

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF060320\
 Data File : BF120522.D
 Acq On : 3 Jun 2020 21:05
 Operator : JU/CG
 Sample : L2839-12
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM6-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	1.952	9	11	18	rBV	603814	701035	21.60%	1.905%
2	2.081	28	33	40	rVB	761280	976282	30.08%	2.653%
3	5.457	602	607	610	rBV	2543618	2318996	71.45%	6.302%
4	6.475	775	780	783	rBV	2425739	2438932	75.15%	6.628%
5	6.663	809	812	818	rBV	63470	60927	1.88%	0.166%
6	6.834	837	841	844	rBV	1040841	800200	24.66%	2.175%
7	6.992	863	868	871	rBV	2666179	2279151	70.22%	6.194%
8	7.398	931	937	940	rBV	1867343	1736731	53.51%	4.720%
9	8.116	1055	1059	1062	rBV	1277346	1007572	31.05%	2.738%
10	9.192	1237	1242	1245	rBV	3769253	3245516	100.00%	8.820%
11	9.586	1305	1309	1312	rVB	572317	505073	15.56%	1.373%
12	9.727	1330	1333	1339	rBV	52878	50177	1.55%	0.136%
13	9.869	1351	1357	1361	rBV	1248582	1131350	34.86%	3.075%
14	9.904	1361	1363	1367	rVB	62469	46822	1.44%	0.127%
15	10.075	1388	1392	1396	rBV2	54617	50194	1.55%	0.136%
16	10.416	1447	1450	1454	rBV	57441	48429	1.49%	0.132%
17	10.663	1487	1492	1495	rBV	1955833	1867963	57.56%	5.077%
18	11.016	1546	1552	1556	rBV5	32671	57064	1.76%	0.155%
19	11.157	1573	1576	1580	rVB	45467	42642	1.31%	0.116%
20	11.216	1580	1586	1590	rBV5	22484	42483	1.31%	0.115%
21	11.304	1598	1601	1605	rBV	69800	62957	1.94%	0.171%
22	11.357	1605	1610	1612	rVV	1249859	1111525	34.25%	3.021%
23	11.380	1612	1614	1617	rVV	780753	572826	17.65%	1.557%
24	11.433	1620	1623	1625	rVB	151732	125730	3.87%	0.342%
25	11.586	1643	1649	1651	rBV2	74087	97692	3.01%	0.265%
26	11.816	1683	1688	1691	rBV2	154056	172601	5.32%	0.469%
27	11.857	1691	1695	1697	rVV2	139562	183311	5.65%	0.498%
28	11.886	1697	1700	1703	rVV	370715	317979	9.80%	0.864%
29	11.933	1706	1708	1711	rVV	56910	54346	1.67%	0.148%
30	11.969	1711	1714	1717	rVV	160374	179277	5.52%	0.487%
31	11.992	1717	1718	1721	rVB	72975	43283	1.33%	0.118%
32	12.151	1742	1745	1747	rBV	67039	59737	1.84%	0.162%
33	12.174	1747	1749	1753	rVB	90059	74097	2.28%	0.201%
34	12.304	1765	1771	1773	rBV3	34223	38671	1.19%	0.105%

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

35	12.410	1785	1789	1792	rBV	66721	72461	2.23%	0.197%
36	12.445	1792	1795	1797	rVV	47177	48259	1.49%	0.131%
37	12.492	1800	1803	1808	rVB2	86024	102988	3.17%	0.280%
38	12.551	1809	1813	1814	rVV	306551	256320	7.90%	0.697%
39	12.568	1814	1816	1820	rVB	1151064	1000008	30.81%	2.718%
40	12.663	1830	1832	1835	rBV	44812	42718	1.32%	0.116%
41	12.798	1849	1855	1861	rBV	957214	1043442	32.15%	2.836%
42	12.845	1861	1863	1869	rVB2	47898	58252	1.79%	0.158%
43	12.916	1872	1875	1876	rBV2	57039	47736	1.47%	0.130%
44	12.945	1876	1880	1883	rVV	3280352	2970013	91.51%	8.072%
45	13.015	1886	1892	1896	rBV2	73483	135065	4.16%	0.367%
46	13.104	1903	1907	1908	rBV4	52029	54274	1.67%	0.147%
47	13.127	1908	1911	1914	rVB	142184	134114	4.13%	0.364%
48	13.174	1914	1919	1920	rBV4	40802	65171	2.01%	0.177%
49	13.192	1920	1922	1925	rVB	88595	66629	2.05%	0.181%
50	13.227	1925	1928	1931	rBV2	104283	121404	3.74%	0.330%
51	13.321	1942	1944	1947	rVB	64501	49965	1.54%	0.136%
52	13.351	1947	1949	1953	rVB2	63375	60467	1.86%	0.164%
53	13.392	1953	1956	1959	rBV4	36118	44583	1.37%	0.121%
54	13.421	1959	1961	1967	rVB6	27993	39318	1.21%	0.107%
55	13.515	1972	1977	1981	rBV5	57130	118610	3.65%	0.322%
56	13.633	1994	1997	2001	rVB	101383	104054	3.21%	0.283%
57	13.745	2012	2016	2018	rBV2	95253	114070	3.51%	0.310%
58	13.774	2018	2021	2023	rVV	96930	101053	3.11%	0.275%
59	13.798	2023	2025	2029	rVV2	85066	132162	4.07%	0.359%
60	13.845	2029	2033	2040	rVB2	267388	387304	11.93%	1.053%
61	13.998	2051	2059	2061	rBV3	1264379	2029652	62.54%	5.516%
62	14.021	2061	2063	2066	rVB	535055	412577	12.71%	1.121%
63	14.104	2074	2077	2079	rBV	73547	60684	1.87%	0.165%
64	14.162	2084	2087	2089	rVV3	71277	76678	2.36%	0.208%
65	14.210	2093	2095	2100	rVB2	41004	62741	1.93%	0.171%
66	14.380	2122	2124	2127	rVB	94849	84144	2.59%	0.229%
67	14.415	2128	2130	2132	rVB	69329	55461	1.71%	0.151%
68	14.468	2136	2139	2144	rVV3	117392	199994	6.16%	0.544%
69	14.527	2147	2149	2153	rVB3	59786	56029	1.73%	0.152%
70	14.768	2187	2190	2193	rBV	114034	128275	3.95%	0.349%
71	14.845	2200	2203	2206	rBV2	130679	143921	4.43%	0.391%

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

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72	15.027	2230	2234	2238	rBV	436531	704025	21.69%	1.913%
73	15.104	2244	2247	2250	rBV	157383	169781	5.23%	0.461%
74	15.139	2250	2253	2259	rVB	96599	148197	4.57%	0.403%
75	15.327	2282	2285	2291	rVB	261458	347626	10.71%	0.945%
76	15.392	2292	2296	2301	rVB	367606	436145	13.44%	1.185%
77	15.457	2302	2307	2310	rVV	625241	857567	26.42%	2.331%
78	15.480	2310	2311	2317	rVB2	188760	178265	5.49%	0.484%
79	15.915	2381	2385	2390	rVB	180358	259013	7.98%	0.704%
80	16.898	2547	2552	2560	rBV2	170802	323232	9.96%	0.878%
81	17.333	2622	2626	2633	rBV	112941	189994	5.85%	0.516%

