

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF060220\
 Data File : BF120486.D
 Acq On : 2 Jun 2020 11:18
 Operator : JU/CG
 Sample : L2792-07
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM107-COMP-2020-0-6

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-BF052820.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	1.916	3	5	11	rVB2	458402	552420	15.58%	1.265%
2	2.016	20	22	30	rVB3	25712	43035	1.21%	0.099%
3	2.099	30	36	44	rVB	1008286	1272040	35.87%	2.912%
4	5.122	541	550	558	rBV2	34766	96226	2.71%	0.220%
5	5.463	604	608	611	rBV	3053264	2827561	79.73%	6.473%
6	6.122	717	720	724	rVB	97256	81054	2.29%	0.186%
7	6.475	775	780	783	rBV	2888512	2917485	82.26%	6.679%
8	6.845	839	843	846	rBV	1278545	1101877	31.07%	2.523%
9	6.998	865	869	872	rBV	3068121	2616722	73.78%	5.991%
10	7.404	933	938	941	rBV	2256793	1971610	55.59%	4.514%
11	8.122	1056	1060	1064	rBV	1505815	1387380	39.12%	3.176%
12	9.198	1239	1243	1246	rBV	4063439	3546537	100.00%	8.120%
13	9.316	1258	1263	1265	rBV2	82018	75984	2.14%	0.174%
14	9.539	1298	1301	1304	rBV2	103569	87208	2.46%	0.200%
15	9.592	1304	1310	1314	rVV	653234	541456	15.27%	1.240%
16	9.628	1314	1316	1319	rVB2	63212	51466	1.45%	0.118%
17	9.739	1330	1335	1339	rVB2	62211	70121	1.98%	0.161%
18	9.881	1352	1359	1362	rBV	1847515	1821533	51.36%	4.170%
19	9.910	1362	1364	1365	rVV	68055	54532	1.54%	0.125%
20	9.928	1365	1367	1370	rVB	87092	65676	1.85%	0.150%
21	10.081	1389	1393	1397	rBV	76952	79170	2.23%	0.181%
22	10.339	1433	1437	1441	rVB	83212	81854	2.31%	0.187%
23	10.428	1449	1452	1456	rVB	86790	74303	2.10%	0.170%
24	10.669	1488	1493	1496	rBV	2132460	1870972	52.75%	4.283%
25	10.745	1503	1506	1511	rBV3	45147	60462	1.70%	0.138%
26	11.028	1549	1554	1558	rVB3	112891	137666	3.88%	0.315%
27	11.169	1575	1578	1582	rVB	48712	43598	1.23%	0.100%
28	11.263	1591	1594	1599	rVB	52406	53252	1.50%	0.122%
29	11.310	1599	1602	1606	rBV	66258	71955	2.03%	0.165%
30	11.369	1606	1612	1614	rVV	1455338	1432532	40.39%	3.280%
31	11.392	1614	1616	1619	rVV	826981	660610	18.63%	1.512%
32	11.439	1622	1624	1627	rVB	240809	177841	5.01%	0.407%
33	11.592	1645	1650	1653	rBV2	66243	100254	2.83%	0.230%
34	11.663	1659	1662	1666	rBV2	32696	47925	1.35%	0.110%

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 Integrator: RTE
 Smoothing : ON
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If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.827	1685	1690	1693	rBV	226527	305696	8.62%	0.700%
36	11.863	1693	1696	1698	rVV3	255648	326631	9.21%	0.748%
37	11.892	1698	1701	1706	rVV2	528545	632210	17.83%	1.447%
38	11.939	1707	1709	1712	rVV	83921	102859	2.90%	0.235%
39	11.980	1712	1716	1718	rVV2	207415	238704	6.73%	0.546%
40	11.998	1718	1719	1722	rVV	91687	85863	2.42%	0.197%
41	12.039	1722	1726	1729	rVV2	61070	120350	3.39%	0.276%
42	12.080	1729	1733	1739	rVV6	70034	141432	3.99%	0.324%
43	12.145	1741	1744	1745	rVV	81202	71054	2.00%	0.163%
44	12.163	1745	1747	1749	rVV	95837	82709	2.33%	0.189%
45	12.186	1749	1751	1754	rVB	67457	53776	1.52%	0.123%
46	12.416	1787	1790	1794	rBV	72709	91166	2.57%	0.209%
47	12.457	1794	1797	1799	rVV2	59352	67572	1.91%	0.155%
48	12.498	1801	1804	1806	rVV2	119119	134295	3.79%	0.307%
49	12.557	1810	1814	1815	rVV2	280672	240715	6.79%	0.551%
50	12.580	1815	1818	1821	rVB	1279419	1053473	29.70%	2.412%
51	12.622	1822	1825	1827	rBV2	41741	42032	1.19%	0.096%
52	12.674	1831	1834	1837	rBV2	60641	58352	1.65%	0.134%
53	12.810	1851	1857	1860	rBV	1019004	1009037	28.45%	2.310%
54	12.857	1863	1865	1869	rVB2	46429	45768	1.29%	0.105%
55	12.951	1877	1881	1884	rBV	3195601	2785700	78.55%	6.378%
56	13.033	1889	1895	1897	rBV3	97998	186292	5.25%	0.427%
57	13.092	1901	1905	1906	rVV3	85355	117760	3.32%	0.270%
58	13.133	1910	1912	1916	rVB	190704	184112	5.19%	0.422%
59	13.204	1922	1924	1926	rVB	123177	94280	2.66%	0.216%
60	13.239	1927	1930	1933	rVB2	100579	89362	2.52%	0.205%
61	13.333	1943	1946	1948	rVB	52244	45340	1.28%	0.104%
62	13.363	1949	1951	1955	rBV	63591	58967	1.66%	0.135%
63	13.639	1996	1998	2003	rVB	78945	85076	2.40%	0.195%
64	13.757	2014	2018	2020	rBV2	88764	109621	3.09%	0.251%
65	13.786	2020	2023	2025	rVV	85653	91035	2.57%	0.208%
66	13.804	2025	2026	2028	rVV	71841	67084	1.89%	0.154%
67	13.857	2031	2035	2041	rVB2	380663	491849	13.87%	1.126%
68	14.004	2053	2060	2063	rBV2	1522958	2193322	61.84%	5.021%
69	14.033	2063	2065	2073	rVV	581903	611619	17.25%	1.400%
70	14.110	2076	2078	2081	rBV	68331	56081	1.58%	0.128%
71	14.392	2118	2126	2129	rBV2	140294	311282	8.78%	0.713%

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Sampling : 1 Min Area: 3 % of largest Peak
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72	14.474	2138	2140	2142	rBV	123789	124809	3.52%	0.286%
73	14.780	2190	2192	2196	rBV2	96613	102691	2.90%	0.235%
74	15.045	2233	2237	2240	rBV	436820	690364	19.47%	1.581%
75	15.115	2246	2249	2252	rBV	233215	255583	7.21%	0.585%
76	15.157	2252	2256	2265	rVB2	250361	431237	12.16%	0.987%
77	15.345	2284	2288	2293	rVB	271938	335048	9.45%	0.767%
78	15.404	2294	2298	2304	rVB	387090	454962	12.83%	1.042%
79	15.468	2304	2309	2312	rBV	959198	1205418	33.99%	2.760%
80	15.933	2384	2388	2394	rVB	353825	513835	14.49%	1.176%
81	16.921	2551	2556	2560	rBV2	156968	309076	8.71%	0.708%
82	17.104	2581	2587	2593	rBV2	320275	699293	19.72%	1.601%

