

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF031820\
 Data File : BF119424.D
 Acq On : 18 Mar 2020 19:30
 Operator : CG/JU
 Sample : L1942-13
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3-SA

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-BF031820.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	2.163	8	13	25	rVB	320112	468153	4.91%	0.478%
2	3.340	210	213	224	rBV	44923	98996	1.04%	0.101%
3	5.087	505	510	517	rVB	127279	169486	1.78%	0.173%
4	5.469	569	575	578	rBV	8642464	9164872	96.11%	9.350%
5	6.475	740	746	749	rBV	7661917	9344821	97.99%	9.534%
6	6.675	777	780	783	rBV	273272	201392	2.11%	0.205%
7	6.857	807	811	814	rBV	2363164	2177792	22.84%	2.222%
8	7.010	833	837	841	rBV	7144655	6969988	73.09%	7.111%
9	7.416	900	906	909	rBV	5653191	5935615	62.24%	6.056%
10	8.139	1024	1029	1032	rBV	3383546	2891588	30.32%	2.950%
11	9.222	1207	1213	1216	rBV	10260515	9536067	100.00%	9.729%
12	9.610	1275	1279	1282	rVB	1955150	1478720	15.51%	1.509%
13	9.757	1301	1304	1307	rBV	142574	107196	1.12%	0.109%
14	9.898	1322	1328	1331	rBV	3702640	3091647	32.42%	3.154%
15	9.927	1331	1333	1337	rVB	203031	156365	1.64%	0.160%
16	10.098	1358	1362	1365	rBV	157470	141304	1.48%	0.144%
17	10.192	1374	1378	1385	rBV	169160	250082	2.62%	0.255%
18	10.445	1417	1421	1426	rVB2	245060	213051	2.23%	0.217%
19	10.686	1456	1462	1465	rBV	6317135	6109061	64.06%	6.233%
20	11.280	1561	1563	1569	rVB	141530	116738	1.22%	0.119%
21	11.386	1577	1581	1583	rBV	3547808	3106421	32.58%	3.169%
22	11.410	1583	1585	1588	rVV	2219714	1631557	17.11%	1.665%
23	11.457	1588	1593	1596	rVB	603642	517473	5.43%	0.528%
24	11.610	1613	1619	1622	rBV	172303	169060	1.77%	0.172%
25	11.886	1663	1666	1668	rBV	206181	173552	1.82%	0.177%
26	11.916	1668	1671	1675	rVV	1108129	959742	10.06%	0.979%
27	11.963	1675	1679	1681	rVV2	155465	179059	1.88%	0.183%
28	11.998	1681	1685	1687	rVV	430888	431476	4.52%	0.440%
29	12.021	1687	1689	1692	rVB	139331	114208	1.20%	0.117%
30	12.180	1713	1716	1718	rBV	164271	149276	1.57%	0.152%
31	12.439	1756	1760	1763	rBV	112293	124597	1.31%	0.127%
32	12.480	1763	1767	1772	rVB3	98144	134585	1.41%	0.137%
33	12.598	1783	1787	1791	rBV	2721950	2384567	25.01%	2.433%
34	12.704	1800	1805	1811	rBV3	432936	560573	5.88%	0.572%

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Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	12.827	1820	1826	1829	rBV	2529803	2482744	26.04%	2.533%
36	12.874	1832	1834	1839	rVB3	126771	146894	1.54%	0.150%
37	12.945	1843	1846	1847	rBV	146101	113759	1.19%	0.116%
38	12.980	1847	1852	1855	rVV	7539643	7968092	83.56%	8.129%
39	13.045	1858	1863	1866	rBV4	146906	245025	2.57%	0.250%
40	13.133	1875	1878	1879	rBV2	103518	115516	1.21%	0.118%
41	13.157	1879	1882	1884	rVB2	317665	295363	3.10%	0.301%
42	13.221	1889	1893	1896	rBV	309107	263413	2.76%	0.269%
43	13.257	1896	1899	1902	rBV2	222674	222373	2.33%	0.227%
44	13.380	1918	1920	1924	rBV2	136294	126858	1.33%	0.129%
45	13.557	1943	1950	1952	rBV3	160516	210527	2.21%	0.215%
46	13.657	1965	1967	1973	rVB2	73486	109346	1.15%	0.112%
47	13.774	1982	1987	1989	rBV2	130487	162498	1.70%	0.166%
48	13.804	1989	1992	1994	rVV	172947	170943	1.79%	0.174%
49	13.827	1994	1996	2000	rVV2	144053	196097	2.06%	0.200%
50	13.874	2000	2004	2010	rVB2	931864	1018115	10.68%	1.039%
51	14.027	2025	2030	2032	rBV2	2341600	2974421	31.19%	3.035%
52	14.051	2032	2034	2039	rVB	956325	797047	8.36%	0.813%
53	14.127	2045	2047	2051	rVB	171886	150090	1.57%	0.153%
54	14.204	2056	2060	2063	rVB2	381450	380176	3.99%	0.388%
55	14.245	2063	2067	2071	rBV2	81422	105648	1.11%	0.108%
56	14.409	2092	2095	2098	rVB	164267	147258	1.54%	0.150%
57	14.509	2108	2112	2114	rBV2	626627	627830	6.58%	0.641%
58	14.557	2118	2120	2124	rVB3	106847	103431	1.08%	0.106%
59	14.815	2158	2164	2171	rVB	663891	746006	7.82%	0.761%
60	15.068	2202	2207	2210	rBV	670299	1143826	11.99%	1.167%
61	15.162	2218	2223	2230	rVB2	665725	876114	9.19%	0.894%
62	15.368	2254	2258	2264	rVB	462816	566310	5.94%	0.578%
63	15.427	2264	2268	2274	rVB	680900	855277	8.97%	0.873%
64	15.498	2274	2280	2283	rBV	1410359	1853120	19.43%	1.891%
65	15.551	2286	2289	2297	rVB	512624	642322	6.74%	0.655%
66	15.992	2358	2364	2368	rBV	333850	533758	5.60%	0.545%
67	16.039	2368	2372	2380	rVB8	104871	232866	2.44%	0.238%
68	16.509	2446	2452	2459	rBV	159681	301910	3.17%	0.308%
69	16.786	2495	2499	2505	rVB4	57429	127282	1.33%	0.130%
70	16.956	2522	2528	2536	rVB2	332215	667697	7.00%	0.681%
71	17.133	2551	2558	2563	rBV3	63783	167534	1.76%	0.171%

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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

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Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.M
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72	17.198	2564	2569	2584	rVB6	79250	231100	2.42%	0.236%
73	17.398	2596	2603	2609	rBV	269229	532040	5.58%	0.543%
74	17.633	2638	2643	2649	rVB	78203	149633	1.57%	0.153%

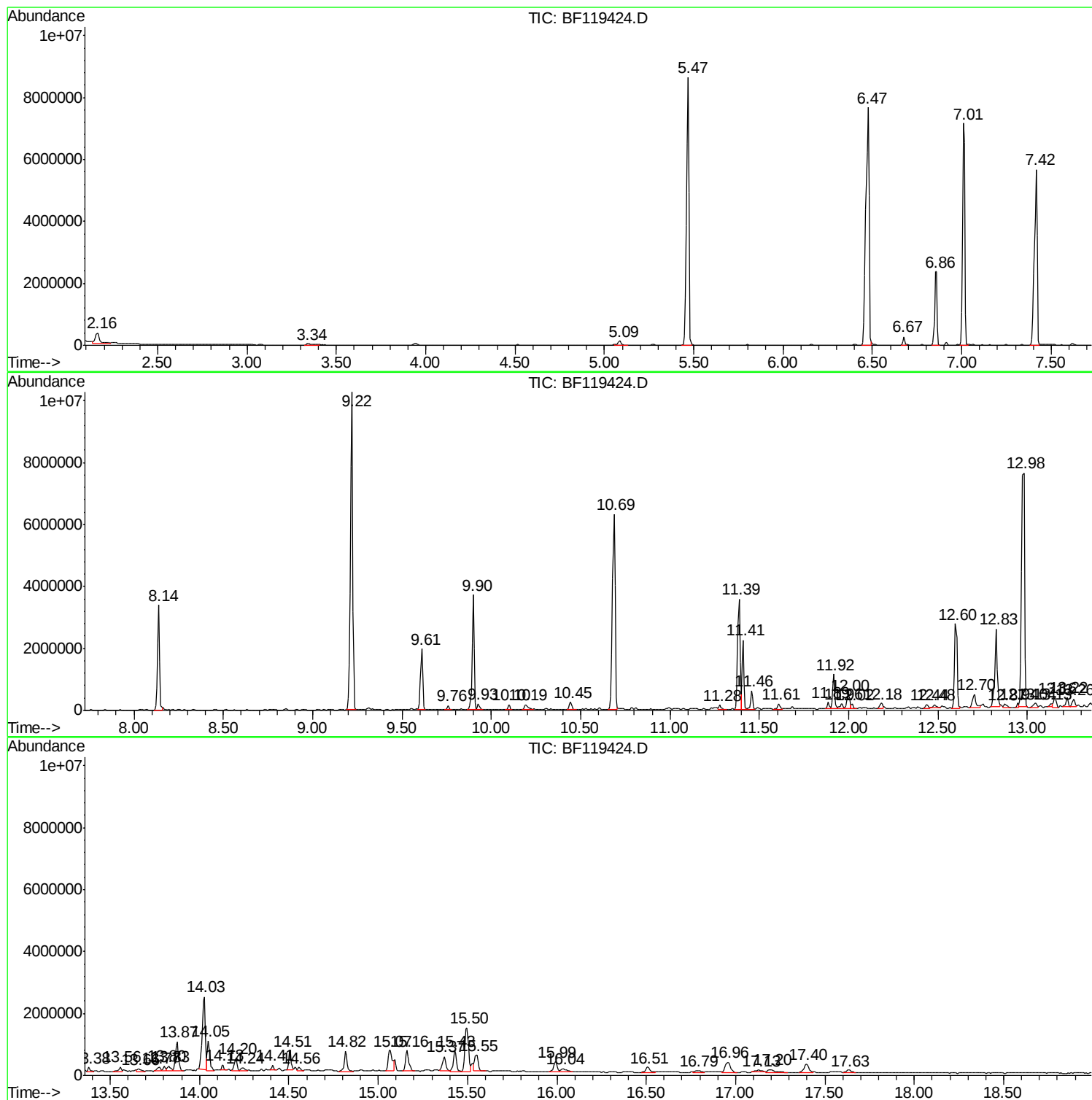
Sum of corrected areas: 98019339

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ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
TP-3-SA