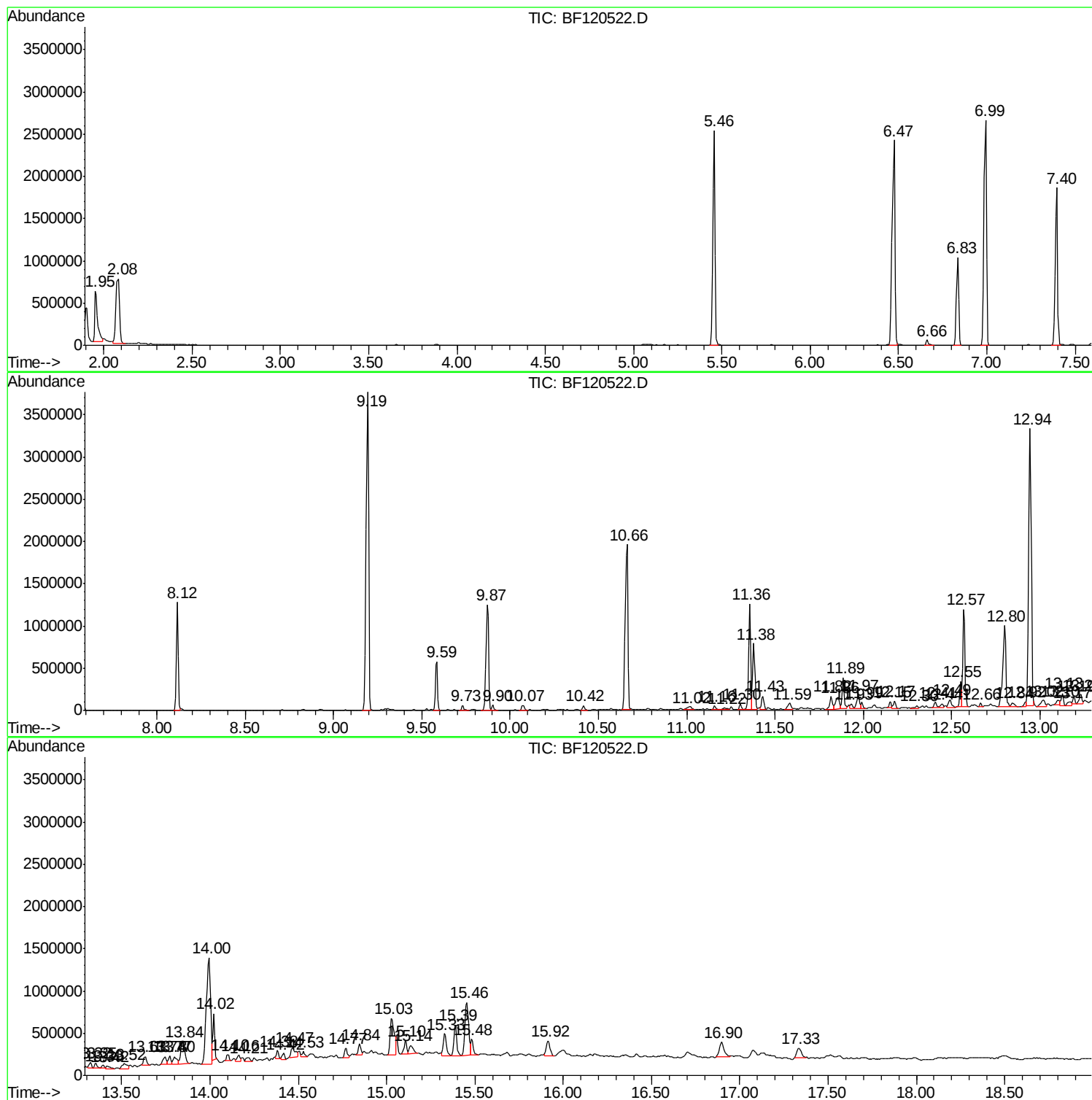
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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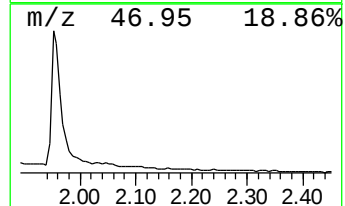
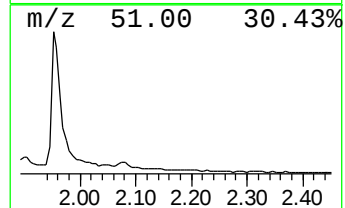
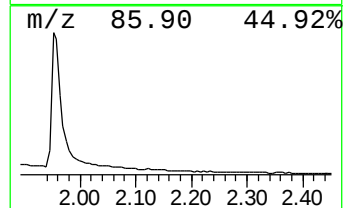
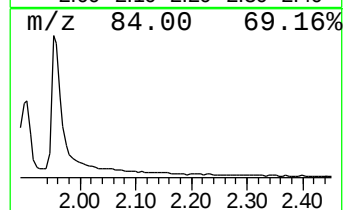
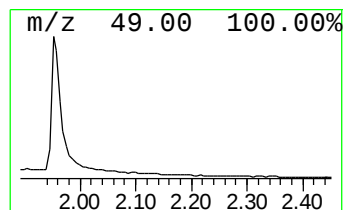
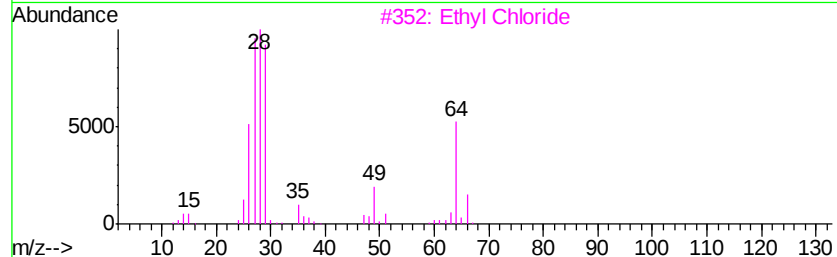
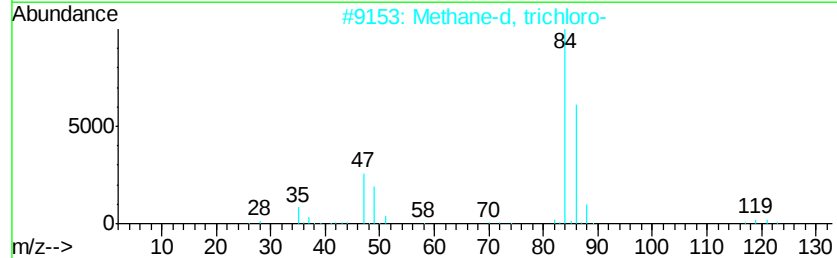
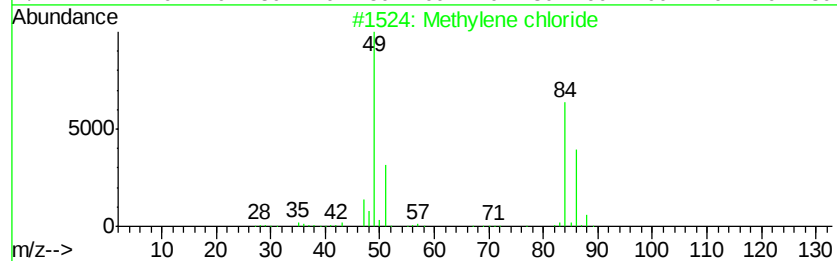
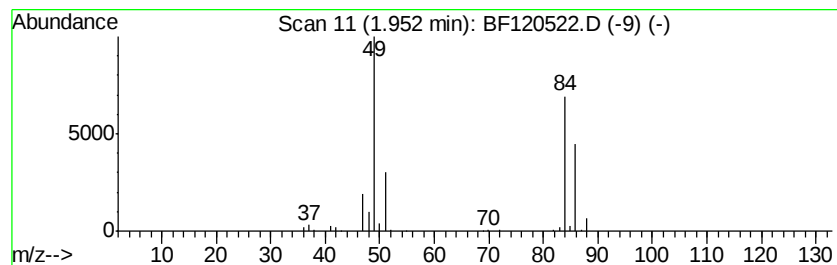
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 1 Methylene chloride Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.95	17.52 ng	701035	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methylene chloride	84	CH2Cl2	000075-09-2	95
2		Methane-d, trichloro-	119	CDCl3	000865-49-6	10
3		Ethyl Chloride	64	C2H5Cl	000075-00-3	4
4		Acetic acid, chloro-, ethyl ester	122	C4H7ClO2	000105-39-5	2
5		2-Chloroethanol	80	C2H5ClO	000107-07-3	1



