

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042121\
 Data File : BF123676.D
 Acq On : 21 Apr 2021 20:06
 Operator : JU/CG
 Sample : M2090-03 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SEDGWICK-COMP-1

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF041521.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF123676.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.046	499	503	510	rBV	176704	198259	7.65%	0.466%
2	5.469	571	575	587	rVB	2654480	2437980	94.04%	5.728%
3	6.481	743	747	758	rBV	3177901	2592532	100.00%	6.091%
4	6.687	778	782	785	rBV4	35036	41300	1.59%	0.097%
5	6.851	806	810	813	rBV	1485087	1245295	48.03%	2.926%
6	7.410	901	905	908	rBV	2131041	1647825	63.56%	3.872%
7	7.451	910	912	916	rVB	54893	41315	1.59%	0.097%
8	8.134	1023	1028	1031	rBV	1912231	1637971	63.18%	3.849%
9	8.157	1031	1032	1037	rVB	149489	86229	3.33%	0.203%
10	8.645	1110	1115	1119	rBV7	19645	37172	1.43%	0.087%
11	8.728	1126	1129	1132	rBV	84364	66784	2.58%	0.157%
12	8.851	1147	1150	1153	rVB	80053	68614	2.65%	0.161%
13	8.951	1163	1167	1171	rBV	73129	76418	2.95%	0.180%
14	9.210	1208	1211	1215	rBV	2866109	2494574	96.22%	5.861%
15	9.298	1223	1226	1227	rBV	83256	75366	2.91%	0.177%
16	9.481	1253	1257	1261	rBV2	53037	77956	3.01%	0.183%
17	9.551	1266	1269	1271	rVV	106836	99177	3.83%	0.233%
18	9.575	1271	1273	1276	rVV	59233	58579	2.26%	0.138%
19	9.604	1276	1278	1282	rVV2	87501	116047	4.48%	0.273%
20	9.669	1286	1289	1294	rVB2	49529	56865	2.19%	0.134%
21	9.751	1300	1303	1309	rVB	130165	118830	4.58%	0.279%
22	9.828	1314	1316	1320	rVB	87565	64178	2.48%	0.151%
23	9.892	1324	1327	1330	rVV	1825380	1511132	58.29%	3.551%
24	9.928	1330	1333	1337	rVB	188402	168515	6.50%	0.396%
25	9.998	1342	1345	1350	rVB4	32026	42030	1.62%	0.099%
26	10.051	1350	1354	1358	rBV6	37940	54647	2.11%	0.128%
27	10.098	1358	1362	1366	rVB	157417	145390	5.61%	0.342%
28	10.139	1366	1369	1376	rVB4	46235	53168	2.05%	0.125%
29	10.210	1378	1381	1384	rBV	39147	42334	1.63%	0.099%
30	10.304	1393	1397	1398	rBV3	36213	39109	1.51%	0.092%
31	10.328	1398	1401	1403	rVB2	93268	79916	3.08%	0.188%
32	10.439	1417	1420	1427	rBV	220957	214797	8.29%	0.505%
33	10.510	1427	1432	1433	rVV	43493	54706	2.11%	0.129%
34	10.528	1433	1435	1437	rVV2	51487	52752	2.03%	0.124%
35	10.551	1437	1439	1441	rVV2	70388	61988	2.39%	0.146%
36	10.598	1445	1447	1451	rVB	76741	68112	2.63%	0.160%

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Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF041521.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	10.681	1456	1461	1466	rBV	1258465	1067548	41.18%	2.508%
38	10.804	1479	1482	1487	rVB3	102018	154735	5.97%	0.364%
39	10.992	1509	1514	1516	rBV3	66590	75064	2.90%	0.176%
40	11.022	1516	1519	1521	rVB3	40844	41300	1.59%	0.097%

41	11.122	1531	1536	1538	rBV3	45356	66560	2.57%	0.156%
42	11.139	1538	1539	1544	rVV	51333	48908	1.89%	0.115%
43	11.186	1544	1547	1550	rVB	71635	62056	2.39%	0.146%
44	11.251	1551	1558	1560	rVV3	101596	147615	5.69%	0.347%
45	11.280	1560	1563	1568	rVB	166364	165138	6.37%	0.388%

46	11.380	1577	1580	1582	rBV	1482470	1347519	51.98%	3.166%
47	11.410	1582	1585	1588	rVV	1759265	1521380	58.68%	3.575%
48	11.457	1590	1593	1596	rVB	501755	420359	16.21%	0.988%
49	11.492	1596	1599	1602	rBV4	43025	49896	1.92%	0.117%
50	11.616	1616	1620	1621	rBV	87220	79814	3.08%	0.188%

51	11.633	1621	1623	1626	rVB4	32639	32545	1.26%	0.076%
52	11.680	1628	1631	1634	rBV	113027	83369	3.22%	0.196%
53	11.716	1634	1637	1643	rVB3	65684	65486	2.53%	0.154%
54	11.780	1644	1648	1651	rBV	517814	456484	17.61%	1.073%
55	11.886	1663	1666	1668	rBV	265311	206772	7.98%	0.486%

56	11.916	1668	1671	1674	rVB	351899	295235	11.39%	0.694%
57	11.963	1676	1679	1682	rBV2	132210	119605	4.61%	0.281%
58	11.998	1682	1685	1687	rVV2	398639	402494	15.53%	0.946%
59	12.022	1687	1689	1694	rVB	210077	161913	6.25%	0.380%
60	12.069	1694	1697	1698	rBV2	32230	36458	1.41%	0.086%

61	12.086	1698	1700	1703	rVB	94122	70395	2.72%	0.165%
62	12.180	1713	1716	1718	rBV	181682	140138	5.41%	0.329%
63	12.204	1718	1720	1724	rVB2	116410	99943	3.86%	0.235%
64	12.333	1740	1742	1744	rVB2	79414	61438	2.37%	0.144%
65	12.363	1744	1747	1749	rBV	111196	105834	4.08%	0.249%

66	12.386	1749	1751	1754	rVB	63387	50280	1.94%	0.118%
67	12.439	1757	1760	1762	rVV	179542	169770	6.55%	0.399%
68	12.469	1762	1765	1767	rVV	1702532	1620635	62.51%	3.808%
69	12.486	1767	1768	1772	rVV	1488077	1047066	40.39%	2.460%
70	12.522	1772	1774	1779	rVB2	137728	146874	5.67%	0.345%

71	12.575	1780	1783	1784	rVV	243378	198569	7.66%	0.467%
72	12.604	1784	1788	1791	rVB	2009526	1906787	73.55%	4.480%
73	12.692	1801	1803	1806	rBV2	54382	49474	1.91%	0.116%
74	12.751	1811	1813	1816	rVV	94828	87337	3.37%	0.205%
75	12.827	1820	1826	1829	rBV	1846804	2052314	79.16%	4.822%

76	12.880	1832	1835	1840	rVB3	90010	131369	5.07%	0.309%
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77	12.945	1844	1846	1847	rBV	96818	72781	2.81%	0.171%
78	12.969	1847	1850	1858	rVB	2104721	2117946	81.69%	4.976%
79	13.051	1859	1864	1867	rBV3	158068	268104	10.34%	0.630%
80	13.139	1876	1879	1880	rBV3	100472	119856	4.62%	0.282%
81	13.157	1880	1882	1885	rVB	329990	264740	10.21%	0.622%
82	13.210	1888	1891	1892	rBV	106229	111206	4.29%	0.261%
83	13.227	1892	1894	1896	rVB	192548	157350	6.07%	0.370%
84	13.263	1896	1900	1902	rBV2	240549	239270	9.23%	0.562%
85	13.351	1913	1915	1918	rVB	108647	97251	3.75%	0.228%
86	13.386	1918	1921	1924	rBV2	137915	131322	5.07%	0.309%
87	13.551	1947	1949	1953	rBV3	123139	141720	5.47%	0.333%
88	13.663	1966	1968	1973	rVB	153199	185532	7.16%	0.436%
89	13.810	1991	1993	1995	rBV2	131327	80725	3.11%	0.190%
90	13.833	1995	1997	2001	rBV2	109058	165456	6.38%	0.389%
91	13.880	2002	2005	2007	rVB2	170084	174006	6.71%	0.409%
92	14.033	2026	2031	2033	rBV2	1655440	2277087	87.83%	5.350%
93	14.057	2033	2035	2038	rVB	884503	699764	26.99%	1.644%
94	14.139	2047	2049	2051	rVB	155216	109518	4.22%	0.257%
95	14.198	2057	2059	2062	rBV2	175976	163371	6.30%	0.384%
96	14.504	2109	2111	2115	rVB3	215621	233252	9.00%	0.548%
97	15.086	2206	2210	2213	rBV	775346	1227937	47.36%	2.885%
98	15.392	2259	2262	2268	rVB	449748	600922	23.18%	1.412%
99	15.451	2269	2272	2278	rVB	671990	816077	31.48%	1.917%
100	15.521	2280	2284	2287	rBV	885264	1131556	43.65%	2.659%