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.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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Instrument :
BNA_F
ClientSampleId :
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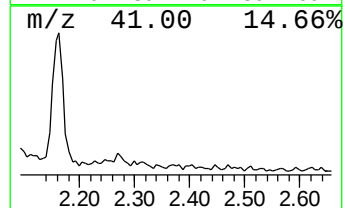
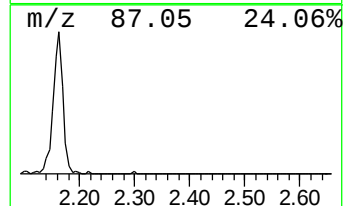
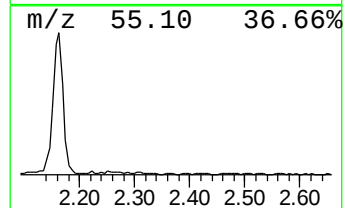
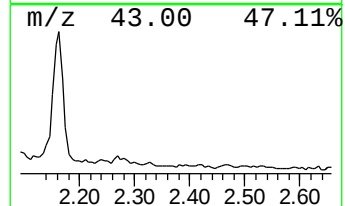
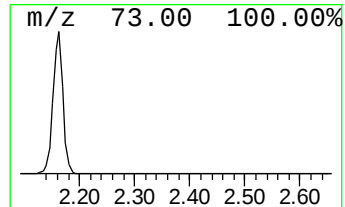
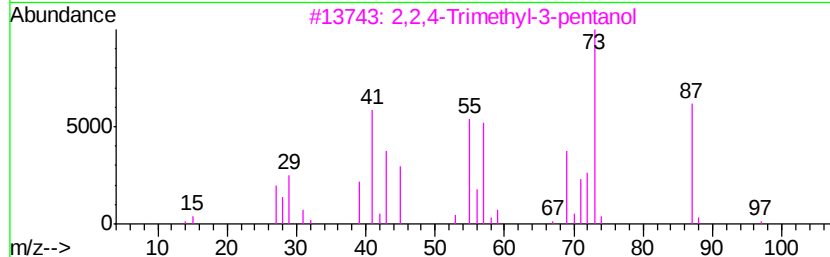
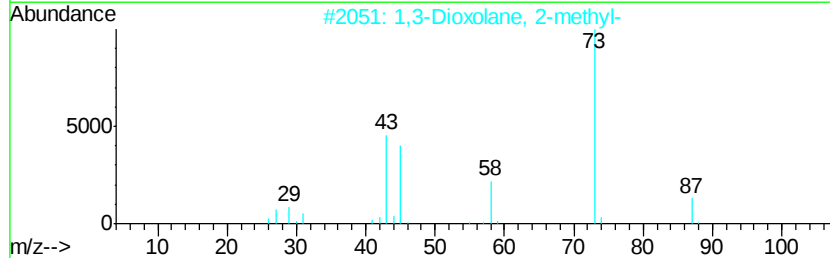
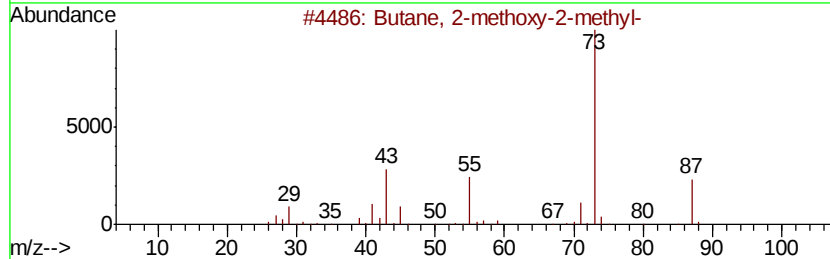
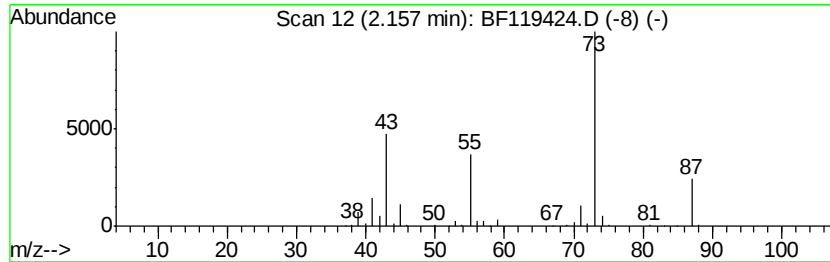
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.16	4.30 ng	468153	1,4-Dichlorobenzene-d4	6.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	47
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
5			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	25



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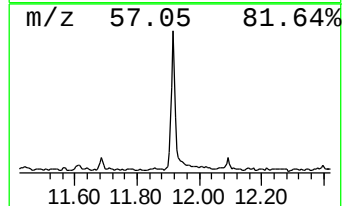
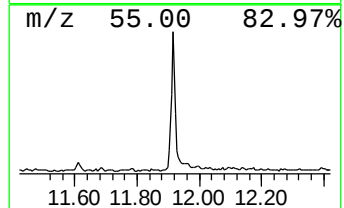
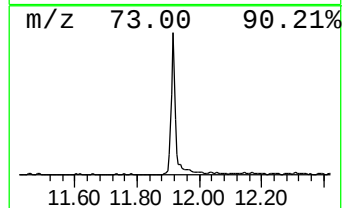
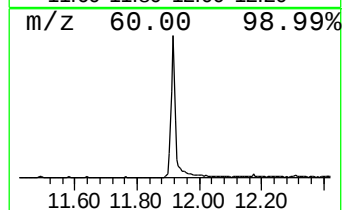
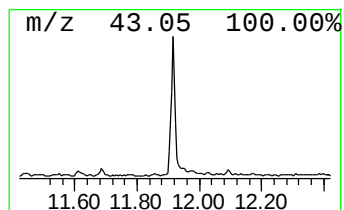
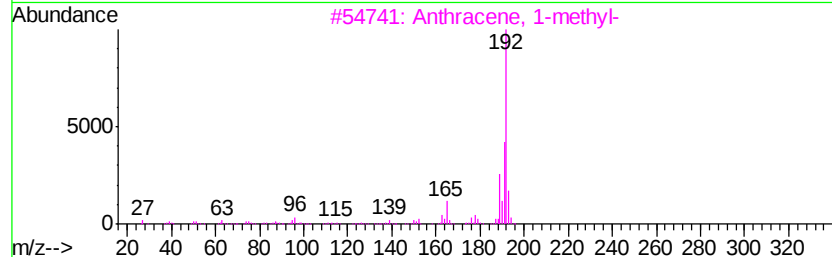
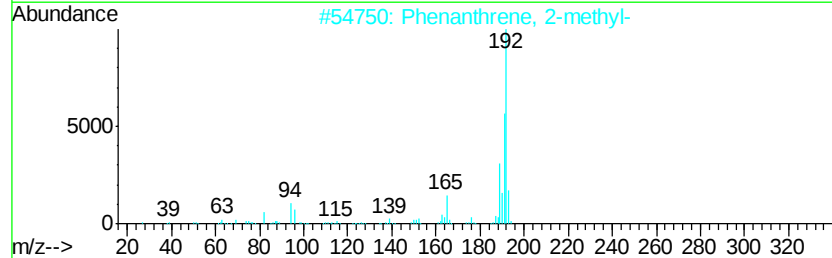
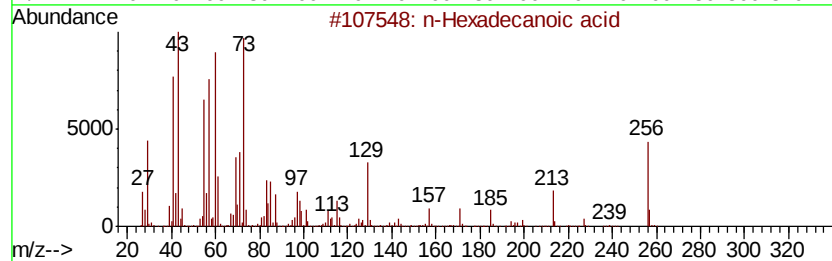
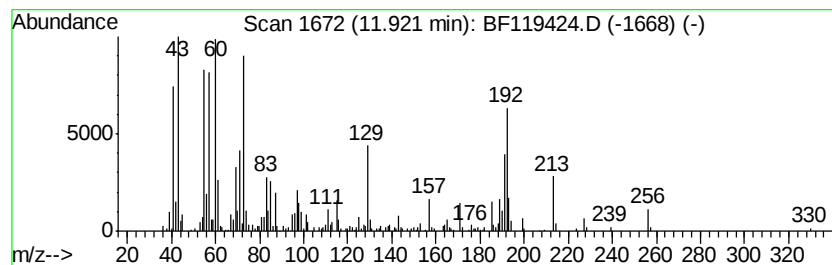
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.92	6.18 ng	959742	Phenanthrene-d10	11.39	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93