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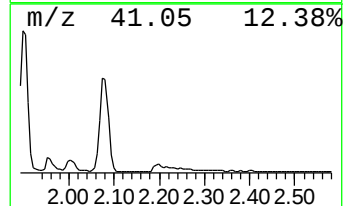
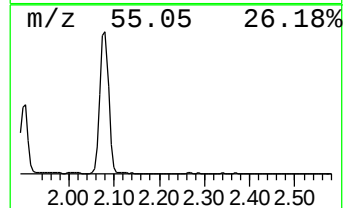
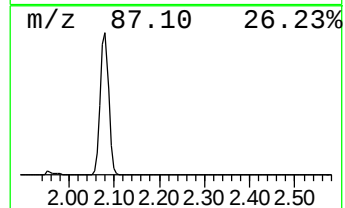
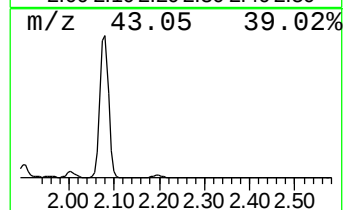
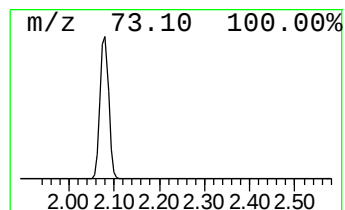
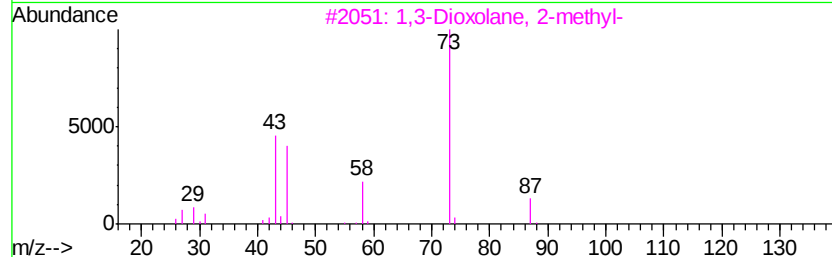
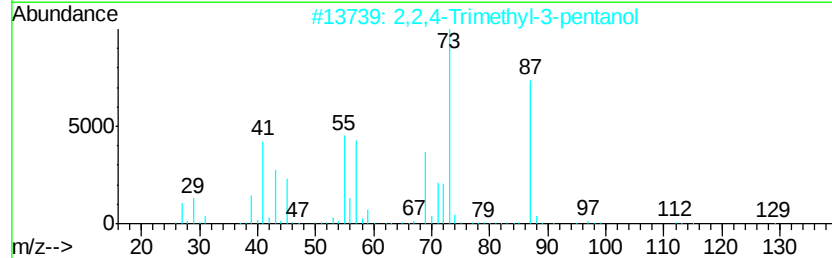
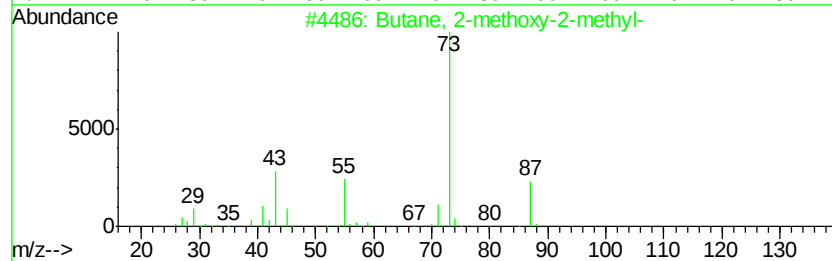
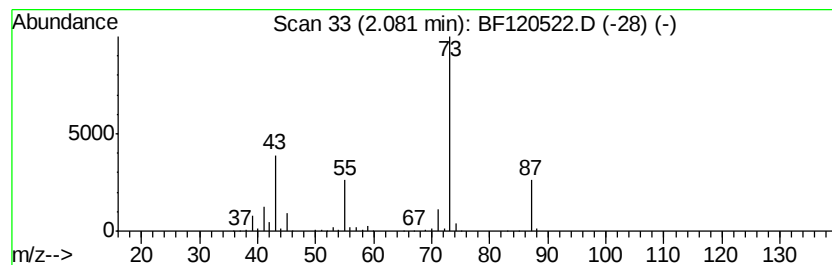
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
2.08	24.40 ng	976282	1,4-Dichlorobenzene-d4	6.83		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		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		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	10



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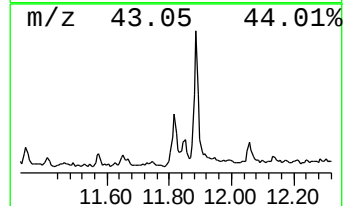
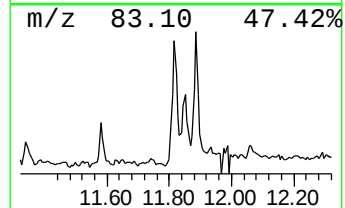
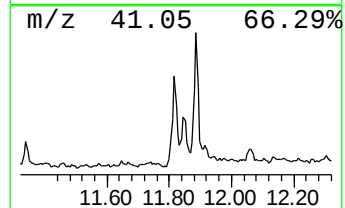
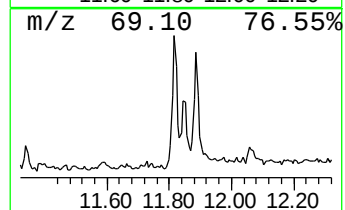
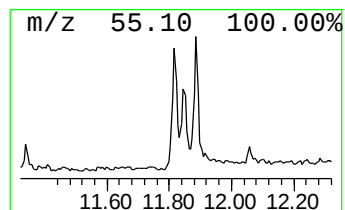
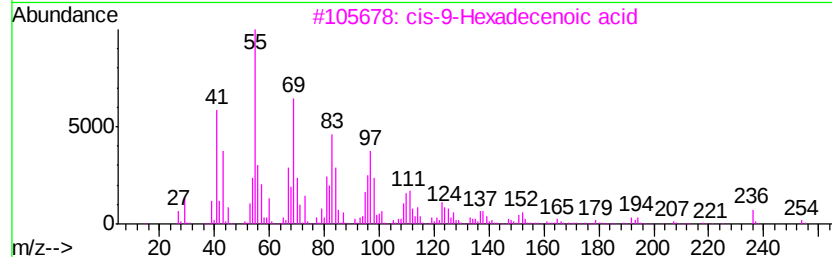
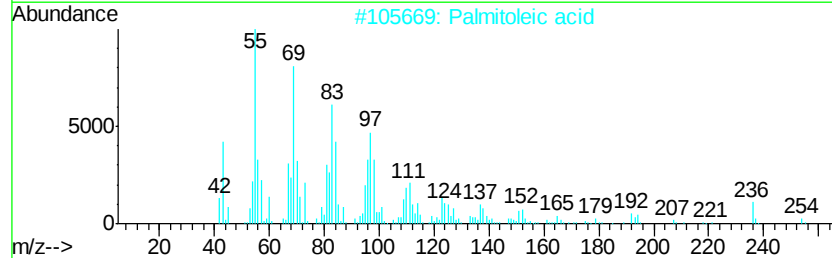
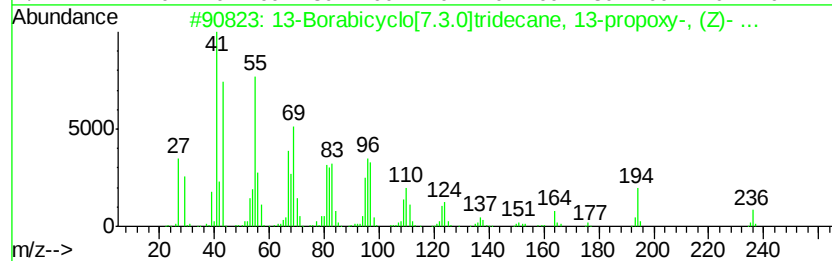
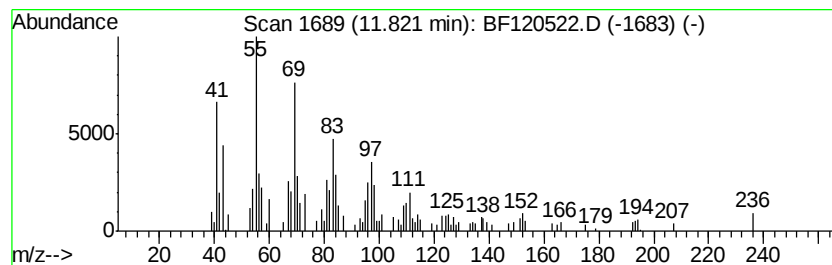
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 13-Borabicyclo[7.3.0]tridec... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.82	3.11 ng	172601	Phenanthrene-d10	11.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	13-Borabicyclo[7.3.0]tridecane, ...	236	C15H29BO	1000156-41-7	93
2	Palmitoleic acid	254	C16H30O2	000373-49-9	87
3	cis-9-Hexadecenoic acid	254	C16H30O2	1000333-19-5	81
4	Hexadecenoic acid, Z-11-	254	C16H30O2	002416-20-8	74
5	Oxacyclotetradecane-2,11-dione, ...	240	C14H24O3	074685-36-2	64