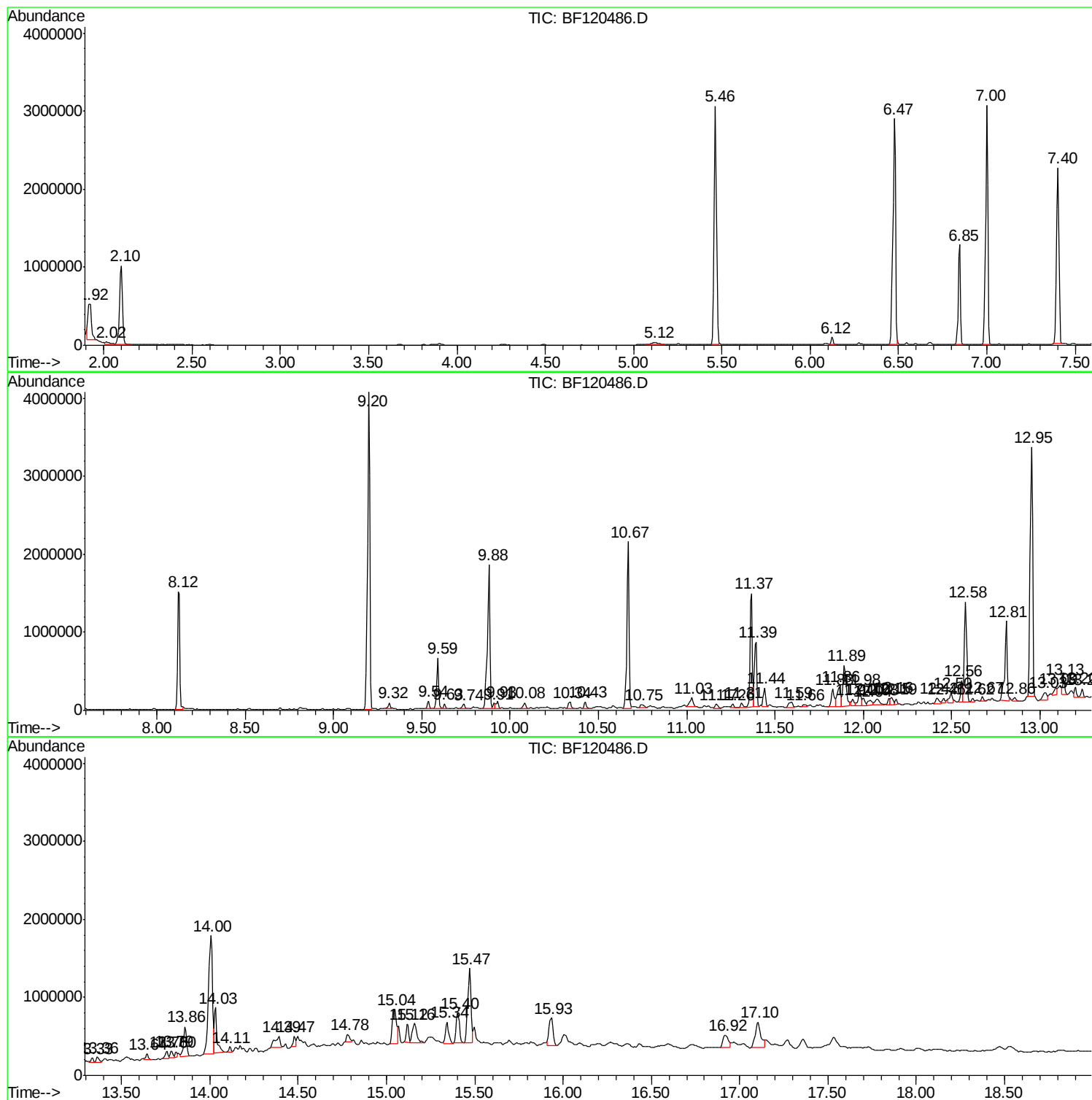
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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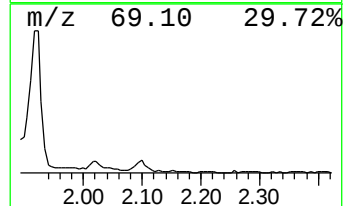
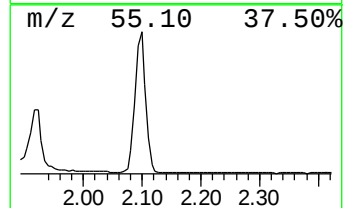
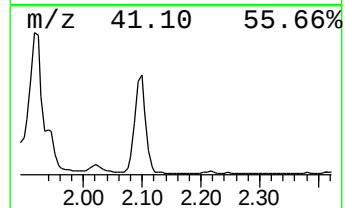
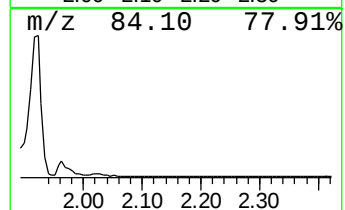
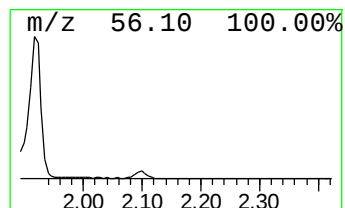
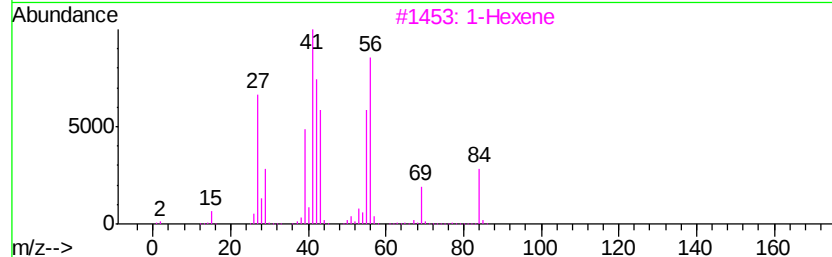
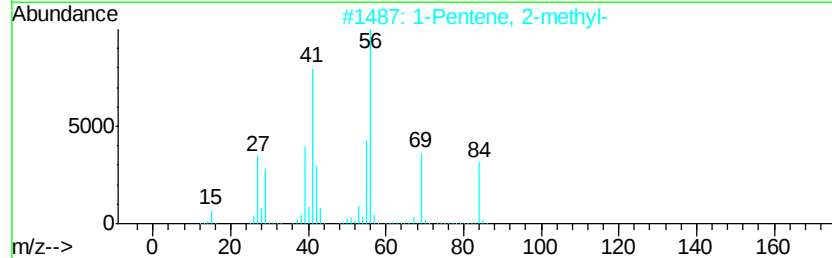
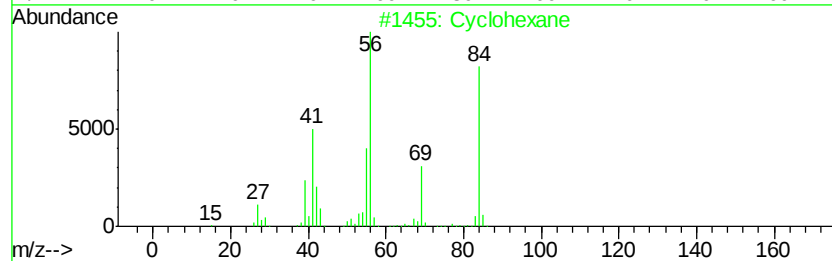
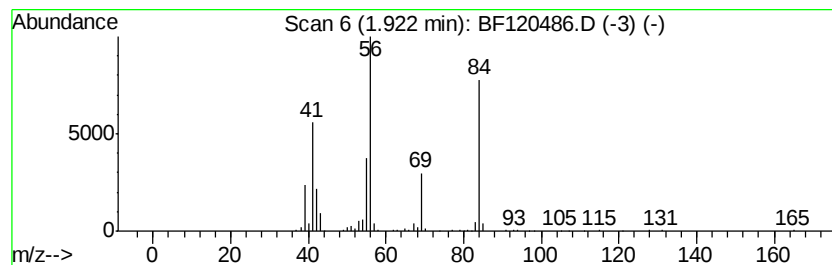
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Cyclohexane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.92	10.03 ng	552420	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane	84	C6H12	000110-82-7	95
2		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	72
3		1-Hexene	84	C6H12	000592-41-6	59
4		2H-Pyran-2,6(3H)-dione, dihydro-...	142	C7H10O3	004160-82-1	53
5		2-Amino-oxazole	84	C3H4N2O	004570-45-0	50



