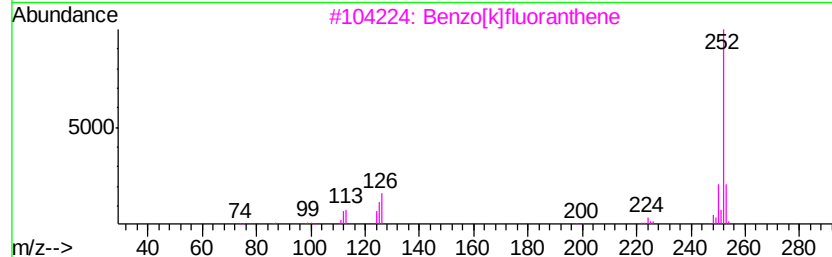
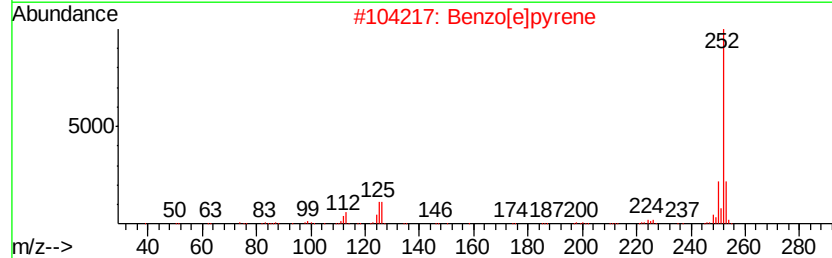
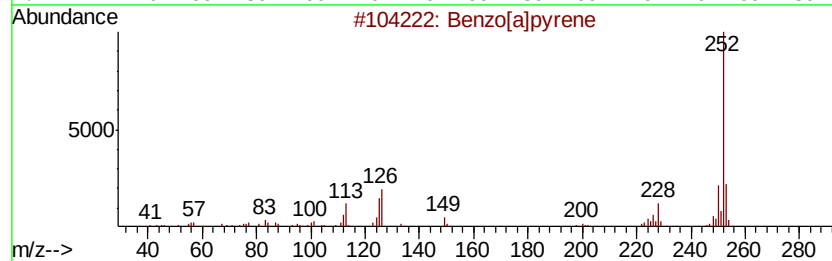
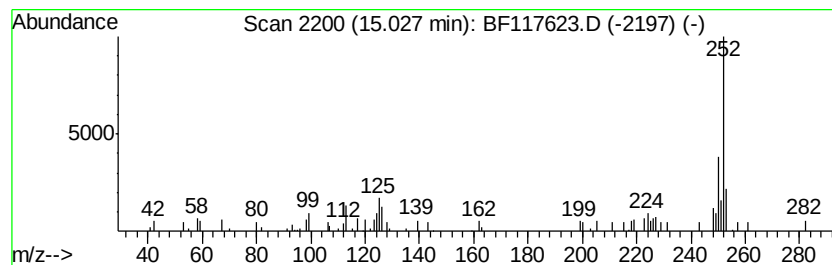
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzo[e]pyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.03	7.03 ng	55777	Perylene-d12	15.33	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[a]pyrene	252	C20H12	000050-32-8	64
2	Benzo[e]pyrene	252	C20H12	000192-97-2	52
3	Benzo[k]fluoranthene	252	C20H12	000207-08-9	52
4	Perylene	252	C20H12	000198-55-0	50
5	12H-[1,3]Benzothiazolo[2,3-b]qui...	252	C14H8N2OS	1000338-32-6	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103019\
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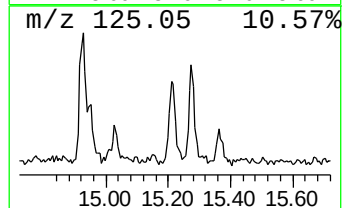
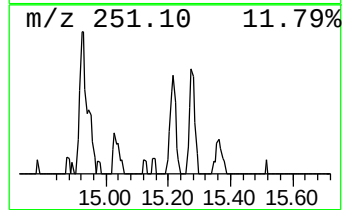
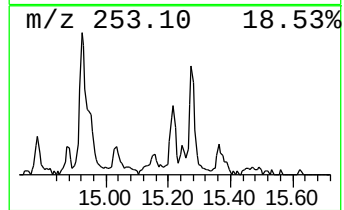
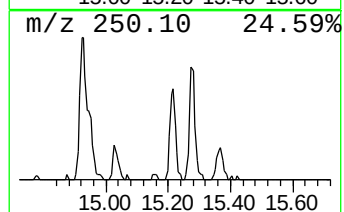
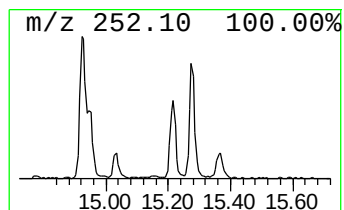