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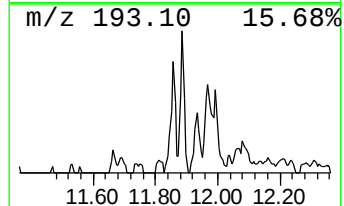
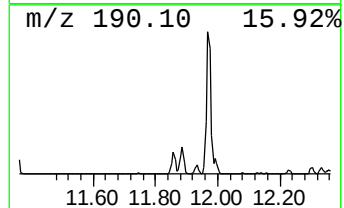
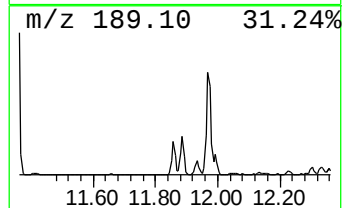
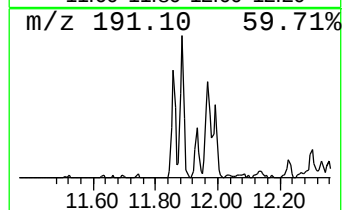
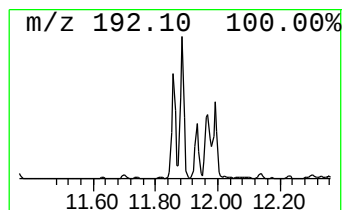
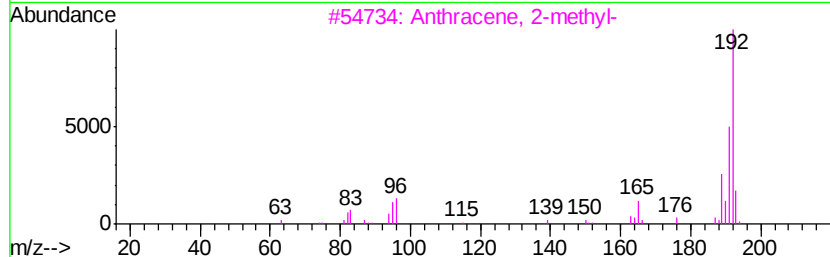
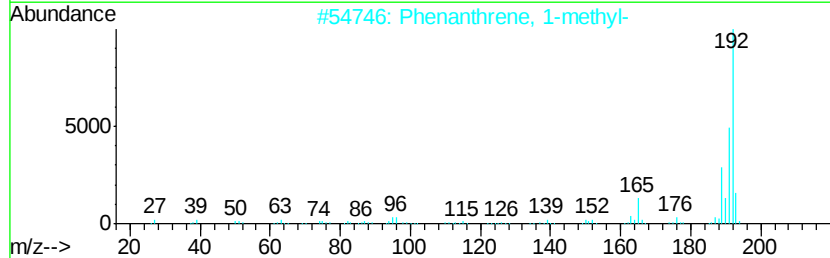
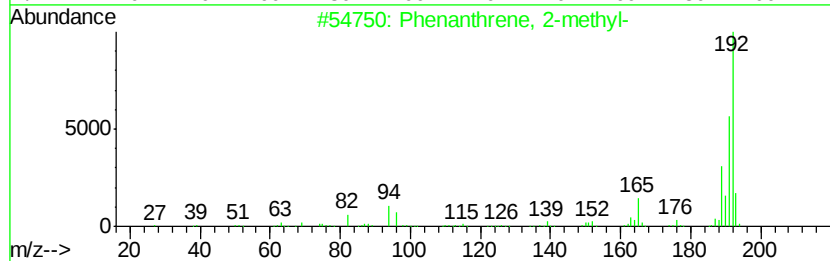
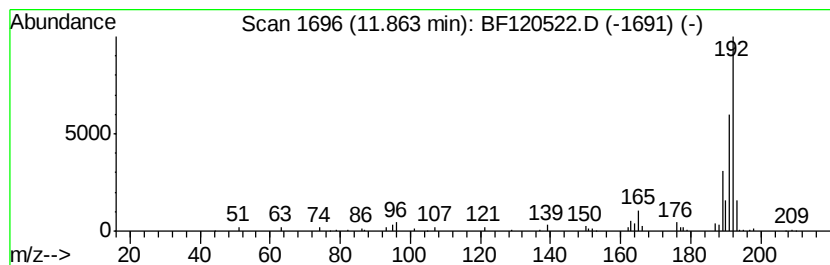
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ClientSampleId :
NM6-COMP-2020-0-6

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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 5 Phenanthrene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.86	3.30 ng	183311	Phenanthrene-d10	11.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060320\
Data File : BF120522.D
Acq On : 3 Jun 2020 21:05
Operator : JU/CG
Sample : L2839-12
Misc :
ALS Vial : 19 Sample Multiplier: 1