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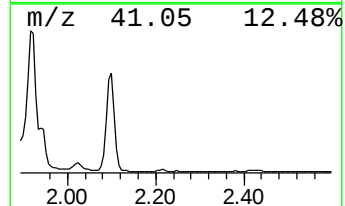
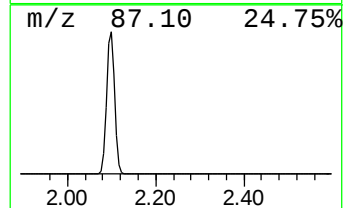
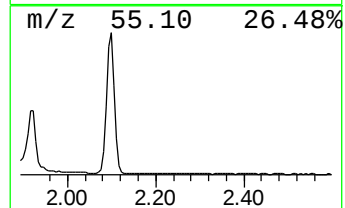
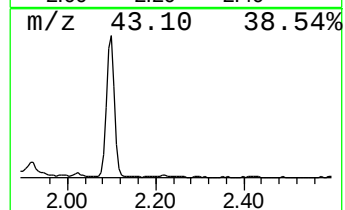
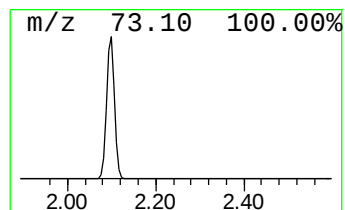
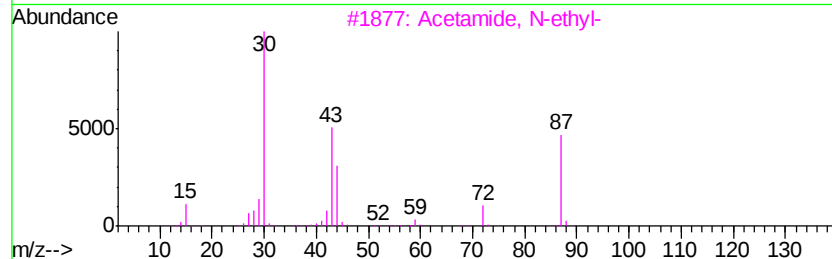
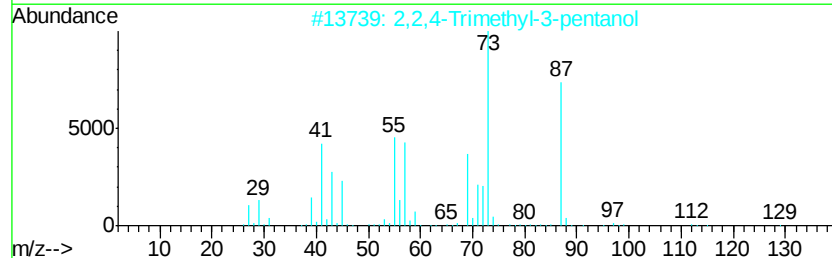
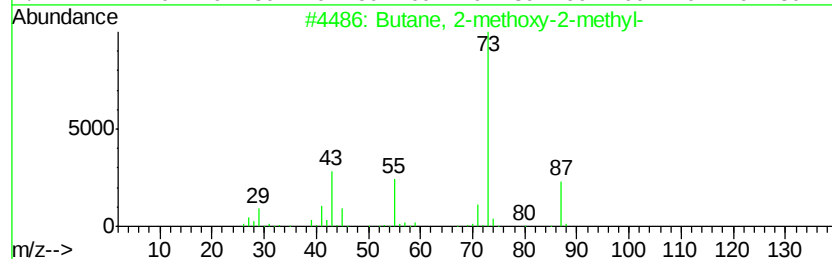
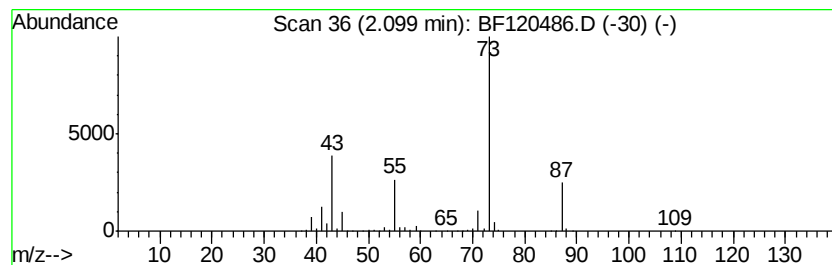
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.10	23.09 ng	1272040	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	27
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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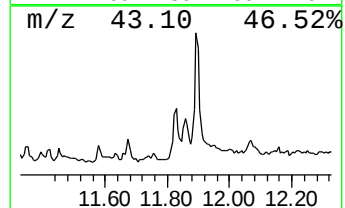
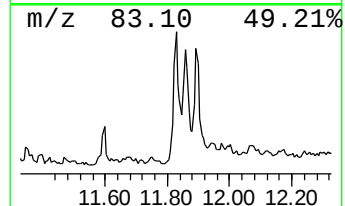
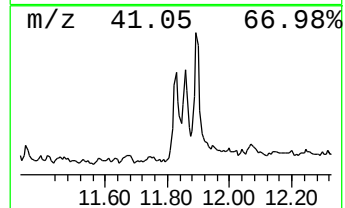
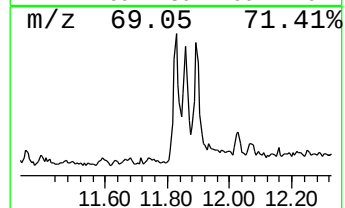
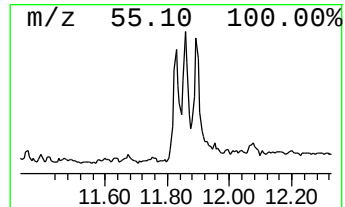
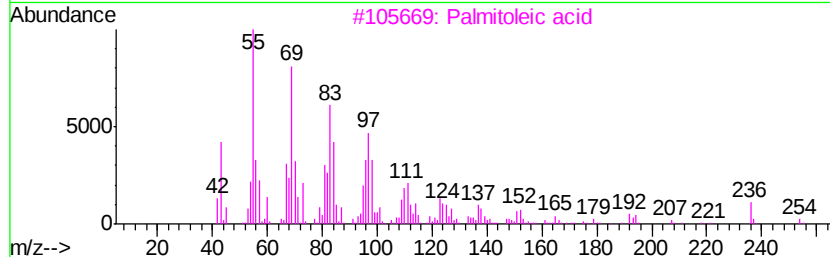
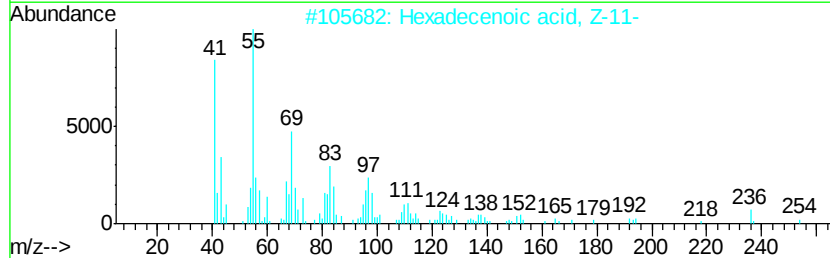
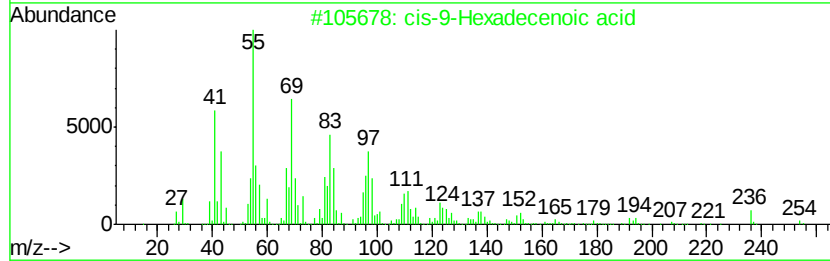
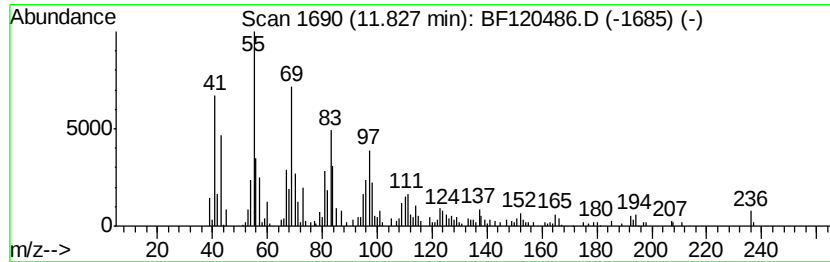
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 cis-9-Hexadecenoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.83	4.27 ng	305696	Phenanthrene-d10	11.37	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	cis-9-Hexadecenoic acid	254	C16H30O2	1000333-19-5	93
2	Hexadecenoic acid, Z-11-	254	C16H30O2	002416-20-8	91
3	Palmitoleic acid	254	C16H30O2	000373-49-9	91
4	Z-7-Tetradecenoic acid	226	C14H26O2	1000130-98-4	87
5	Oxacyclotridecan-2-one	198	C12H22O2	000947-05-7	78