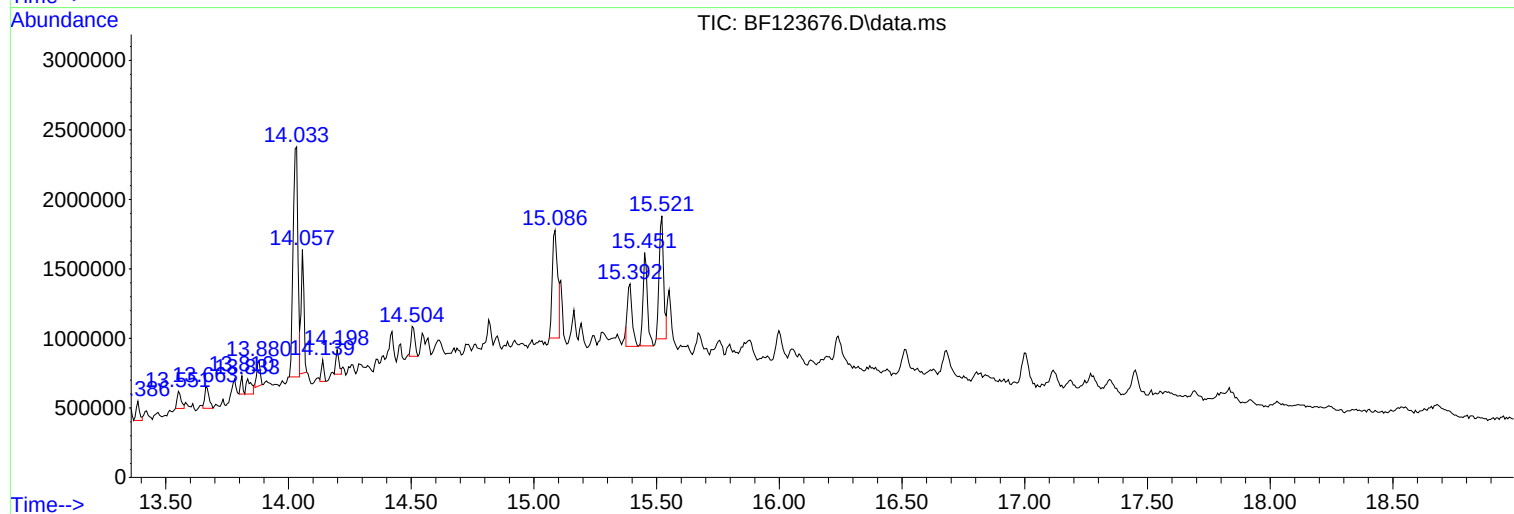
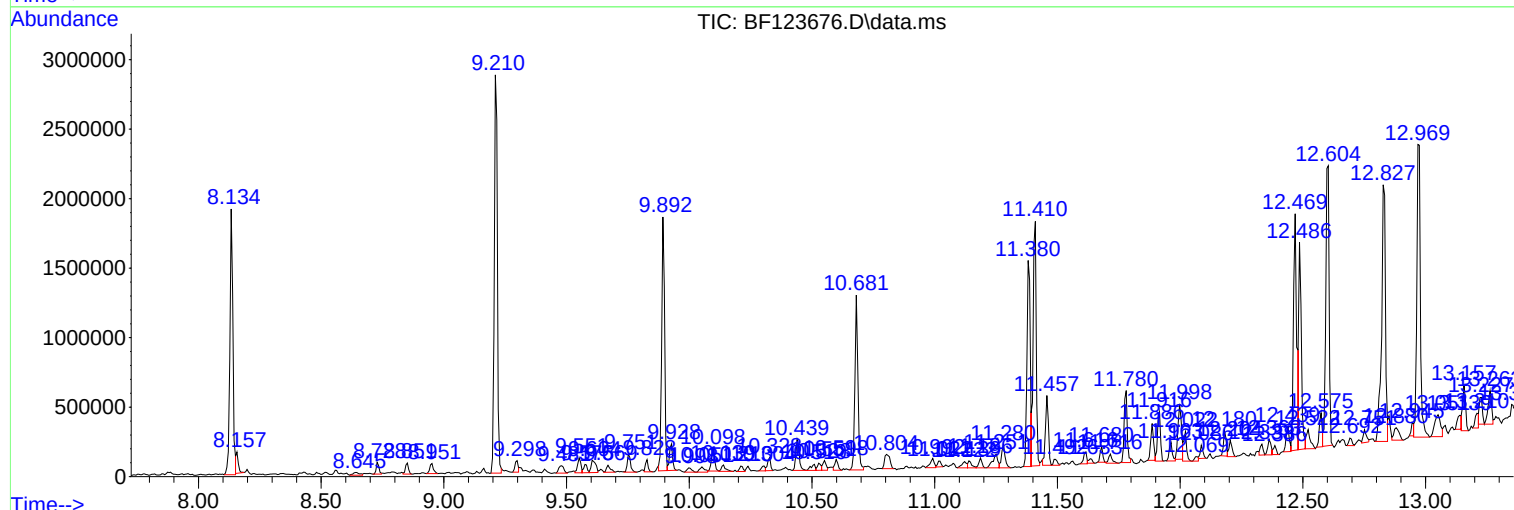
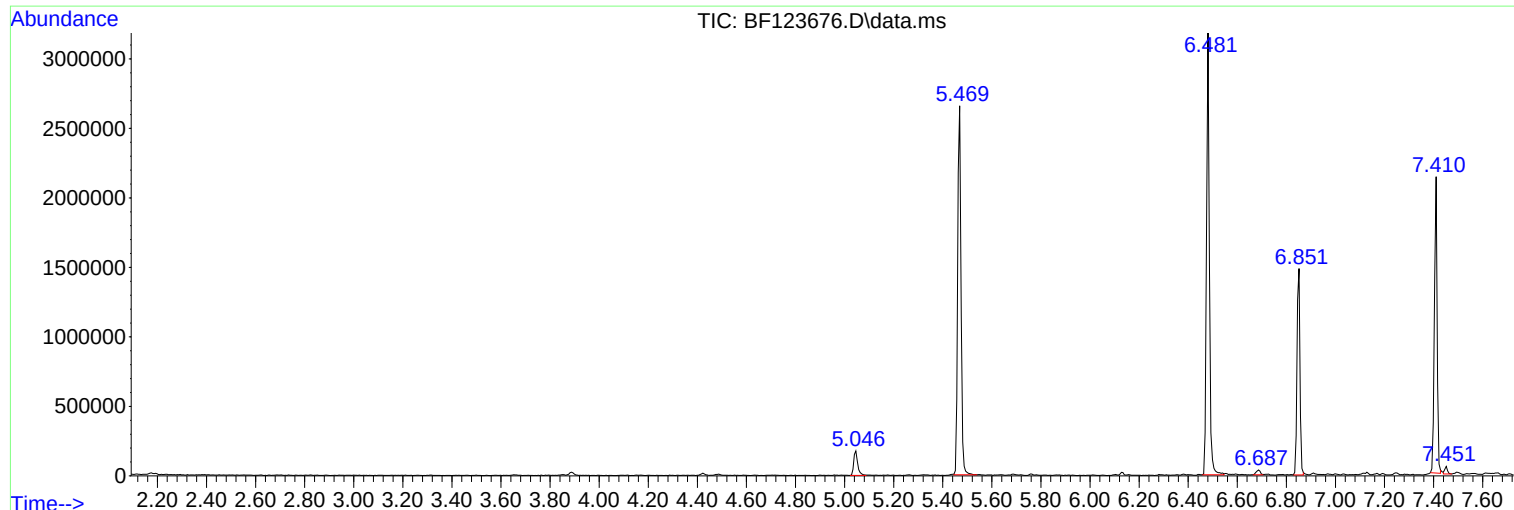
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF041521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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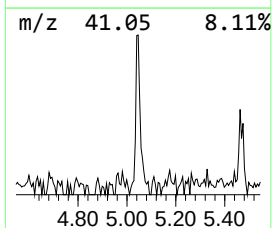
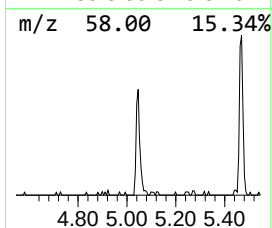
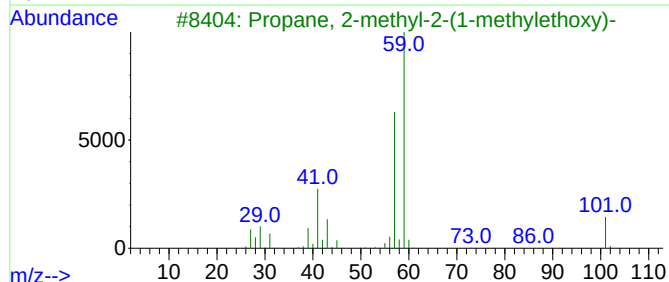
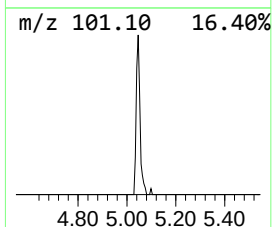
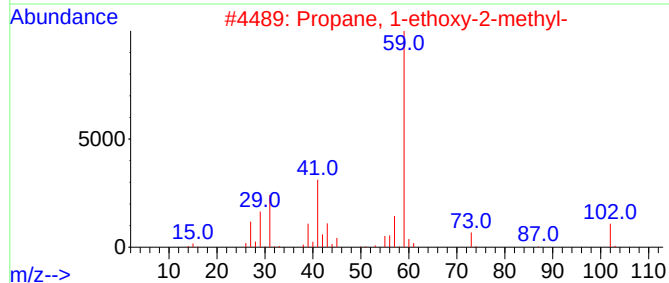
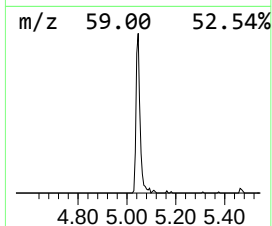
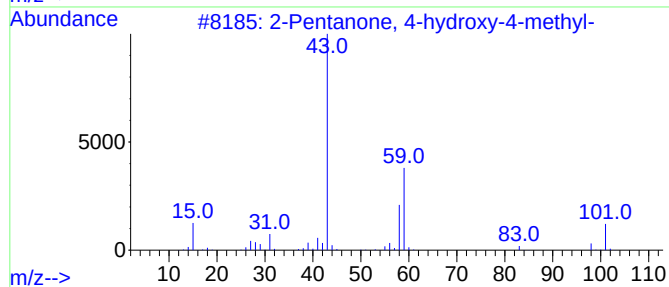
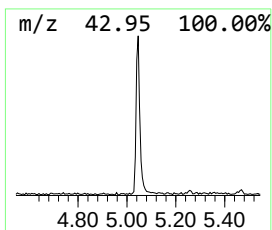
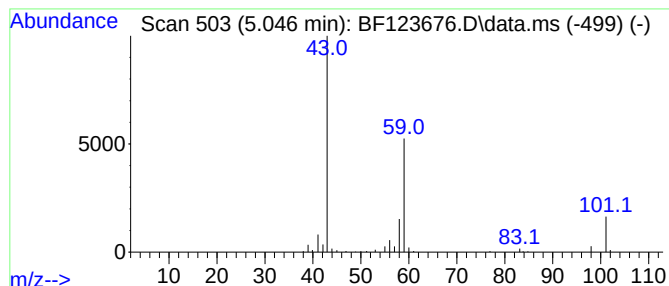
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF041521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.046	3.18 ng	198259	1,4-Dichlorobenzene-d4	6.851

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53		
2	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	37		
3	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36		
4	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28		
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25		



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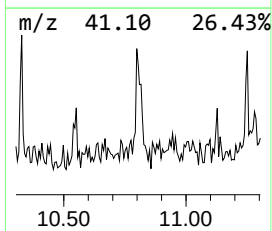
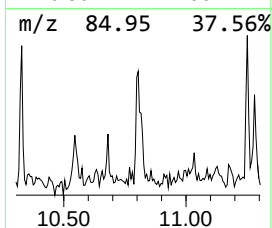
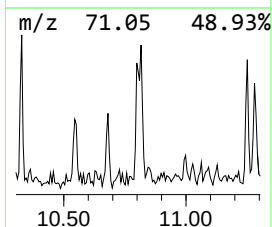
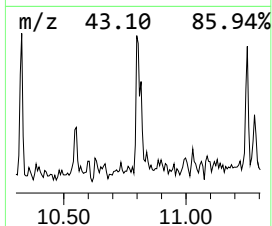
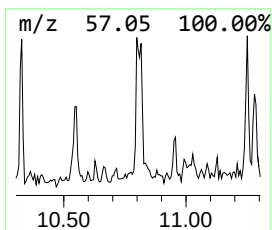
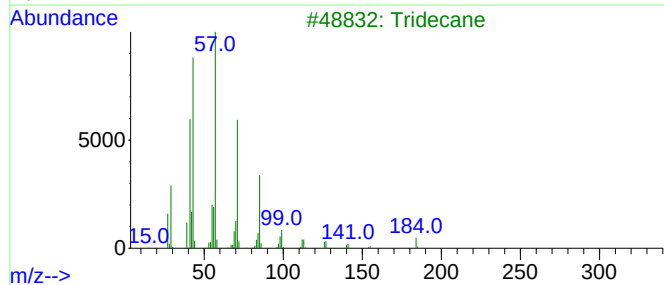
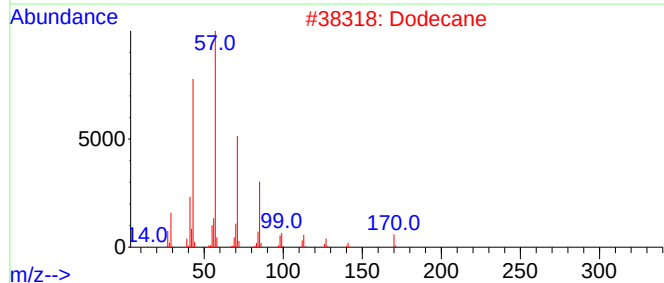
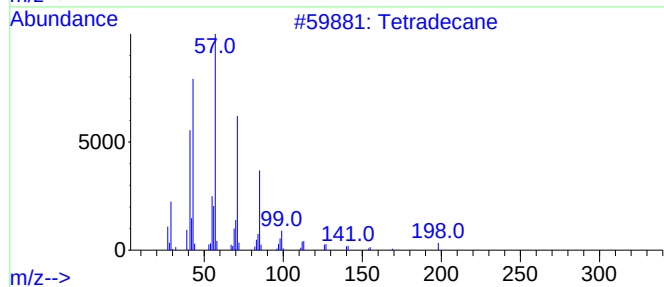
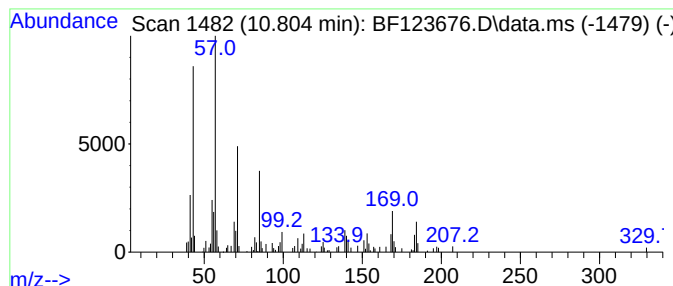
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Tetradecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.804	2.30 ng	154735	Phenanthrene-d10	11.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	70
2			Dodecane	170	C12H26	000112-40-3	70
3			Tridecane	184	C13H28	000629-50-5	68
4			Dodecane, 3-methyl-	184	C13H28	017312-57-1	62
5			Sulfurous acid, butyl octadecyl ...	390	C22H46O3S	1000309-18-5	59



