

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
 Data File : BF120252.D
 Acq On : 30 Apr 2020 21:07
 Operator : MA/SJ
 Sample : L2452-05
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 227

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-BF040820.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	3.493	233	239	247	rVB	92685	157403	3.58%	0.404%
2	5.198	526	529	531	rBV	93018	91268	2.08%	0.234%
3	5.263	537	540	562	rBV	3079617	3554408	80.85%	9.130%
4	6.316	714	719	735	rBV	3225847	3524235	80.16%	9.053%
5	6.651	772	776	779	rBV	844181	676367	15.39%	1.737%
6	6.810	798	803	806	rBV	3169649	2865551	65.18%	7.361%
7	7.222	868	873	876	rBV	2668084	2304684	52.42%	5.920%
8	7.933	991	994	997	rBV	1153786	869513	19.78%	2.233%
9	9.022	1174	1179	1182	rBV	4657163	4153779	94.48%	10.670%
10	9.169	1201	1204	1216	rVB	106894	116120	2.64%	0.298%
11	9.416	1242	1246	1249	rBV	522076	391010	8.89%	1.004%
12	9.692	1289	1293	1296	rBV	1234536	971852	22.11%	2.496%
13	9.722	1296	1298	1302	rVB	55356	47519	1.08%	0.122%
14	10.239	1383	1386	1391	rVB2	77999	64474	1.47%	0.166%
15	10.486	1423	1428	1431	rBV	3077255	2633936	59.91%	6.766%
16	11.174	1542	1545	1548	rBV	1139841	1042775	23.72%	2.679%
17	11.198	1548	1549	1553	rVV	260288	184857	4.20%	0.475%
18	11.251	1555	1558	1561	rVB	122098	96813	2.20%	0.249%
19	11.680	1628	1631	1633	rBV	72983	67863	1.54%	0.174%
20	11.721	1633	1638	1642	rVV2	795107	852533	19.39%	2.190%
21	11.751	1642	1643	1647	rVV2	97620	93227	2.12%	0.239%
22	11.792	1647	1650	1652	rVV	199458	202177	4.60%	0.519%
23	11.810	1652	1653	1657	rVB	59261	46755	1.06%	0.120%
24	11.974	1679	1681	1684	rBV	53845	47327	1.08%	0.122%
25	12.227	1721	1724	1728	rBV	73221	80004	1.82%	0.206%
26	12.268	1728	1731	1736	rVB4	50021	74523	1.70%	0.191%
27	12.316	1736	1739	1743	rBV3	72966	88166	2.01%	0.226%
28	12.392	1748	1752	1755	rBV	779608	761757	17.33%	1.957%
29	12.439	1755	1760	1766	rVB6	37085	71627	1.63%	0.184%
30	12.510	1769	1772	1775	rBV	403908	423453	9.63%	1.088%
31	12.616	1785	1790	1793	rBV	807991	833521	18.96%	2.141%
32	12.668	1797	1799	1804	rVB2	46211	63907	1.45%	0.164%
33	12.739	1807	1811	1812	rBV2	62252	54700	1.24%	0.141%
34	12.774	1812	1817	1820	rVV	4598416	4396235	100.00%	11.292%

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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

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

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35	12.839	1823	1828	1832	rBV3	87734	124047	2.82%	0.319%
36	12.921	1840	1842	1844	rBV2	76957	89266	2.03%	0.229%
37	12.951	1844	1847	1850	rVB	164766	167107	3.80%	0.429%
38	13.015	1855	1858	1860	rVB	141825	124264	2.83%	0.319%
39	13.051	1860	1864	1867	rBV2	121290	129522	2.95%	0.333%
40	13.145	1877	1880	1882	rBV	75191	68196	1.55%	0.175%
41	13.174	1882	1885	1888	rBV	82308	74452	1.69%	0.191%
42	13.257	1895	1899	1903	rVB7	56358	76715	1.75%	0.197%
43	13.457	1931	1933	1938	rVB	55814	58264	1.33%	0.150%
44	13.563	1948	1951	1954	rBV2	72764	91912	2.09%	0.236%
45	13.674	1966	1970	1977	rBV2	377076	490528	11.16%	1.260%
46	13.815	1987	1994	1997	rBV2	1202531	1577417	35.88%	4.052%
47	13.845	1997	1999	2001	rVB	280014	215936	4.91%	0.555%
48	14.204	2056	2060	2063	rBV	108421	134276	3.05%	0.345%
49	14.239	2063	2066	2070	rVB2	66404	69752	1.59%	0.179%
50	14.304	2073	2077	2079	rBV2	116111	146006	3.32%	0.375%
51	14.598	2124	2127	2132	rVB2	117045	125263	2.85%	0.322%
52	14.827	2162	2166	2176	rBV	427213	726413	16.52%	1.866%
53	14.909	2176	2180	2187	rVB3	195512	279883	6.37%	0.719%
54	15.109	2210	2214	2220	rBV	236183	284402	6.47%	0.731%
55	15.162	2220	2223	2228	rVB	278321	313707	7.14%	0.806%
56	15.227	2229	2234	2237	rVV	726010	896363	20.39%	2.302%
57	15.257	2237	2239	2245	rVB4	172090	222696	5.07%	0.572%
58	15.656	2304	2307	2313	rVB	101368	139160	3.17%	0.357%
59	16.568	2458	2462	2471	rVB2	125368	241111	5.48%	0.619%
60	16.974	2527	2531	2538	rVB	86071	159691	3.63%	0.410%