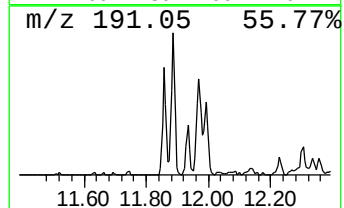
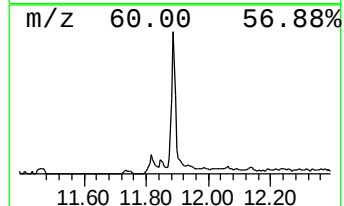
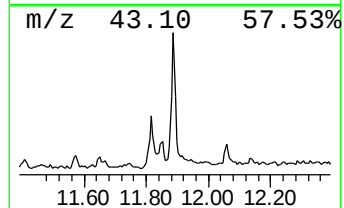
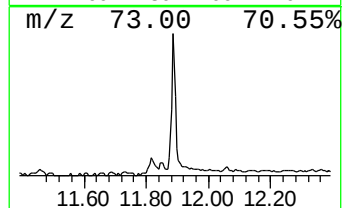
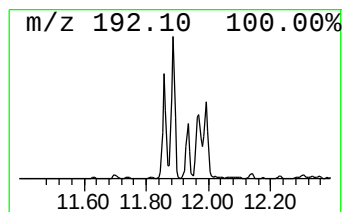
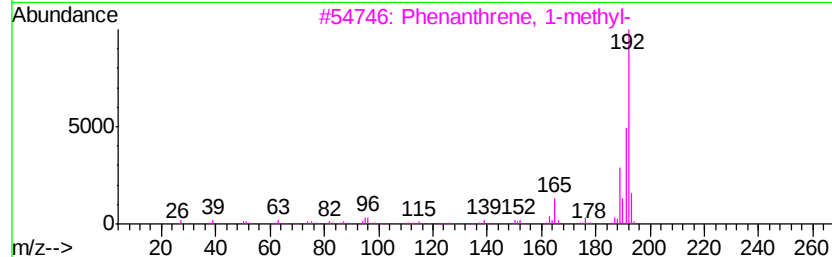
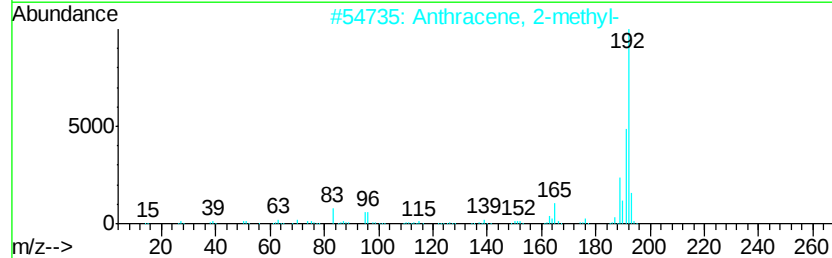
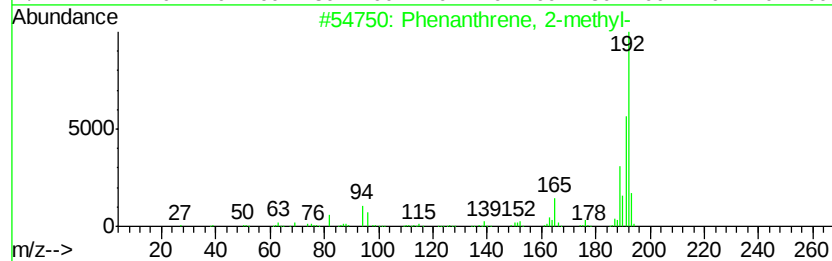
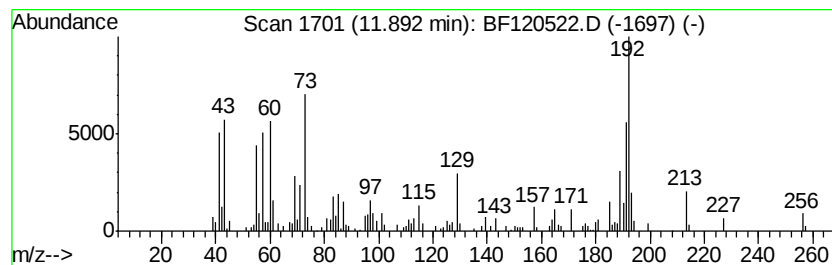
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.89	5.72 ng	317979	Phenanthrene-d10	11.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
4	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	92
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060320\
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Sample : L2839-12
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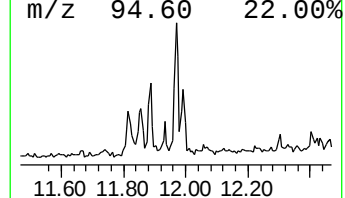
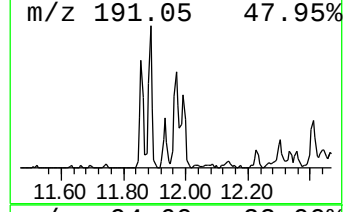
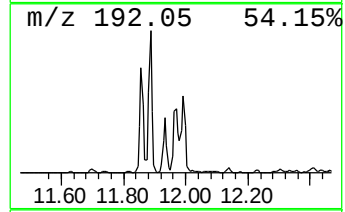
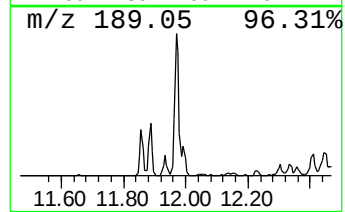
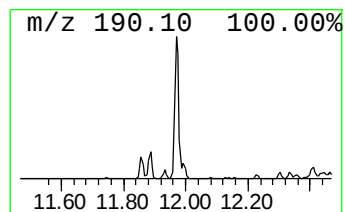
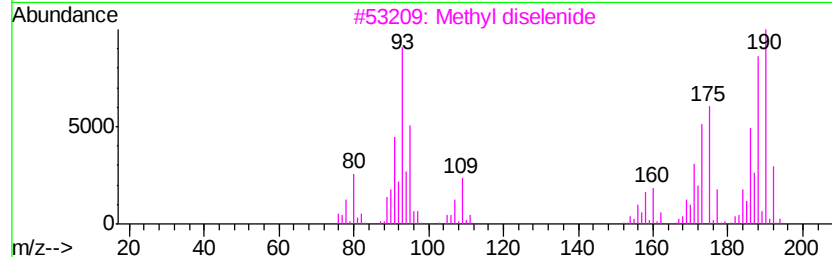
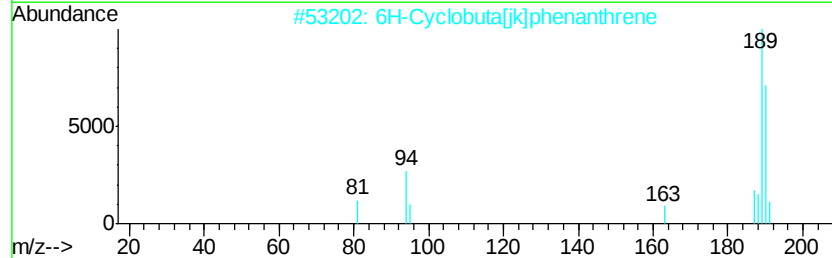
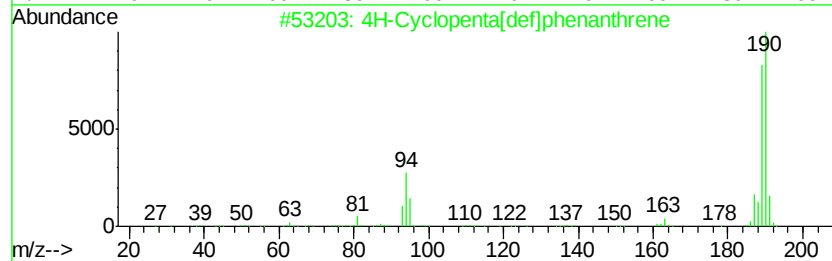
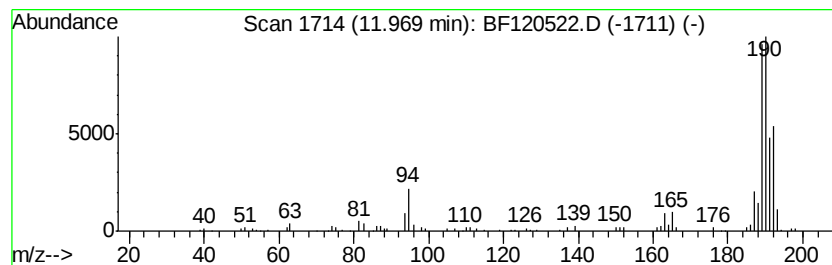
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.97	3.23 ng	179277	Phenanthrene-d10		11.36
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2	6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	58
3	Methyl diselenide	190	C2H6Se2	007101-31-7	41
4	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	35
5	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060320\
Data File : BF120522.D
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ClientSampleId :
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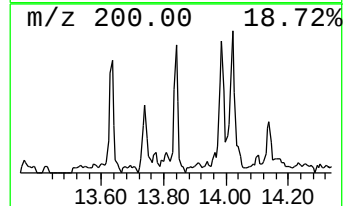
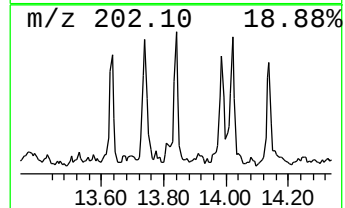
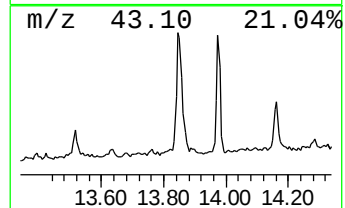
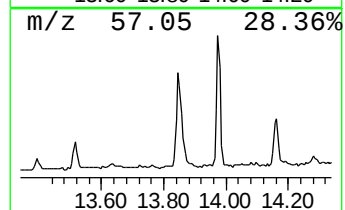
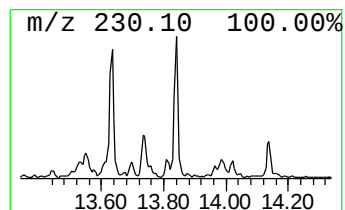
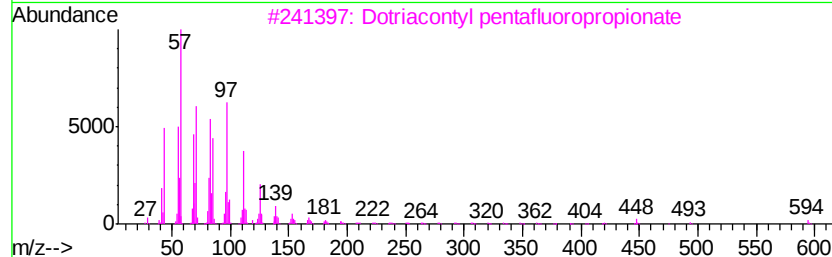
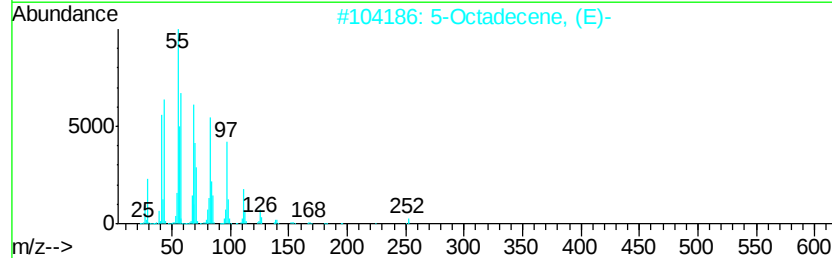
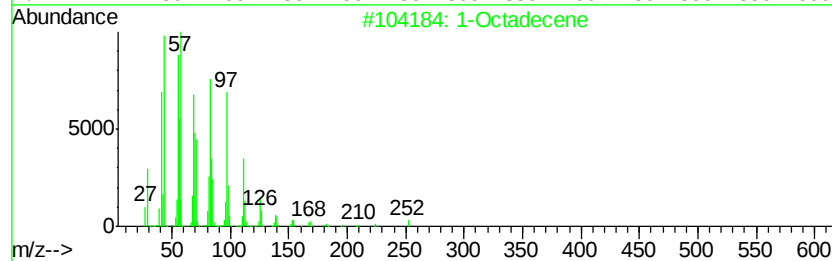
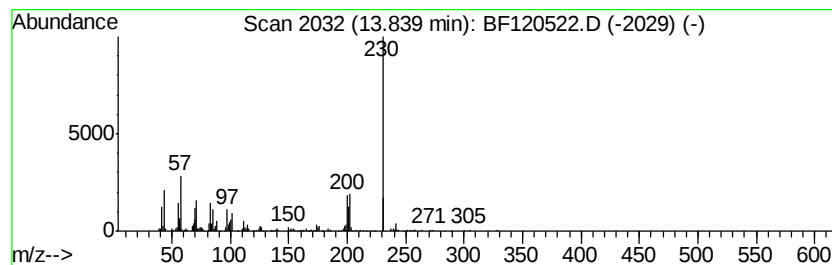
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1-Octadecene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.84	3.82 ng	387304	Chrysene-d12	14.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Octadecene	252	C18H36	000112-88-9	94
2		5-Octadecene, (E)-	252	C18H36	007206-21-5	86
3		Dotriacontyl pentafluoropropionate	612	C35H65F5O2	1000351-81-4	64
4		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	64
5		Octacosyl heptafluorobutyrate	606	C32H57F7O2	1000351-83-6	60