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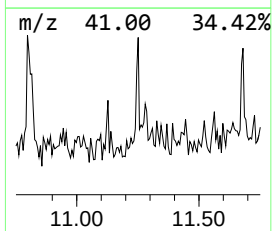
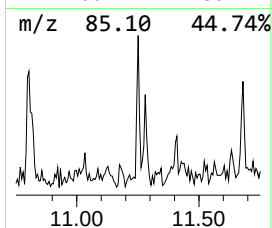
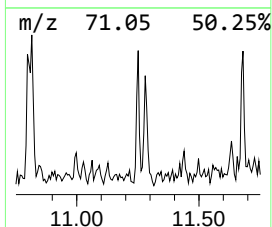
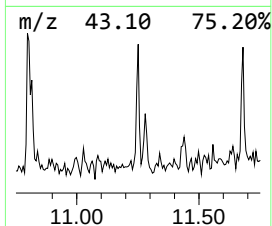
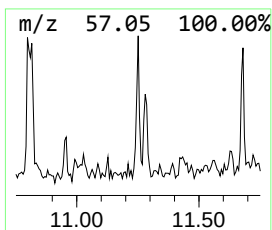
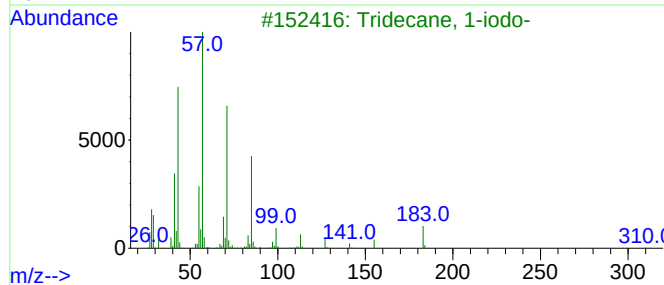
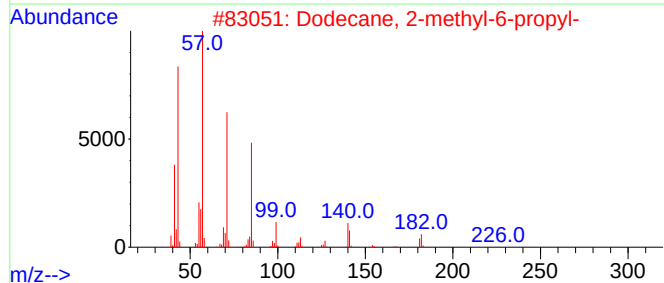
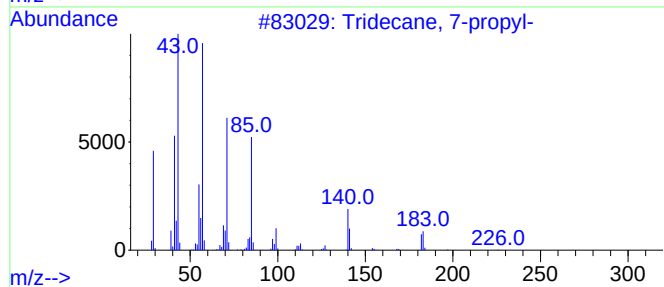
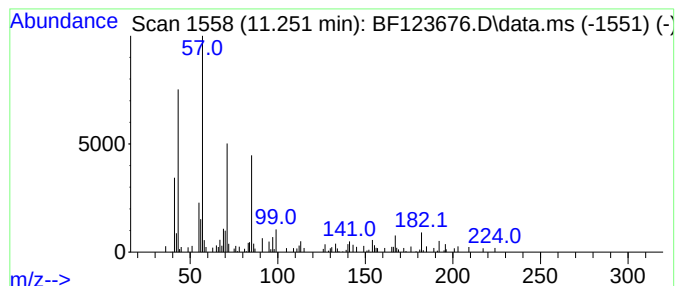
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Tridecane, 7-propyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.251	2.19 ng	147615	Phenanthrene-d10	11.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 7-propyl-	226	C16H34	055045-09-5	80
2			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	64
3			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	64
4			Pentadecane	212	C15H32	000629-62-9	64
5			Nonadecane, 9-methyl-	282	C20H42	013287-24-6	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042121\
Data File : BF123676.D
Acq On : 21 Apr 2021 20:06
Operator : JU/CG
Sample : M2090-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

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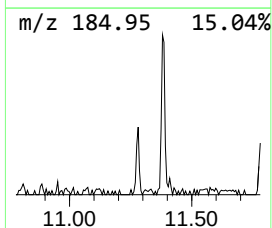
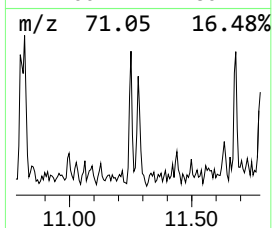
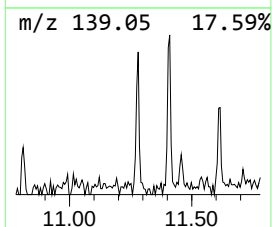
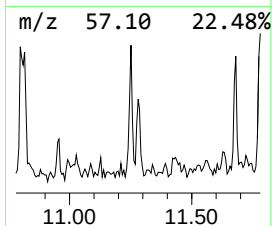
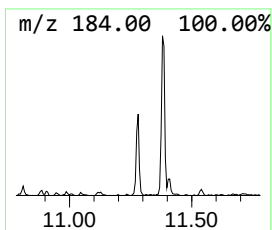
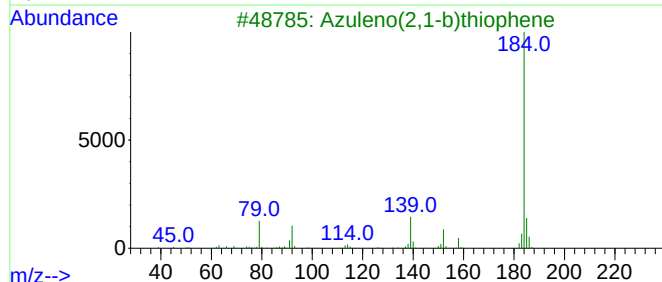
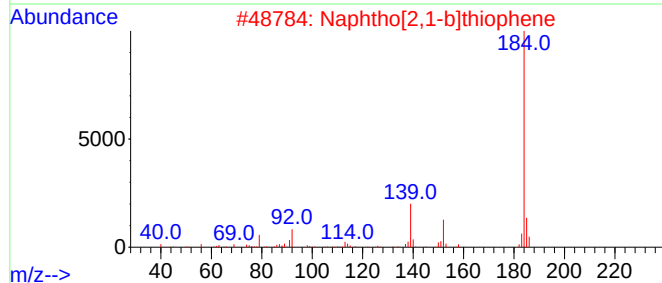
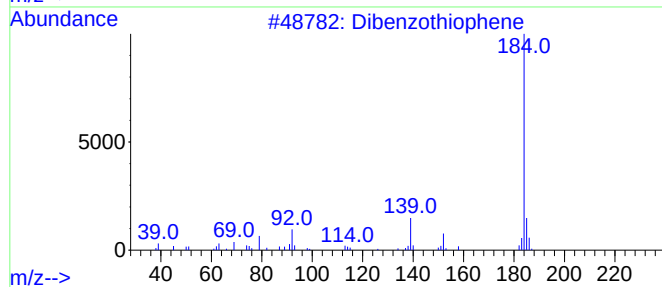
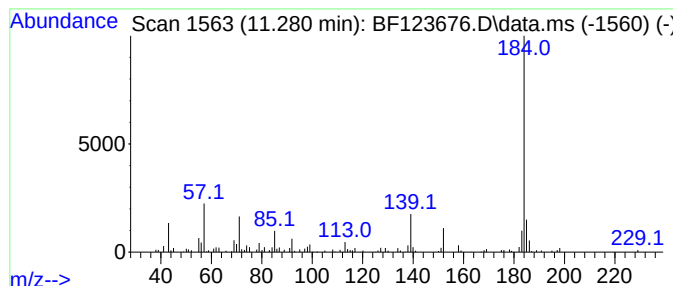
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Dibenzothiophene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.281	2.45 ng	165138	Phenanthrene-d10	11.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzothiophene	184	C12H8S	000132-65-0	93
2			Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	90
3			Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	87
4			Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	83
5			Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042121\
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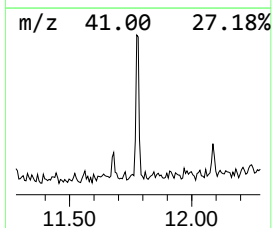
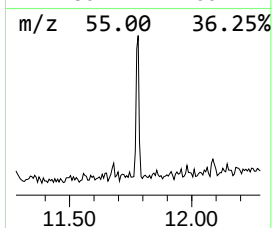
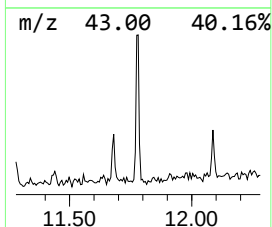
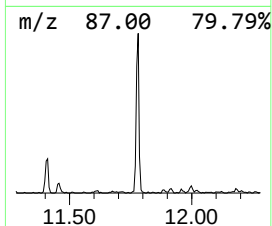
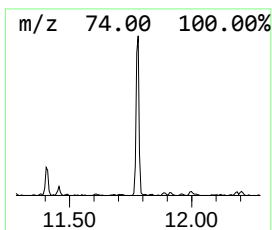
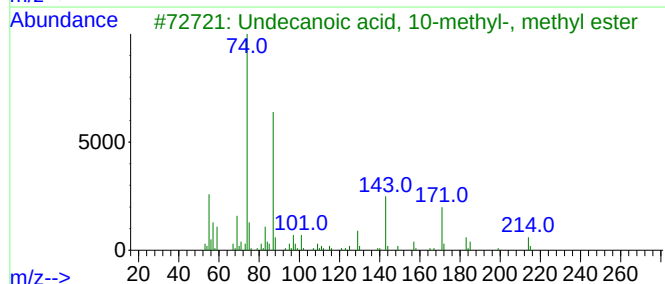
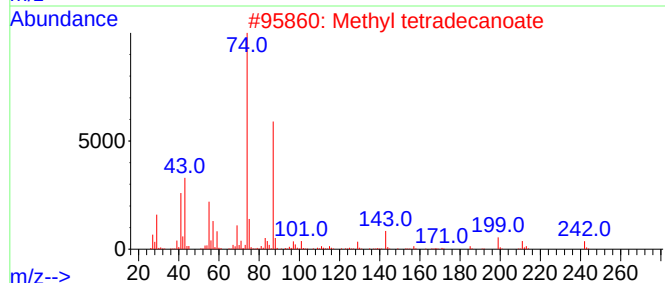
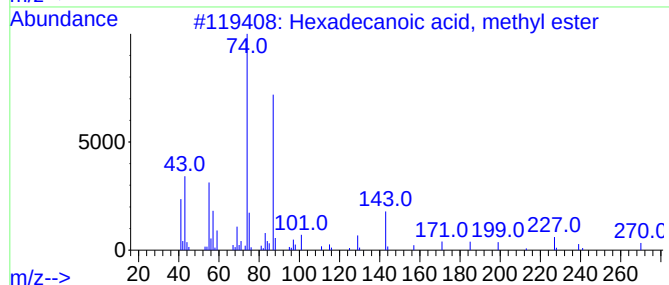
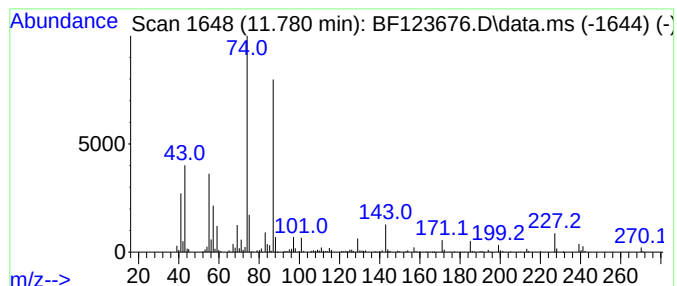
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Hexadecanoic acid, methyl e... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.781	6.78 ng	456484	Phenanthrene-d10	11.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	95
2			Methyl tetradecanoate	242	C15H30O2	000124-10-7	87
3			Undecanoic acid, 10-methyl-, met...	214	C13H26O2	005129-56-6	86
4			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	86
5			Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042121\
Data File : BF123676.D
Acq On : 21 Apr 2021 20:06
Operator : JU/CG
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ALS Vial : 18 Sample Multiplier: 1