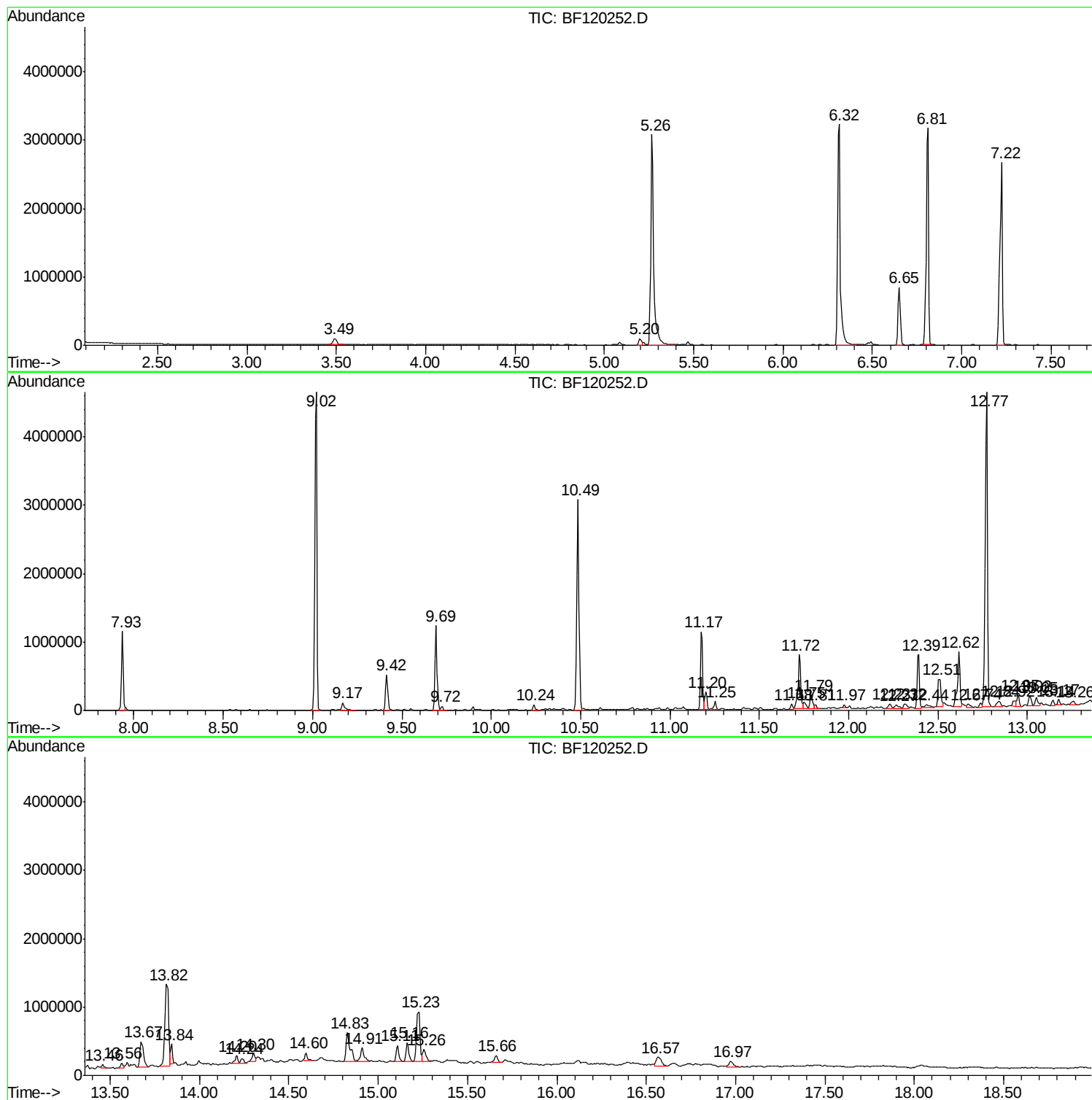
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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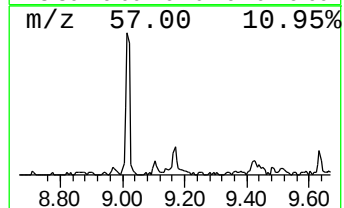
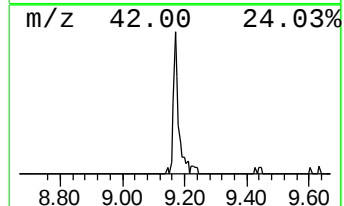
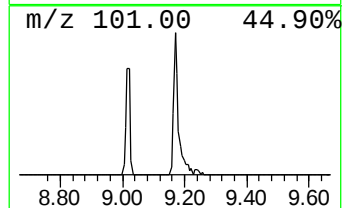
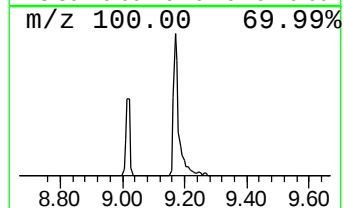
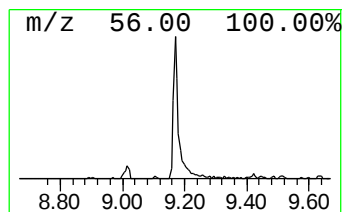
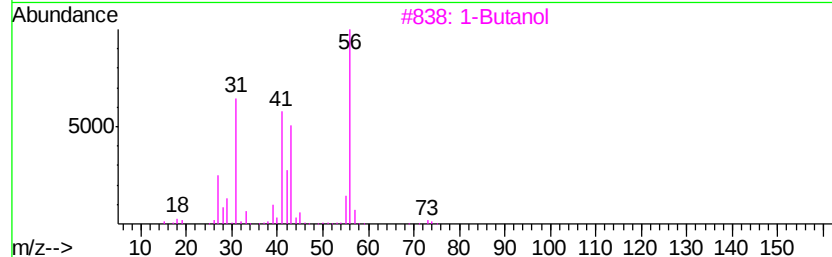
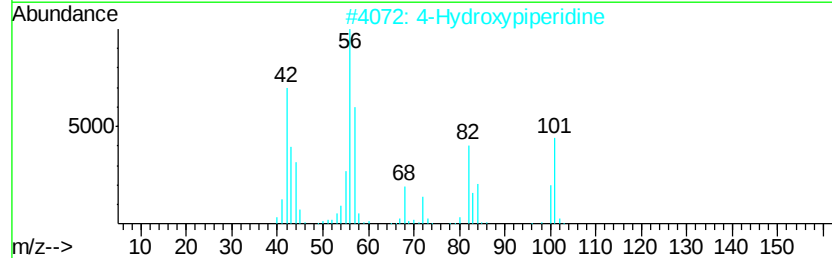
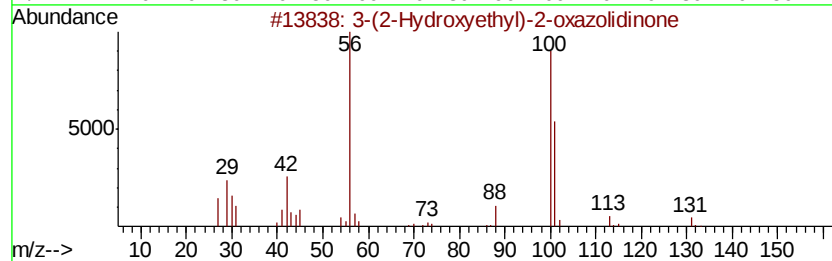
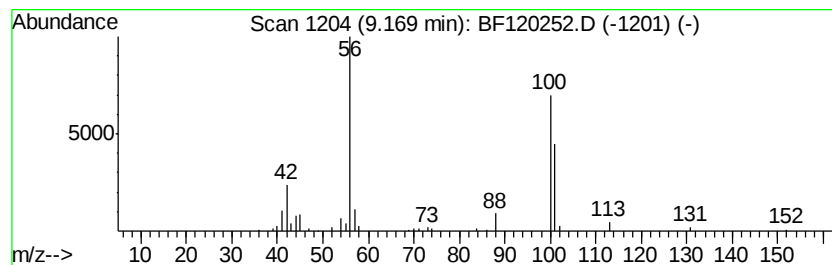
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 3-(2-Hydroxyethyl)-2-oxazol... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.17	2.39 ng	116120	Acenaphthene-d10	9.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-(2-Hydroxyethyl)-2-oxazolidinone	131	C5H9NO3	003356-88-5	83
2		4-Hydroxypiperidine	101	C5H11NO	005382-16-1	47
3		1-Butanol	74	C4H10O	000071-36-3	38
4		Formic acid, hexyl ester	130	C7H14O2	000629-33-4	23
5		Imidazole, 5-fluoro-4-methyl-	100	C4H5FN2	041367-01-5	14



