

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052225\
 Data File : BF142510.D
 Acq On : 22 May 2025 15:07
 Operator : RC/JU
 Sample : Q2075-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS-10

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-BF052025.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF142510.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.110	487	496	506	rVB	131287	198901	10.61%	0.897%
2	5.510	556	564	569	rBV	1364883	1785150	95.22%	8.048%
3	6.516	727	735	743	rBV	1375595	1874773	100.00%	8.452%
4	6.728	766	771	778	rBV2	33210	46801	2.50%	0.211%
5	6.898	794	800	805	rVV	553928	690908	36.85%	3.115%
6	7.286	861	866	871	rBV4	11194	20638	1.10%	0.093%
7	7.457	885	895	899	rBV	853265	1142835	60.96%	5.153%
8	7.486	899	900	904	rVV	35365	35319	1.88%	0.159%
9	7.710	933	938	944	rVB2	34766	55541	2.96%	0.250%
10	8.180	1005	1018	1024	rBV	718443	995565	53.10%	4.489%
11	8.233	1024	1027	1030	rVB2	18718	23918	1.28%	0.108%
12	8.392	1049	1054	1062	rVB6	16723	28262	1.51%	0.127%
13	8.469	1062	1067	1072	rBV5	17486	30476	1.63%	0.137%
14	8.592	1085	1088	1095	rVB	17511	23623	1.26%	0.107%
15	8.680	1100	1103	1107	rBV4	14236	19093	1.02%	0.086%
16	8.769	1113	1118	1121	rBV	54805	70213	3.75%	0.317%
17	8.998	1152	1157	1161	rBV3	16473	27143	1.45%	0.122%
18	9.198	1188	1191	1195	rVV2	15472	20478	1.09%	0.092%
19	9.251	1195	1200	1205	rVV	1384060	1750396	93.37%	7.892%
20	9.333	1209	1214	1219	rBV	65716	90158	4.81%	0.406%
21	9.592	1251	1258	1260	rBV2	23444	42851	2.29%	0.193%
22	9.651	1264	1268	1271	rVV3	30235	50754	2.71%	0.229%
23	9.798	1289	1293	1297	rBV	35412	48802	2.60%	0.220%
24	9.863	1299	1304	1308	rVB	65170	82005	4.37%	0.370%
25	9.933	1308	1316	1321	rBV	750645	961994	51.31%	4.337%
26	10.139	1346	1351	1355	rVB	40179	60777	3.24%	0.274%
27	10.180	1355	1358	1363	rVB5	14760	20225	1.08%	0.091%
28	10.251	1367	1370	1372	rBV3	14394	19777	1.05%	0.089%
29	10.363	1385	1389	1393	rVB	67188	84552	4.51%	0.381%
30	10.480	1404	1409	1415	rBV	70908	98001	5.23%	0.442%
31	10.580	1422	1426	1432	rVB	26890	43513	2.32%	0.196%
32	10.645	1432	1437	1444	rVB	159145	207207	11.05%	0.934%
33	10.721	1444	1450	1455	rBV	833588	1053283	56.18%	4.749%
34	10.839	1465	1470	1478	rVB	77569	130276	6.95%	0.587%
35	11.027	1498	1502	1505	rBV3	20470	30025	1.60%	0.135%
36	11.227	1532	1536	1540	rVB	31961	40182	2.14%	0.181%

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Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052025.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	11.286	1540	1546	1548	rVV2	53648	82540	4.40%	0.372%
38	11.316	1548	1551	1557	rVB	59875	81823	4.36%	0.369%
39	11.421	1564	1569	1571	rBV	607329	802915	42.83%	3.620%
40	11.445	1571	1573	1578	rVV	627829	714495	38.11%	3.221%
41	11.498	1578	1582	1585	rVB	69865	91565	4.88%	0.413%
42	11.651	1602	1608	1614	rBV2	39841	72945	3.89%	0.329%
43	11.710	1614	1618	1622	rVV2	46645	65690	3.50%	0.296%
44	11.751	1622	1625	1631	rVB3	16228	24436	1.30%	0.110%
45	11.945	1650	1658	1664	rBV2	193241	364434	19.44%	1.643%
46	12.039	1669	1674	1681	rVB2	109632	211533	11.28%	0.954%
47	12.121	1684	1688	1691	rBV	30250	38345	2.05%	0.173%
48	12.221	1698	1705	1706	rBV	46161	68163	3.64%	0.307%
49	12.239	1706	1708	1713	rVB	51107	60566	3.23%	0.273%
50	12.474	1745	1748	1751	rBV	26664	31980	1.71%	0.144%
51	12.510	1751	1754	1759	rVB2	34420	45615	2.43%	0.206%
52	12.563	1759	1763	1768	rVB	27947	39283	2.10%	0.177%
53	12.639	1771	1776	1780	rBV	767368	947690	50.55%	4.273%
54	12.680	1780	1783	1788	rVB3	33966	48496	2.59%	0.219%
55	12.733	1788	1792	1795	rBV	36540	49046	2.62%	0.221%
56	12.868	1807	1815	1820	rBV	561045	836752	44.63%	3.773%
57	12.910	1820	1822	1828	rVB2	31608	44869	2.39%	0.202%
58	13.010	1831	1839	1846	rVB	999617	1392481	74.27%	6.278%
59	13.080	1846	1851	1856	rBV2	45049	90816	4.84%	0.409%
60	13.192	1863	1870	1874	rVB	72640	132420	7.06%	0.597%
61	13.257	1878	1881	1884	rVV	52847	72914	3.89%	0.329%
62	13.298	1884	1888	1895	rVB2	57159	88988	4.75%	0.401%
63	13.392	1901	1904	1906	rBV	18843	20291	1.08%	0.091%
64	13.421	1906	1909	1914	rVB3	20431	24293	1.30%	0.110%
65	13.698	1952	1956	1961	rVB	49857	72264	3.85%	0.326%
66	13.810	1970	1975	1978	rBV2	56221	92946	4.96%	0.419%
67	13.839	1978	1980	1982	rVV	21671	22798	1.22%	0.103%
68	13.904	1987	1991	2001	rVB2	97372	158515	8.46%	0.715%
69	14.062	2010	2018	2021	rBV2	660070	1156663	61.70%	5.215%
70	14.204	2039	2042	2048	rBV2	31467	55996	2.99%	0.252%
71	14.445	2080	2083	2087	rVV	38073	50229	2.68%	0.226%
72	14.486	2087	2090	2093	rVB2	31660	37829	2.02%	0.171%
73	14.574	2102	2105	2107	rBV2	20080	25505	1.36%	0.115%
74	15.115	2191	2197	2206	rBV	266137	619342	33.04%	2.792%
75	15.221	2212	2215	2220	rVB	36221	49775	2.65%	0.224%
76	15.427	2245	2250	2255	rVB	129319	208568	11.12%	0.940%

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Instrument :
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ClientSampleId :
SS-10

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

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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	15.486	2255	2260	2266	rVV	152085	231671	12.36%	1.044%
78	15.556	2266	2272	2284	rVB2	335887	624560	33.31%	2.816%
79	17.050	2520	2526	2535	rVB2	75671	171453	9.15%	0.773%
80	17.298	2564	2568	2572	rBV2	11149	21455	1.14%	0.097%
81	17.503	2596	2603	2616	rVB	55656	130760	6.97%	0.590%