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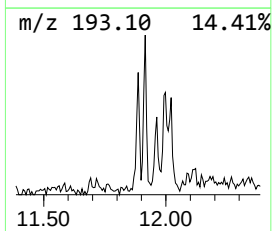
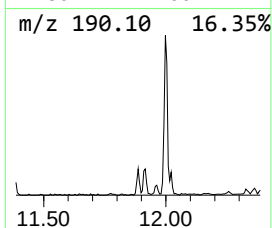
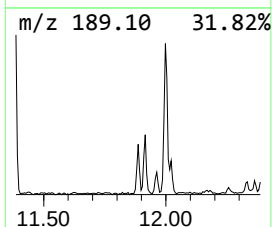
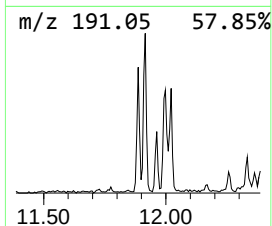
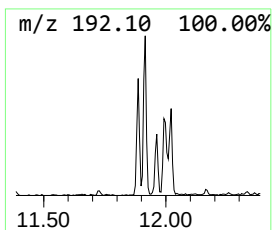
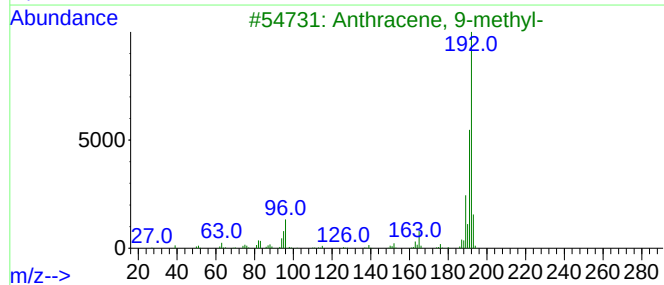
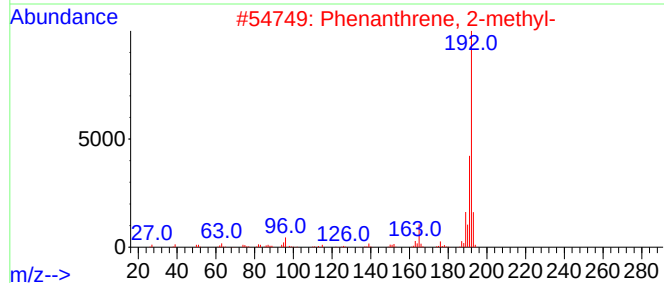
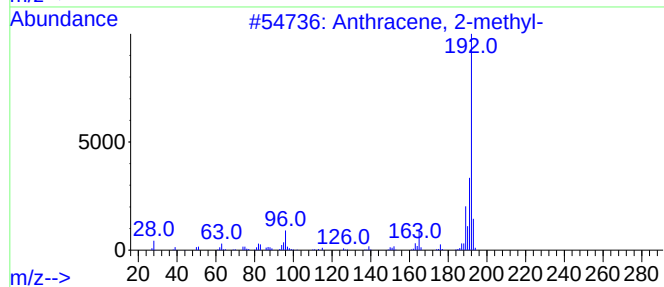
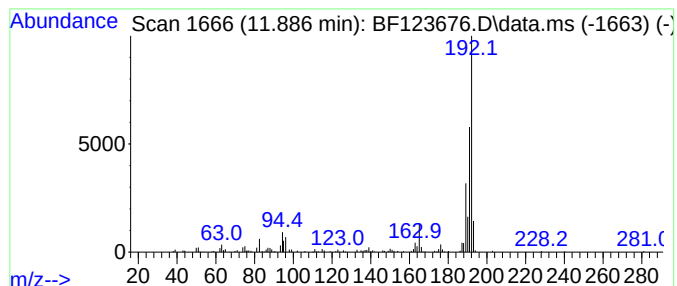
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Anthracene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.886	3.07 ng	206772	Phenanthrene-d10	11.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
5			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042121\
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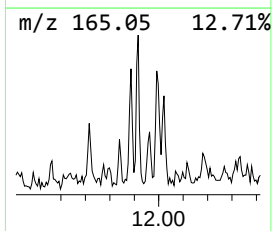
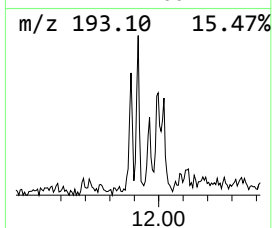
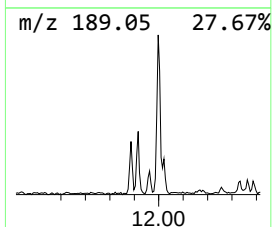
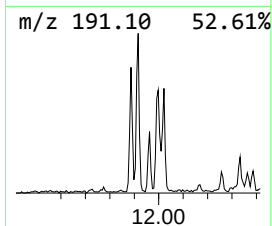
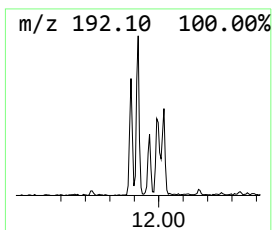
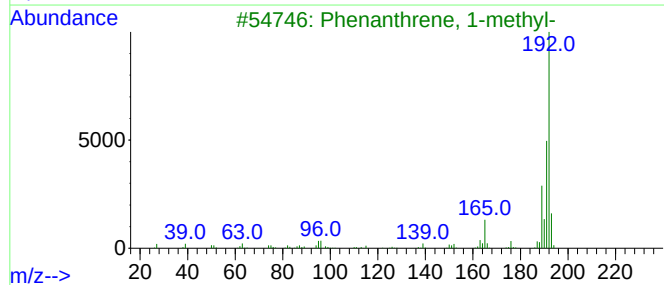
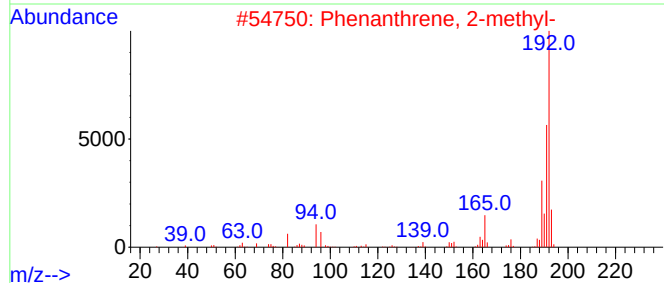
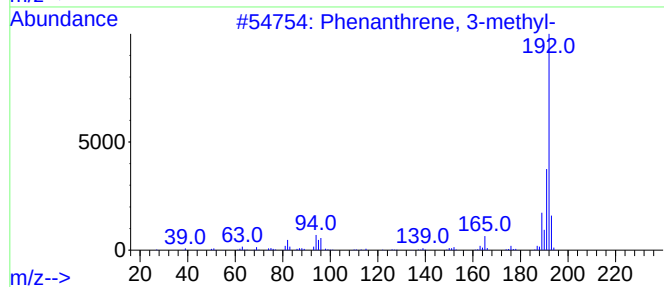
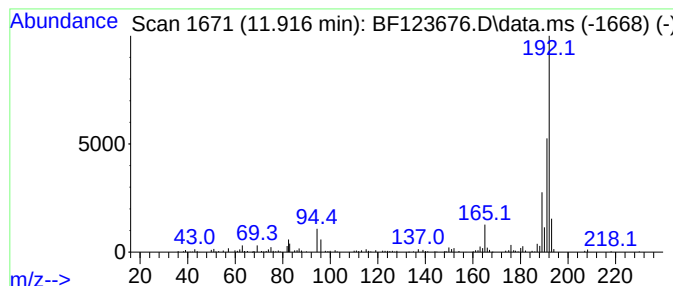
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenanthrene, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.916	4.38 ng	295235	Phenanthrene-d10	11.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	96
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
4			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	94
5			1H-Indene, 1-phenyl-	192	C15H12	001961-96-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042121\
Data File : BF123676.D
Acq On : 21 Apr 2021 20:06
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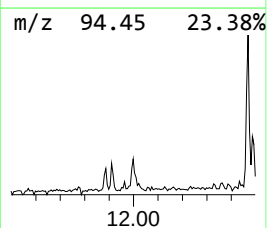
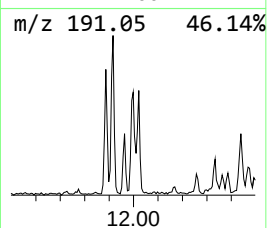
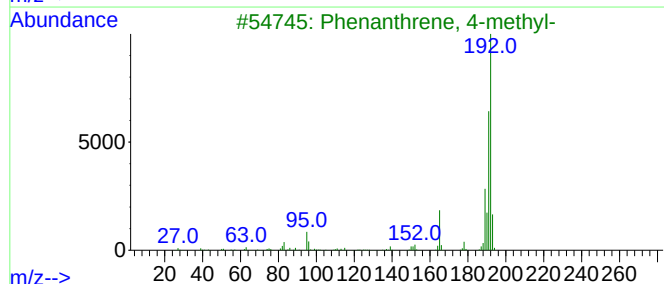
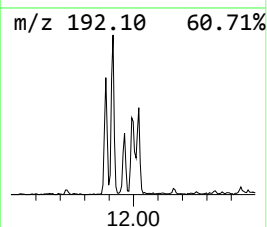
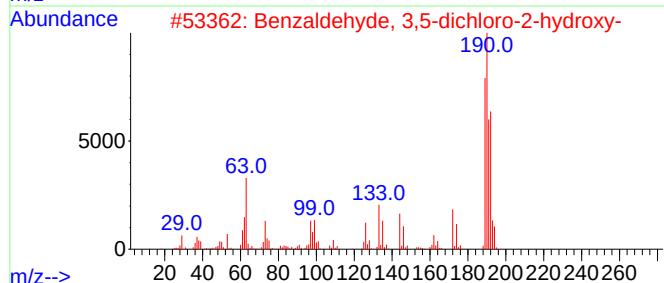
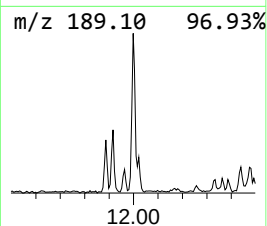
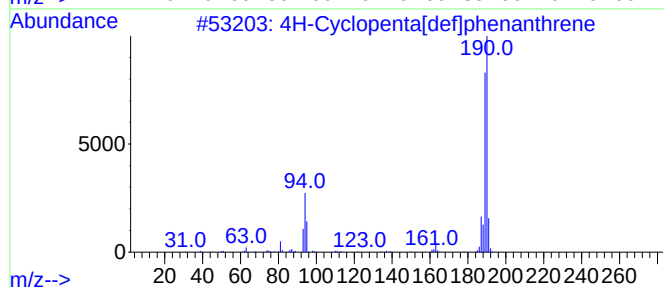
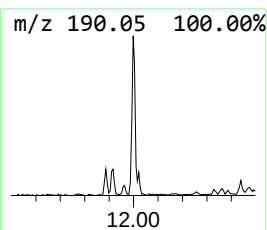
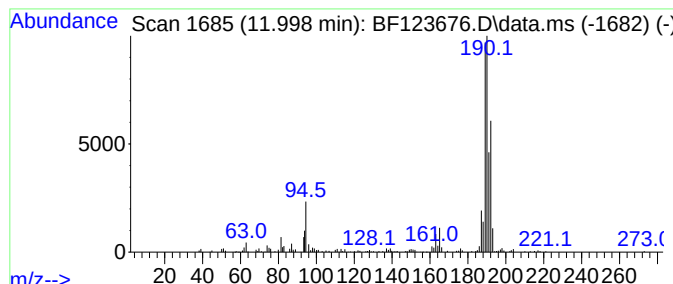
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.998	5.97 ng	402494	Phenanthrene-d10	11.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2			Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	50
3			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	35
4			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	30
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042121\
Data File : BF123676.D
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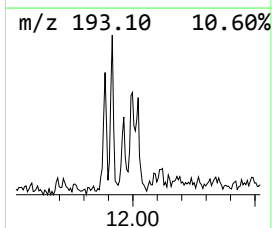
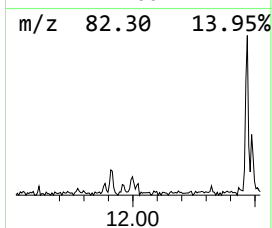
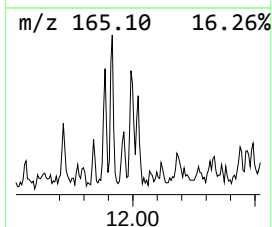
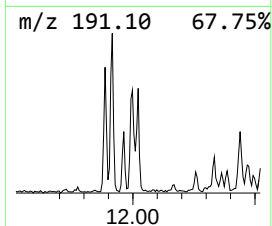
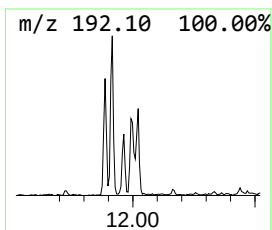
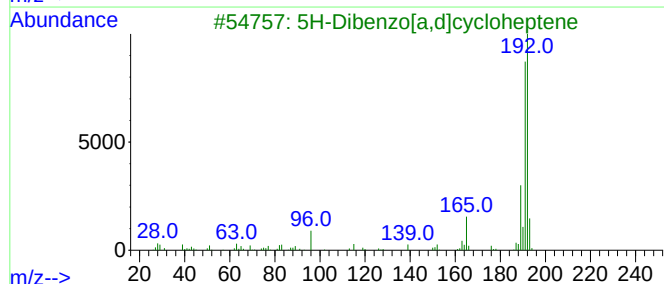
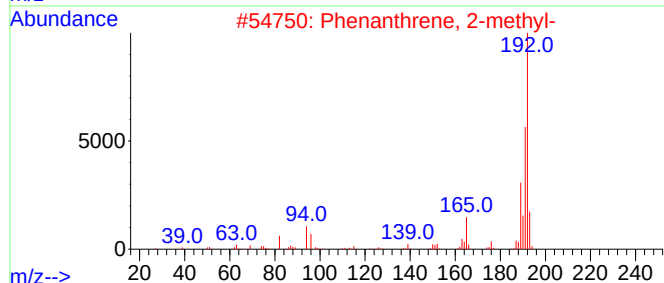
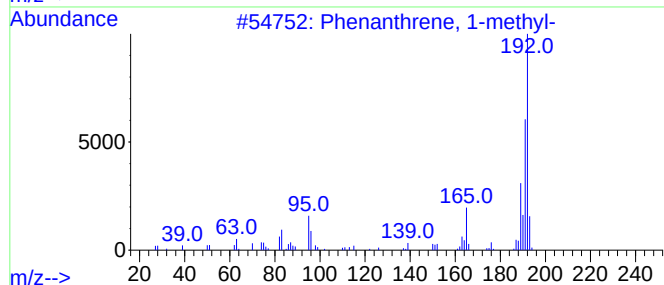
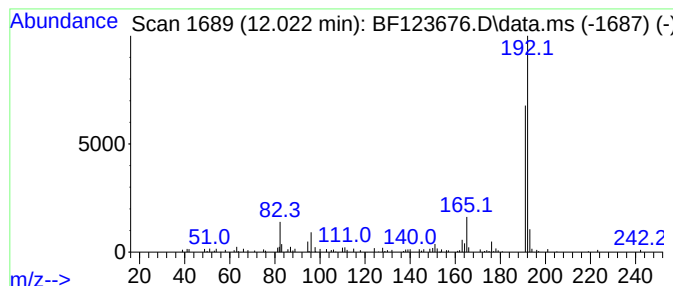
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenanthrene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.022	2.40 ng	161913	Phenanthrene-d10	11.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90
3			5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	90
4			Anthracene, 9-methyl-	192	C15H12	000779-02-2	90
5			Anthracene, 2-methyl-	192	C15H12	000613-12-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042121\
Data File : BF123676.D
Acq On : 21 Apr 2021 20:06
Operator : JU/CG
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ALS Vial : 18 Sample Multiplier: 1