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060320\
Data File : BF120522.D
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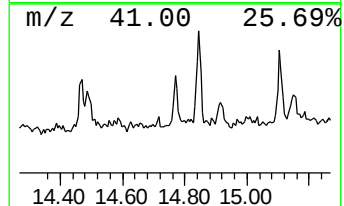
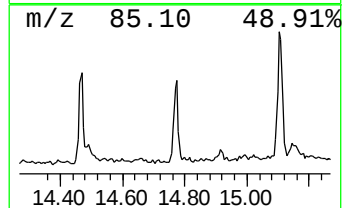
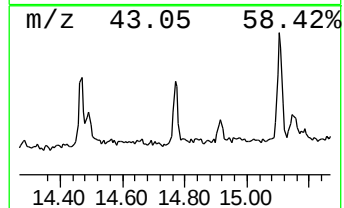
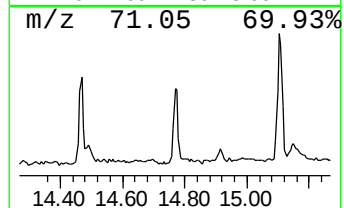
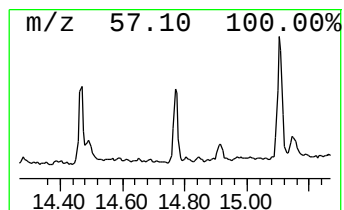
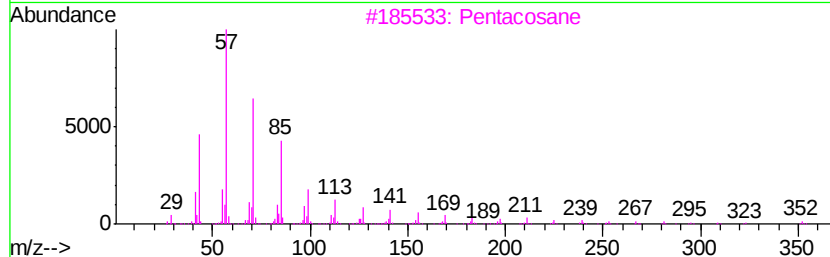
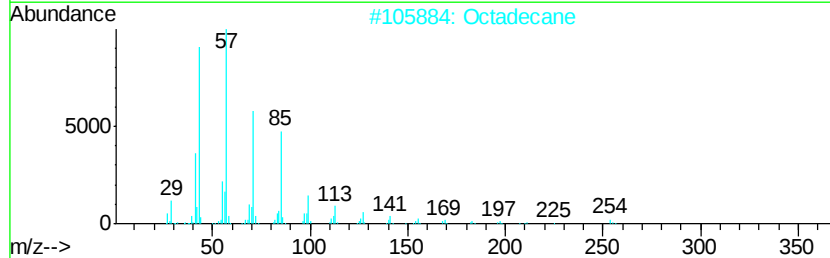
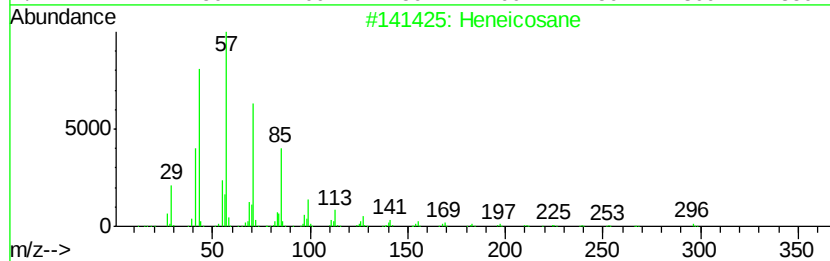
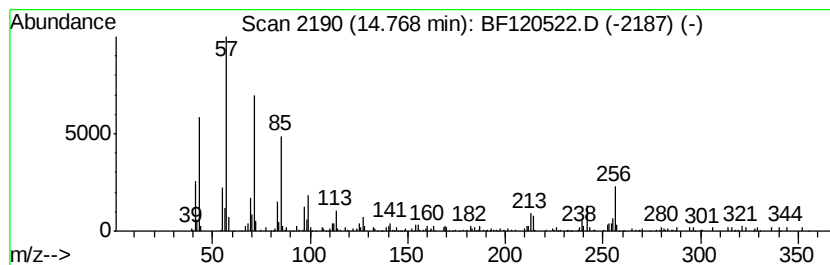
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heneicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.77	2.99 ng	128275	Perylene-d12	15.46