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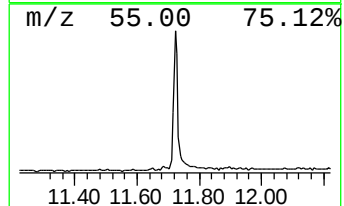
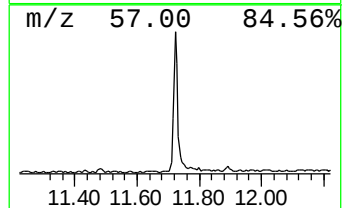
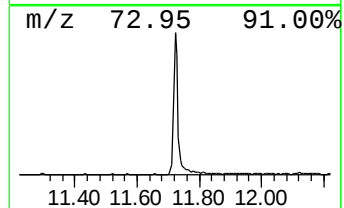
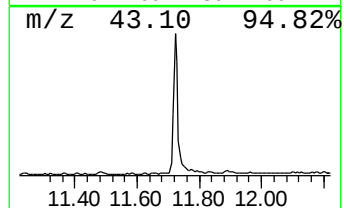
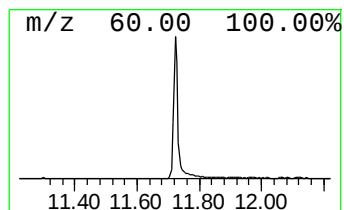
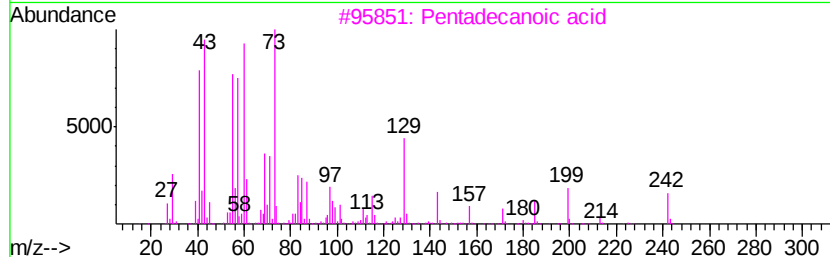
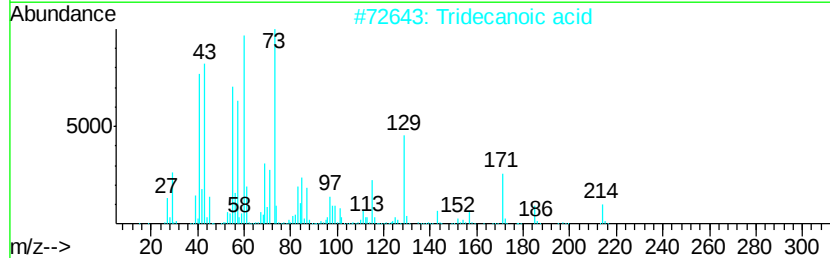
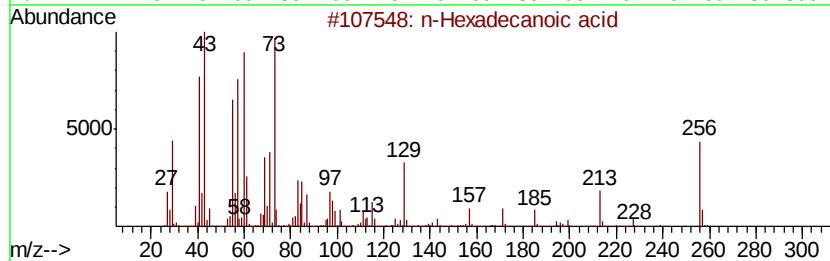
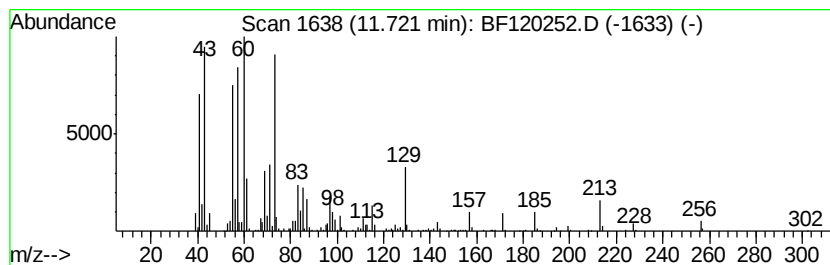
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.72	16.35 ng	852533	Phenanthrene-d10	11.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tridecanoic acid	214	C13H26O2	000638-53-9	93
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	87
5		Undecanoic acid	186	C11H22O2	000112-37-8	80



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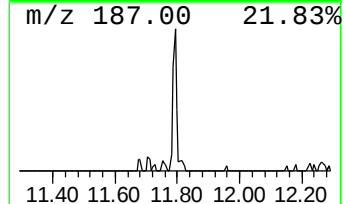
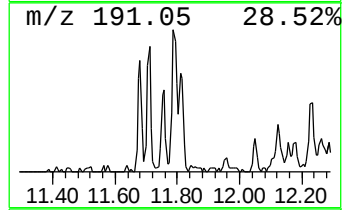
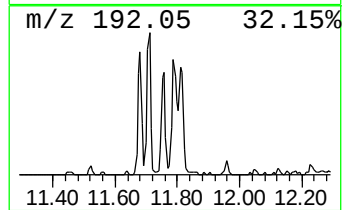
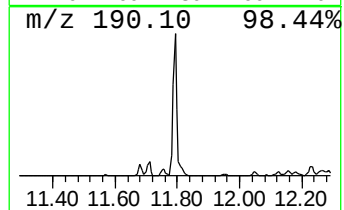
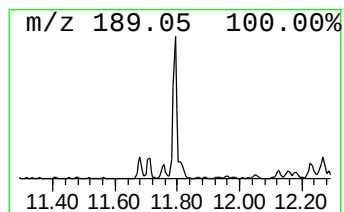
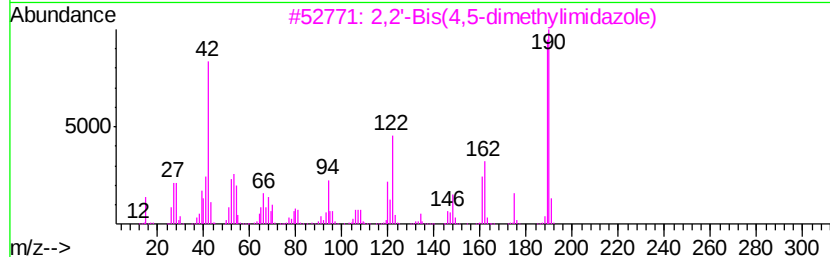
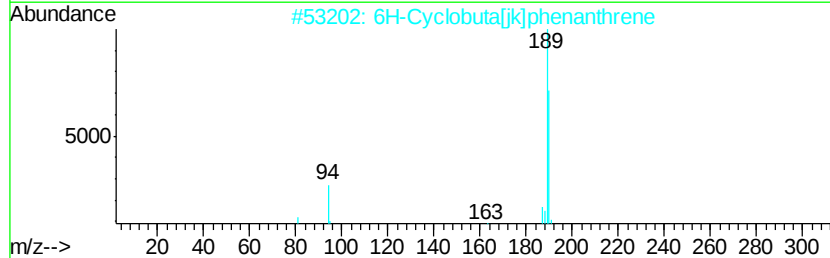
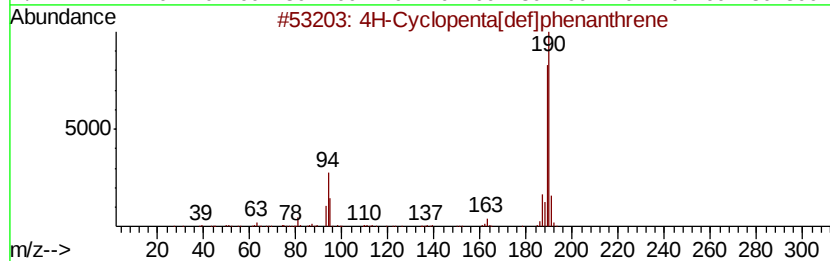
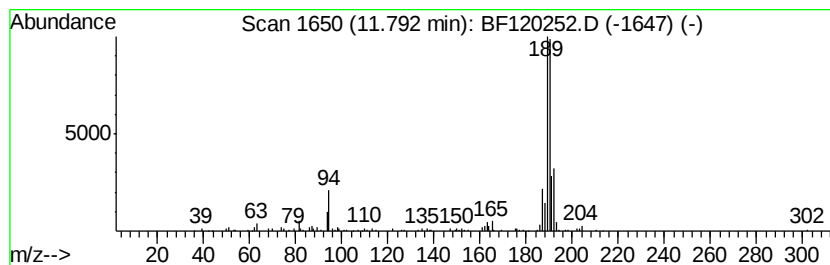
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.79	3.88 ng	202177	Phenanthrene-d10	11.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	89
2		6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	59
3		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	42
4		1-(Phenylethynyl)-1H-1,2,3-benzo...	219	C14H9N3	209912-23-2	27
5		Pyrimidine, 2-chloro-5-phenyl-	190	C10H7ClN2	022536-62-5	25