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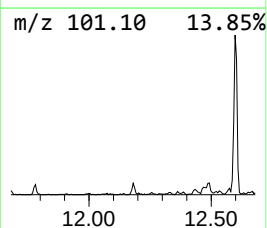
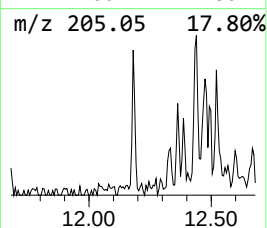
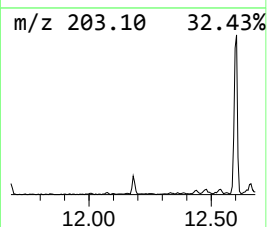
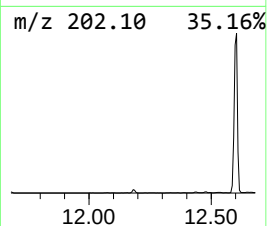
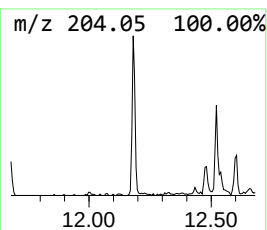
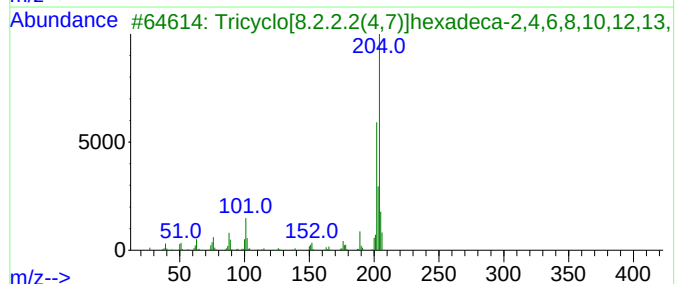
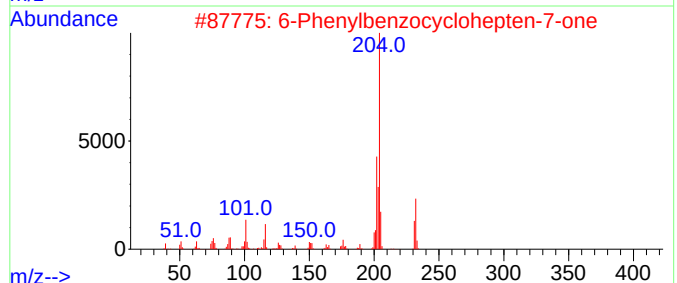
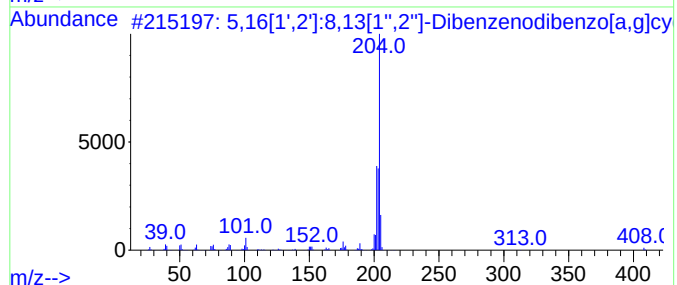
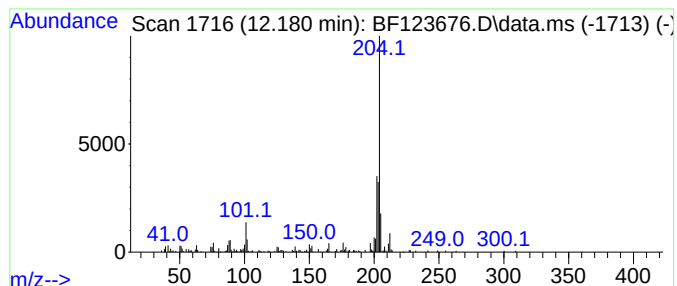
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.180	2.08 ng	140138	Phenanthrene-d10	11.381

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	86	
2	6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	86	
3	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	81	
4	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	81	
5	9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	58	



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042121\
Data File : BF123676.D
Acq On : 21 Apr 2021 20:06
Operator : JU/CG
Sample : M2090-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SEDGWICK-COMP-1

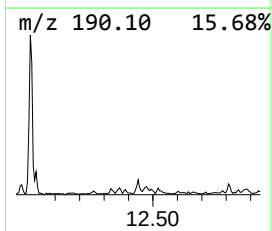
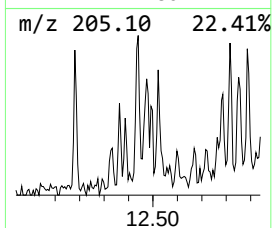
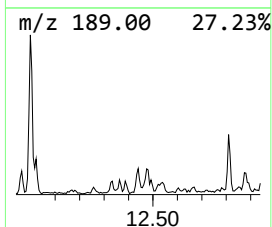
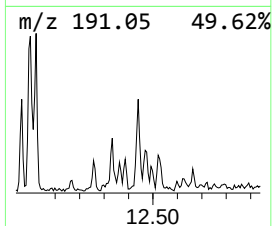
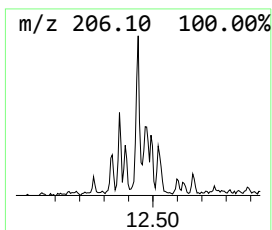
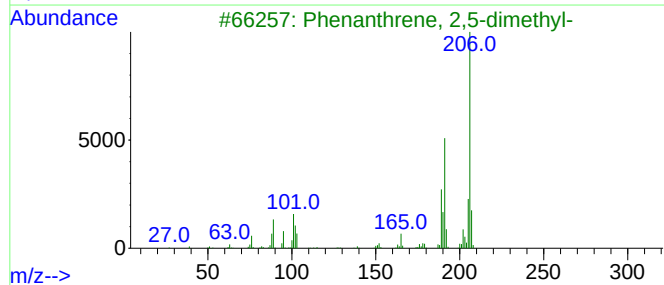
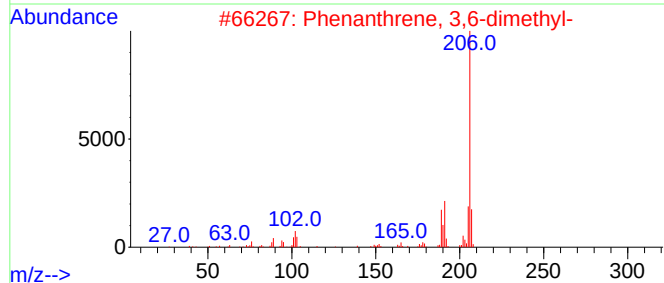
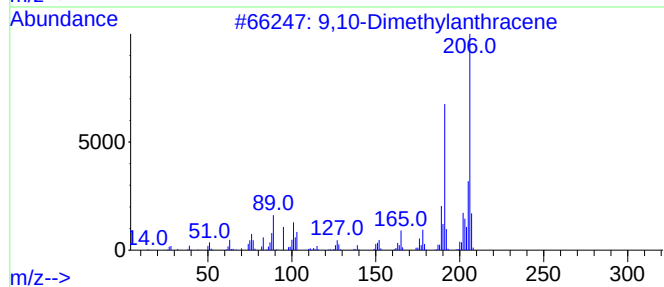
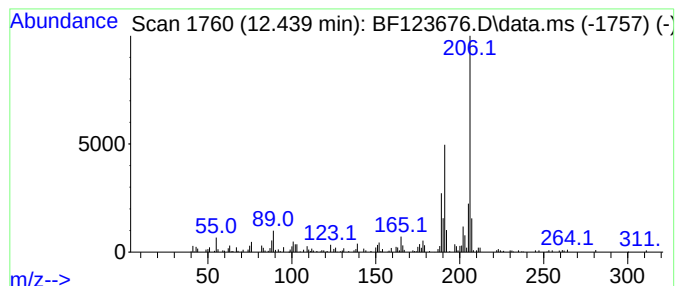
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 9,10-Dimethylantracene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.439	2.52 ng	169770	Phenanthrene-d10	11.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Dimethylantracene	206	C16H14	000781-43-1	91
2			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	91
3			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	90
4			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	81
5			1H-Indene, 2-methyl-3-phenyl-	206	C16H14	035099-60-6	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042121\
Data File : BF123676.D
Acq On : 21 Apr 2021 20:06
Operator : JU/CG
Sample : M2090-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