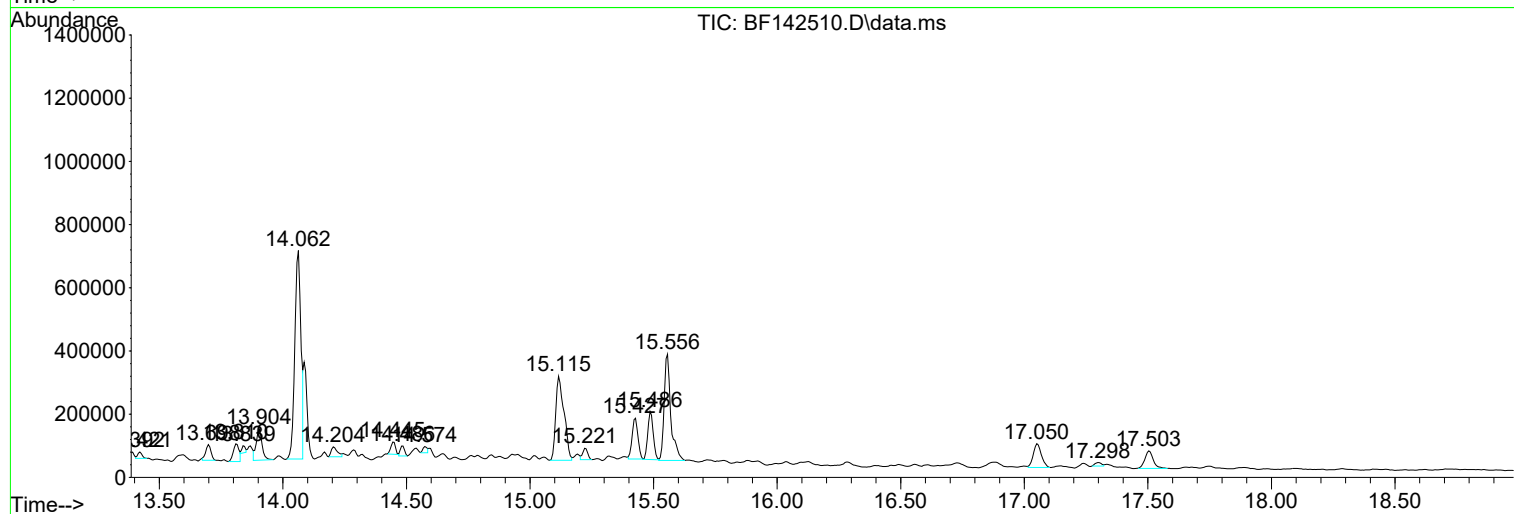
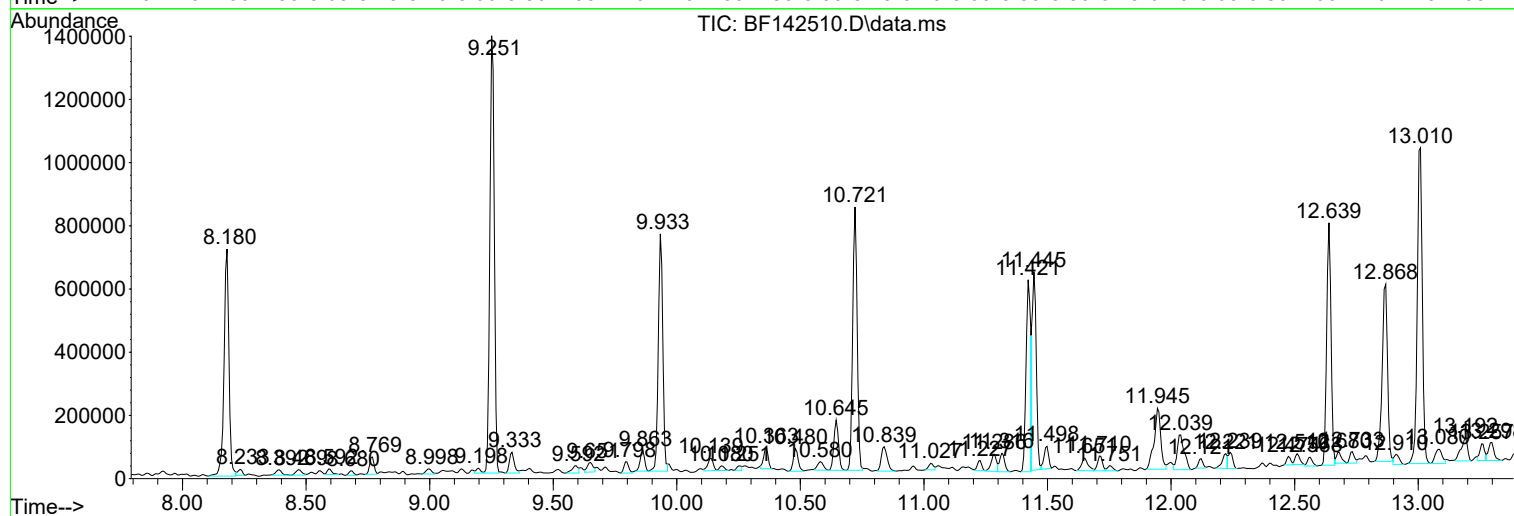
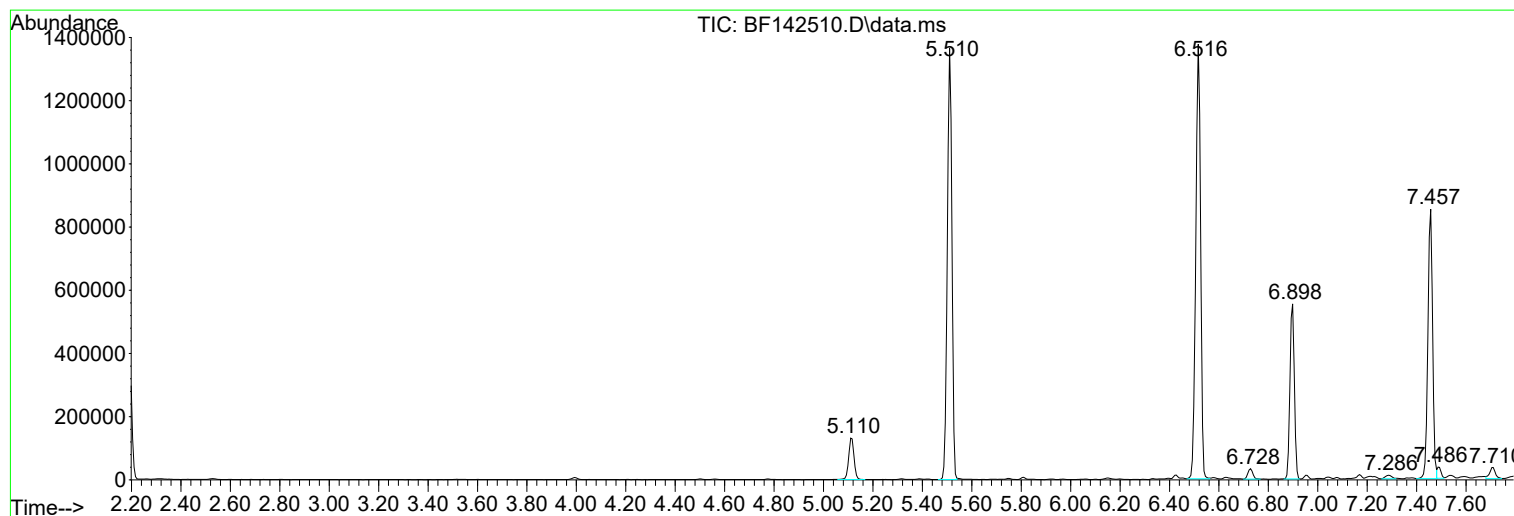
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Misc :
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Instrument :
BNA_F
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052225\
Data File : BF142510.D
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Instrument :
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SS-10