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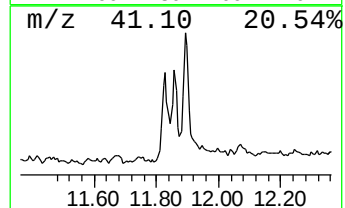
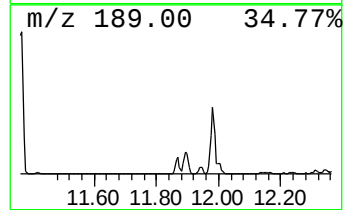
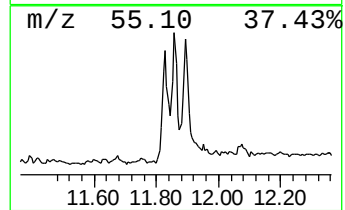
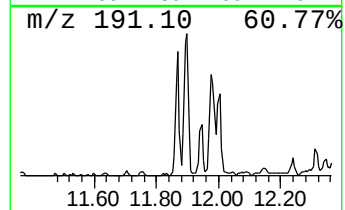
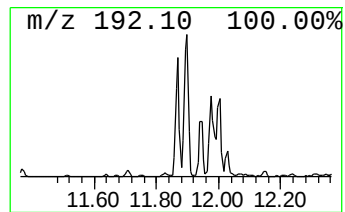
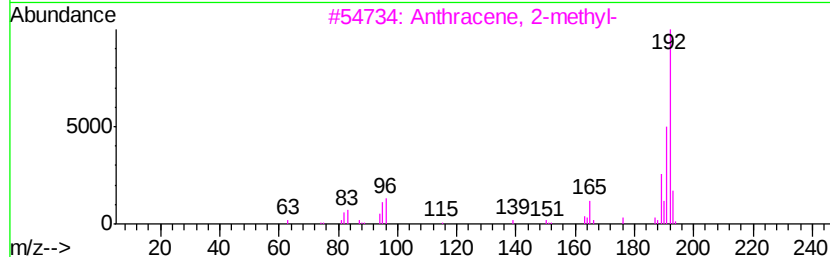
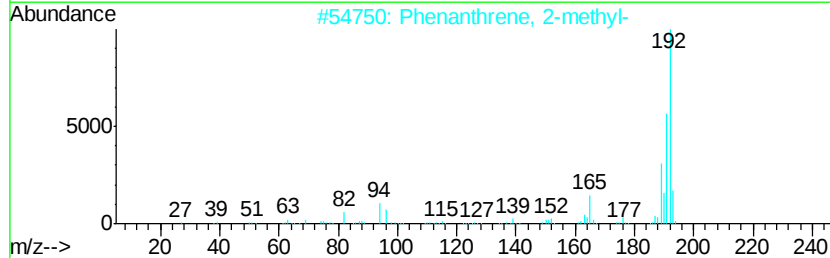
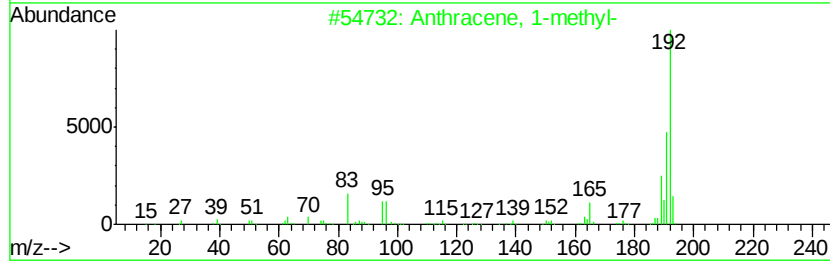
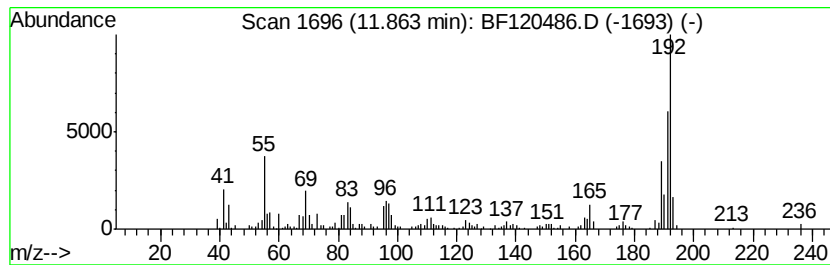
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Anthracene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.86	4.56 ng	326631	Phenanthrene-d10	11.37

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 1-methyl-	192	C15H12	000610-48-0	96
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060220\
Data File : BF120486.D
Acq On : 2 Jun 2020 11:18
Operator : JU/CG
Sample : L2792-07
Misc :
ALS Vial : 3 Sample Multiplier: 1

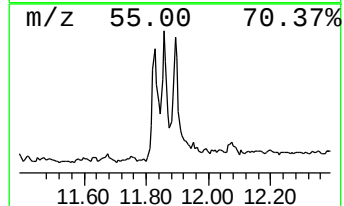
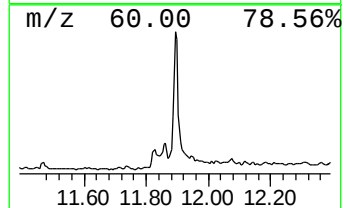
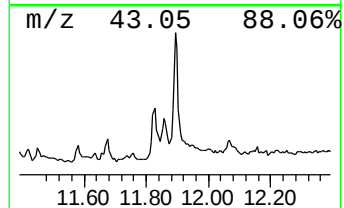
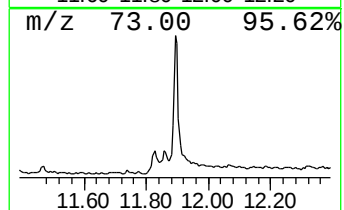
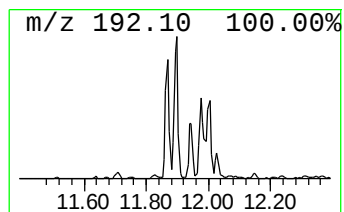
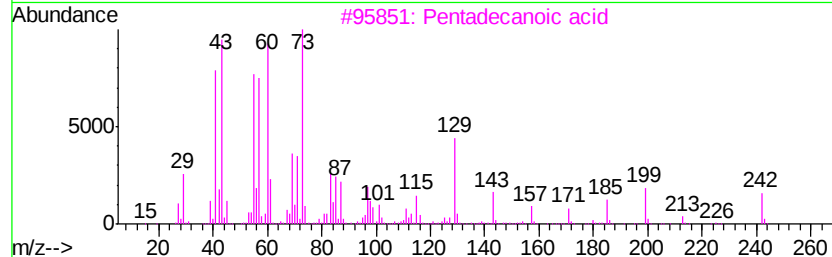
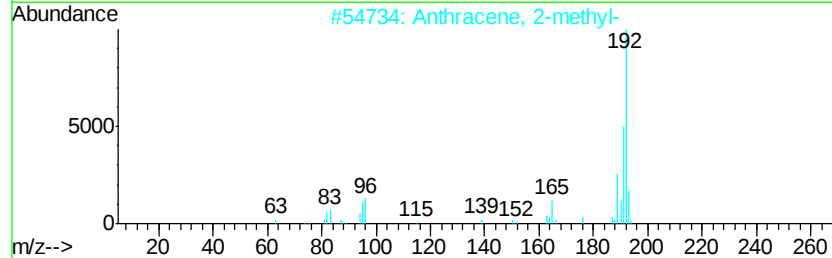
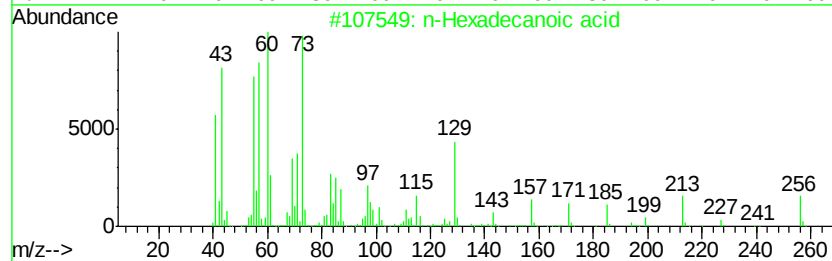
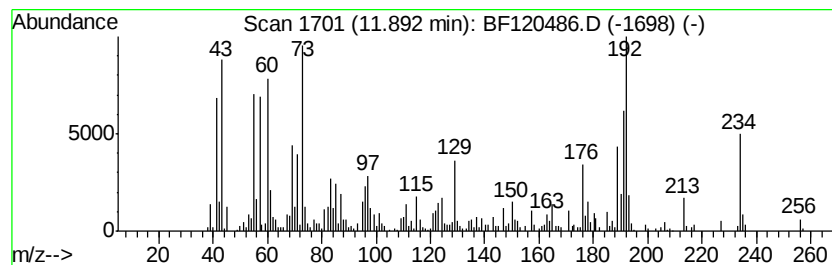
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ClientSampleId :
NM107-COMP-2020-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
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 5

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.89	8.83 ng	632210	Phenanthrene-d10		11.37
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	46
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	38
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	38
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060220\
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 Acq On : 2 Jun 2020 11:18
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 Sample : L2792-07
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