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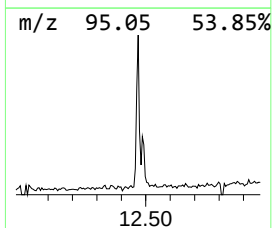
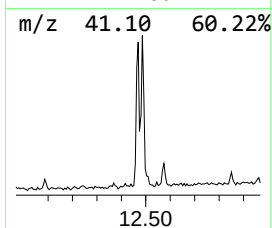
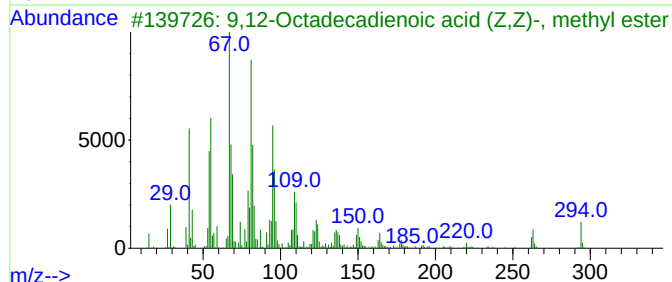
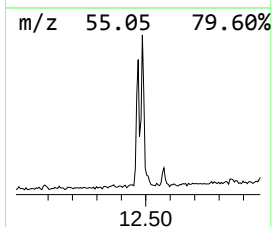
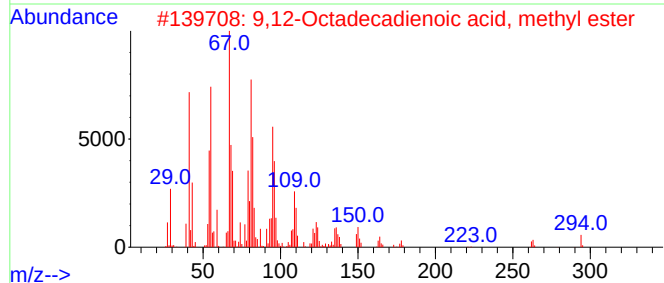
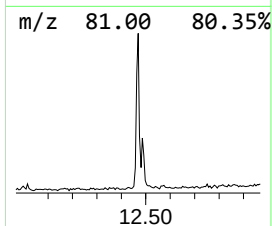
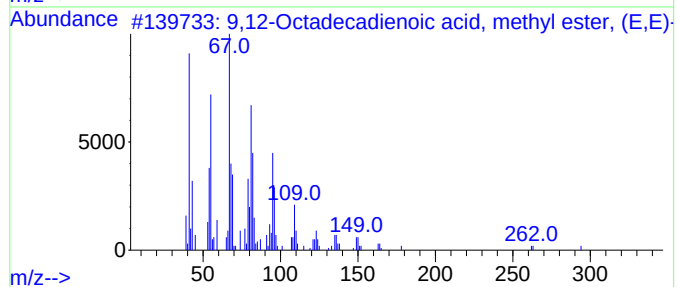
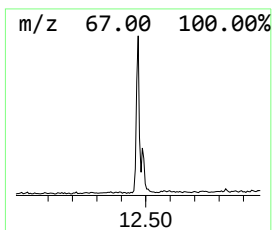
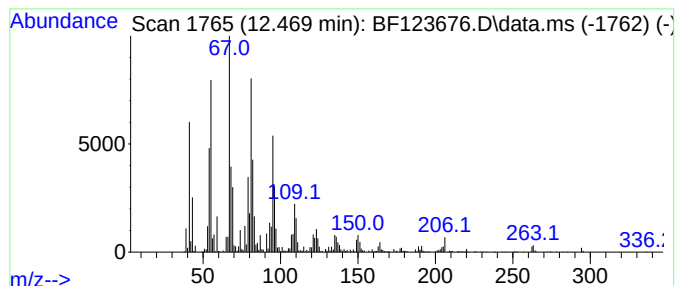
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.469	24.05 ng	1620640	Phenanthrene-d10	11.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99		
2	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99		
3	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99		
4	10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99		
5	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042121\
Data File : BF123676.D
Acq On : 21 Apr 2021 20:06
Operator : JU/CG
Sample : M2090-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

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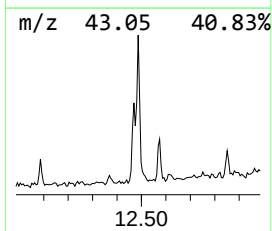
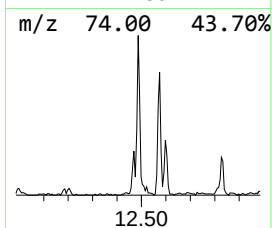
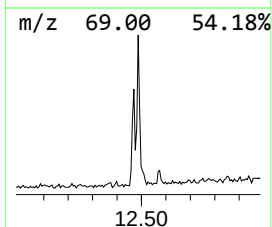
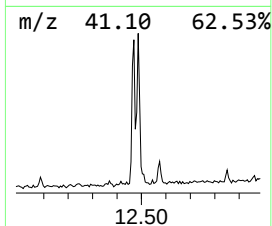
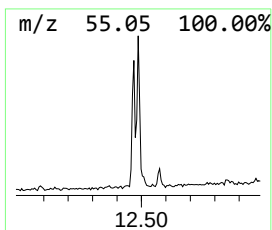
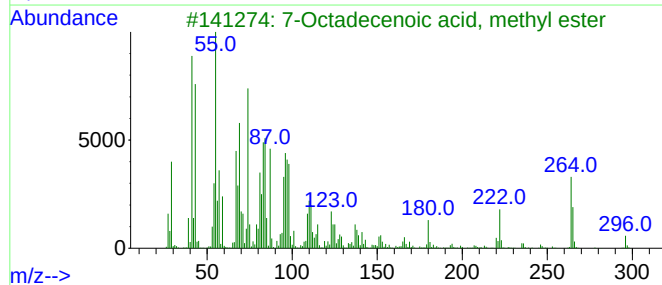
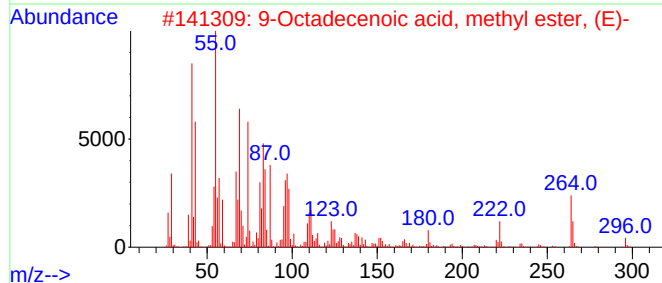
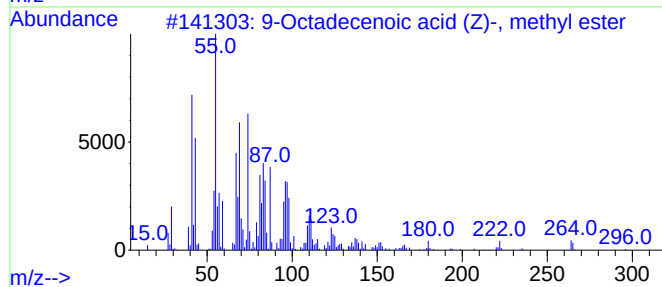
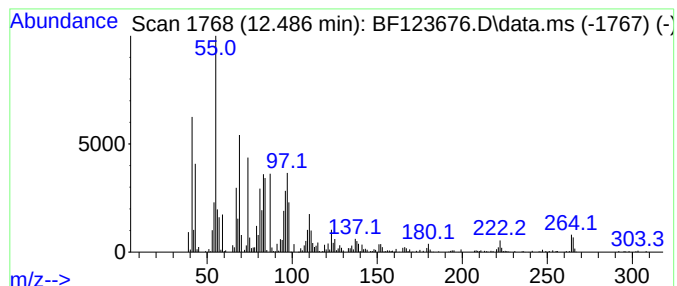
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 9-Octadecenoic acid (Z)-, m... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.486	15.54 ng	1047070	Phenanthrene-d10	11.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	98
2			9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	91
3			7-Octadecenoic acid, methyl ester	296	C19H36O2	057396-98-2	91
4			12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	91
5			10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042121\
Data File : BF123676.D
Acq On : 21 Apr 2021 20:06
Operator : JU/CG
Sample : M2090-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

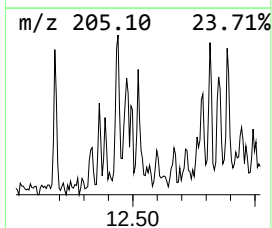
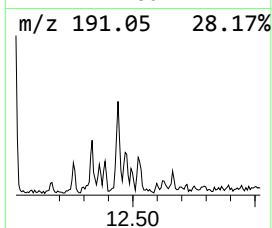
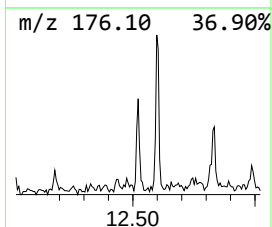
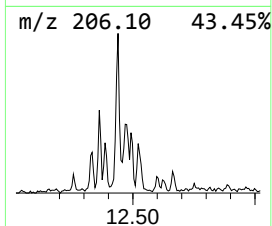
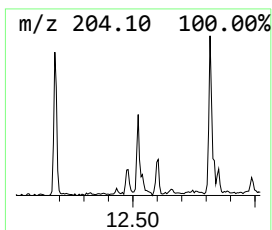
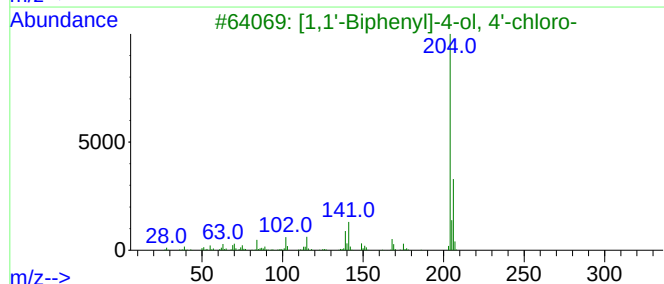
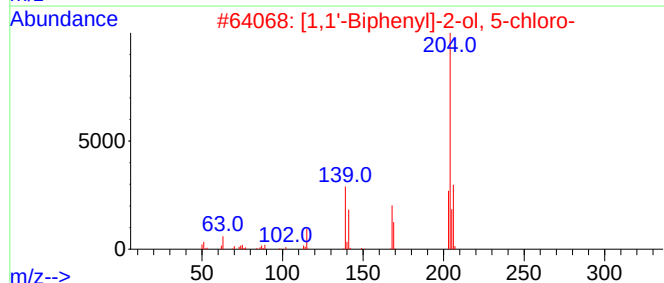
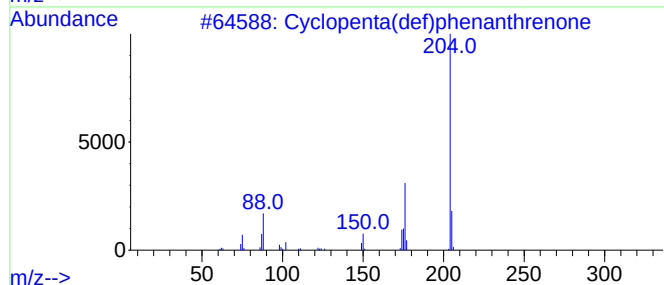
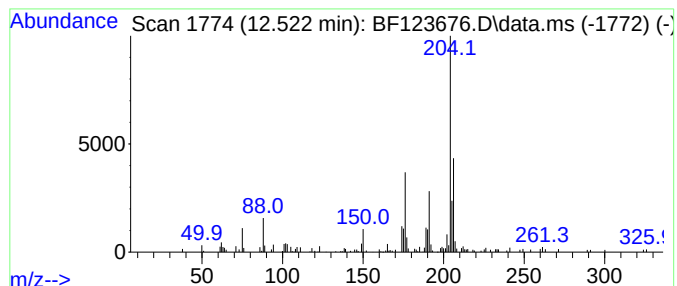
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF041521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Cyclopenta(def)phenanthrenone Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD			R.T.
12.522	2.18 ng	146874	Phenanthrene-d10			11.381
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone		204	C15H8O	005737-13-3	90
2	[1,1'-Biphenyl]-2-ol, 5-chloro-		204	C12H9ClO	000607-12-5	47
3	[1,1'-Biphenyl]-4-ol, 4'-chloro-		204	C12H9ClO	028034-99-3	43
4	Benzene, 1-chloro-4-phenoxy-		204	C12H9ClO	007005-72-3	43
5	2-Chloro-4-phenylphenol		204	C12H9ClO	000092-04-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042121\
Data File : BF123676.D
Acq On : 21 Apr 2021 20:06
Operator : JU/CG
Sample : M2090-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