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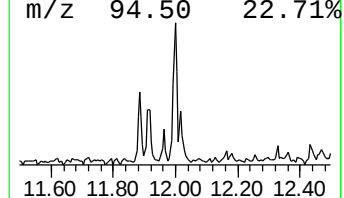
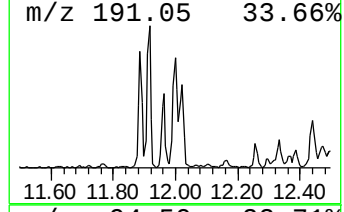
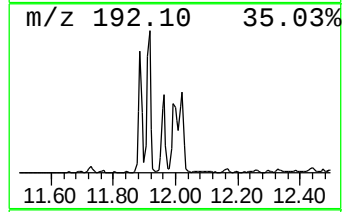
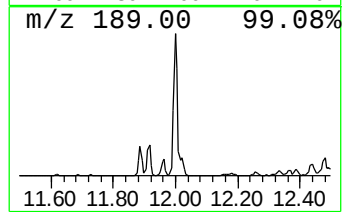
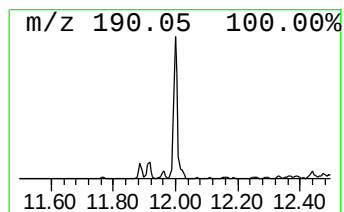
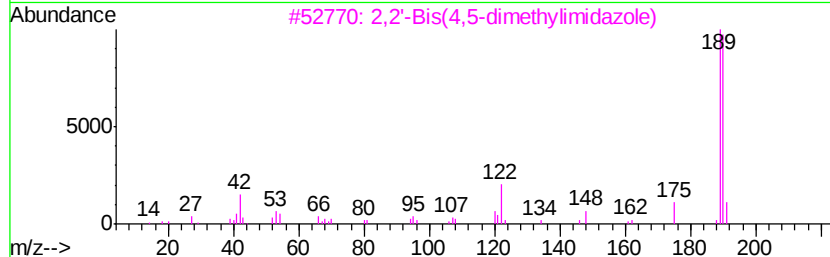
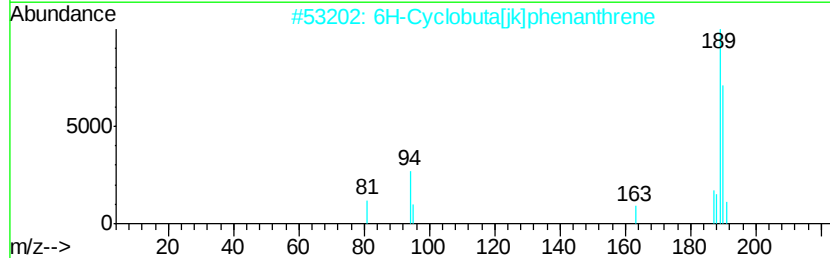
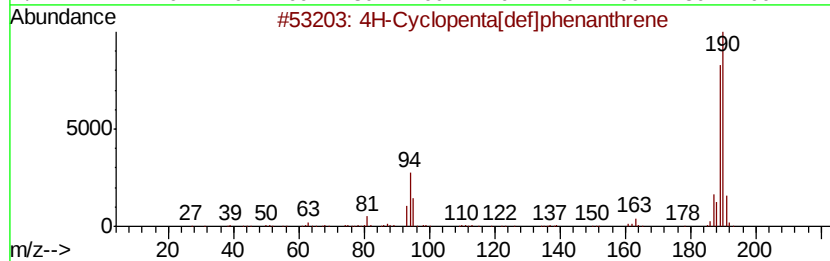
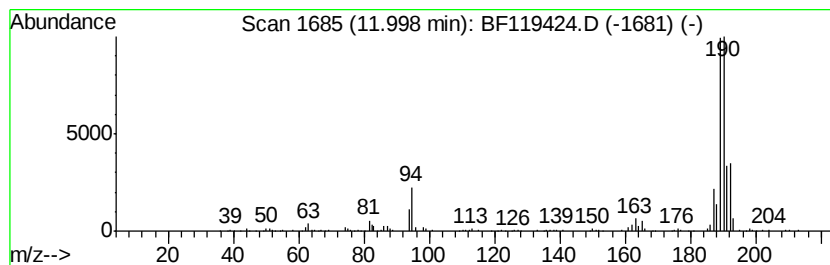
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 4H-Cyclopenta[def]phenanthrene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.00	2.78 ng	431476	Phenanthrene-d10	11.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2		6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	56
3		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	40
4		Megastigmatrienone	190	C13H18O	038818-55-2	25
5		2-Naphthaldehyde, 1,2,3,4-tetra...	190	C12H14O2	002472-02-8	18



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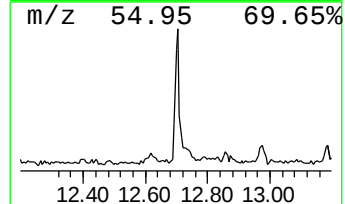
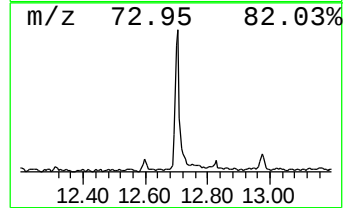
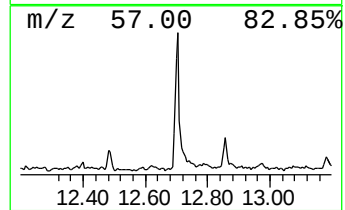
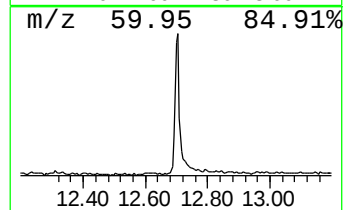
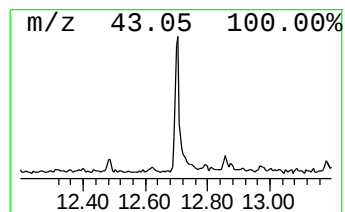
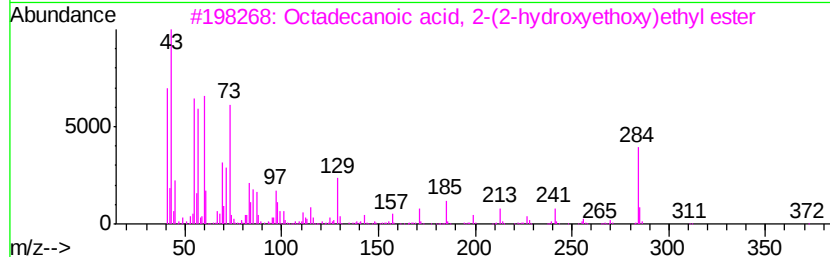
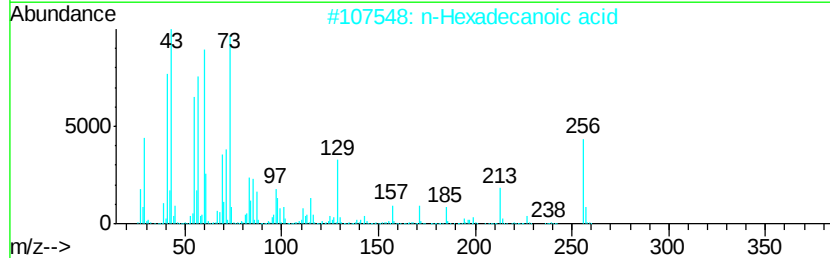
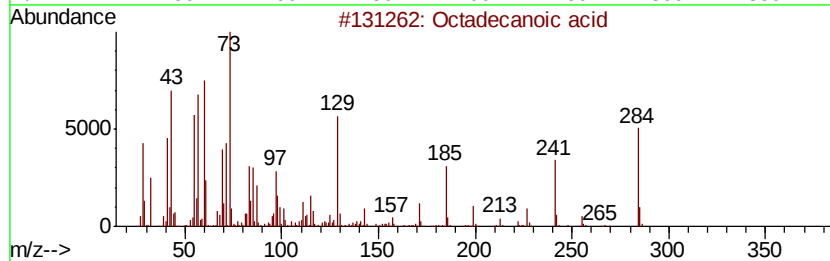
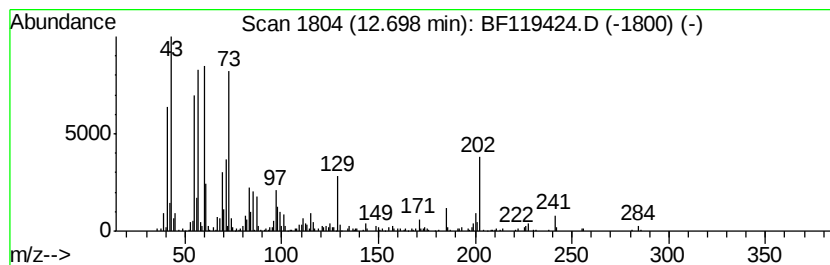
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	3.61 ng	560573	Phenanthrene-d10	11.39

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	98
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
3		Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	87
4		Tridecanoic acid	214	C13H26O2	000638-53-9	81
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031820\
 Data File : BF119424.D
 Acq On : 18 Mar 2020 19:30
 Operator : CG/JU
 Sample : L1942-13
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3-SA

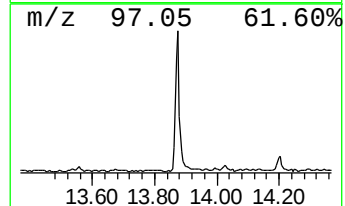
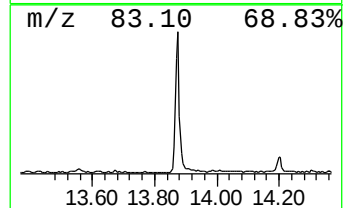
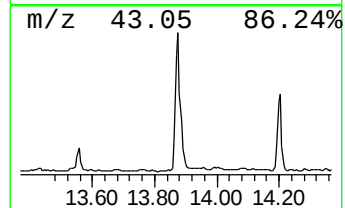
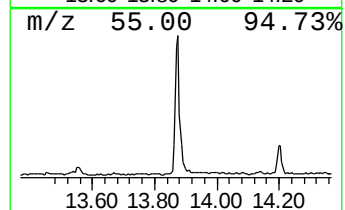
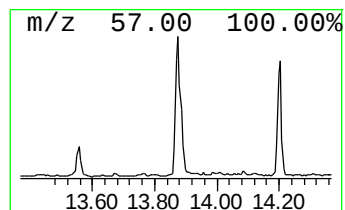
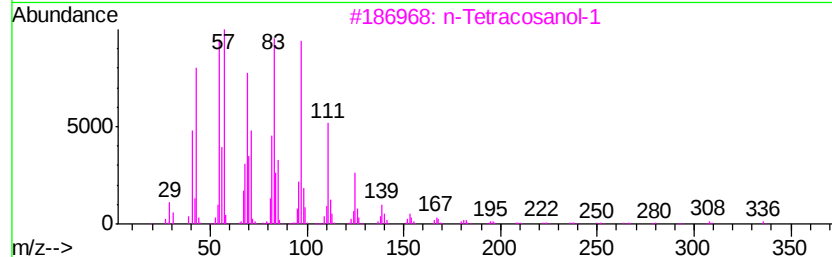
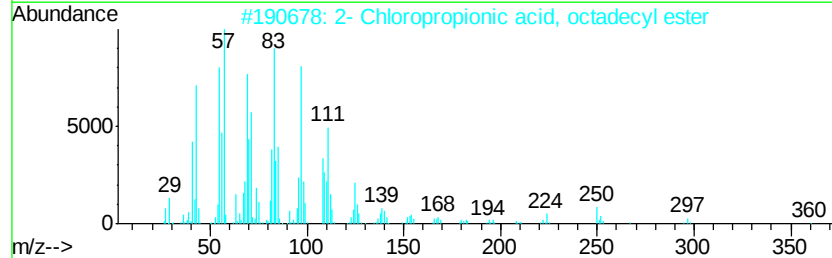
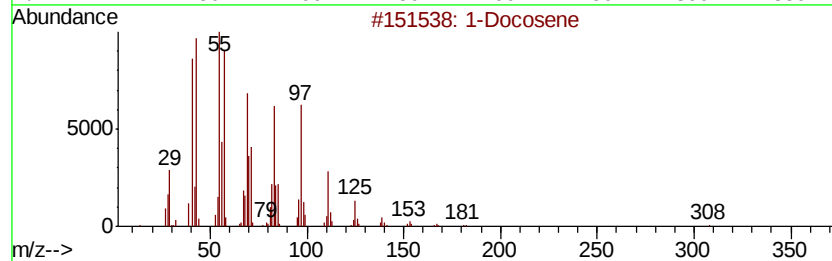
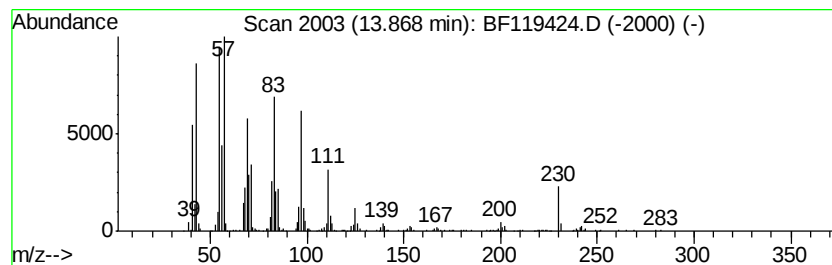
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 6 1-Docosene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	6.85 ng	1018120	Chrysene-d12	14.03