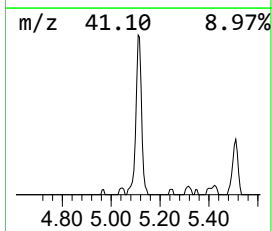
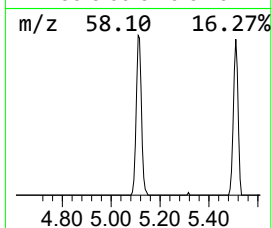
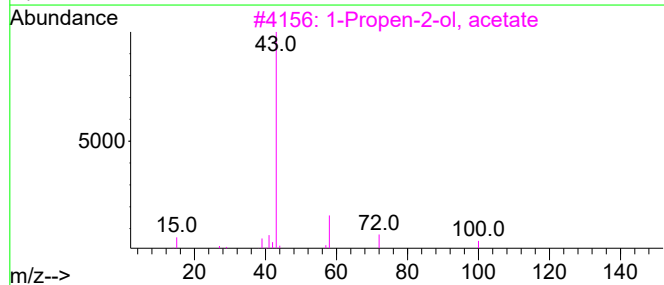
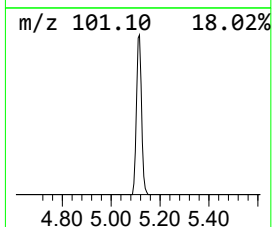
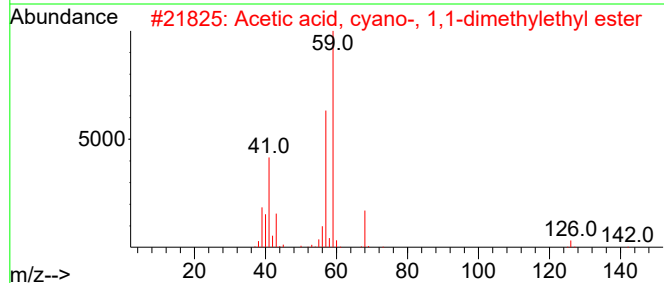
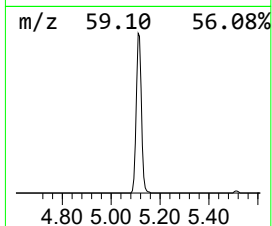
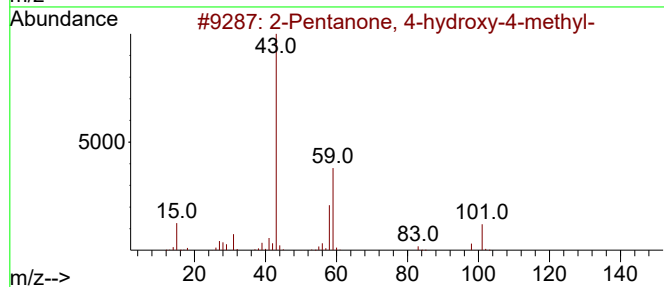
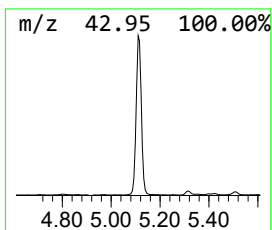
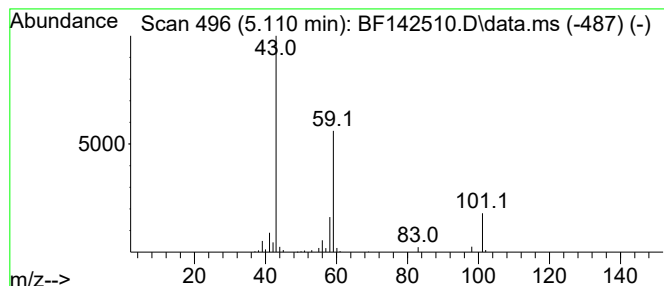
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.110	5.76 ng	198901	1,4-Dichlorobenzene-d4	6.898

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59		
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25		
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12		
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9		
5	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	9		



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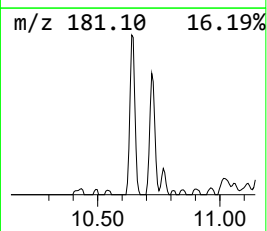
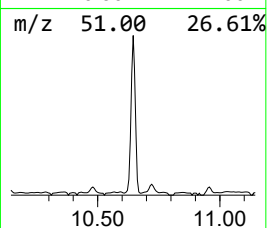
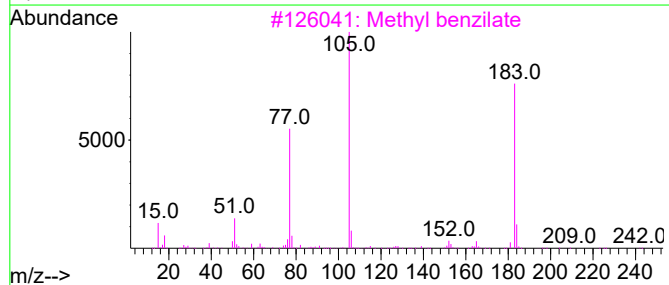
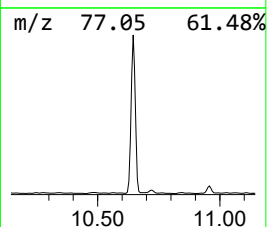
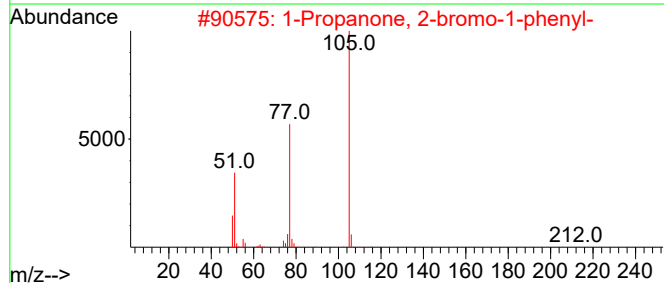
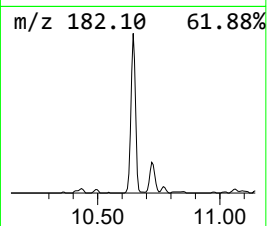
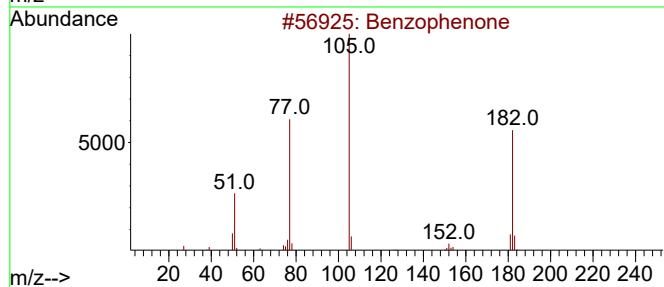
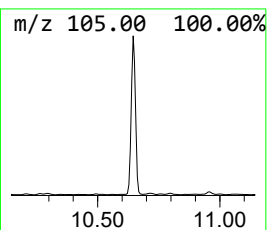
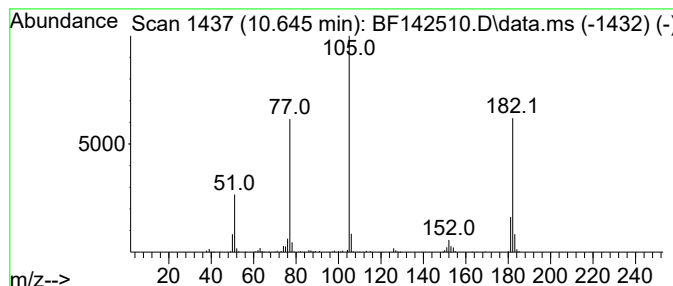
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.645	4.31 ng	207207	Acenaphthene-d10	9.933