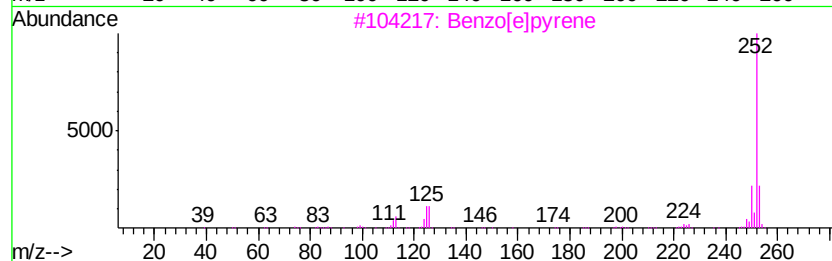
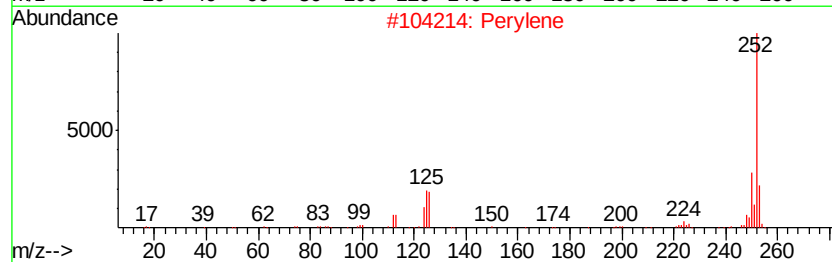
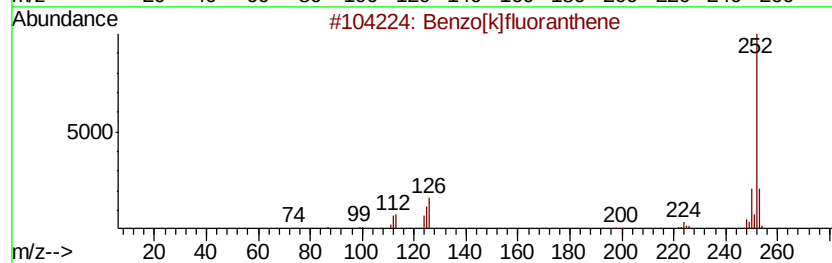
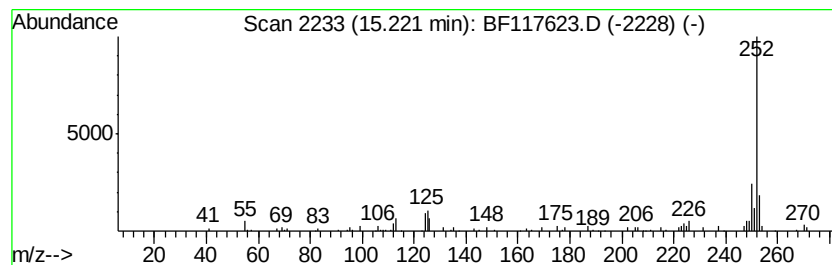
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Perylene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.22	9.70 ng	76883	Perylene-d12	15.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[k]fluoranthene	252	C20H12	000207-08-9	96
2	Perylene	252	C20H12	000198-55-0	95
3	Benzo[e]pyrene	252	C20H12	000192-97-2	95
4	Benzo[a]pyrene	252	C20H12	000050-32-8	93
5	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103019\
Data File : BF117623.D
Acq On : 30 Oct 2019 14:39
Operator : JU
Sample : K5542-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

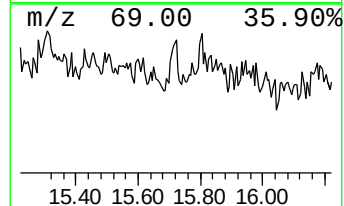
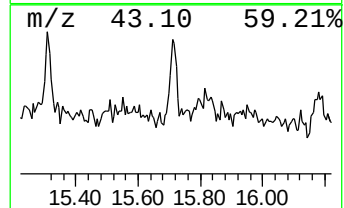
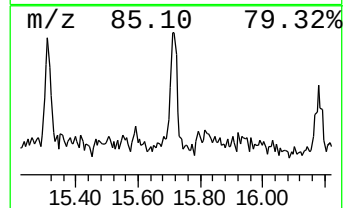
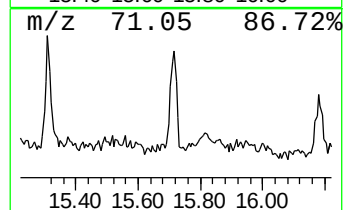