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene		308	C22H44	001599-67-3	94
2	2- Chloropropionic acid, octadec...		360	C21H41ClO2	088104-31-8	91
3	n-Tetracosanol-1		354	C24H50O	000506-51-4	91
4	Behenic alcohol		326	C22H46O	000661-19-8	91
5	1-Octadecanol		270	C18H38O	000112-92-5	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031820\
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Misc :
ALS Vial : 9 Sample Multiplier: 1

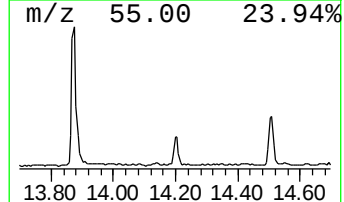
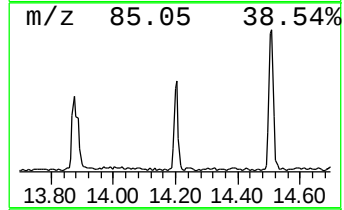
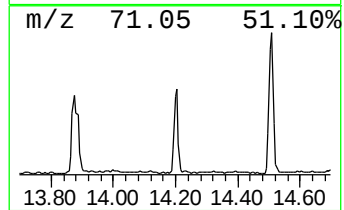
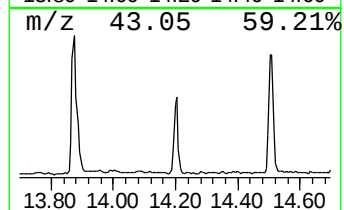
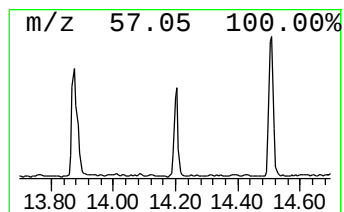
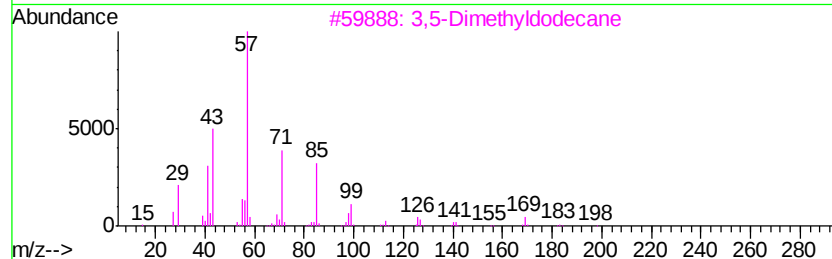
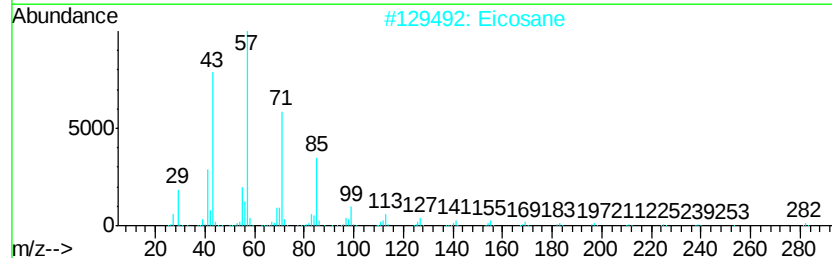
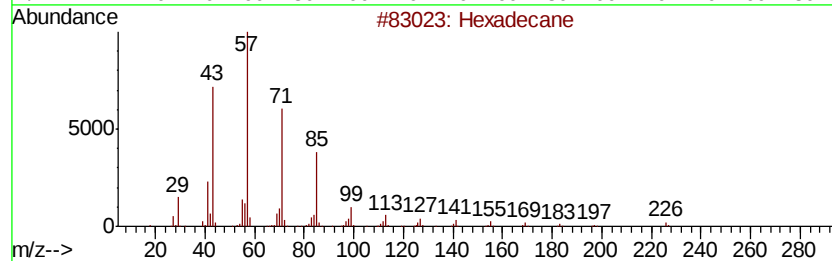
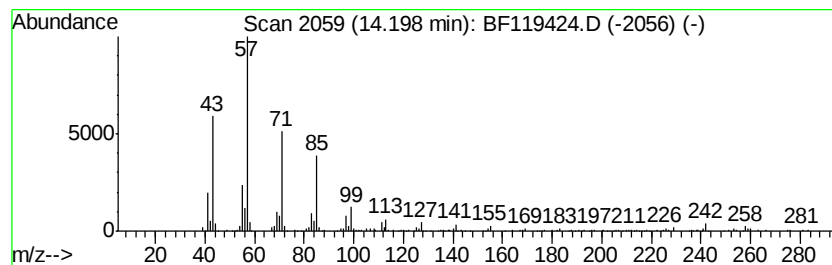
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.
14.20	2.56 ng	380176	Chrysene-d12		14.03
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	97
2	Eicosane	282	C20H42	000112-95-8	76
3	3,5-Dimethyldodecane	198	C14H30	107770-99-0	68
4	Pentadecane	212	C15H32	000629-62-9	64
5	Undecane, 3,8-dimethyl-	184	C13H28	017301-30-3	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031820\
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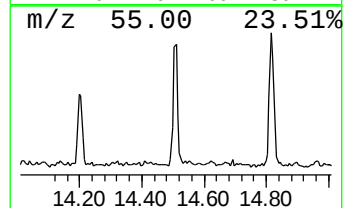
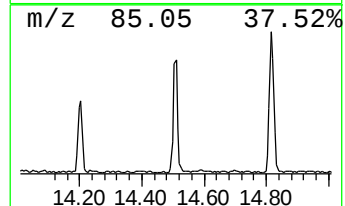
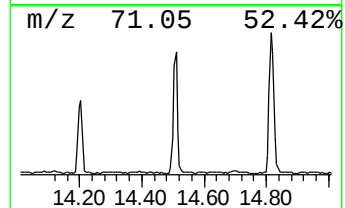
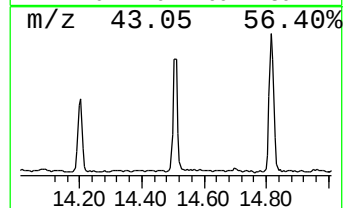
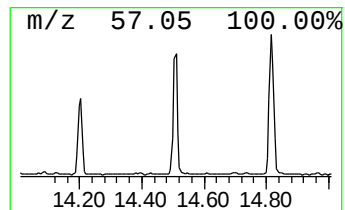
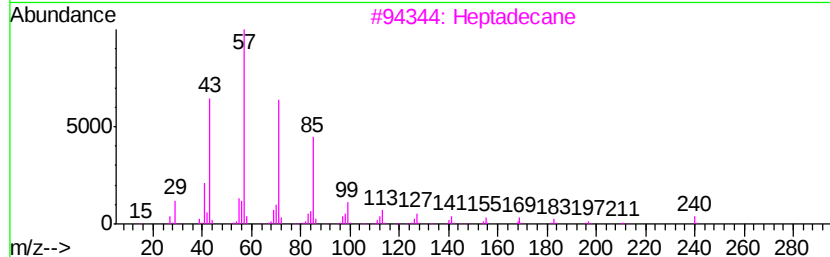
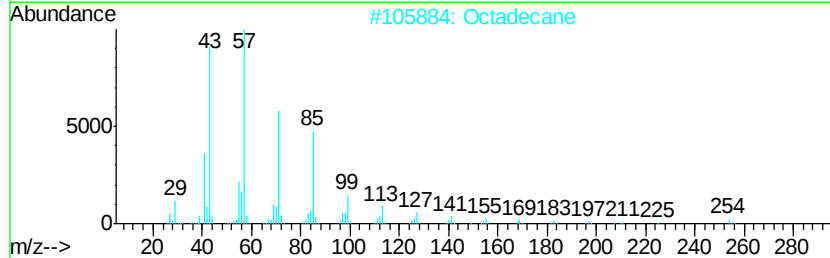
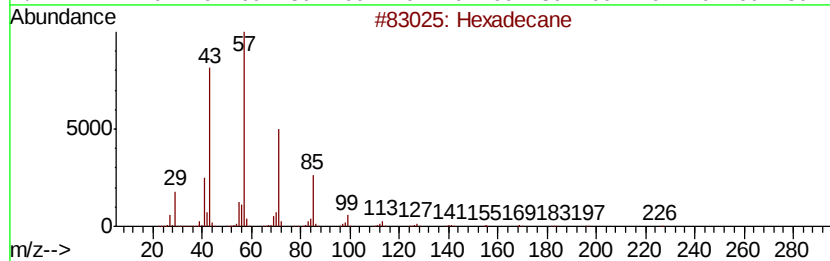
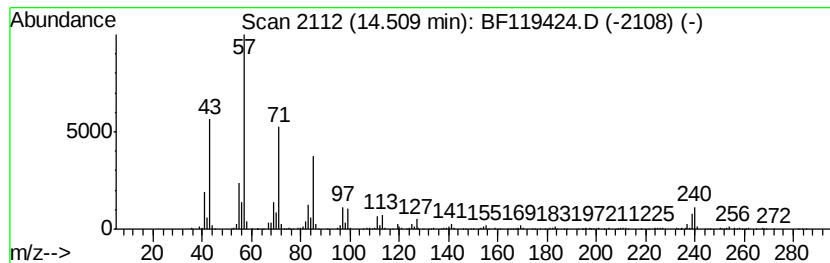
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.51	4.22 ng	627830	Chrysene-d12	14.03	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	93
2	Octadecane	254	C18H38	000593-45-3	87
3	Heptadecane	240	C17H36	000629-78-7	83
4	Heneicosane	296	C21H44	000629-94-7	81
5	Tetratetracontane	619	C44H90	007098-22-8	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031820\
 Data File : BF119424.D
 Acq On : 18 Mar 2020 19:30
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Instrument :