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			1-Propanone, 2-bromo-1-phenyl-	212	C9H9BrO	002114-00-3	47
3			Methyl benzilate	242	C15H14O3	000076-89-1	47
4			acetic acid, 2-bromo-, 2-oxo-2-p...	256	C10H9BrO3	1000401-50-0	47
5			Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	46



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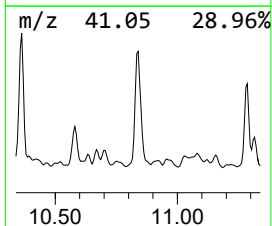
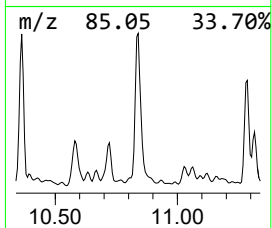
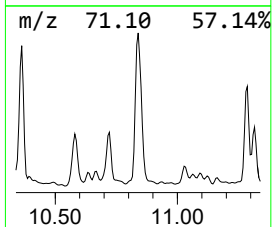
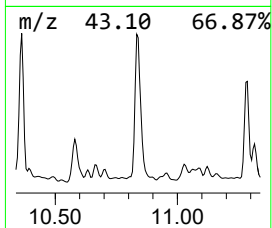
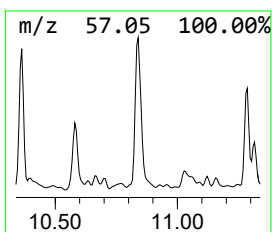
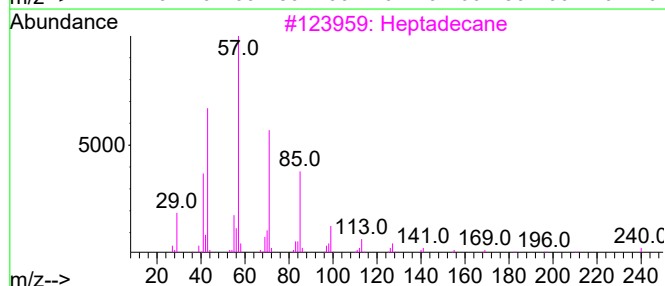
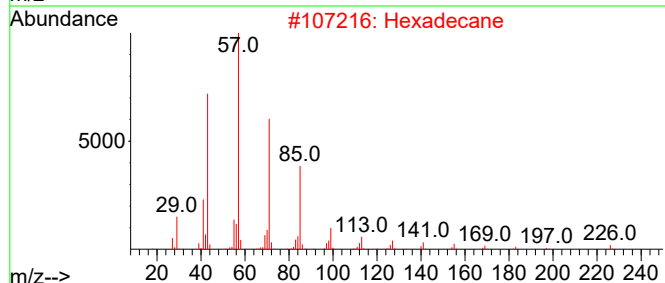
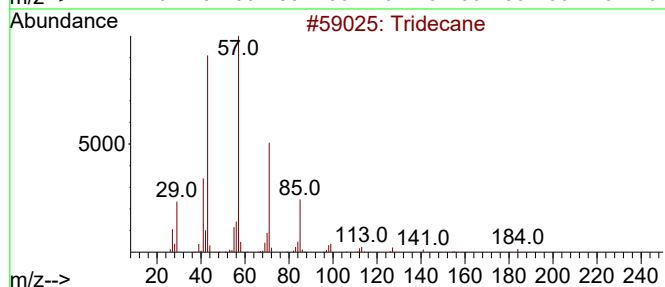
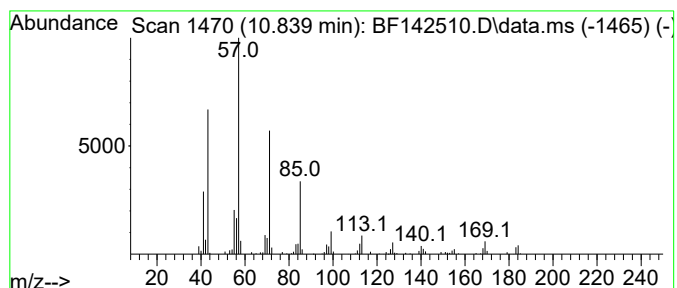
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Tridecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.839	3.25 ng	130276	Phenanthrene-d10	11.421

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	87
2			Hexadecane	226	C16H34	000544-76-3	86
3			Heptadecane	240	C17H36	000629-78-7	86
4			Tridecane, 3-methyl-	198	C14H30	006418-41-3	86
5			Eicosane	282	C20H42	000112-95-8	86