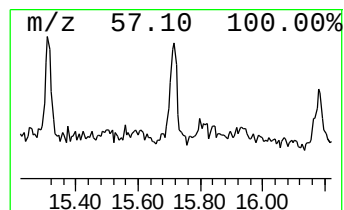
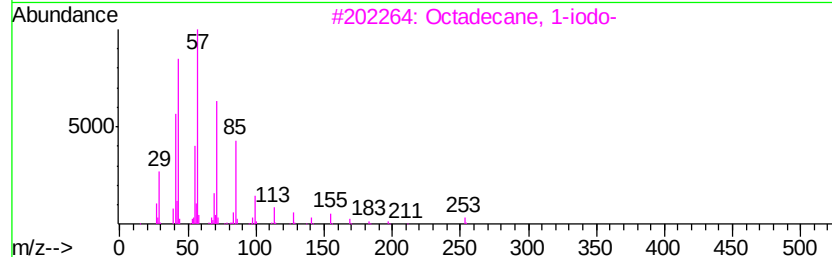
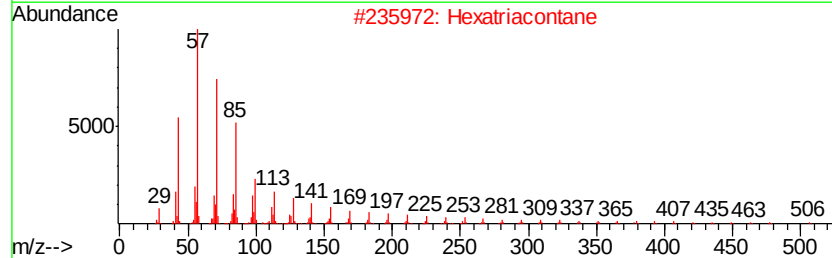
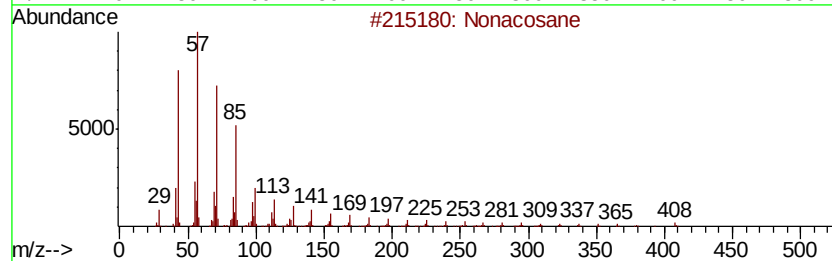
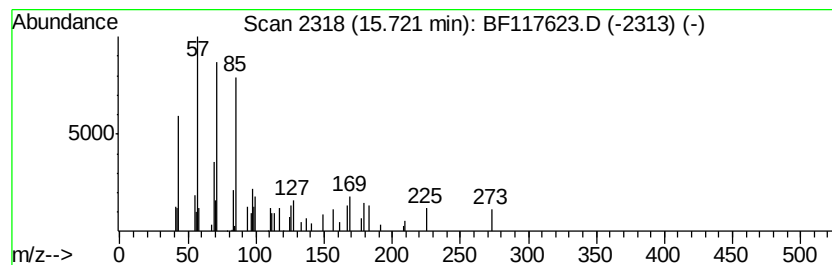
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Nonacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.72	4.03 ng	31924	Perylene-d12	15.33
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Nonacosane	408 C29H60	000630-03-5	53
2	Hexatriacontane	507 C36H74	000630-06-8	53
3	Octadecane, 1-iodo-	380 C18H37I	000629-93-6	50
4	Nonane, 1-iodo-	254 C9H19I	004282-42-2	47
5	Nonane, 5-methyl-5-propyl-	184 C13H28	017312-75-3	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103019\
Data File : BF117623.D
Acq On : 30 Oct 2019 14:39
Operator : JU
Sample : K5542-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
198-90-(0-2)

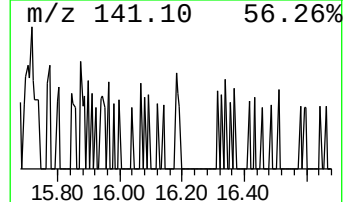
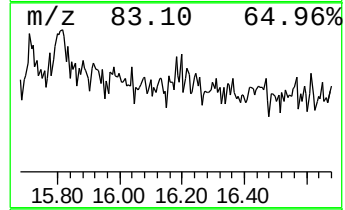