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 ClientSampleId :
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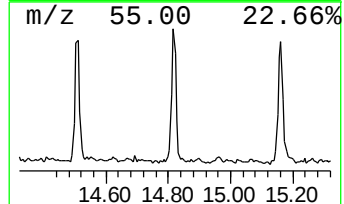
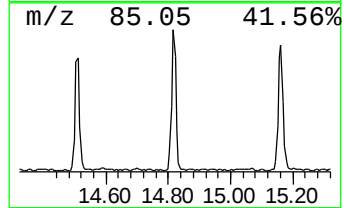
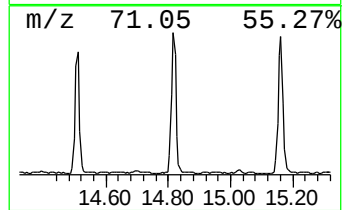
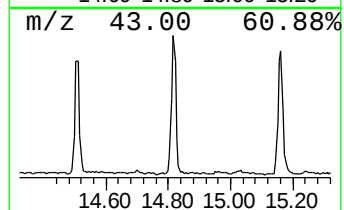
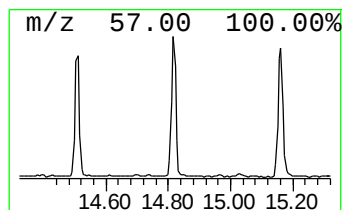
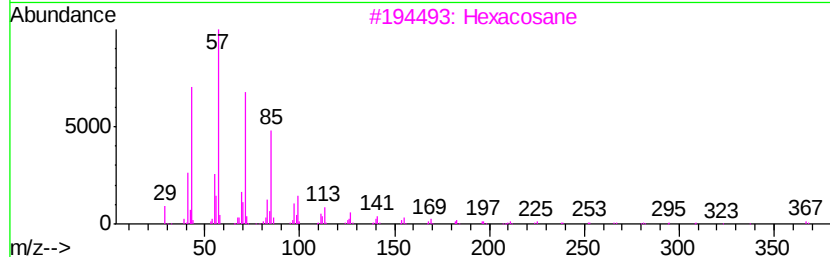
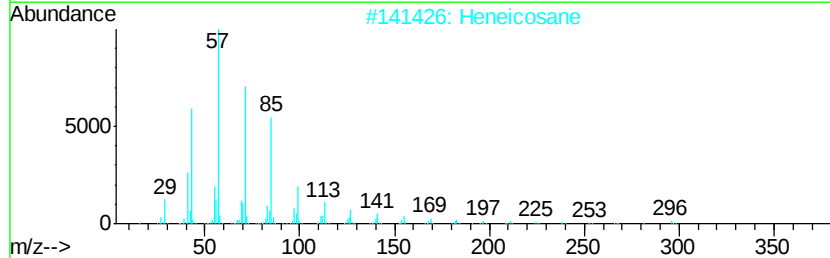
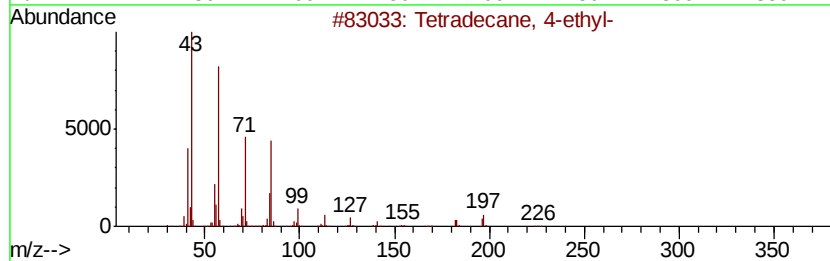
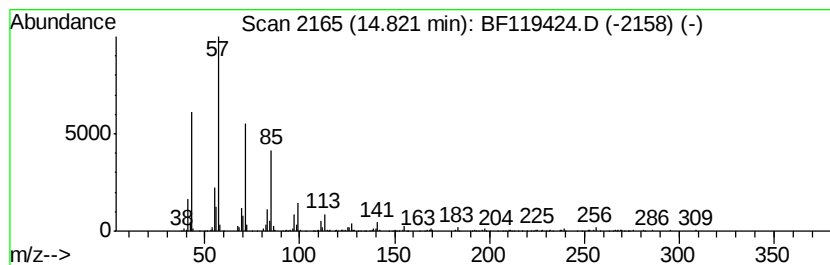
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 Tetradecane, 4-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	8.05 ng	746006	Perylene-d12	15.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane, 4-ethyl-	226	C16H34	055045-14-2	91
2		Heneicosane	296	C21H44	000629-94-7	90
3		Hexacosane	366	C26H54	000630-01-3	90
4		Tetracosane	338	C24H50	000646-31-1	90
5		Hexadecane	226	C16H34	000544-76-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031820\
 Data File : BF119424.D
 Acq On : 18 Mar 2020 19:30
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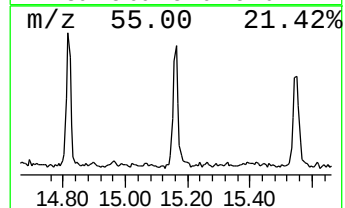
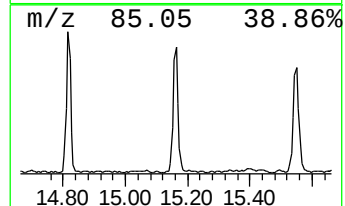
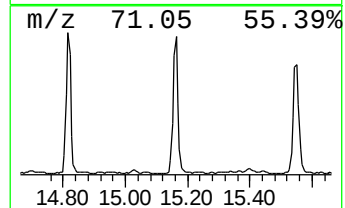
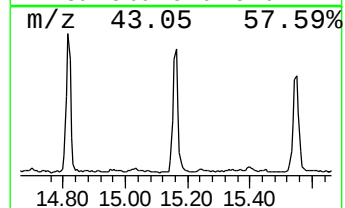
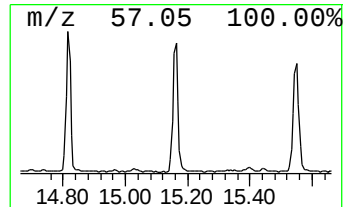
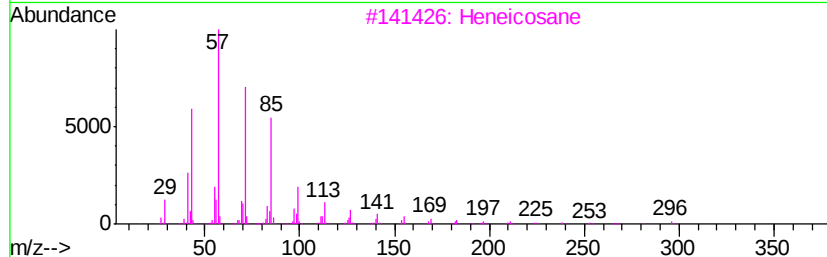
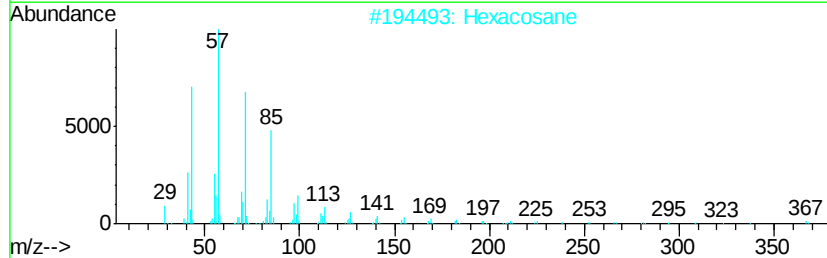
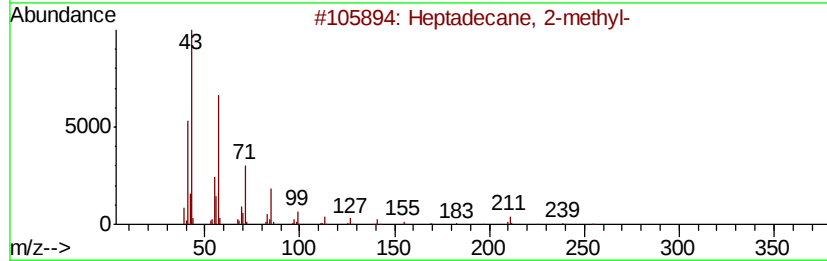
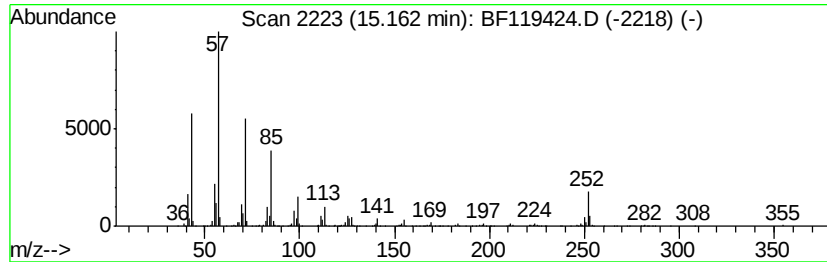
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 10 Heptadecane, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.16	9.46 ng	876114	Perylene-d12	15.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 2-methyl-	254	C18H38	001560-89-0	81
2		Hexacosane	366	C26H54	000630-01-3	81
3		Heneicosane	296	C21H44	000629-94-7	76
4		Heptacosane	380	C27H56	000593-49-7	76
5		2-methyloctacosane	408	C29H60	1000376-72-8	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031820\
Data File : BF119424.D
Acq On : 18 Mar 2020 19:30
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Misc :
ALS Vial : 9 Sample Multiplier: 1