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	90
2	Octadecane		254	C18H38	000593-45-3	83
3	Pentacosane		352	C25H52	000629-99-2	78
4	2-Bromo dodecane		248	C12H25Br	013187-99-0	58
5	2-methyloctacosane		408	C29H60	1000376-72-8	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060320\
Data File : BF120522.D
Acq On : 3 Jun 2020 21:05
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Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
NM6-COMP-2020-0-6

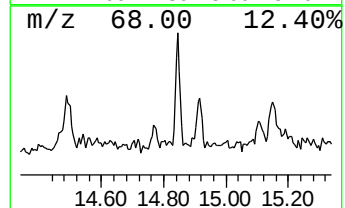
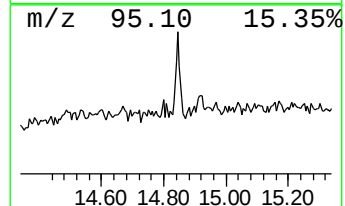
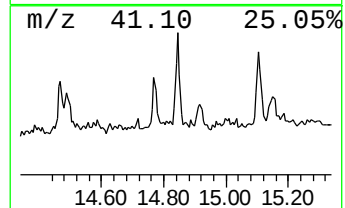
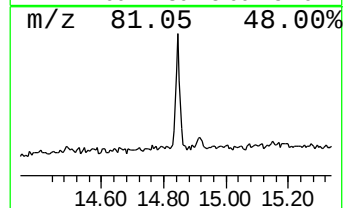
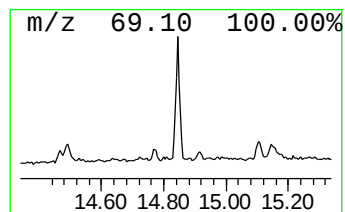
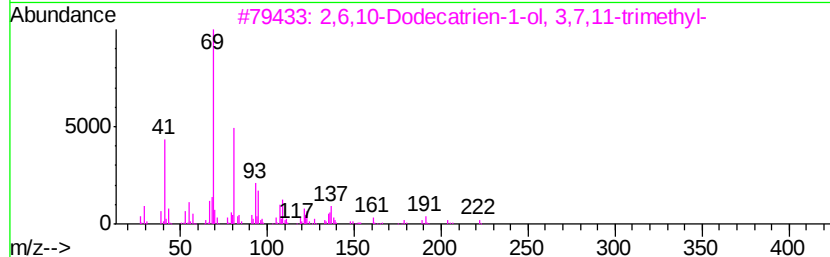
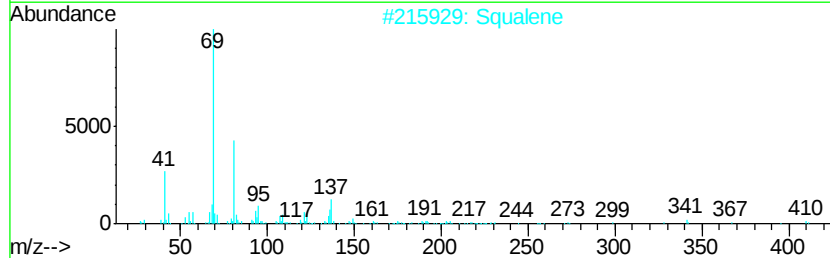
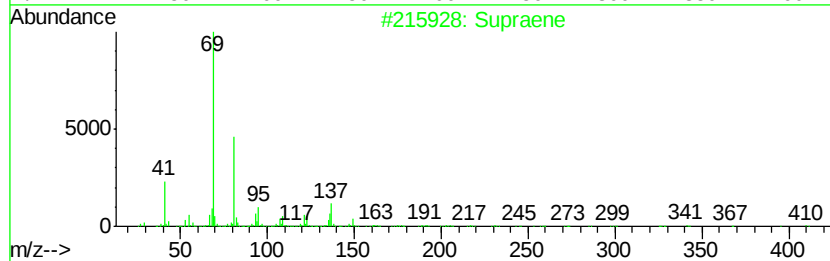
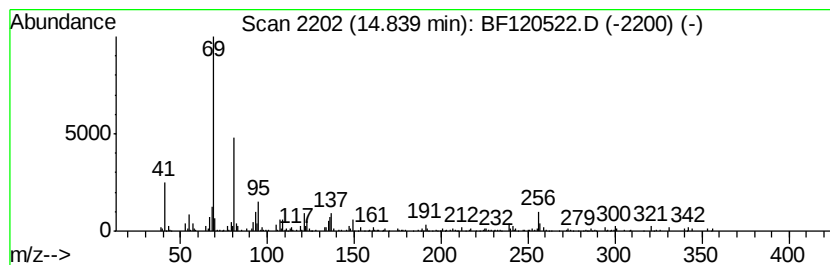
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Supraene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.84	3.36 ng	143921	Perylene-d12	15.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Supraene	410	C30H50	007683-64-9	87
2		Squalene	410	C30H50	000111-02-4	83
3		2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	72
4		2,6,10-Dodecatrienoic acid, 3,7,...	250	C16H26O2	004176-79-8	64
5		4,8,12-Tetradecatrienitrile, 5...	245	C17H27N	005981-31-7	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060320\
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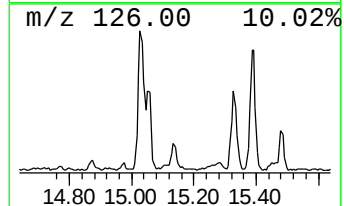
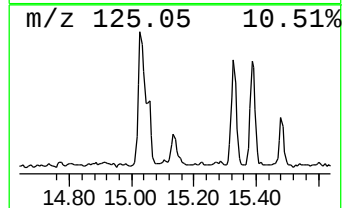
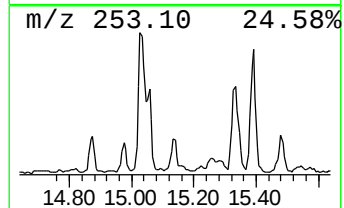
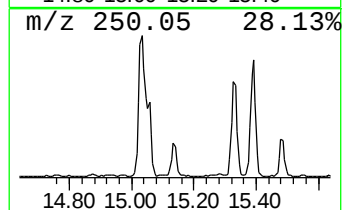
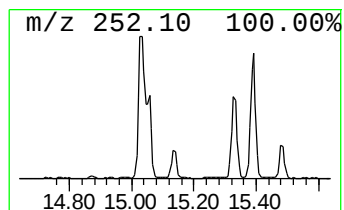
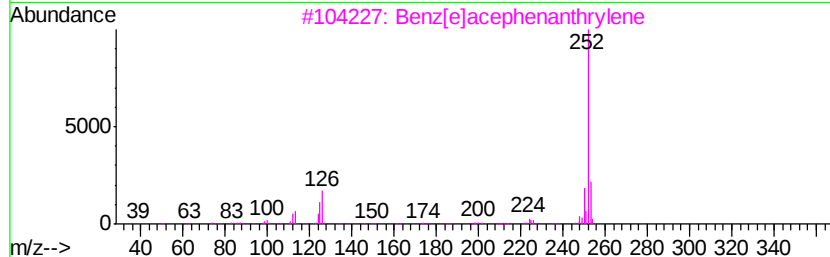
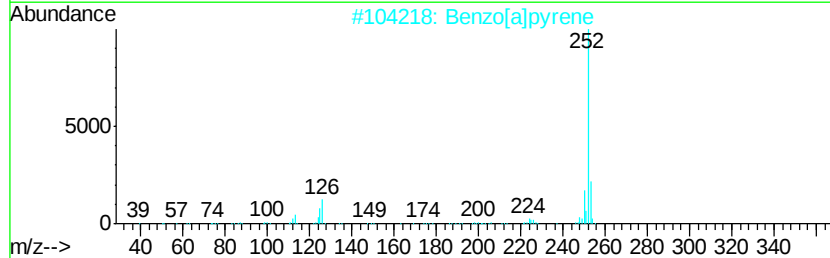
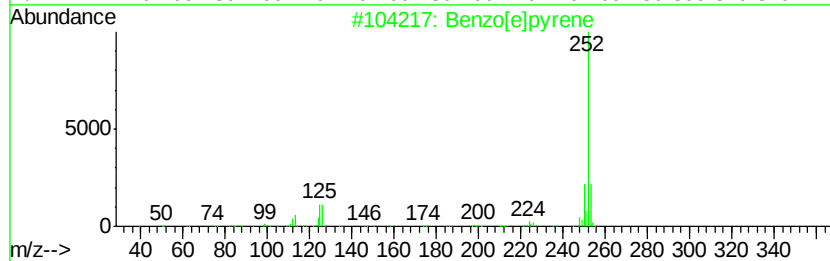
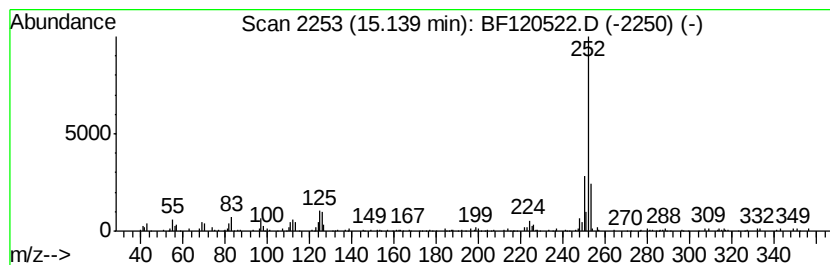
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
15.14	3.46 ng	148197	Perylene-d12	15.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	93
2		Benzo[a]pyrene	252	C20H12	000050-32-8	90
3		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	90
4		Benzo[k]fluoranthene	252	C20H12	000207-08-9	90
5		Perylene	252	C20H12	000198-55-0	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060320\
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Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
NM6-COMP-2020-0-6