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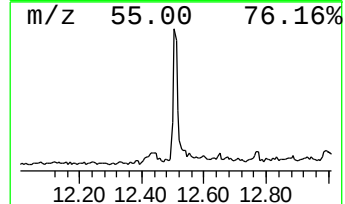
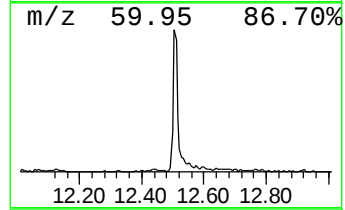
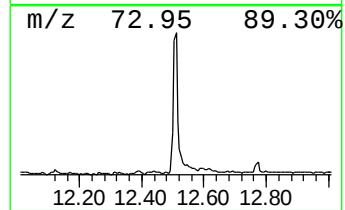
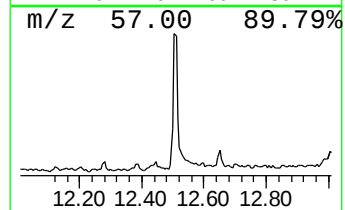
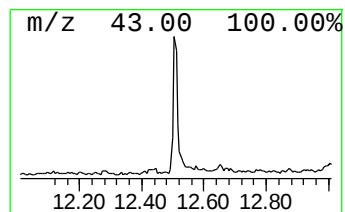
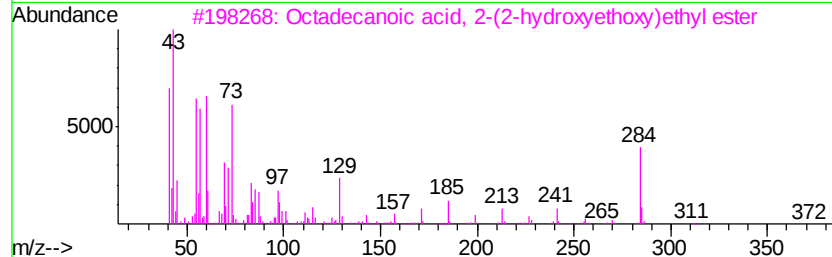
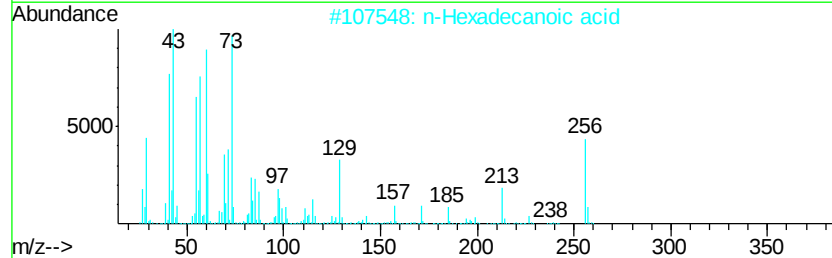
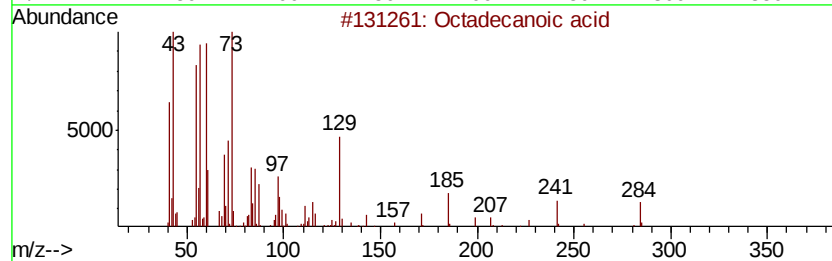
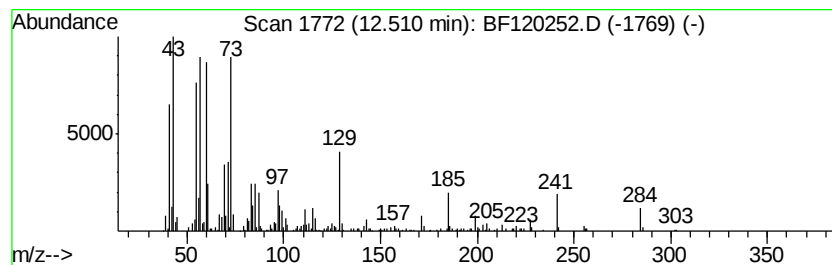
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.51	5.37 ng	423453	Chrysene-d12	13.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	89
3		Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	87
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
5		Tridecanoic acid	214	C13H26O2	000638-53-9	74