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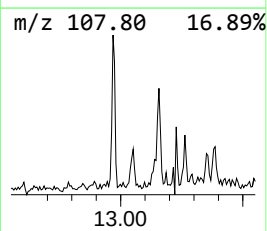
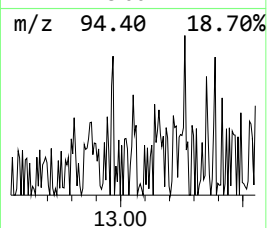
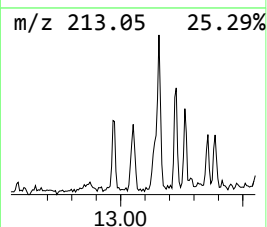
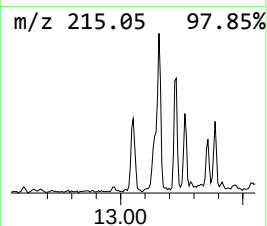
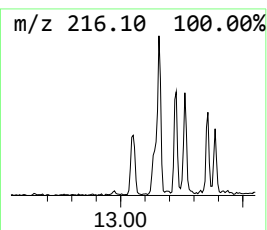
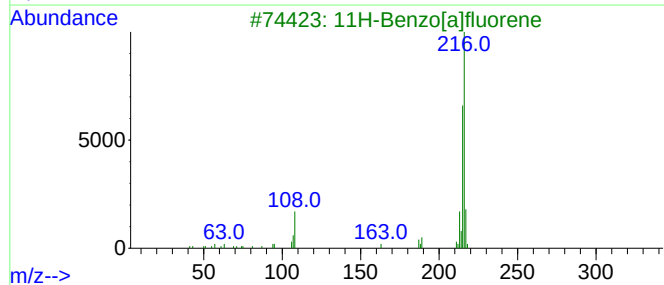
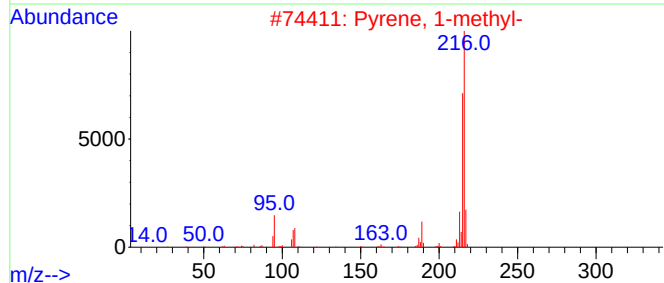
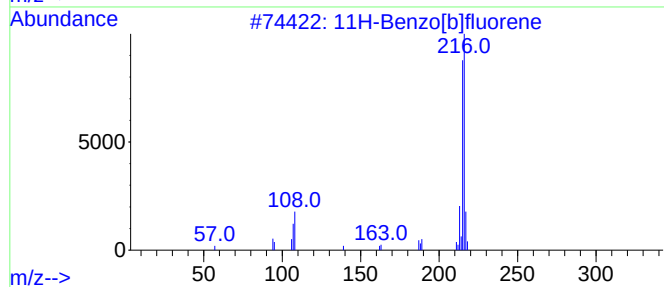
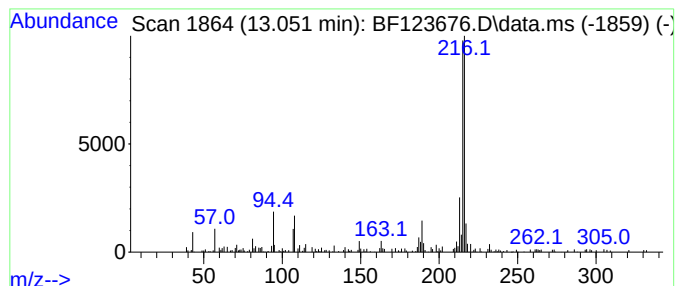
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 11H-Benzo[b]fluorene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.051	2.35 ng	268104	Chrysene-d12	14.033

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	95
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	76
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	72
5			Pyrene, 2-methyl-	216	C17H12	003442-78-2	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042121\
Data File : BF123676.D
Acq On : 21 Apr 2021 20:06
Operator : JU/CG
Sample : M2090-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SEDGWICK-COMP-1

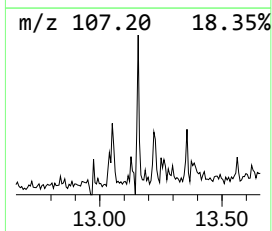
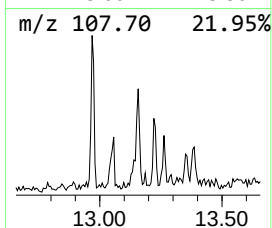
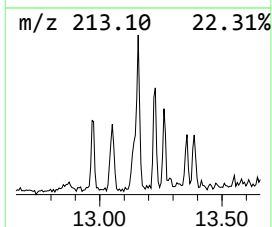
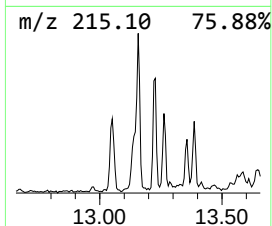
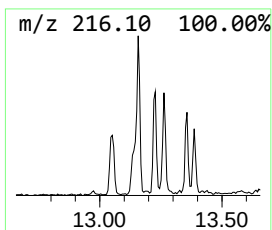
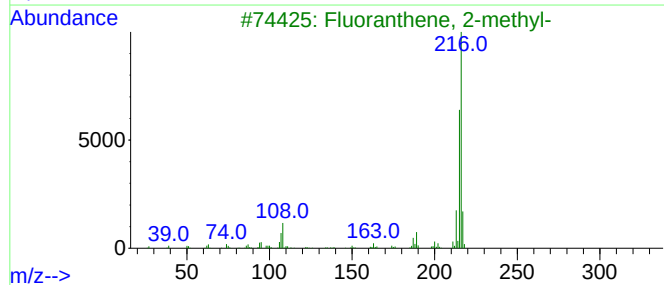
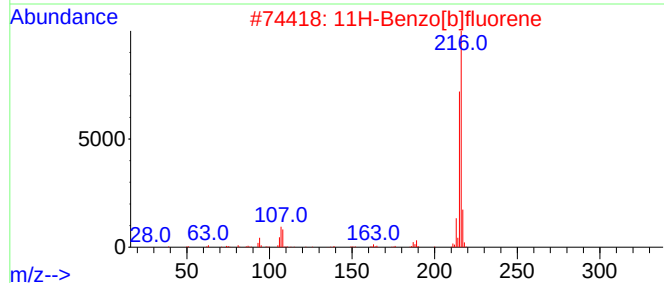
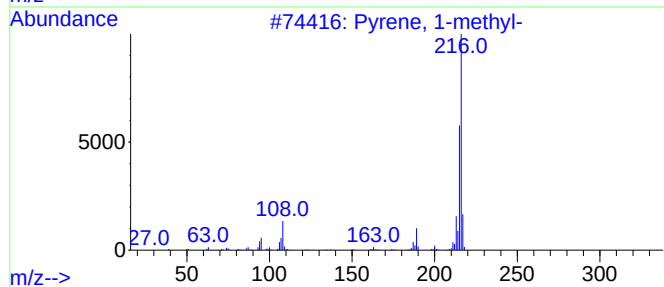
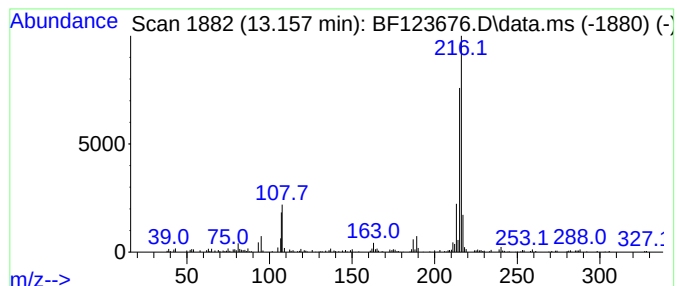
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Pyrene, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.157	2.33 ng	264740	Chrysene-d12	14.033

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
3			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	76
4			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	76
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042121\
Data File : BF123676.D
Acq On : 21 Apr 2021 20:06
Operator : JU/CG
Sample : M2090-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SEDGWICK-COMP-1

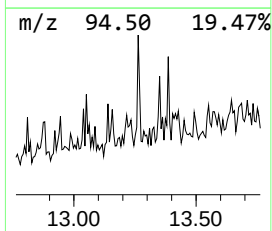
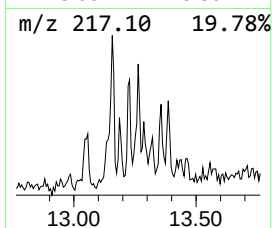
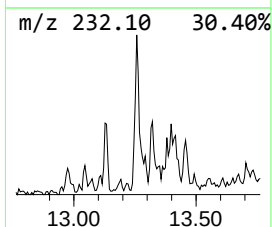
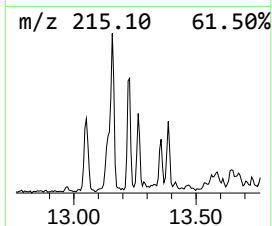
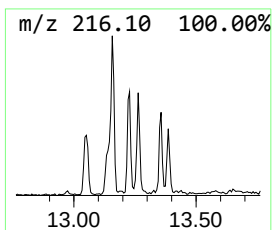
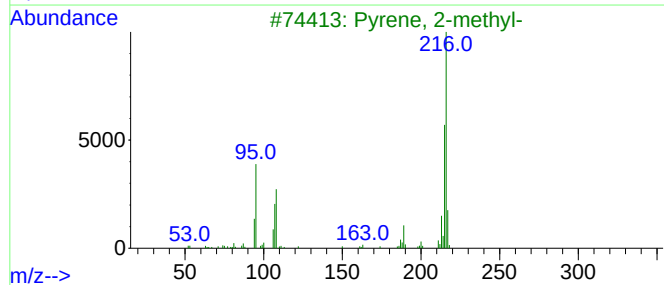
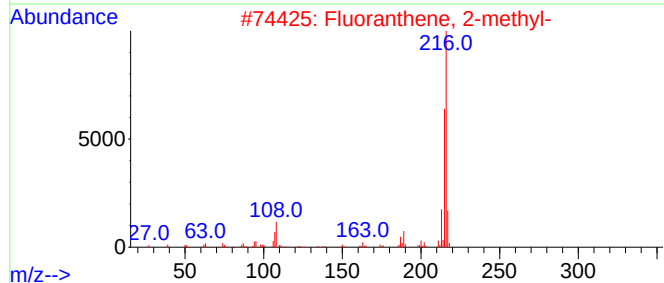
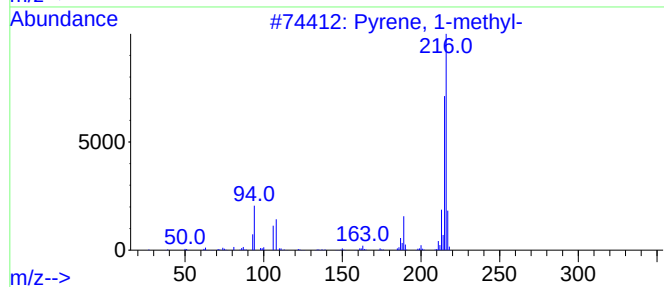
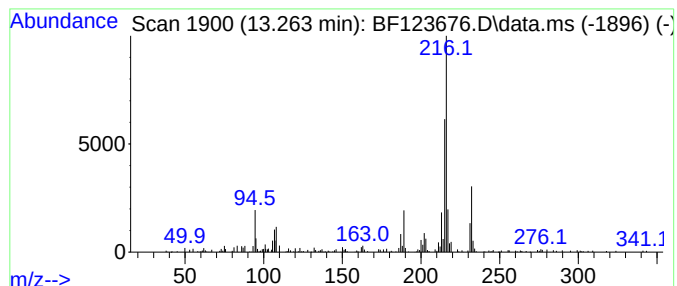
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Fluoranthene, 2-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.263	2.10 ng	239270	Chrysene-d12	14.033