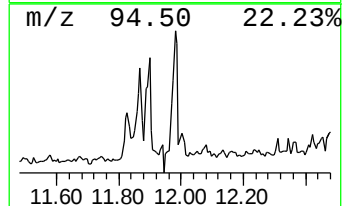
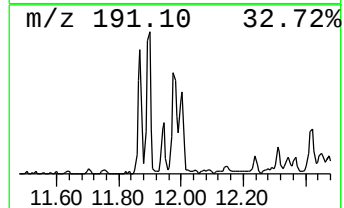
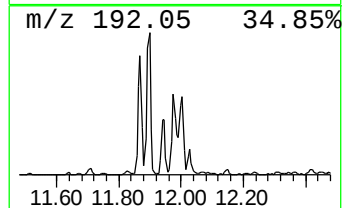
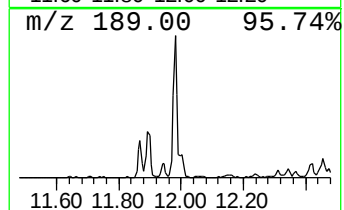
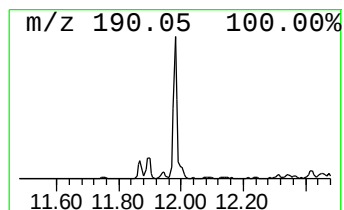
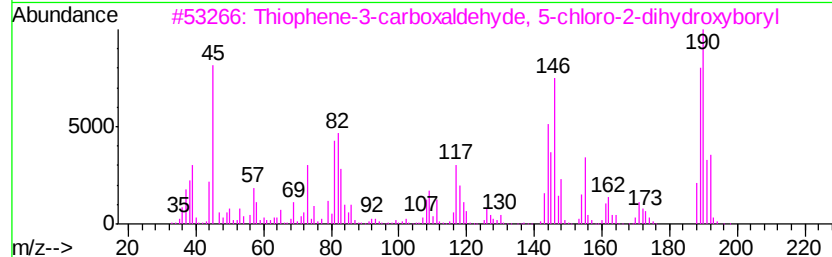
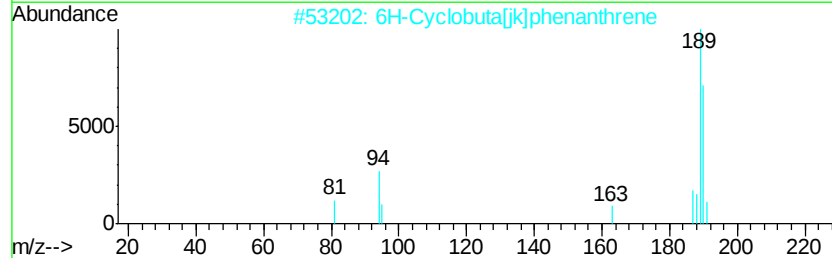
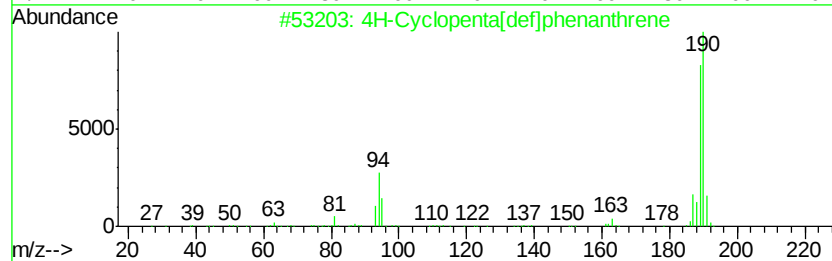
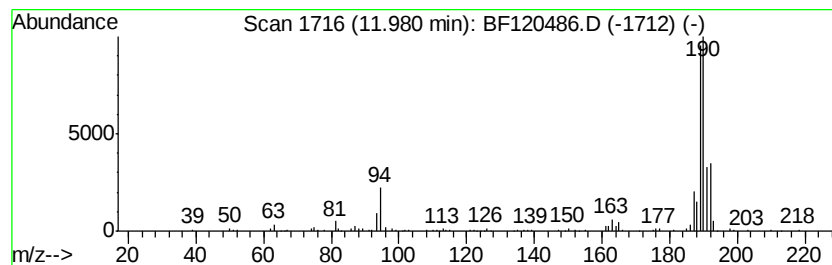
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.98	3.33 ng	238704	Phenanthrene-d10	11.37	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	81
2	6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	72
3	Thiophene-3-carboxaldehyde, 5-chloro-2-dihydroxyboryl	190	C5H4BClO3S	036155-87-0	52
4	2,3,5,6-Tetramethylterephthalaldehyde	190	C12H14O2	007072-01-7	50
5	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060220\
Data File : BF120486.D
Acq On : 2 Jun 2020 11:18
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Sample : L2792-07
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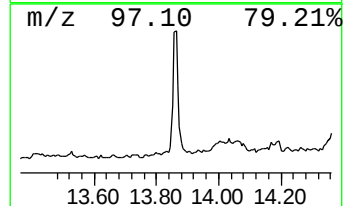
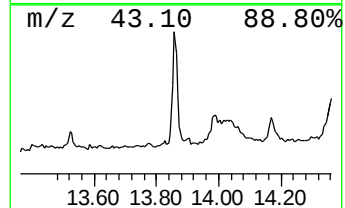
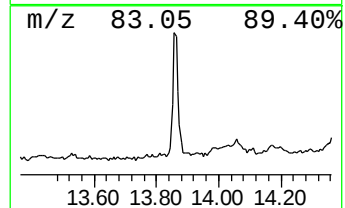
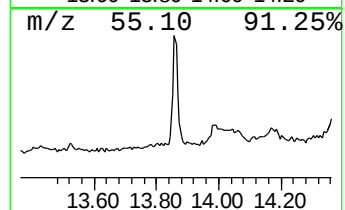
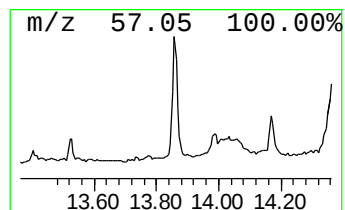
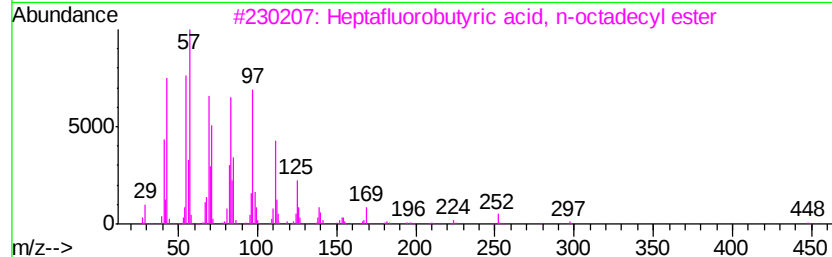
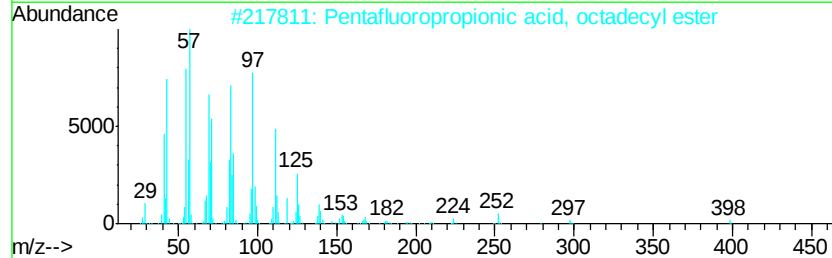
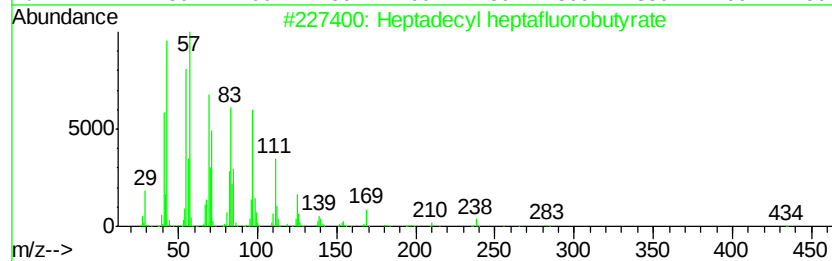
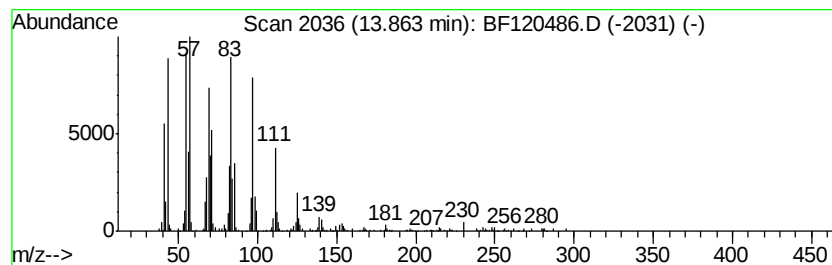
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptadecyl heptafluorobutyrate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.86	4.48 ng	491849	Chrysene-d12	14.01		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	93
2		Pentafluoropropionic acid, octadecyl ester	416	C21H37F5O2	959261-25-7	90
3		Heptafluorobutyric acid, n-octadecyl ester	466	C22H37F7O2	000400-57-7	90
4		Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	90
5		Pentafluoropropionic acid, heptadecyl ester	402	C20H35F5O2	959218-78-1	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060220\
Data File : BF120486.D
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Sample : L2792-07
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Instrument :
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NM107-COMP-2020-0-6