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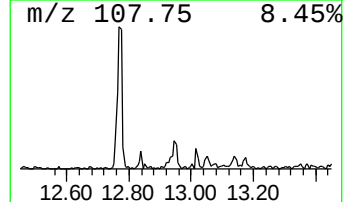
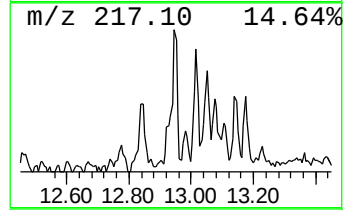
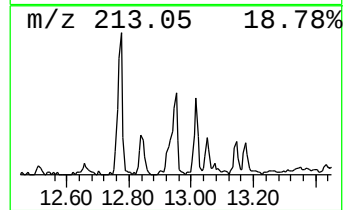
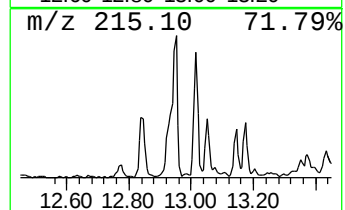
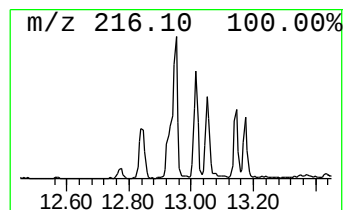
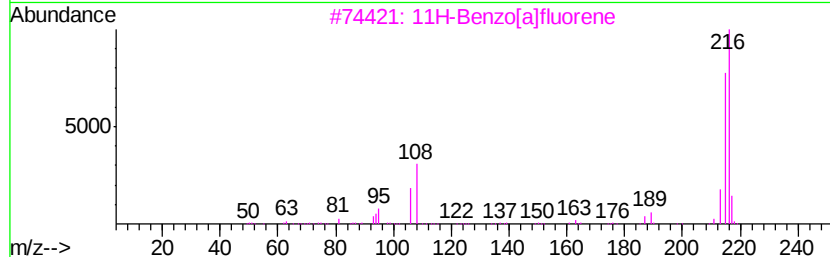
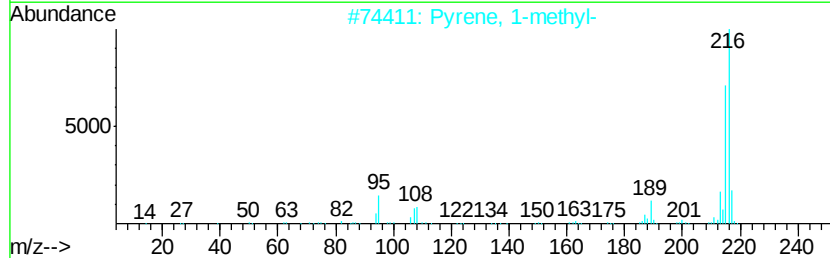
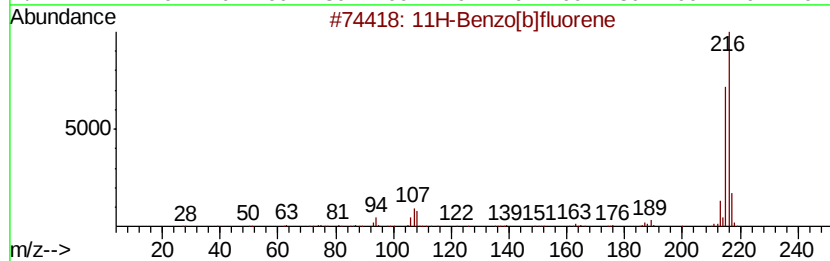
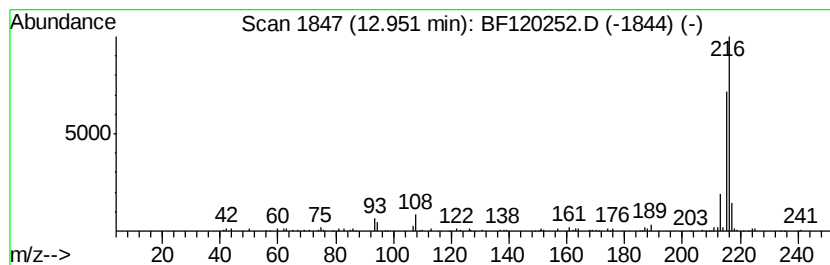
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 11H-Benzo[b]fluorene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.95	2.12 ng	167107	Chrysene-d12	13.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
3		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
4		.beta.-Carboline, 1,2,3,4-tetra...	260	C14H16N2O3	029573-11-3	64
5		Pyrene, 4-methyl-	216	C17H12	003353-12-6	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
Data File : BF120252.D
Acq On : 30 Apr 2020 21:07
Operator : MA/SJ
Sample : L2452-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
227

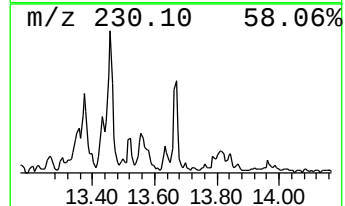
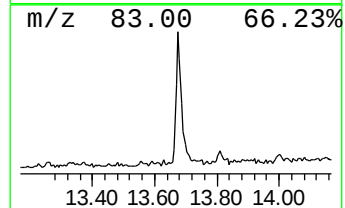
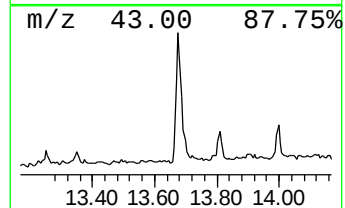
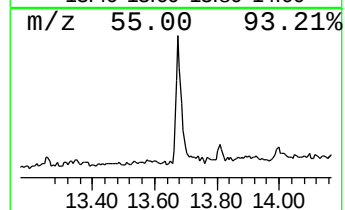
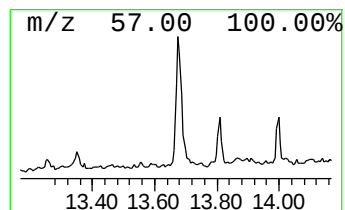
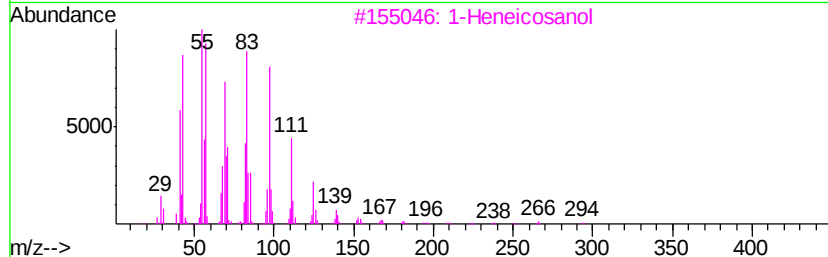
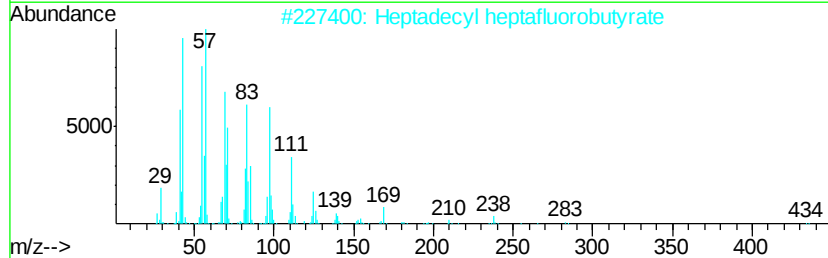
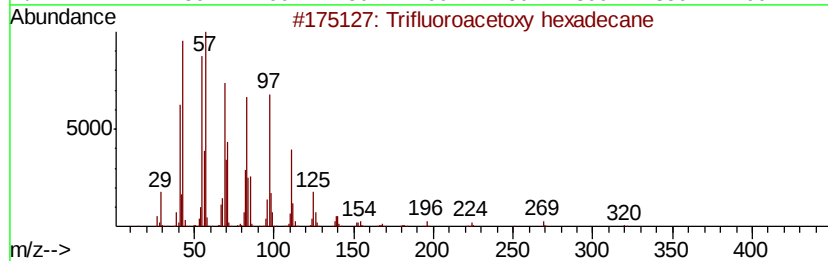
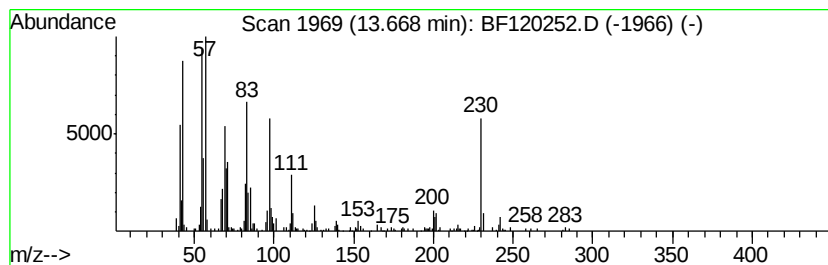
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Trifluoroacetoxy hexadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.67	6.22 ng	490528	Chrysene-d12	13.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
2		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
3		1-Heneicosanol	312	C21H44O	015594-90-8	93
4		Behenic alcohol	326	C22H46O	000661-19-8	93
5		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
Data File : BF120252.D
Acq On : 30 Apr 2020 21:07
Operator : MA/SJ
Sample : L2452-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