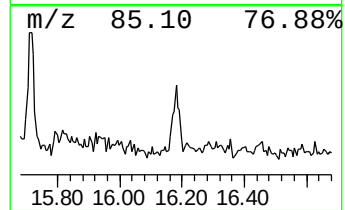
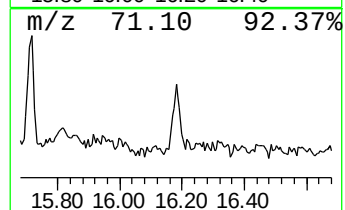
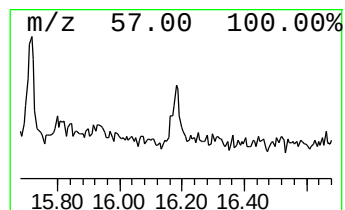
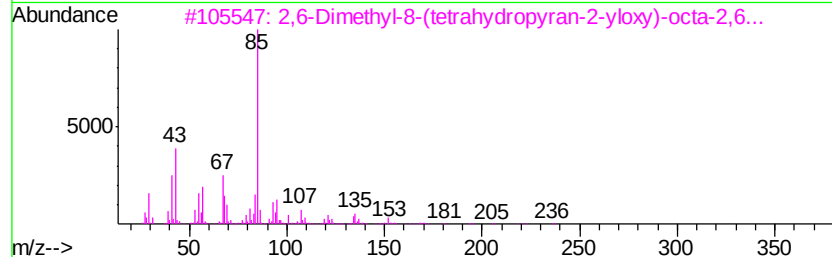
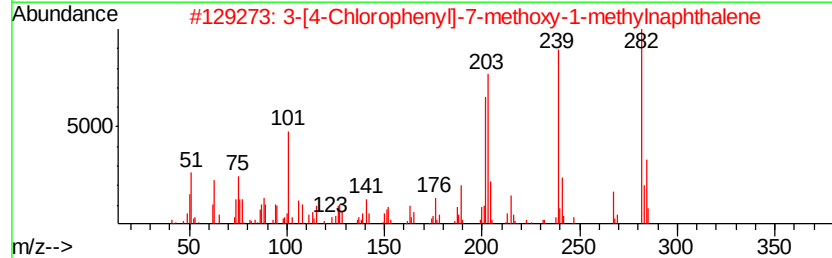
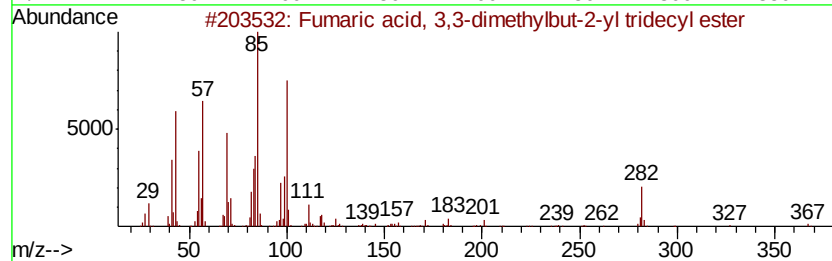
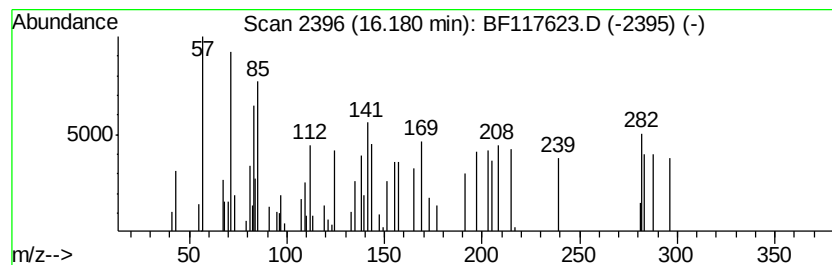
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown16.18 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.18	2.19 ng	17385	Perylene-d12	15.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Fumaric acid, 3,3-dimethylbut-2-...	382	C23H42O4	1000348-71-2	14
2		3-[4-Chlorophenyl]-7-methoxy-1-m...	282	C18H15ClO	055614-38-5	10
3		2,6-Dimethyl-8-(tetrahydropyran-...	254	C15H26O3	038290-53-8	10
4		10-Methylnonadecane	282	C20H42	056862-62-5	10
5		Cholest-22-ene-21-ol, 3,5-dehydr...	498	C33H54O3	1000124-60-2	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF103019\
 Data File : BF117623.D
 Acq On : 30 Oct 2019 14:39
 Operator : JU
 Sample : K5542-07
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 198-90-(0-2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101819.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
2-Pentanone, 4-hy...	4.92	11.4	ng	122432	1	6.73	215640	20.0
unknown6.50	6.50	72.3	ng	778984	1	6.73	215640	20.0
Anthracene, 2-met...	11.78	6.9	ng	81793	4	11.25	238446	20.0
6H-Cyclobuta[jk]p...	11.86	3.9	ng	46094	4	11.25	238446	20.0
Behenic alcohol	13.73	2.9	ng	60877	5	13.89	427632	20.0
unknown14.27	14.27	3.2	ng	69214	5	13.89	427632	20.0
Pentacosane	14.64	2.6	ng	20962	6	15.33	158585	20.0
Supraene	14.71	5.3	ng	42401	6	15.33	158585	20.0
1,1'-Binaphthalene	14.77	2.0	ng	15940	6	15.33	158585	20.0
Benzo[e]pyrene	15.03	7.0	ng	55777	6	15.33	158585	20.0
Perylene	15.22	9.7	ng	76883	6	15.33	158585	20.0
Nonacosane	15.72	4.0	ng	31924	6	15.33	158585	20.0
unknown16.18	16.18	2.2	ng	17385	6	15.33	158585	20.0