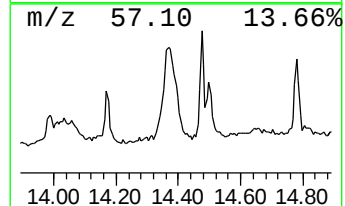
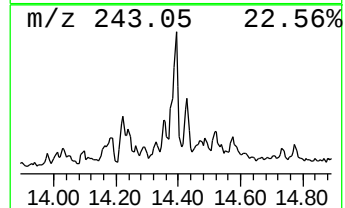
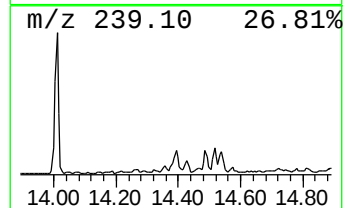
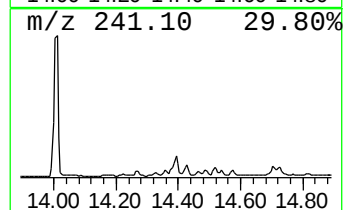
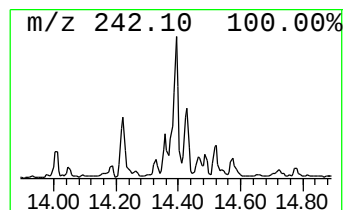
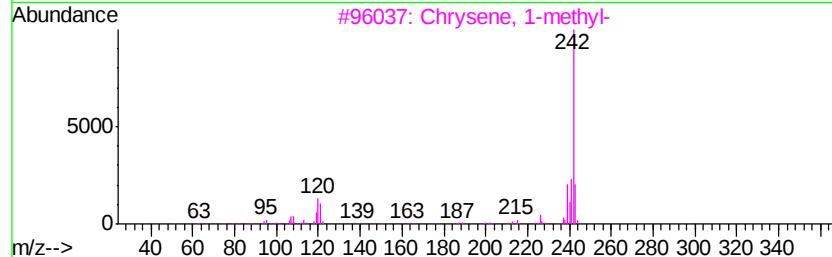
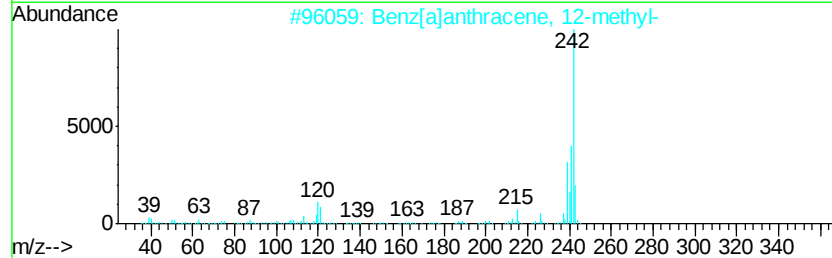
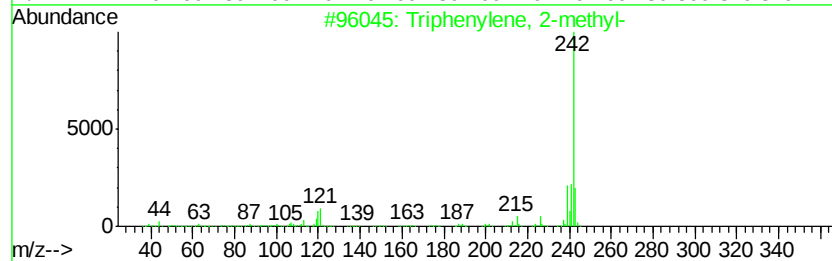
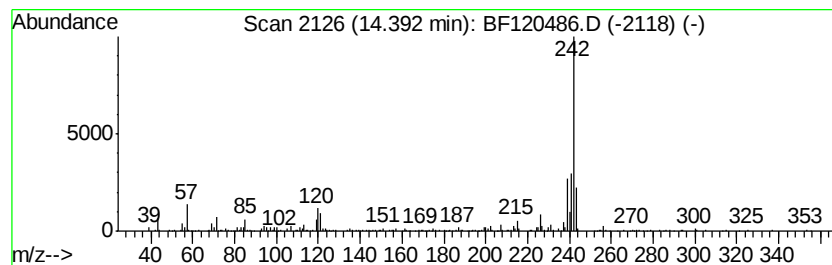
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.39	2.84 ng	311282	Chrysene-d12	14.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Triphenylene, 2-methyl-	242	C19H14	001705-84-6	96
2		Benz[a]anthracene, 12-methyl-	242	C19H14	002422-79-9	96
3		Chrysene, 1-methyl-	242	C19H14	003351-28-8	94
4		Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	94
5		Chrysene, 6-methyl-	242	C19H14	001705-85-7	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060220\
Data File : BF120486.D
Acq On : 2 Jun 2020 11:18
Operator : JU/CG
Sample : L2792-07
Misc :
ALS Vial : 3 Sample Multiplier: 1

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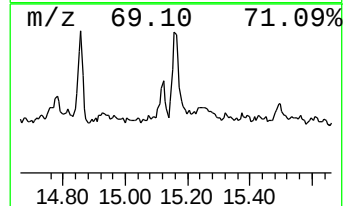
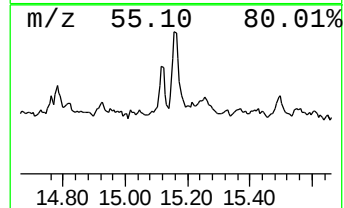
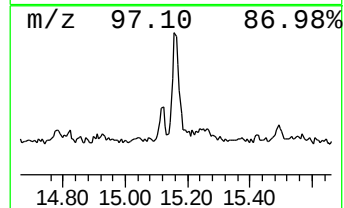
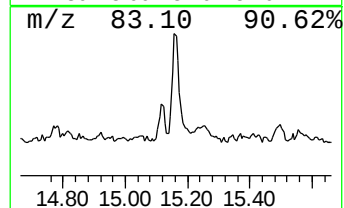
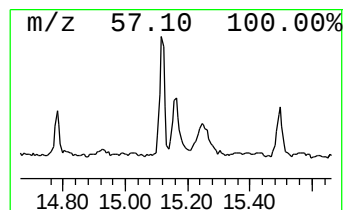
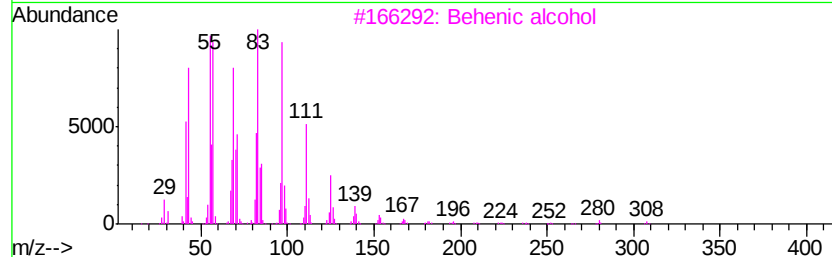
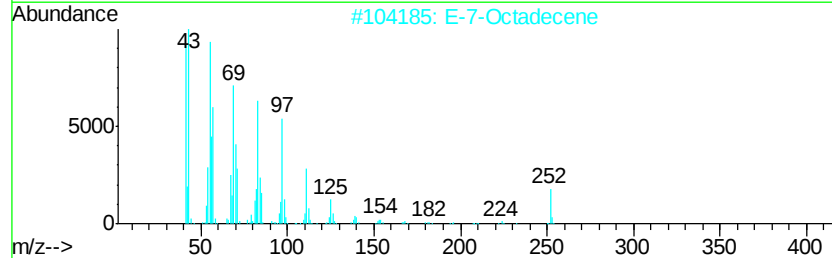
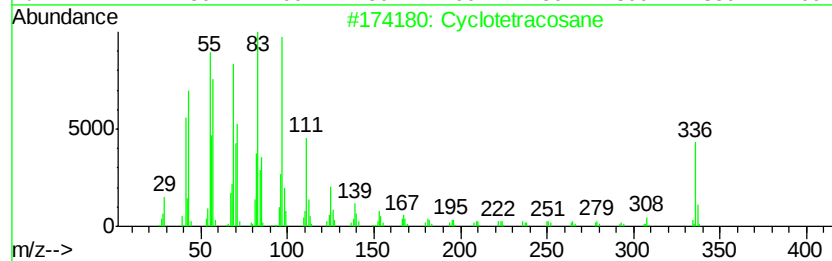
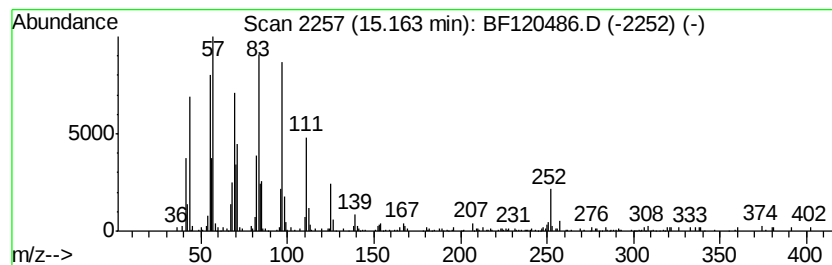
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Cyclotetracosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.16	7.15 ng	431237	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclotetracosane	336	C24H48	000297-03-0	96
2		E-7-Octadecene	252	C18H36	1000130-92-0	94
3		Behenic alcohol	326	C22H46O	000661-19-8	91
4		n-Tetracosanol-1	354	C24H50O	000506-51-4	91
5		1-Heneicosanol	312	C21H44O	015594-90-8	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060220\
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Acq On : 2 Jun 2020 11:18
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Misc :
ALS Vial : 3 Sample Multiplier: 1