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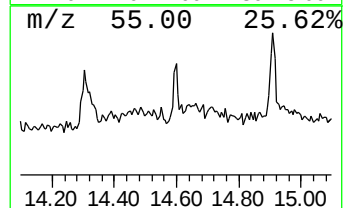
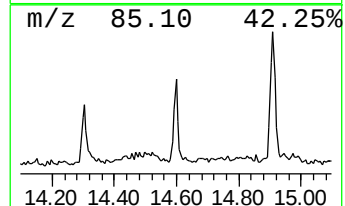
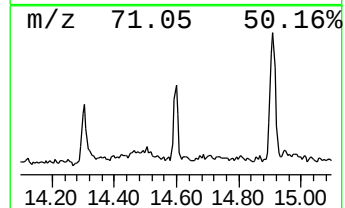
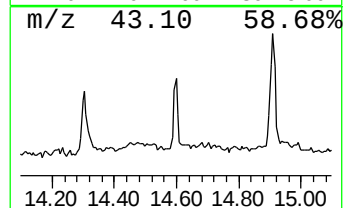
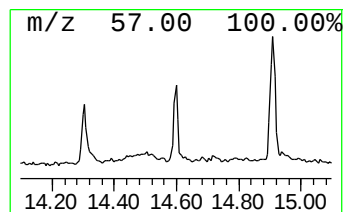
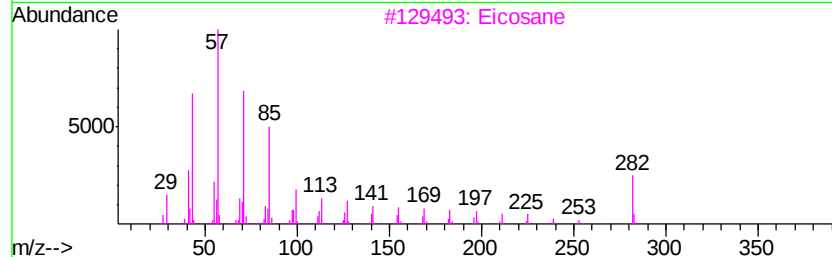
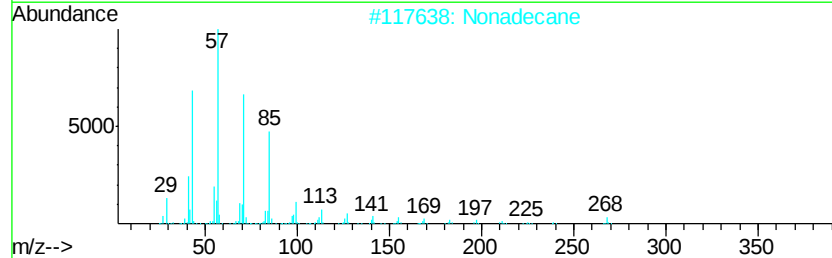
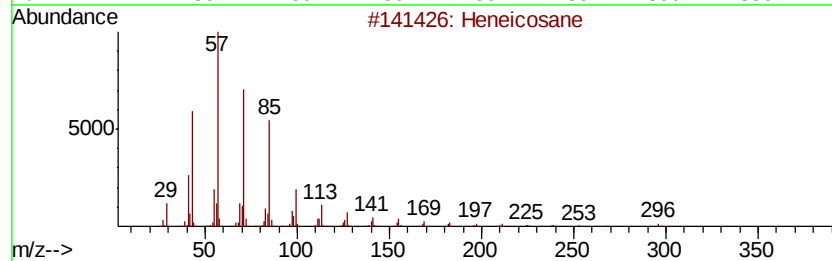
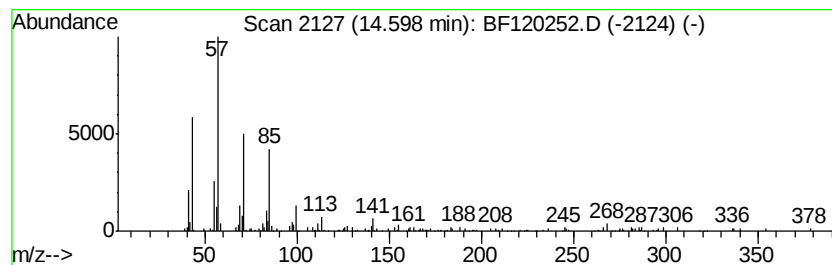
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heneicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.60	2.79 ng	125263	Perylene-d12	15.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	93
2			Nonadecane	268	C19H40	000629-92-5	90
3			Eicosane	282	C20H42	000112-95-8	87
4			Hexacosane	366	C26H54	000630-01-3	80
5			Eicosane, 10-methyl-	296	C21H44	054833-23-7	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
Data File : BF120252.D
Acq On : 30 Apr 2020 21:07
Operator : MA/SJ
Sample : L2452-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