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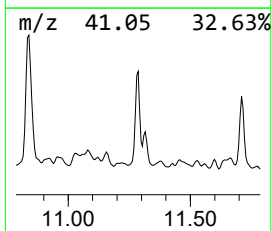
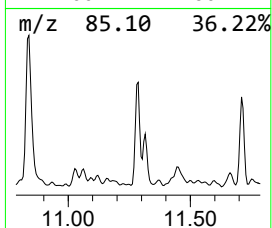
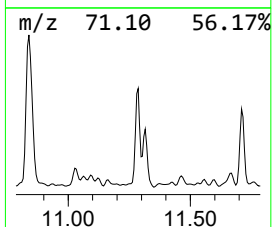
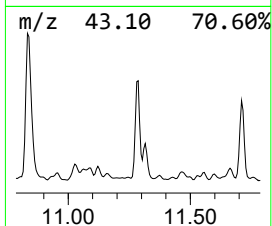
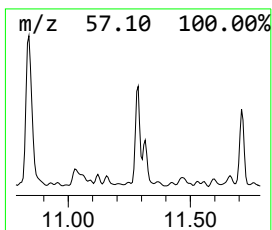
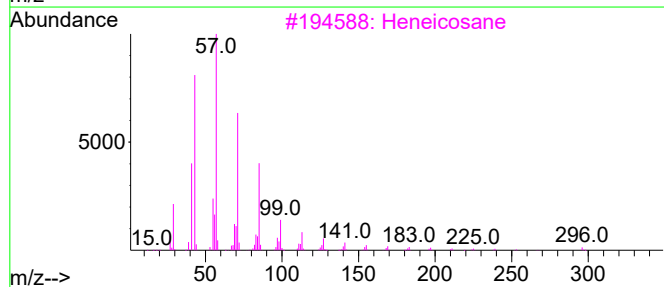
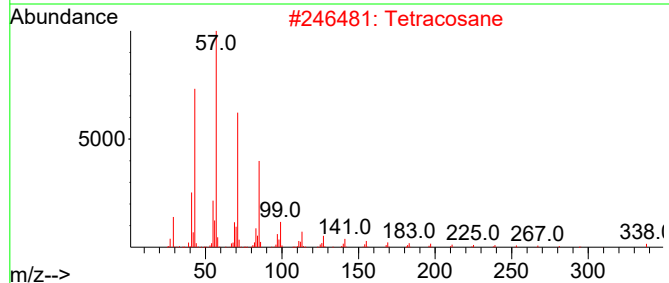
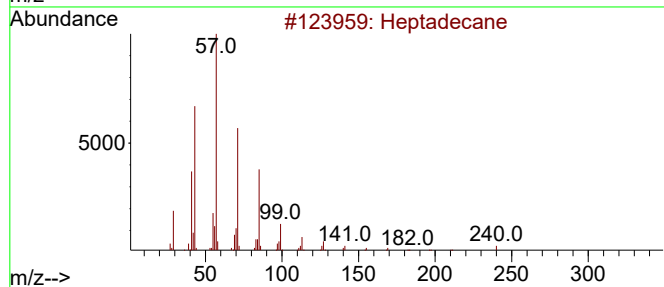
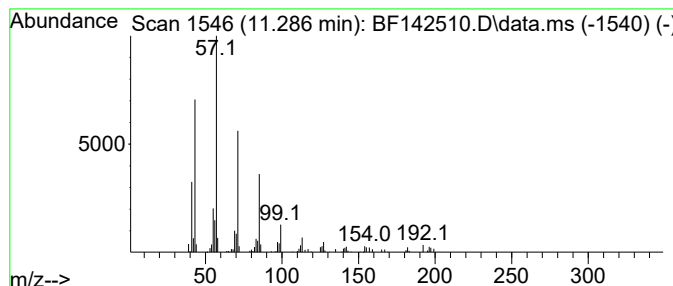
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Heptadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.286	2.06 ng	82540	Phenanthrene-d10	11.421

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	87
2			Tetracosane	338	C24H50	000646-31-1	87
3			Heneicosane	296	C21H44	000629-94-7	86
4			Hexadecane	226	C16H34	000544-76-3	86
5			Pentadecane	212	C15H32	000629-62-9	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052225\
Data File : BF142510.D
Acq On : 22 May 2025 15:07
Operator : RC/JU
Sample : Q2075-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SS-10

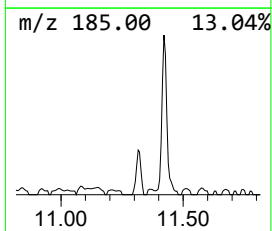
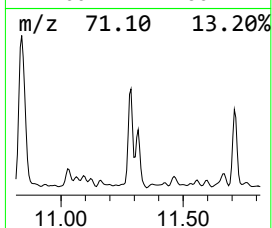
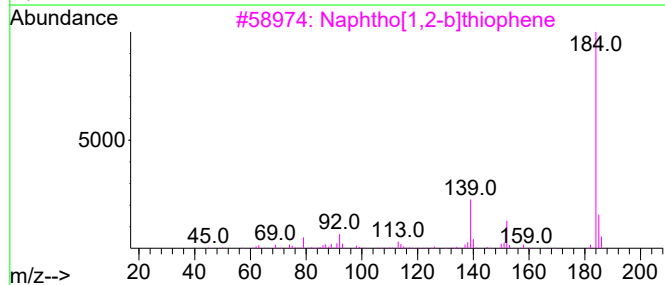
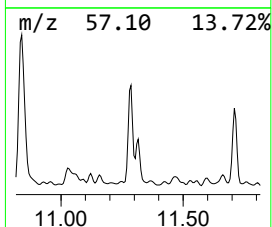
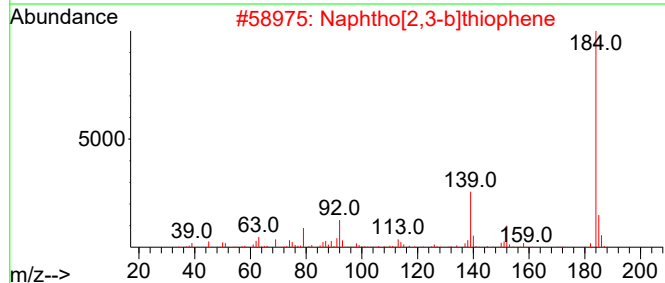
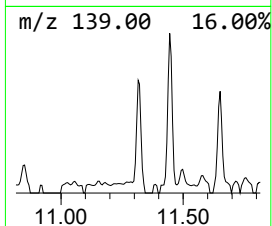
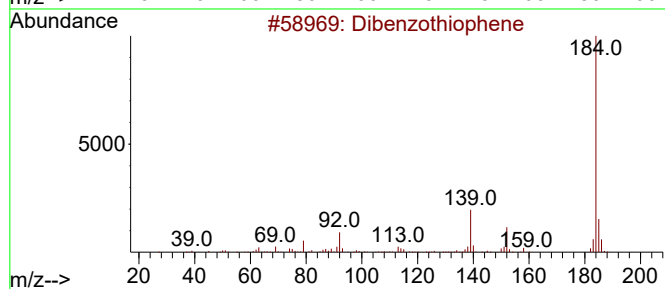
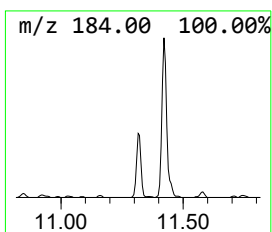
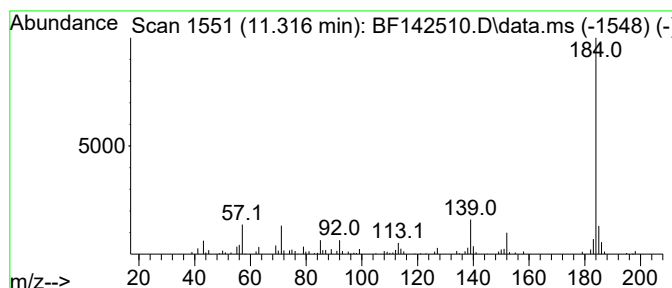
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Dibenzothiophene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.316	2.04 ng	81823	Phenanthrene-d10	11.421

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dibenzothiophene	184	C12H8S	000132-65-0	95
2		Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	95
3		Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	93
4		Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	91
5		Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052225\
Data File : BF142510.D
Acq On : 22 May 2025 15:07
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Sample : Q2075-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