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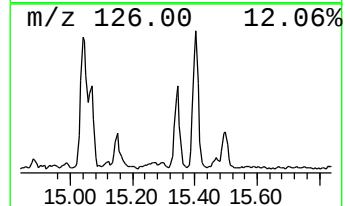
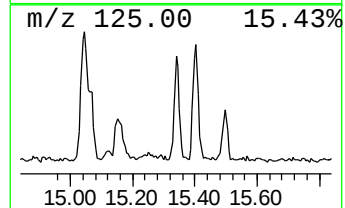
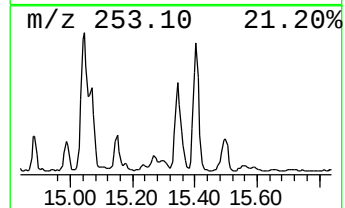
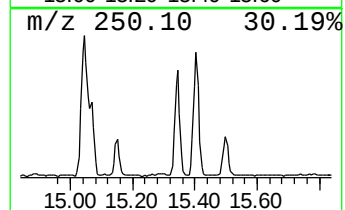
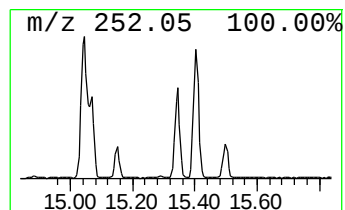
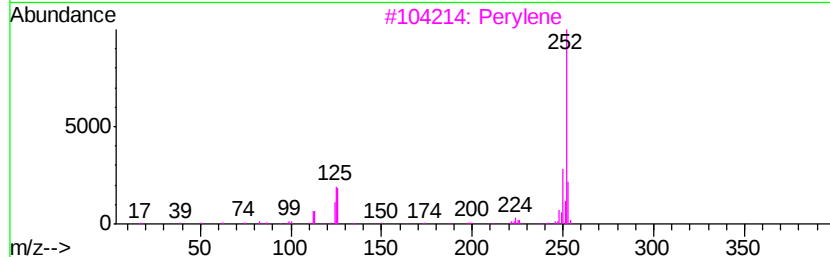
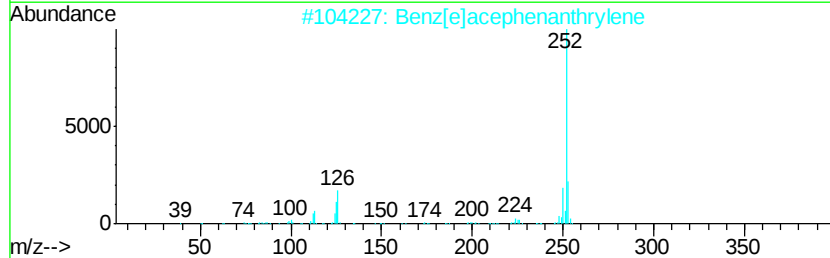
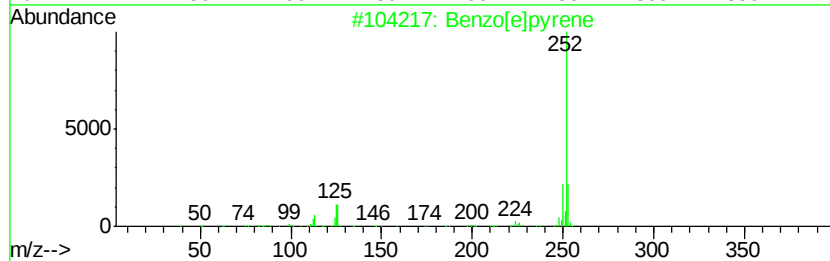
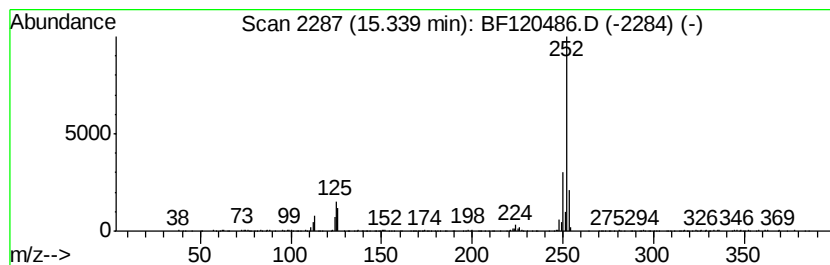
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzo[e]pyrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.34	5.56 ng	335048	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	96
3		Perylene	252	C20H12	000198-55-0	95
4		Benzo[il]fluoranthene	252	C20H12	000205-82-3	94
5		Benzo[a]pyrene	252	C20H12	000050-32-8	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060220\
Data File : BF120486.D
Acq On : 2 Jun 2020 11:18
Operator : JU/CG
Sample : L2792-07
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
NM107-COMP-2020-0-6