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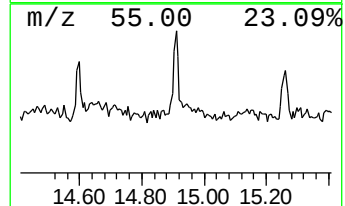
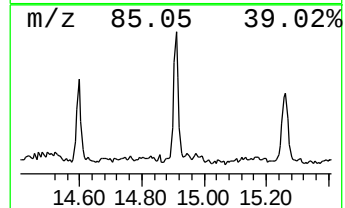
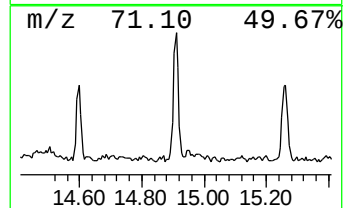
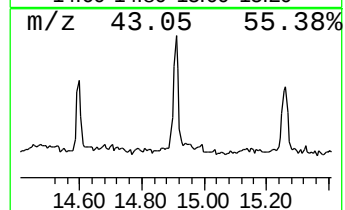
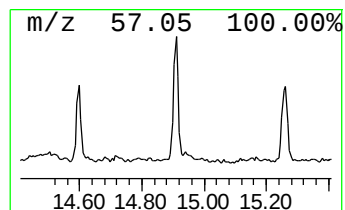
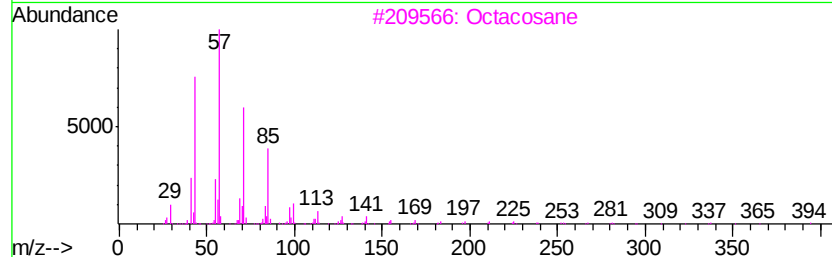
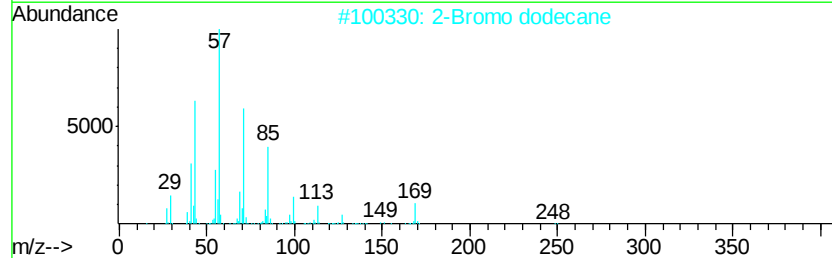
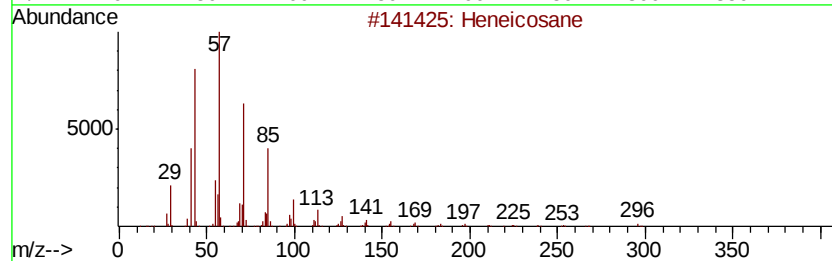
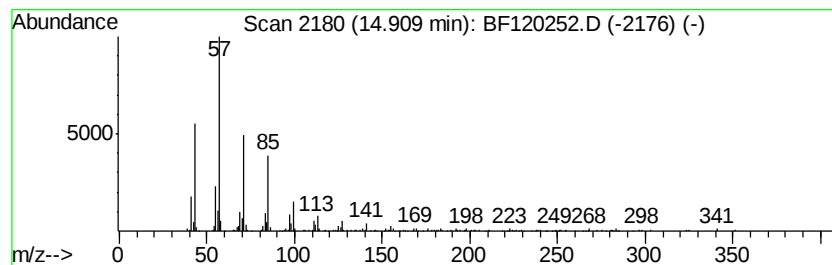
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 2-Bromo dodecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.91	6.24 ng	279883	Perylene-d12	15.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	96
2		2-Bromo dodecane	248	C12H25Br	013187-99-0	91
3		Octacosane	394	C28H58	000630-02-4	90
4		Tridecane, 6-propyl-	226	C16H34	055045-10-8	87
5		Nonadecane	268	C19H40	000629-92-5	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
Data File : BF120252.D
Acq On : 30 Apr 2020 21:07
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Sample : L2452-05
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ALS Vial : 21 Sample Multiplier: 1

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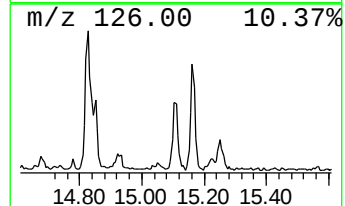
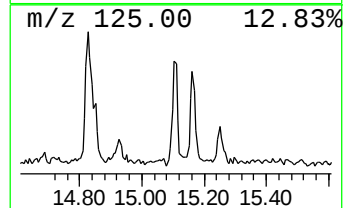
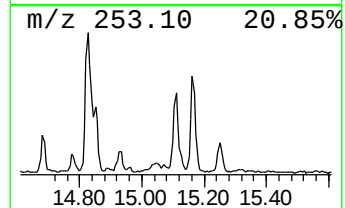
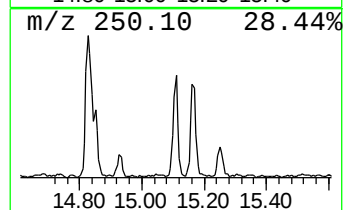
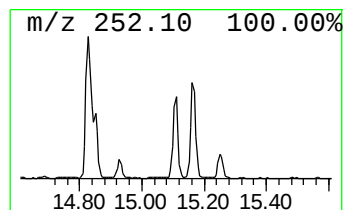
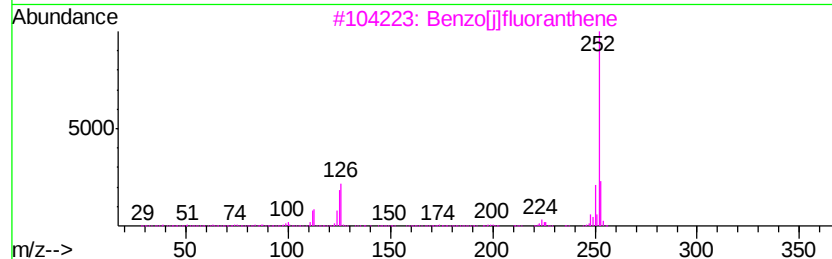
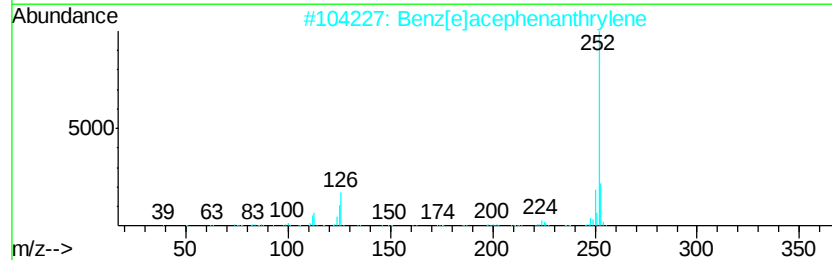
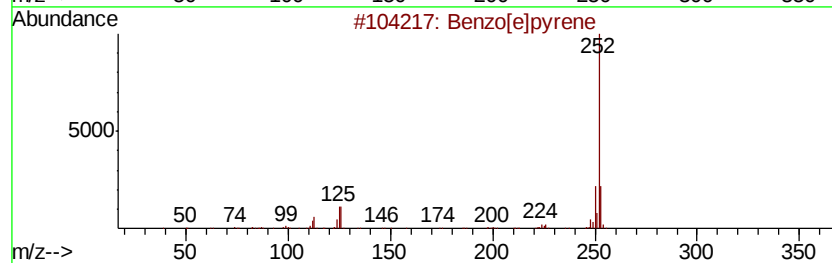
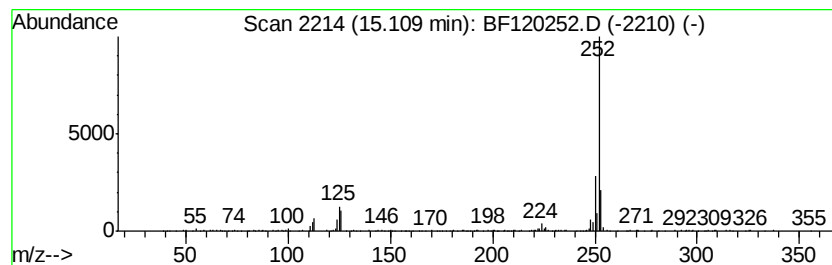
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.11	6.35 ng	284402	Perylene-d12	15.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]bpvrene	252	C20H12	000192-97-2	96
2		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	95
3		Benzo[i]fluoranthene	252	C20H12	000205-82-3	95
4		Benzo[a]bpvrene	252	C20H12	000050-32-8	93
5		Perylene	252	C20H12	000198-55-0	91