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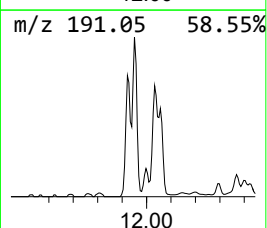
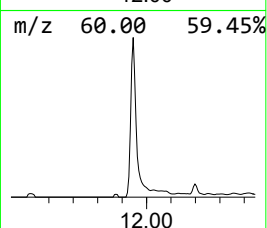
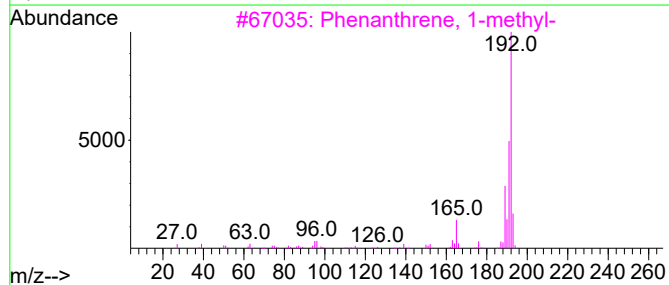
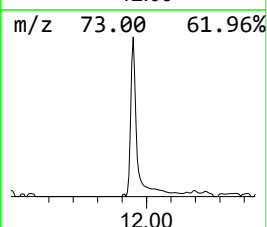
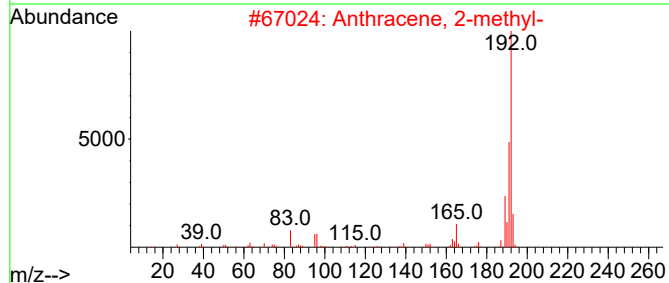
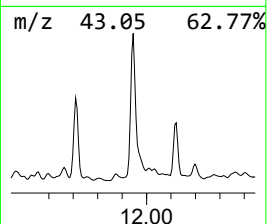
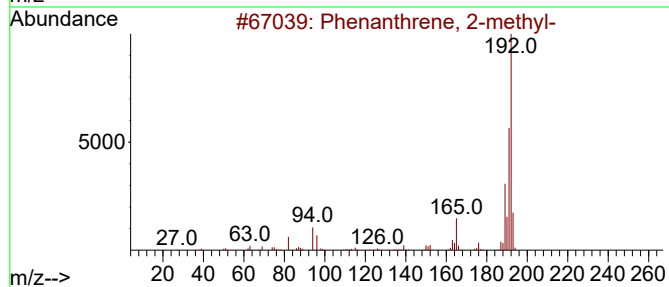
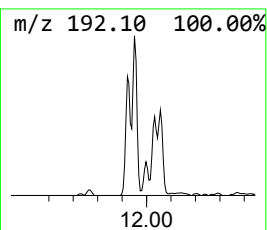
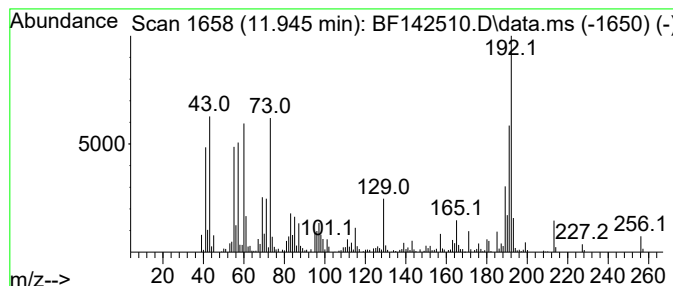
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Phenanthrene, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.945	9.08 ng	364434	Phenanthrene-d10	11.421

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
4			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	95
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052225\
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Acq On : 22 May 2025 15:07
Operator : RC/JU
Sample : Q2075-01
Misc :
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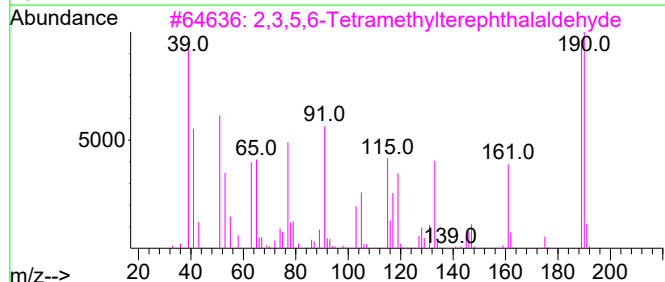
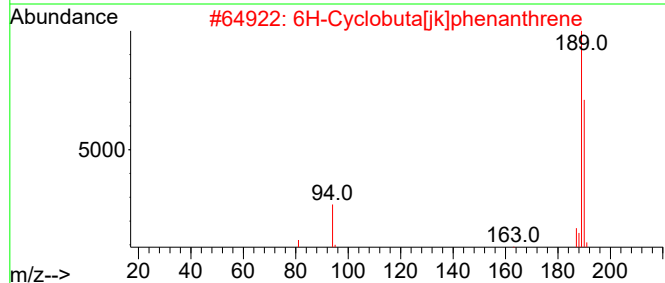
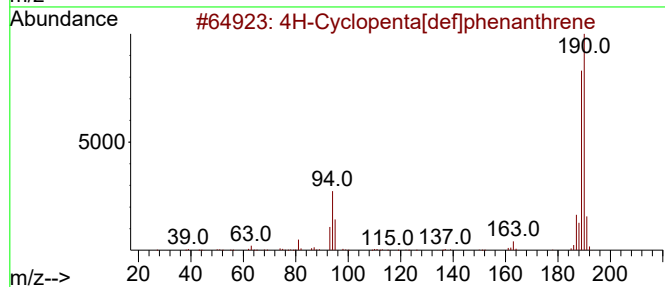
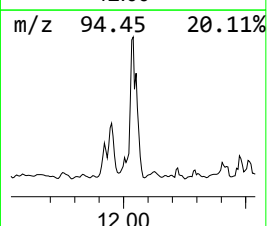
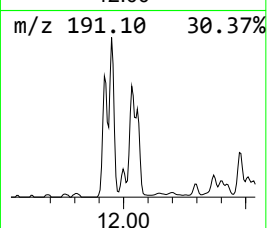
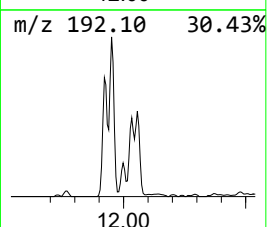
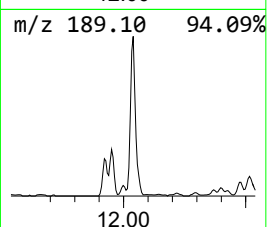
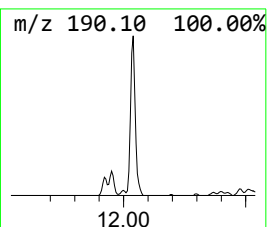
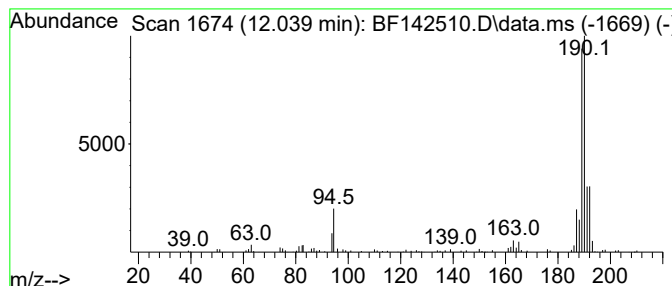
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.039	5.27 ng	211533	Phenanthrene-d10	11.421	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	72
3	2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	50
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45
5	1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052225\
Data File : BF142510.D
Acq On : 22 May 2025 15:07
Operator : RC/JU
Sample : Q2075-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