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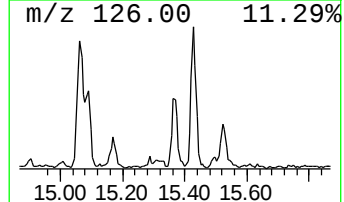
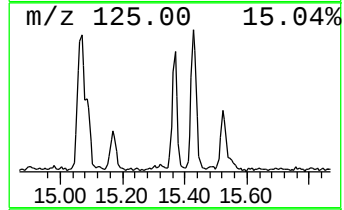
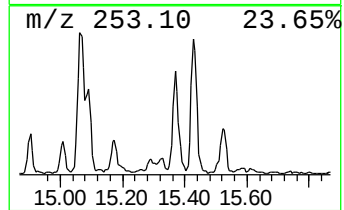
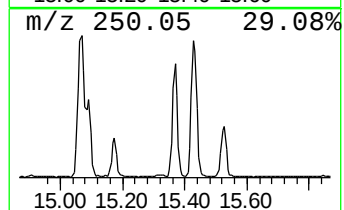
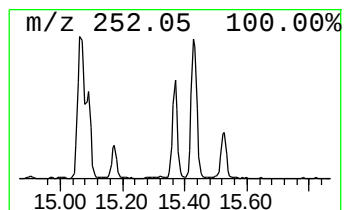
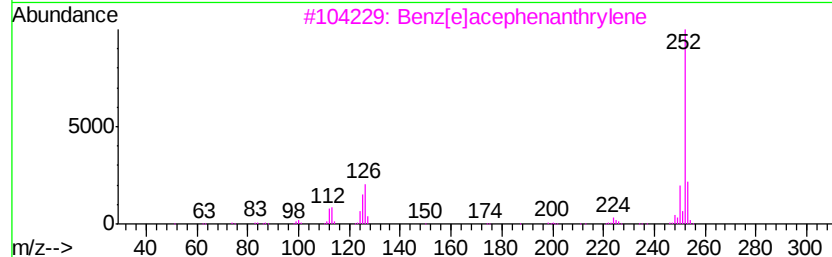
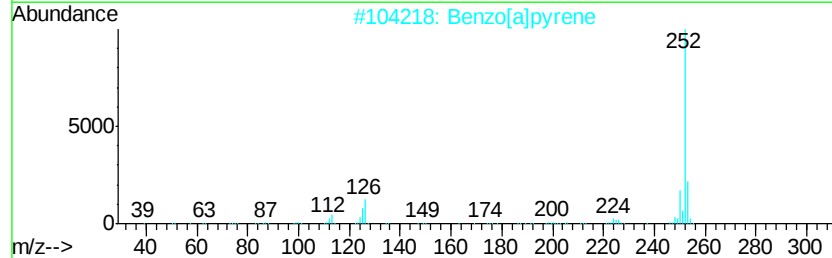
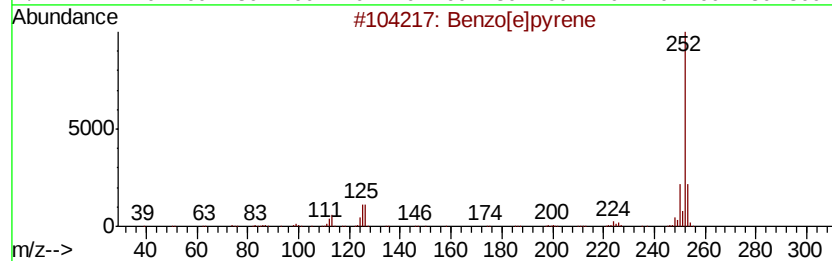
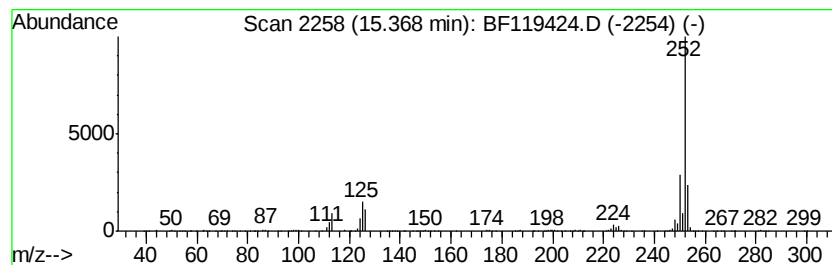
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Benzo[e]pyrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.37	6.11 ng	566310	Perylene-d12	15.50

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[a]pyrene	252	C20H12	000050-32-8	97
3			Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	95
4			Benzo[i]lfluoranthene	252	C20H12	000205-82-3	94
5			Perylene	252	C20H12	000198-55-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031820\
Data File : BF119424.D
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Sample : L1942-13
Misc :
ALS Vial : 9 Sample Multiplier: 1

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ClientSampleId :
TP-3-SA

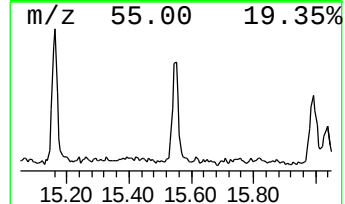
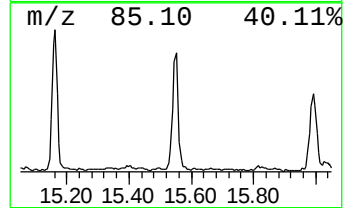
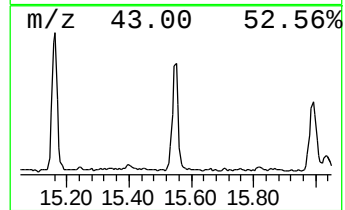
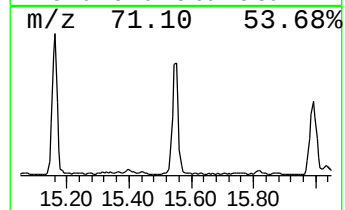
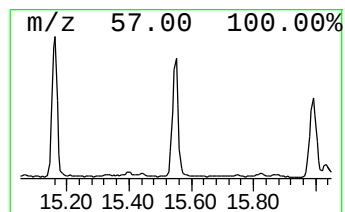
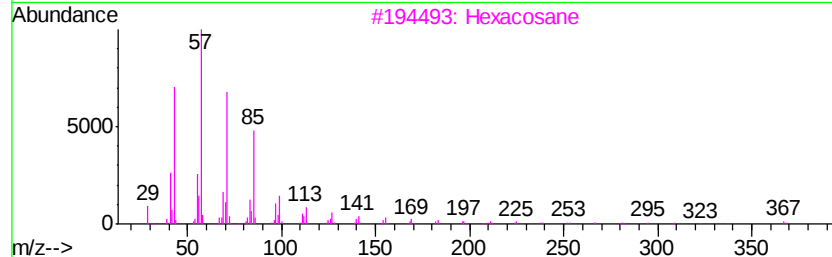
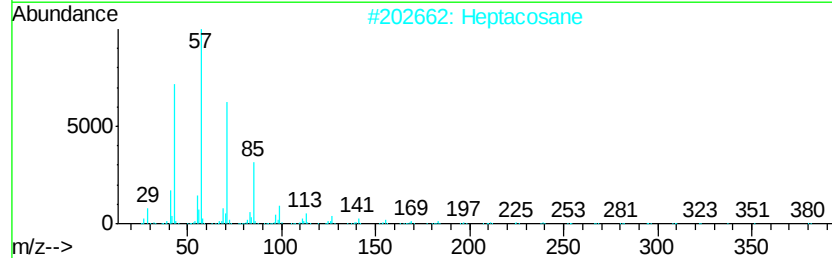
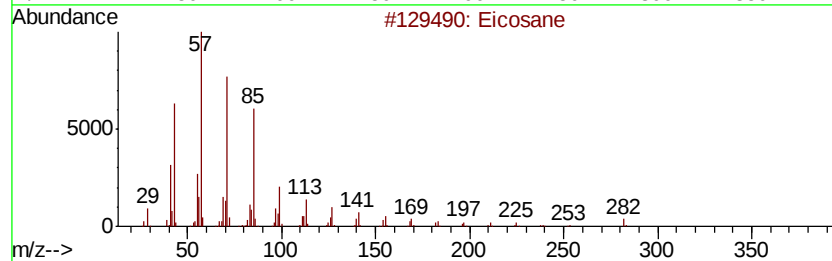
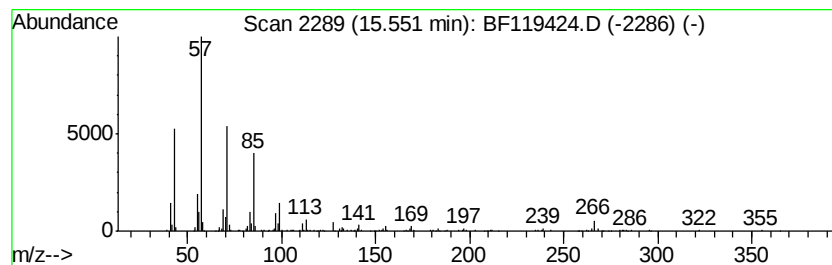
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Eicosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.55	6.93 ng	642322	Perylene-d12	15.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	93
2		Heptacosane	380	C27H56	000593-49-7	87
3		Hexacosane	366	C26H54	000630-01-3	86
4		1-Iodoundecane	282	C11H23I	004282-44-4	81
5		Heptadecane	240	C17H36	000629-78-7	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031820\
Data File : BF119424.D
Acq On : 18 Mar 2020 19:30
Operator : CG/JU
Sample : L1942-13
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-3-SA