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Data File : BF120252.D
Acq On : 30 Apr 2020 21:07
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Sample : L2452-05
Misc :
ALS Vial : 21 Sample Multiplier: 1

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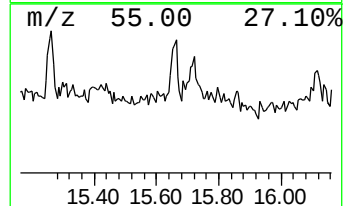
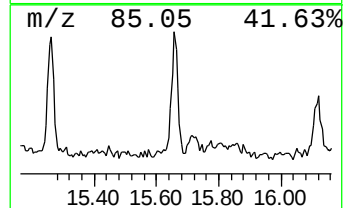
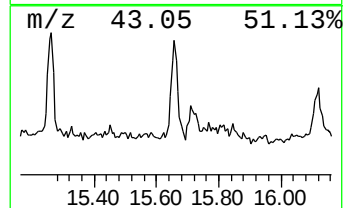
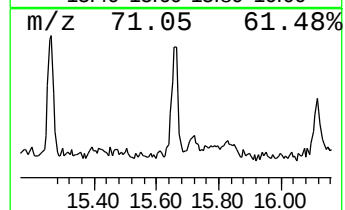
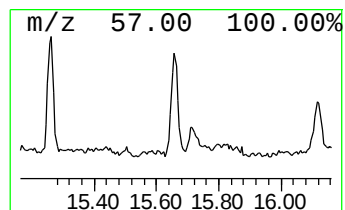
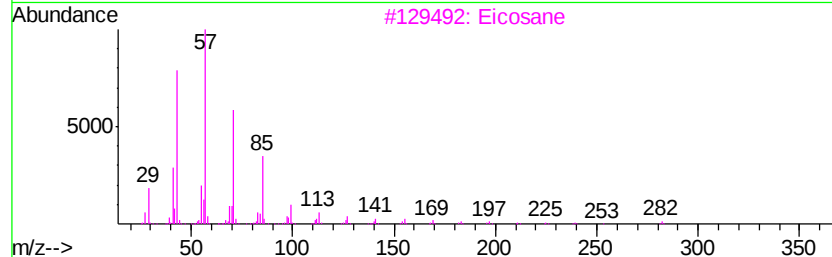
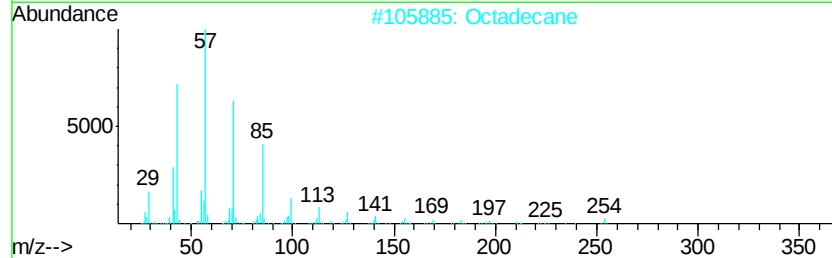
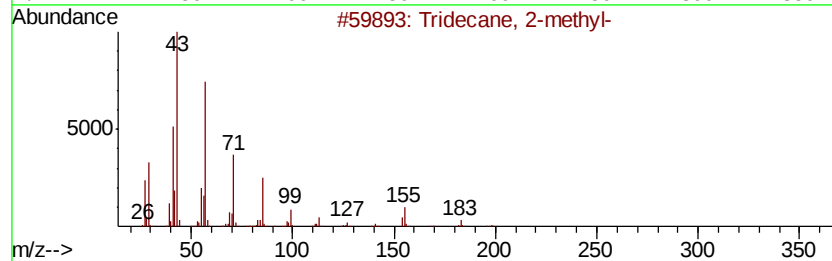
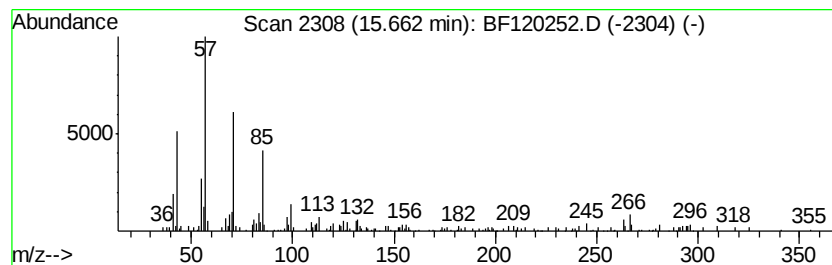
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tridecane, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.66	3.10 ng	139160	Perylene-d12	15.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 2-methyl-	198	C14H30	001560-96-9	86
2		Octadecane	254	C18H38	000593-45-3	86
3		Eicosane	282	C20H42	000112-95-8	86
4		Docosane	310	C22H46	000629-97-0	86
5		Dodecane, 4,9-dipropyl-	254	C18H38	003054-63-5	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF043020\
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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
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n-Hexadecanoic acid	11.72	16.4	ng	852533	4	11.17	1042780	20.0
4H-Cyclopenta[def...	11.79	3.9	ng	202177	4	11.17	1042780	20.0
Octadecanoic acid	12.51	5.4	ng	423453	5	13.82	1577420	20.0
11H-Benzo[b]fluorene	12.95	2.1	ng	167107	5	13.82	1577420	20.0
Trifluoroacetoxy ...	13.67	6.2	ng	490528	5	13.82	1577420	20.0
Heneicosane	14.60	2.8	ng	125263	6	15.23	896363	20.0
2-Bromo dodecane	14.91	6.2	ng	279883	6	15.23	896363	20.0
Benzo[e]pyrene	15.11	6.3	ng	284402	6	15.23	896363	20.0
Tridecane, 2-methyl-	15.66	3.1	ng	139160	6	15.23	896363	20.0