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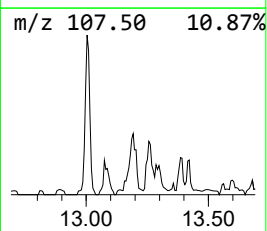
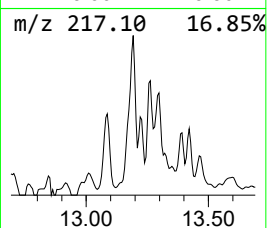
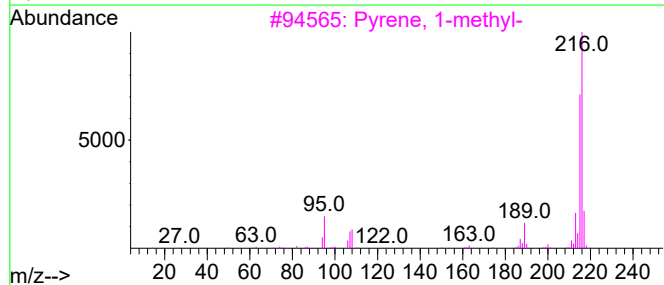
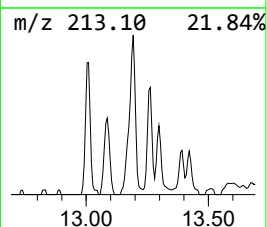
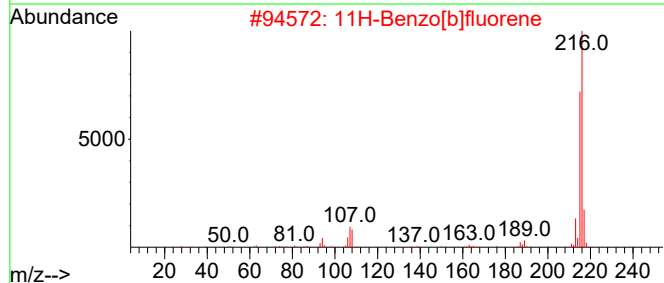
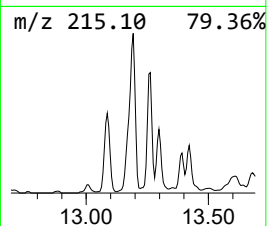
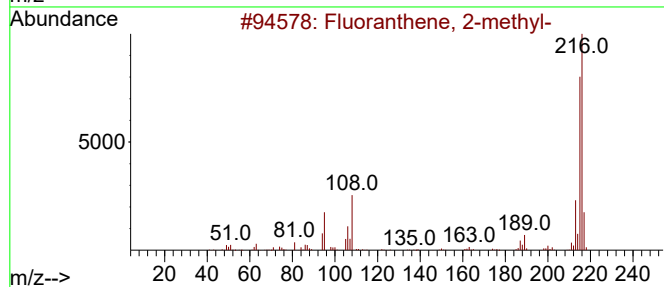
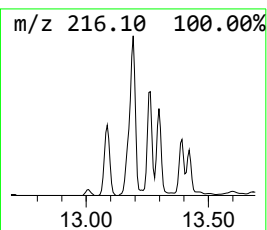
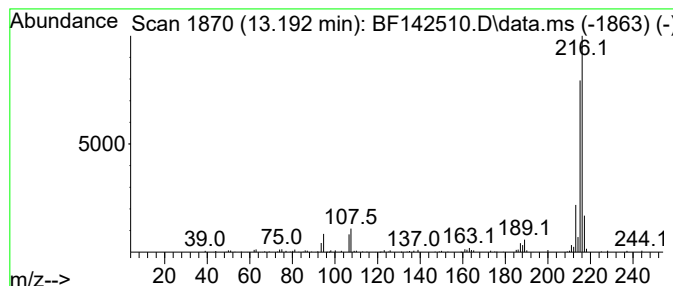
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Fluoranthene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.192	2.29 ng	132420	Chrysene-d12	14.062

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	95
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
3			Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
4			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
5			Pyrene, 2-methyl-	216	C17H12	003442-78-2	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052225\
Data File : BF142510.D
Acq On : 22 May 2025 15:07
Operator : RC/JU
Sample : Q2075-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

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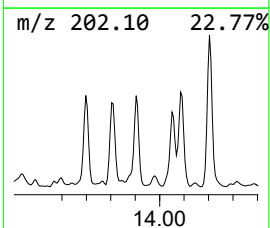
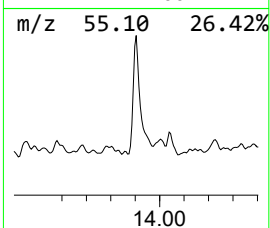
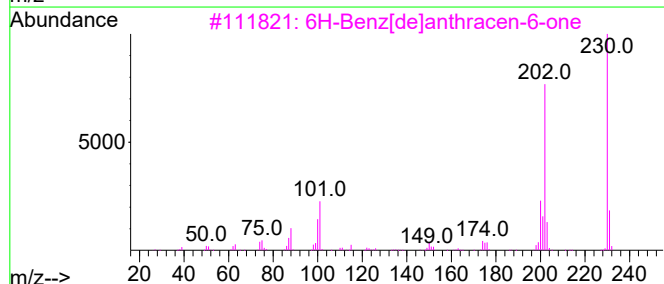
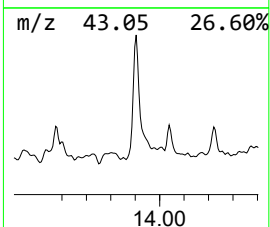
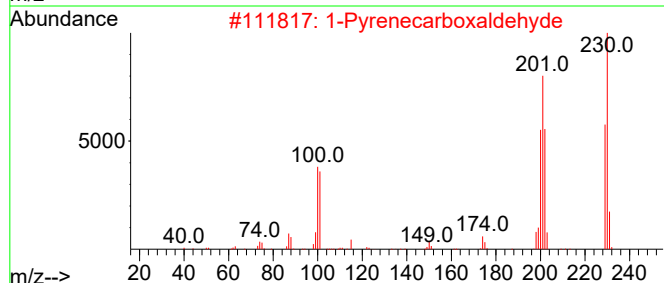
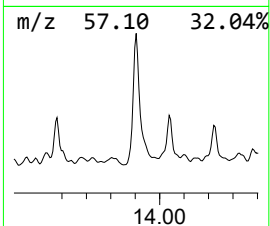
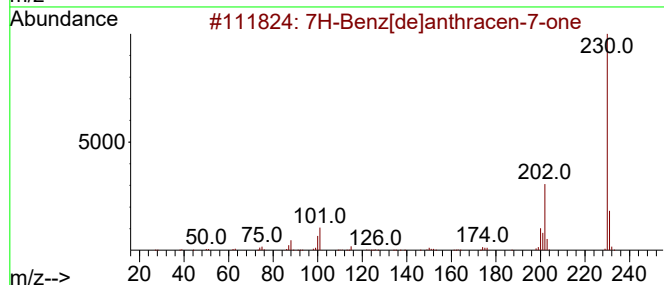
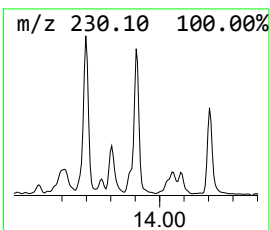
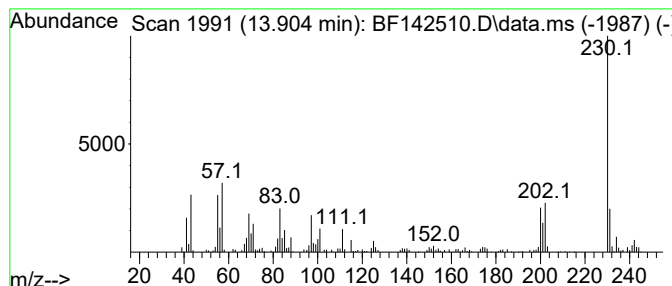
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 7H-Benz[de]anthracen-7-one Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.904	2.74 ng	158515	Chrysene-d12	14.062