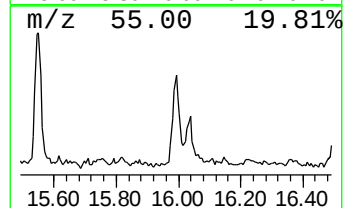
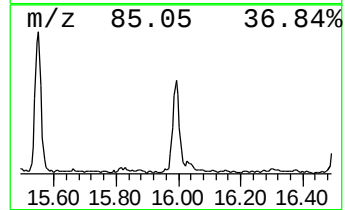
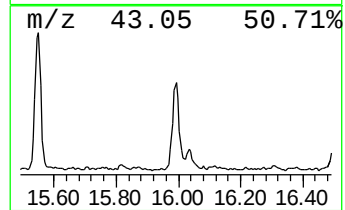
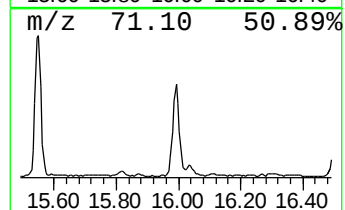
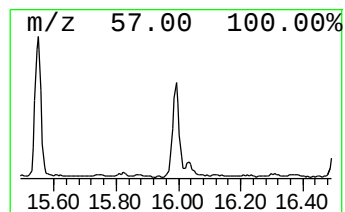
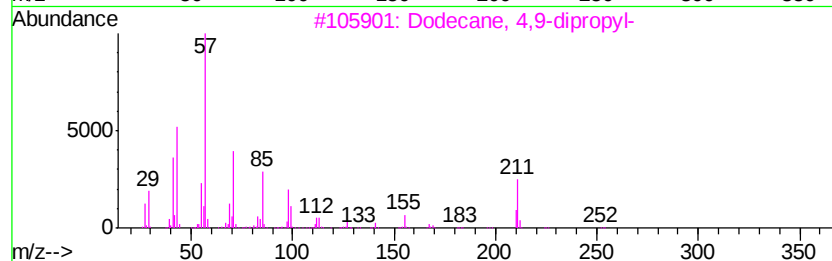
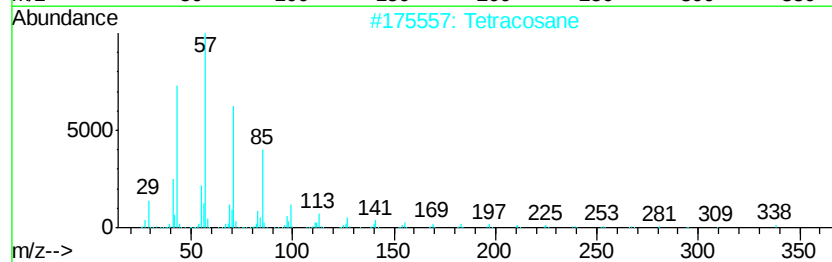
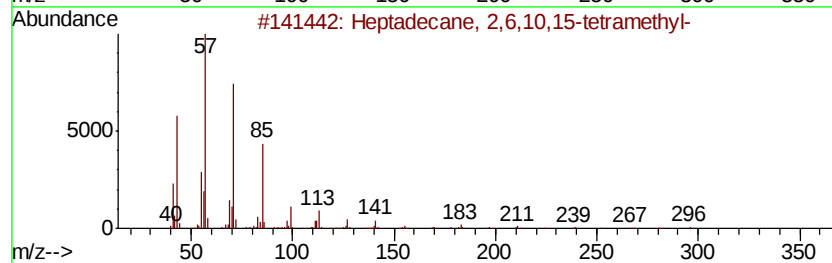
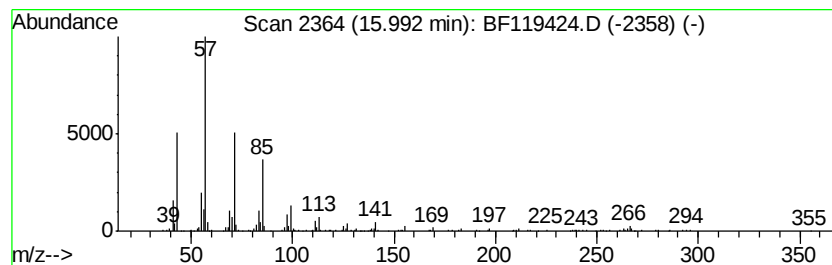
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heptadecane, 2,6,10,15-tetr... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.99	5.76 ng	533758	Perylene-d12	15.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
2		Tetracosane	338	C24H50	000646-31-1	90
3		Dodecane, 4,9-dipropyl-	254	C18H38	003054-63-5	90
4		Hexadecane, 2-methyl-	240	C17H36	001560-92-5	90
5		Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031820\
Data File : BF119424.D
Acq On : 18 Mar 2020 19:30
Operator : CG/JU
Sample : L1942-13
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-3-SA

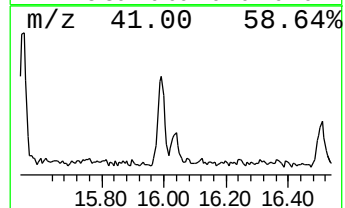
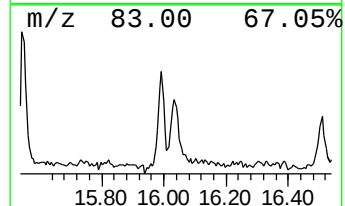
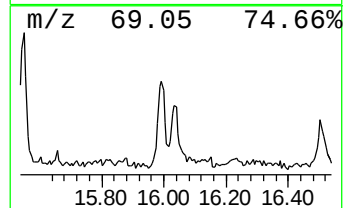
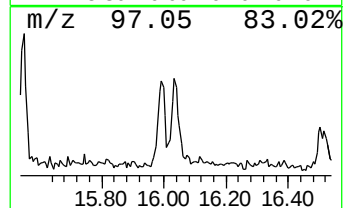
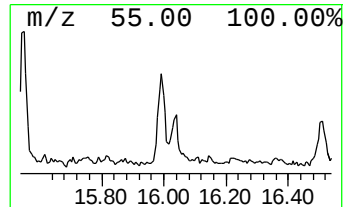
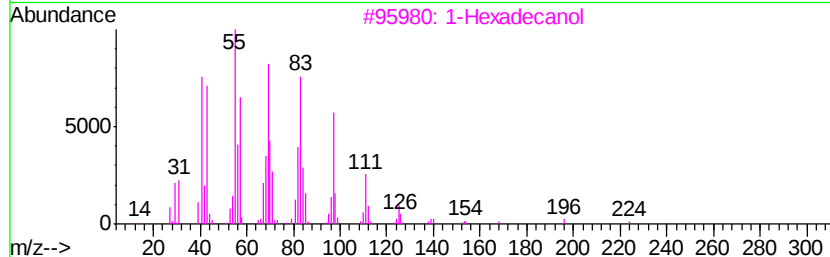
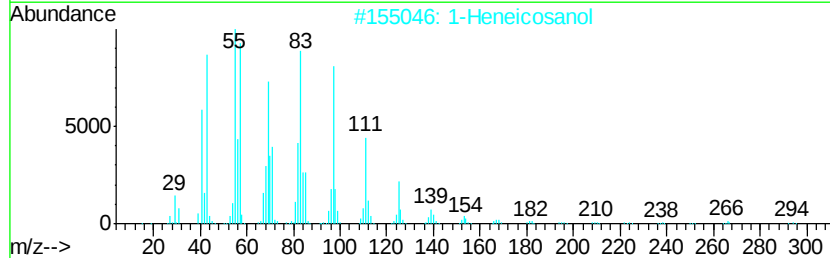
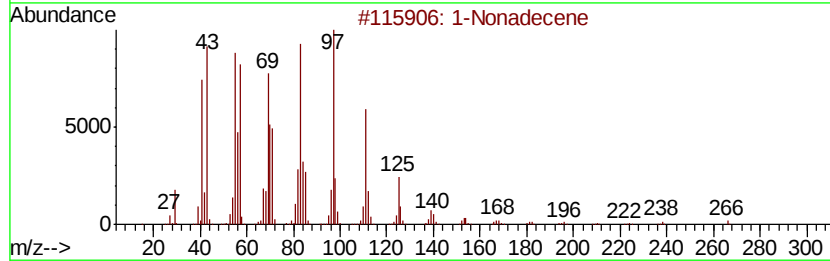
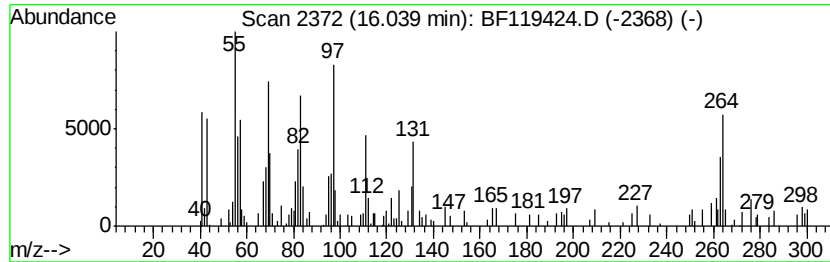
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown16.04 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.04	2.51 ng	232866	Perylene-d12	15.50

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Nonadecene	266	C19H38	018435-45-5	49
2			1-Heneicosanol	312	C21H44O	015594-90-8	49
3			1-Hexadecanol	242	C16H34O	036653-82-4	46
4			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	46
5			Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	45



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031820\
Data File : BF119424.D
Acq On : 18 Mar 2020 19:30
Operator : CG/JU
Sample : L1942-13
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-3-SA