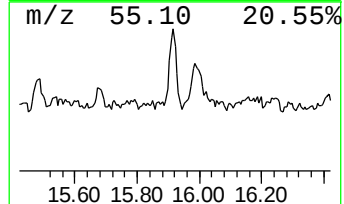
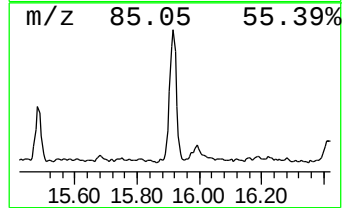
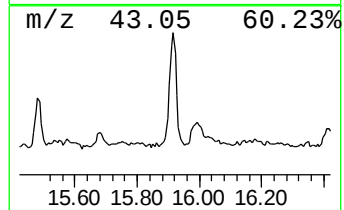
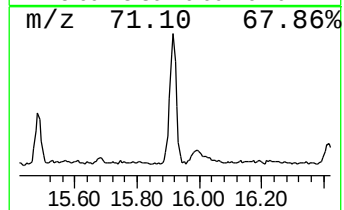
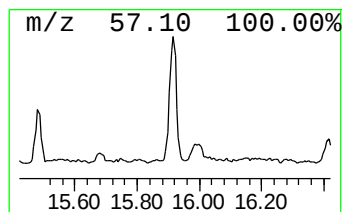
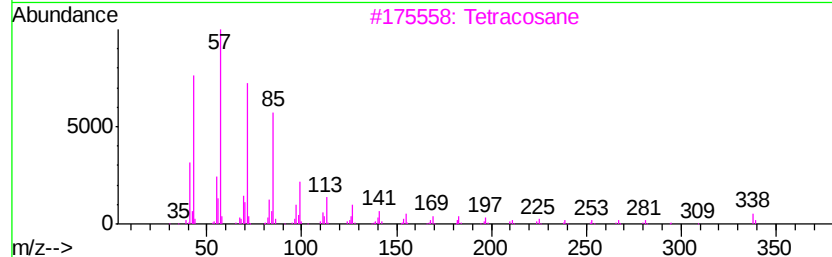
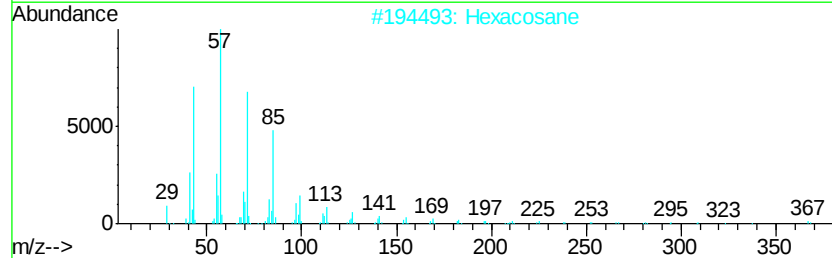
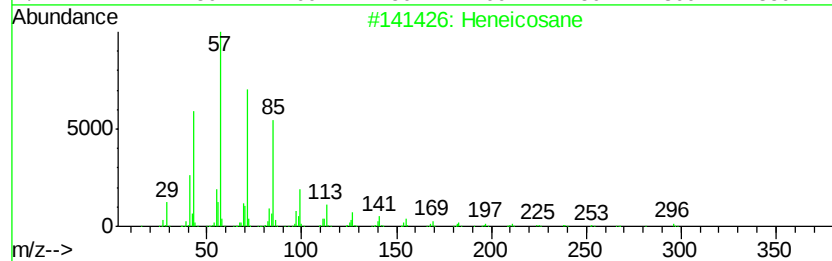
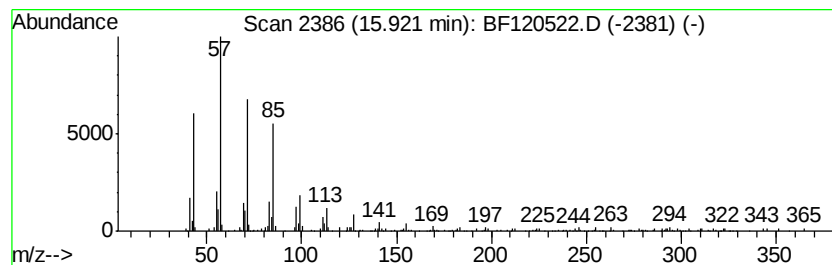
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hexacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.92	6.04 ng	259013	Perylene-d12	15.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Hexacosane	366	C26H54	000630-01-3	91
3		Tetracosane	338	C24H50	000646-31-1	91
4		Nonadecane	268	C19H40	000629-92-5	87
5		Octadecane	254	C18H38	000593-45-3	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF060320\
 Data File : BF120522.D
 Acq On : 3 Jun 2020 21:05
 Operator : JU/CG
 Sample : L2839-12
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM6-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
Methylene chloride	1.95	17.5	ng	701035	1	6.83	800200	20.0
Butane, 2-methoxy...	2.08	24.4	ng	976282	1	6.83	800200	20.0
13-Borabicyclo[7....	11.82	3.1	ng	172601	4	11.36	1111530	20.0
Phenanthrene, 2-m...	11.86	3.3	ng	183311	4	11.36	1111530	20.0
Anthracene, 2-met...	11.89	5.7	ng	317979	4	11.36	1111530	20.0
4H-Cyclopenta[def...	11.97	3.2	ng	179277	4	11.36	1111530	20.0
1-Octadecene	13.84	3.8	ng	387304	5	14.00	2029650	20.0
Heneicosane	14.77	3.0	ng	128275	6	15.46	857567	20.0
Supraene	14.84	3.4	ng	143921	6	15.46	857567	20.0
Benzo[e]pyrene	15.14	3.5	ng	148197	6	15.46	857567	20.0
Hexacosane	15.92	6.0	ng	259013	6	15.46	857567	20.0