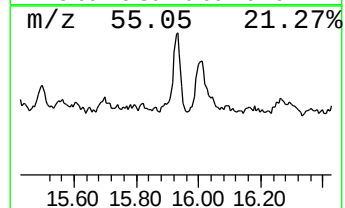
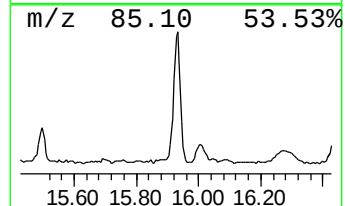
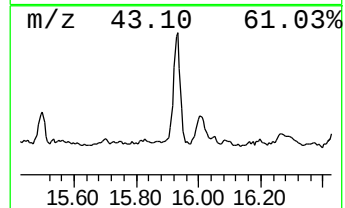
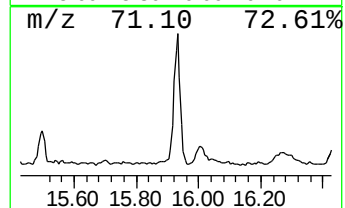
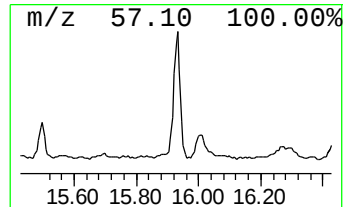
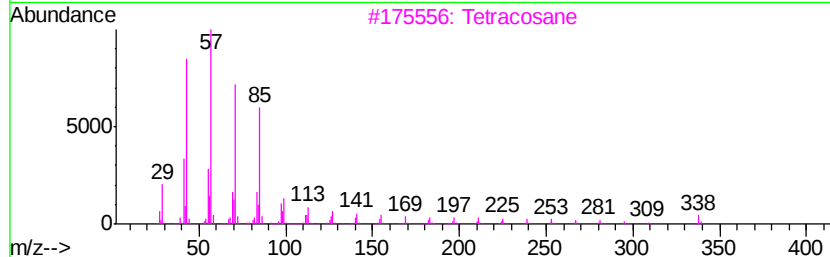
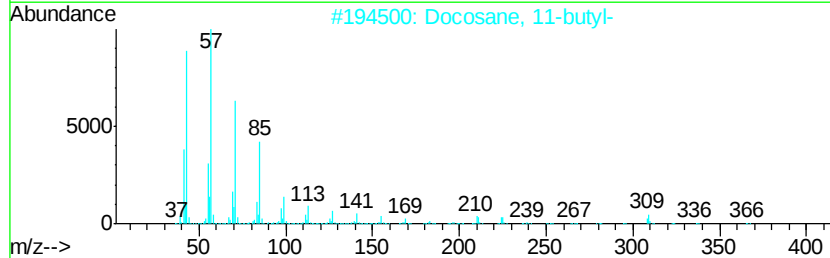
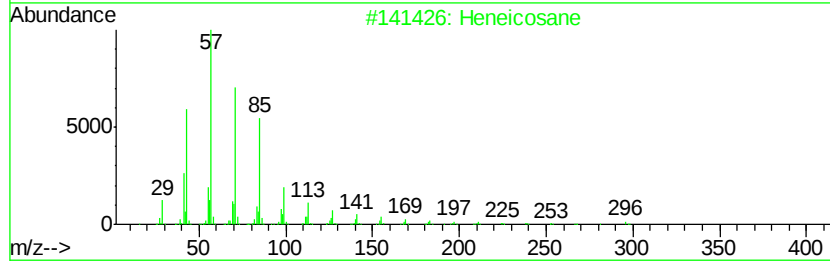
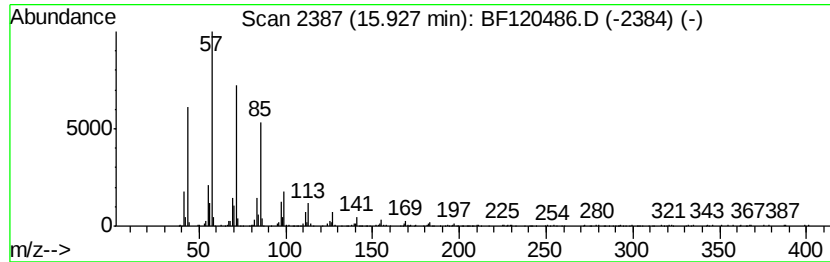
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.93	8.53 ng	513835	Perylene-d12	15.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	97
2			Docosane, 11-butyl-	366	C26H54	013475-76-8	94
3			Tetracosane	338	C24H50	000646-31-1	94
4			Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	94
5			Pentacosane	352	C25H52	000629-99-2	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060220\
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Sample : L2792-07
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
NM107-COMP-2020-0-6

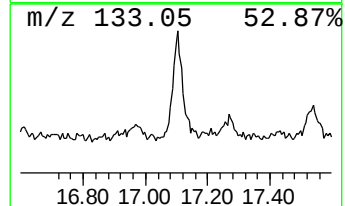
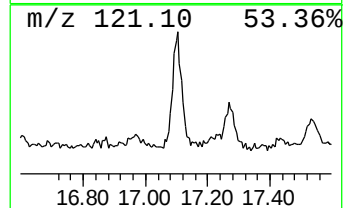
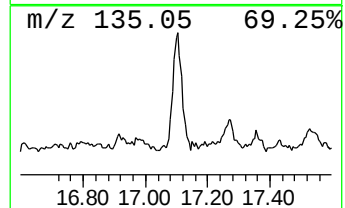
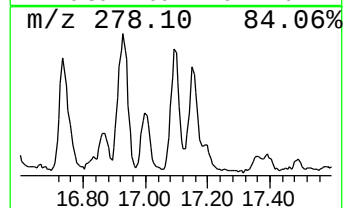
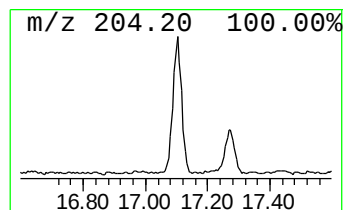
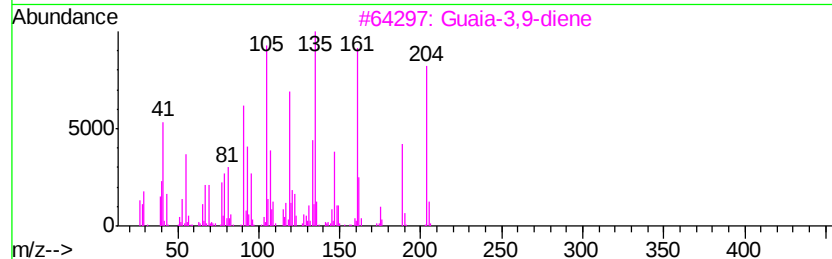
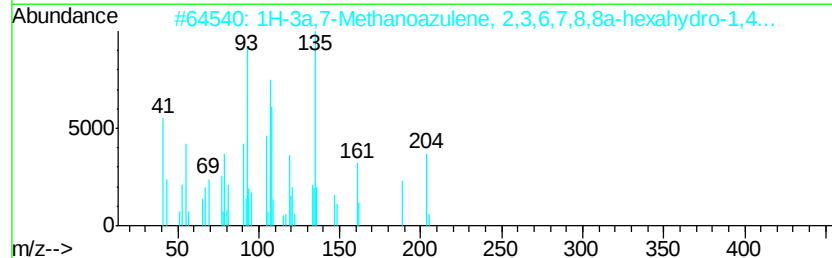
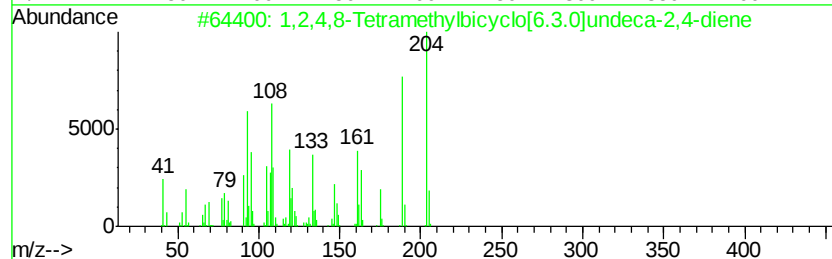
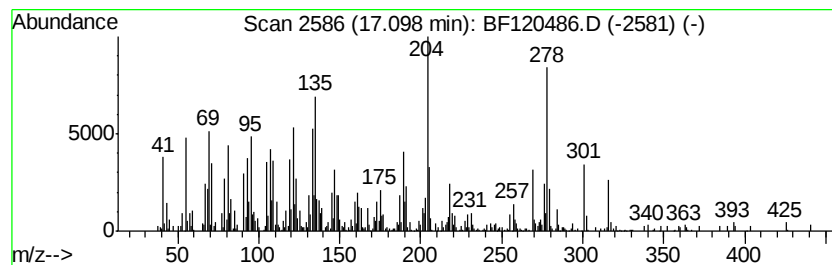
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1,2,4,8-Tetramethylbicyclo[...] Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.10	11.60 ng	699293	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2,4,8-Tetramethylbicyclo[6.3.0]undeca-2,4-diene	204	C15H24	137235-51-9	78
2		1H-3a,7-Methanoazulene, 2,3,6,7,8a-hexahydro-1,4-diene	204	C15H24	000560-32-7	46
3		Guaia-3,9-diene	204	C15H24	000489-83-8	46
4		1,4-Dimethyl-8-isopropylidenetriene	204	C15H24	1000140-07-7	46
5		Farnesyl bromide	284	C15H25Br	006874-67-5	45



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Instrument :
 BNA_F
 ClientSampleId :
 NM107-COMP-2020-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF052820.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
Cyclohexane	1.92	10.0	ng	552420	1	6.85	1101880	20.0
Butane, 2-methoxy...	2.10	23.1	ng	1272040	1	6.85	1101880	20.0
cis-9-Hexadecenoic...	11.83	4.3	ng	305696	4	11.37	1432530	20.0
Anthracene, 1-met...	11.86	4.6	ng	326631	4	11.37	1432530	20.0
n-Hexadecanoic acid	11.89	8.8	ng	632210	4	11.37	1432530	20.0
4H-Cyclopenta[def...	11.98	3.3	ng	238704	4	11.37	1432530	20.0
Heptadecyl hepta...	13.86	4.5	ng	491849	5	14.01	2193320	20.0
Triphenylene, 2-m...	14.39	2.8	ng	311282	5	14.01	2193320	20.0
Cyclotetracosane	15.16	7.2	ng	431237	6	15.47	1205420	20.0
Benzo[e]pyrene	15.34	5.6	ng	335048	6	15.47	1205420	20.0
Heneicosane	15.93	8.5	ng	513835	6	15.47	1205420	20.0
1,2,4,8-Tetrameth...	17.10	11.6	ng	699293	6	15.47	1205420	20.0