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	96
2			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	93
3			Pyrene, 2-methyl-	216	C17H12	003442-78-2	90
4			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	64
5			Pyrene, 4-methyl-	216	C17H12	003353-12-6	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042121\
Data File : BF123676.D
Acq On : 21 Apr 2021 20:06
Operator : JU/CG
Sample : M2090-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SEDGWICK-COMP-1

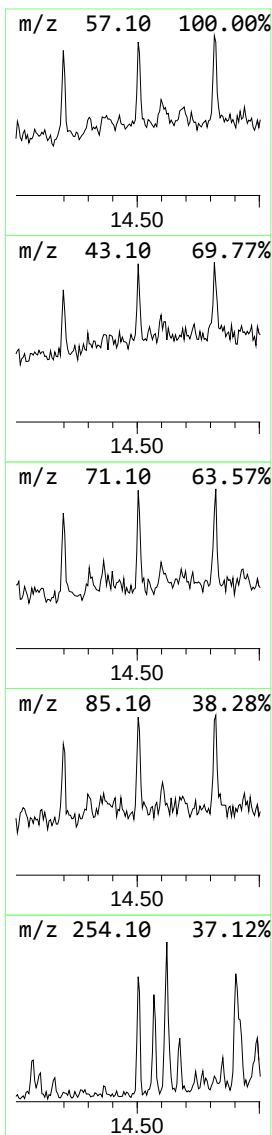
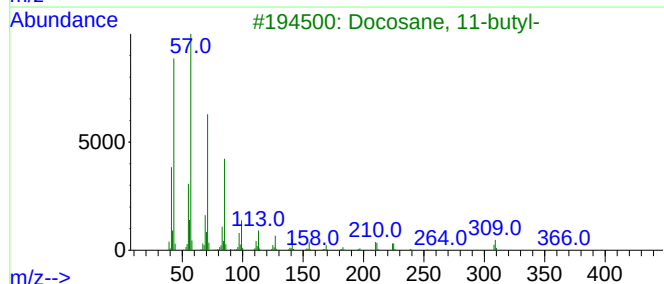
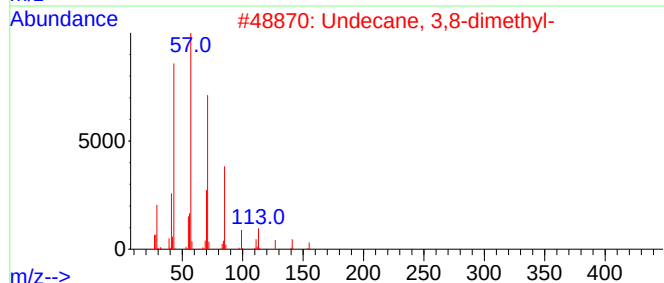
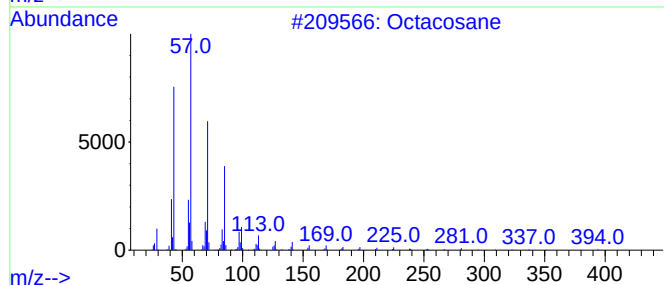
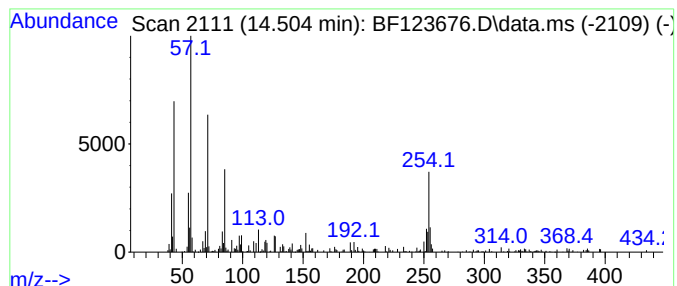
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Octacosane Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.504	2.05 ng	233252	Chrysene-d12	14.033

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	58
2			Undecane, 3,8-dimethyl-	184	C13H28	017301-30-3	52
3			Docosane, 11-butyl-	366	C26H54	013475-76-8	52
4			Nonadecane, 2,3-dimethyl-	296	C21H44	075163-99-4	49
5			Hexadecane	226	C16H34	000544-76-3	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042121\
Data File : BF123676.D
Acq On : 21 Apr 2021 20:06
Operator : JU/CG
Sample : M2090-03 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SEDGWICK-COMP-1

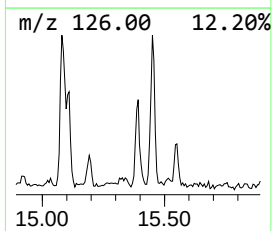
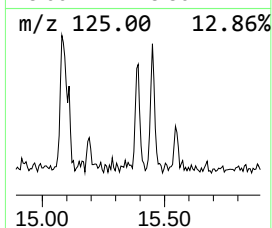
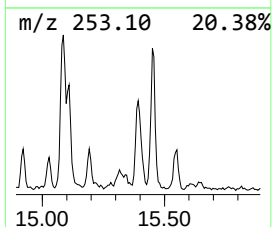
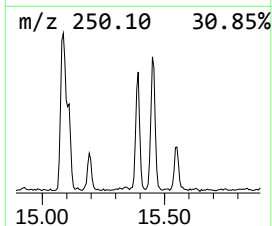
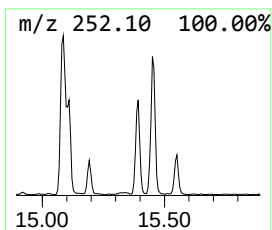
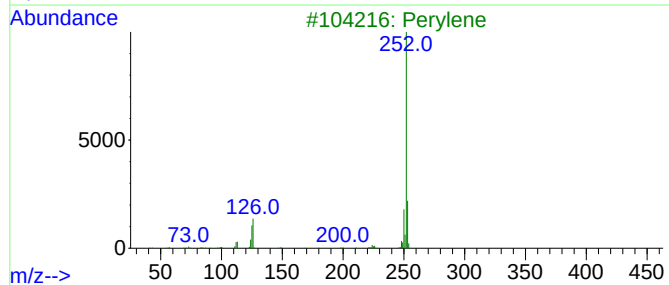
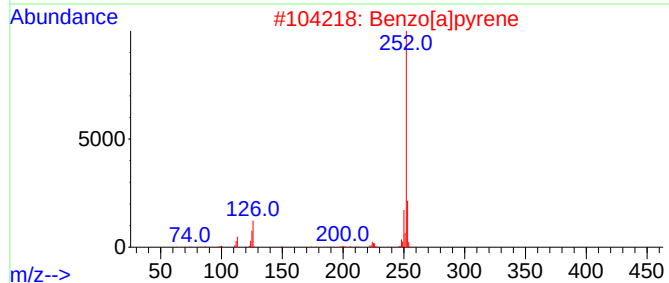
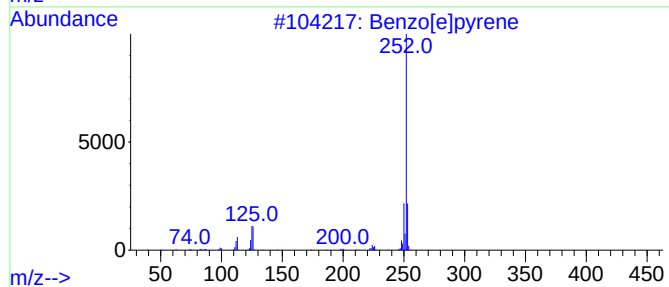
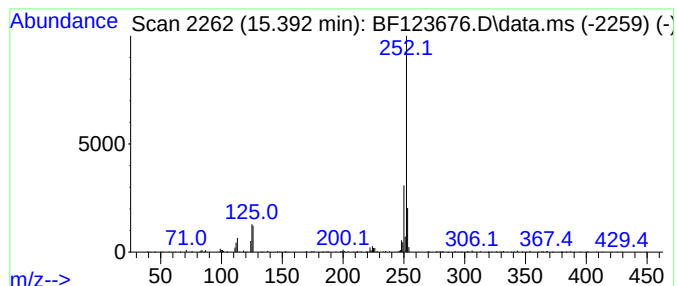
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.392	10.62 ng	600922	Perylene-d12	15.521

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			Perylene	252	C20H12	000198-55-0	96
4			Benz[e]acephenanthrylene	252	C20H12	000205-99-2	96
5			Benzo[j]fluoranthene	252	C20H12	000205-82-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042121\
 Data File : BF123676.D
 Acq On : 21 Apr 2021 20:06
 Operator : JU/CG
 Sample : M2090-03 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SEDGWICK-COMP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF041521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.046	3.2	ng	198259	1	6.851	1245300	20.0
Tetradecane	10.804	2.3	ng	154735	4	11.381	1347520	20.0
Tridecane, 7-pr...	11.251	2.2	ng	147615	4	11.381	1347520	20.0
Dibenzothiophene	11.281	2.5	ng	165138	4	11.381	1347520	20.0
Hexadecanoic ac...	11.781	6.8	ng	456484	4	11.381	1347520	20.0
Anthracene, 2-m...	11.886	3.1	ng	206772	4	11.381	1347520	20.0
Phenanthrene, 3...	11.916	4.4	ng	295235	4	11.381	1347520	20.0
4H-Cyclopenta[d...	11.998	6.0	ng	402494	4	11.381	1347520	20.0
Phenanthrene, 1...	12.022	2.4	ng	161913	4	11.381	1347520	20.0
5,16[1',2']:8,1...	12.180	2.1	ng	140138	4	11.381	1347520	20.0
9,10-Dimethylan...	12.439	2.5	ng	169770	4	11.381	1347520	20.0
9,12-Octadecadi...	12.469	24.1	ng	1620640	4	11.381	1347520	20.0
9-Octadecenoic ...	12.486	15.5	ng	1047070	4	11.381	1347520	20.0
Cyclopenta(def)...	12.522	2.2	ng	146874	4	11.381	1347520	20.0
11H-Benzo[b]flu...	13.051	2.4	ng	268104	5	14.033	2277090	20.0
Pyrene, 1-methyl-	13.157	2.3	ng	264740	5	14.033	2277090	20.0
Fluoranthene, 2...	13.263	2.1	ng	239270	5	14.033	2277090	20.0
Octacosane	14.504	2.0	ng	233252	5	14.033	2277090	20.0
Benzo[e]pyrene	15.392	10.6	ng	600922	6	15.521	1131560	20.0