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	83
2			1-Pyrenecarboxaldehyde	230	C17H10O	003029-19-4	59
3			6H-Benz[de]anthracen-6-one	230	C17H10O	080252-14-8	58
4			o-Terphenyl	230	C18H14	000084-15-1	53
5			Pyrene, 1,3-dimethyl-	230	C18H14	064401-21-4	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052225\
Data File : BF142510.D
Acq On : 22 May 2025 15:07
Operator : RC/JU
Sample : Q2075-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

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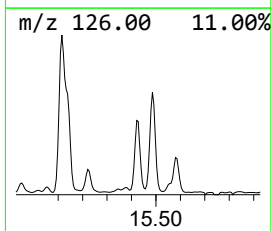
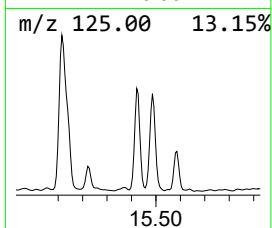
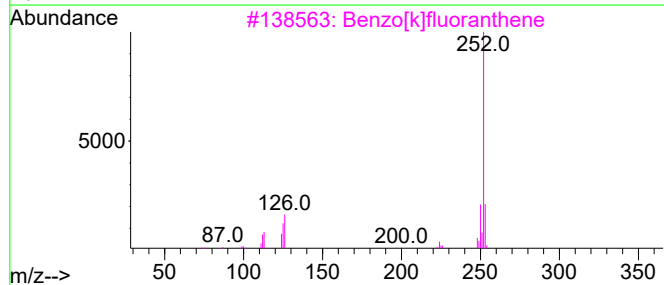
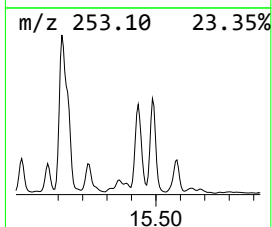
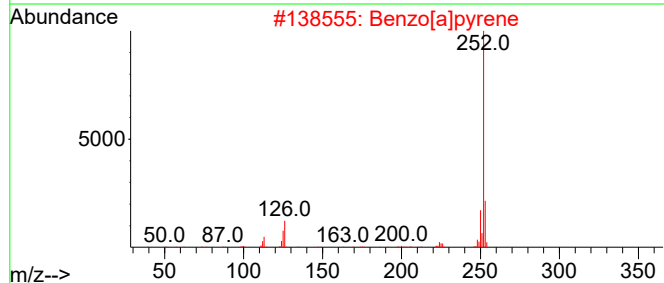
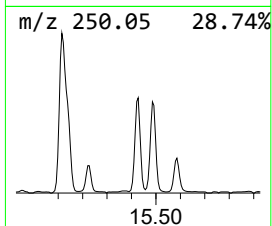
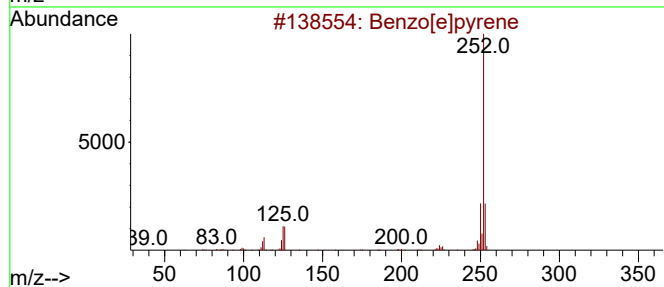
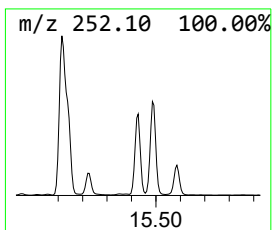
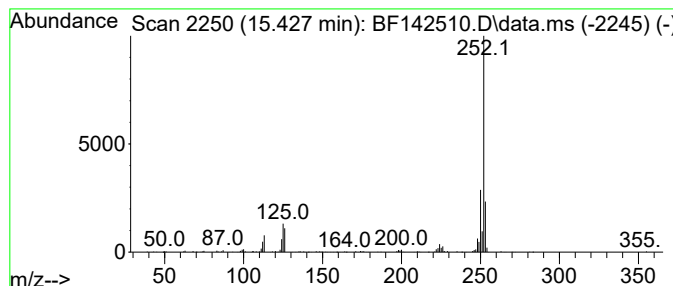
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.427	6.68 ng	208568	Perylene-d12	15.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	99
2			Benzo[a]pyrene	252	C20H12	000050-32-8	96
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
4			Benzo[j]fluoranthene	252	C20H12	000205-82-3	95
5			Benzo[b]fluoranthene	252	C20H12	000205-99-2	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052225\
Data File : BF142510.D
Acq On : 22 May 2025 15:07
Operator : RC/JU
Sample : Q2075-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SS-10

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.110	5.8	ng	198901	1	6.898	690908	20.0
Benzophenone	10.645	4.3	ng	207207	3	9.933	961994	20.0
Tridecane	10.839	3.3	ng	130276	4	11.421	802915	20.0
Heptadecane	11.286	2.1	ng	82540	4	11.421	802915	20.0
Dibenzothiophene	11.316	2.0	ng	81823	4	11.421	802915	20.0
Phenanthrene, 2...	11.945	9.1	ng	364434	4	11.421	802915	20.0
4H-Cyclopenta[d...	12.039	5.3	ng	211533	4	11.421	802915	20.0
Fluoranthene, 2...	13.192	2.3	ng	132420	5	14.062	1156660	20.0
7H-Benz[de]anth...	13.904	2.7	ng	158515	5	14.062	1156660	20.0
Benzo[e]pyrene	15.427	6.7	ng	208568	6	15.557	624560	20.0