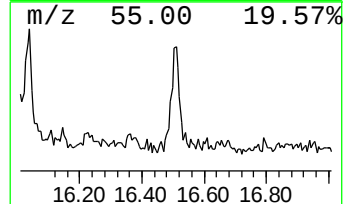
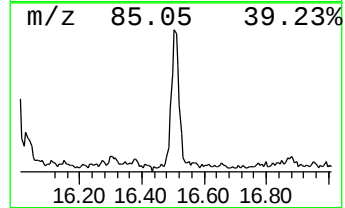
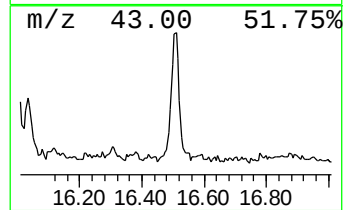
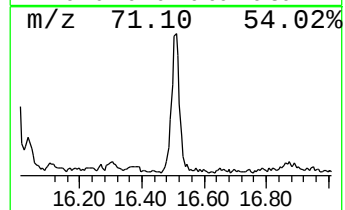
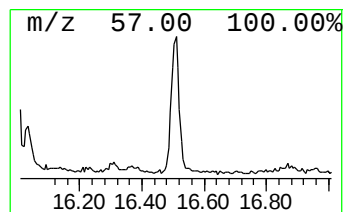
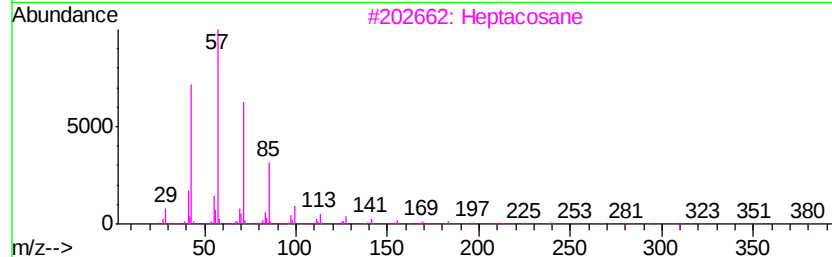
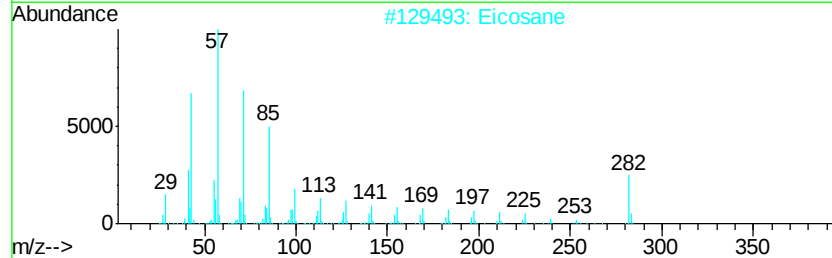
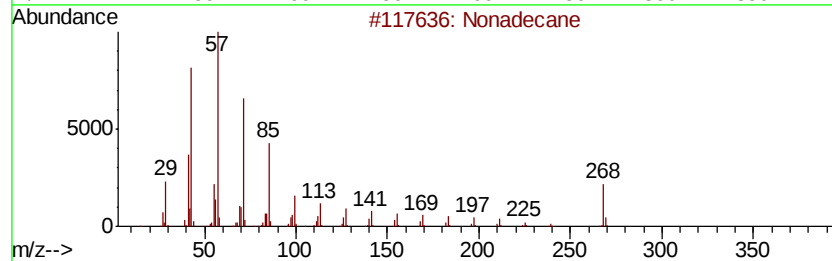
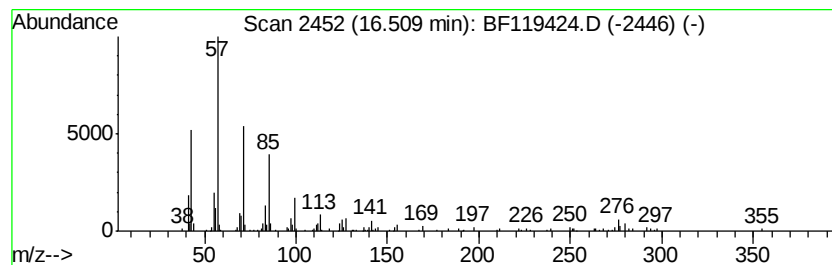
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Nonadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.51	3.26 ng	301910	Perylene-d12	15.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	87
2		Eicosane	282	C20H42	000112-95-8	83
3		Heptacosane	380	C27H56	000593-49-7	83
4		Octacosane	394	C28H58	000630-02-4	81
5		Hexadecane	226	C16H34	000544-76-3	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031820\
Data File : BF119424.D
Acq On : 18 Mar 2020 19:30
Operator : CG/JU
Sample : L1942-13
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-3-SA

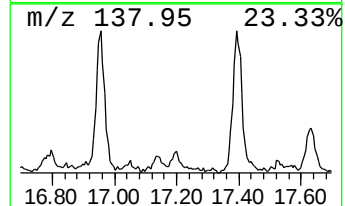
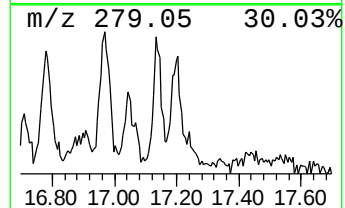
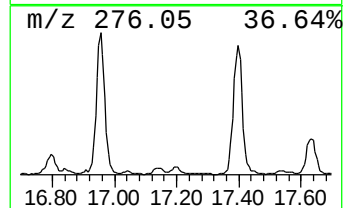
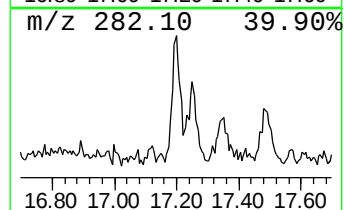
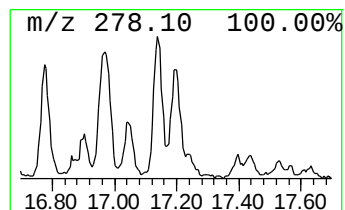
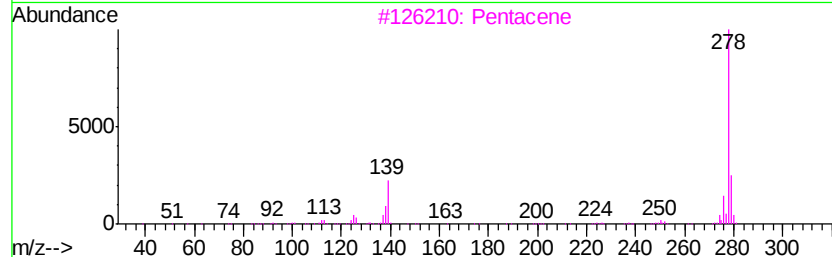
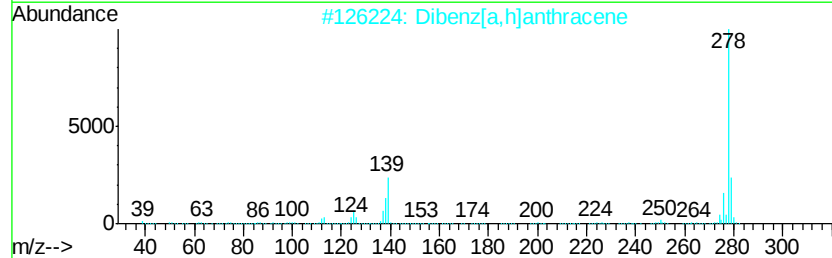
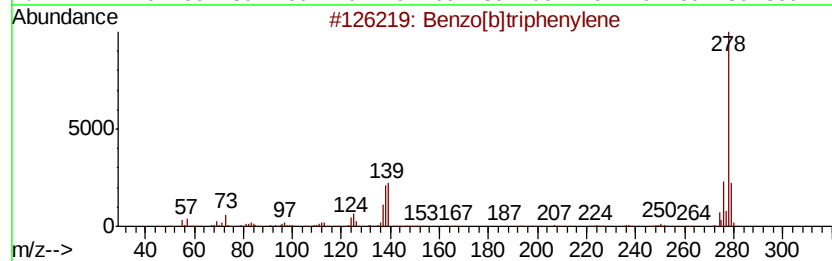
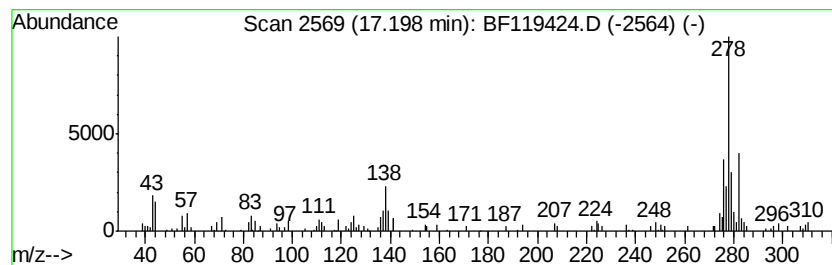
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Benzo[b]triphenylene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.20	2.49 ng	231100	Perylene-d12	15.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]triphenylene	278	C22H14	000215-58-7	55
2		Dibenz[a,h]anthracene	278	C22H14	000053-70-3	49
3		Pentacene	278	C22H14	000135-48-8	46
4		Benzo[a]naphthacene	278	C22H14	000226-88-0	43
5		1-Butaneboronic acid, cyclic dip...	278	C18H19B02	024371-99-1	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF031820\
 Data File : BF119424.D
 Acq On : 18 Mar 2020 19:30
 Operator : CG/JU
 Sample : L1942-13
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3-SA

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.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
Butane, 2-methoxy...	2.16	4.3	ng	468153	1	6.86	2177790	20.0
n-Hexadecanoic acid	11.92	6.2	ng	959742	4	11.39	3106420	20.0
4H-Cyclopenta[def...	12.00	2.8	ng	431476	4	11.39	3106420	20.0
Octadecanoic acid	12.70	3.6	ng	560573	4	11.39	3106420	20.0
1-Docosene	13.87	6.8	ng	1018120	5	14.03	2974420	20.0
Hexadecane	14.20	2.6	ng	380176	5	14.03	2974420	20.0
Octadecane	14.51	4.2	ng	627830	5	14.03	2974420	20.0
Tetradecane, 4-et...	14.82	8.1	ng	746006	6	15.50	1853120	20.0
Heptadecane, 2-me...	15.16	9.5	ng	876114	6	15.50	1853120	20.0
Benzo[e]pyrene	15.37	6.1	ng	566310	6	15.50	1853120	20.0
Eicosane	15.55	6.9	ng	642322	6	15.50	1853120	20.0
Heptadecane, 2,6,...	15.99	5.8	ng	533758	6	15.50	1853120	20.0
unknown16.04	16.04	2.5	ng	232866	6	15.50	1853120	20.0
Nonadecane	16.51	3.3	ng	301910	6	15.50	1853120	20.0
Benzo[b]triphenylene	17.20	2.5	ng	231100	6	15.50	1853120	20.0