

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102125\
 Data File : BF144024.D
 Acq On : 21 Oct 2025 20:56
 Operator : RC/JU
 Sample : Q3403-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH-18

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-BF100625.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF144024.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.487	384	390	396	rBV	12899	22610	1.19%	0.133%
2	5.093	484	493	508	rVB	61691	111042	5.87%	0.655%
3	5.498	556	562	582	rBV	1412424	1876828	99.14%	11.062%
4	6.504	727	733	760	rBV	1410673	1893083	100.00%	11.158%
5	6.881	792	797	802	rBV	531871	649451	34.31%	3.828%
6	7.439	886	892	903	rBV	867381	1117061	59.01%	6.584%
7	7.698	932	936	943	rBV2	10400	19321	1.02%	0.114%
8	8.163	1010	1015	1032	rBV	631808	812113	42.90%	4.787%
9	9.239	1192	1198	1208	rBV	1364442	1788339	94.47%	10.541%
10	9.575	1251	1255	1258	rBV	14943	20552	1.09%	0.121%
11	9.781	1285	1290	1294	rVB	33883	41574	2.20%	0.245%
12	9.922	1306	1314	1327	rBV	596962	791250	41.80%	4.664%
13	10.633	1430	1435	1442	rBV	189815	241646	12.76%	1.424%
14	10.710	1442	1448	1459	rBV	625831	851709	44.99%	5.020%
15	10.833	1464	1469	1476	rVB6	12157	20954	1.11%	0.124%
16	11.016	1495	1500	1503	rBV3	12176	19824	1.05%	0.117%
17	11.410	1562	1567	1577	rBV2	521849	810403	42.81%	4.777%
18	11.745	1620	1624	1629	rVB	16526	22119	1.17%	0.130%
19	11.939	1648	1657	1663	rBV3	193192	382232	20.19%	2.253%
20	12.022	1668	1671	1674	rBV2	63168	85478	4.52%	0.504%
21	12.210	1698	1703	1705	rBV2	21249	32140	1.70%	0.189%
22	12.233	1705	1707	1714	rVB	21472	31152	1.65%	0.184%
23	12.357	1722	1728	1731	rBV2	19070	38355	2.03%	0.226%
24	12.392	1731	1734	1736	rVV2	21585	24304	1.28%	0.143%
25	12.463	1742	1746	1749	rBV	69037	97066	5.13%	0.572%
26	12.498	1749	1752	1758	rVV2	41557	76144	4.02%	0.449%
27	12.551	1758	1761	1769	rVB5	28929	49569	2.62%	0.292%
28	12.627	1769	1774	1778	rBV	123050	155733	8.23%	0.918%
29	12.721	1787	1790	1794	rBV2	26181	32309	1.71%	0.190%
30	12.857	1808	1813	1819	rBV	168845	249298	13.17%	1.469%
31	12.910	1819	1822	1826	rVB2	29346	39141	2.07%	0.231%
32	12.998	1832	1837	1845	rVB	1092601	1411115	74.54%	8.317%
33	13.074	1845	1850	1857	rBV3	40742	80902	4.27%	0.477%
34	13.180	1861	1868	1872	rVB2	43632	98622	5.21%	0.581%
35	13.227	1873	1876	1878	rBV4	14166	19916	1.05%	0.117%
36	13.292	1883	1887	1894	rVB	50900	78495	4.15%	0.463%

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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-BF100625.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	13.380	1898	1902	1905	rVV	41920	54477	2.88%	0.321%
38	13.416	1905	1908	1911	rVB	35620	43594	2.30%	0.257%
39	13.569	1931	1934	1936	rBV3	14598	20676	1.09%	0.122%
40	13.698	1949	1956	1961	rVB2	26950	53911	2.85%	0.318%
41	13.804	1969	1974	1981	rBV5	27306	70902	3.75%	0.418%
42	13.904	1987	1991	2001	rVB3	67558	131788	6.96%	0.777%
43	14.057	2011	2017	2028	rBV2	568551	963314	50.89%	5.678%
44	14.216	2041	2044	2053	rVB7	30017	56677	2.99%	0.334%
45	14.445	2080	2083	2086	rVB	34233	41346	2.18%	0.244%
46	14.480	2087	2089	2092	rVB	25597	24326	1.28%	0.143%
47	14.521	2093	2096	2101	rBV5	26610	51146	2.70%	0.301%
48	14.815	2143	2146	2153	rVB3	46526	84443	4.46%	0.498%
49	14.910	2159	2162	2166	rBV	68989	84027	4.44%	0.495%
50	15.110	2191	2196	2203	rBV	60162	127142	6.72%	0.749%
51	15.415	2244	2248	2254	rVB	43635	66133	3.49%	0.390%
52	15.480	2255	2259	2264	rVV	61532	93391	4.93%	0.550%
53	15.545	2264	2270	2292	rVB	506417	906567	47.89%	5.344%

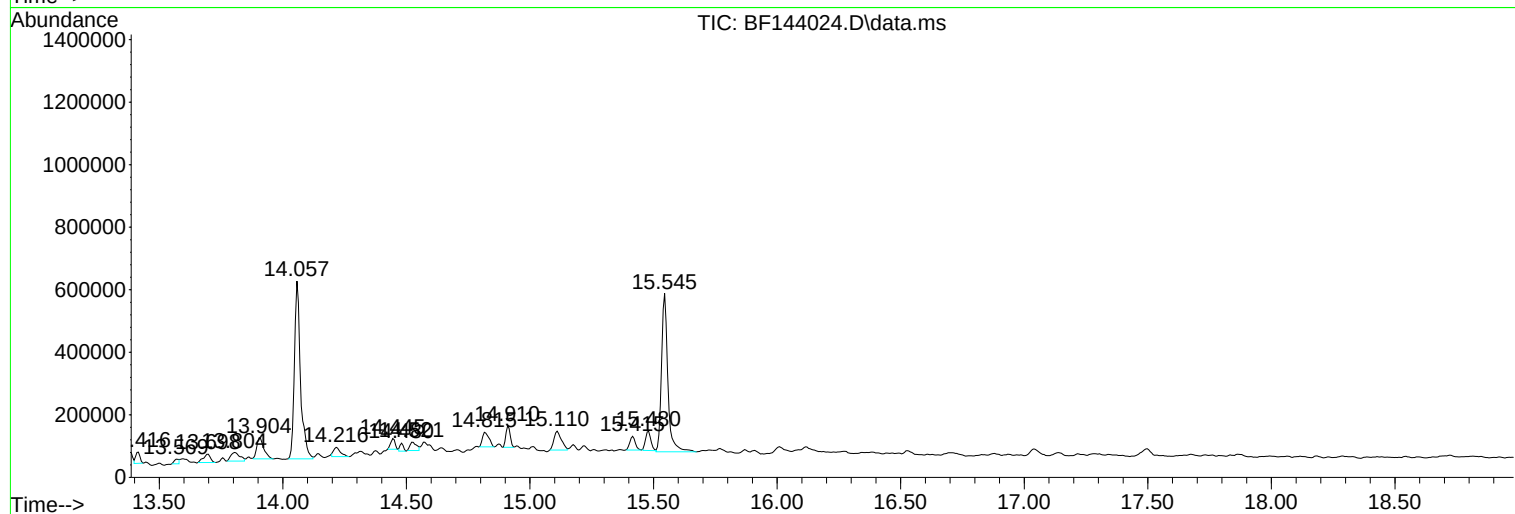
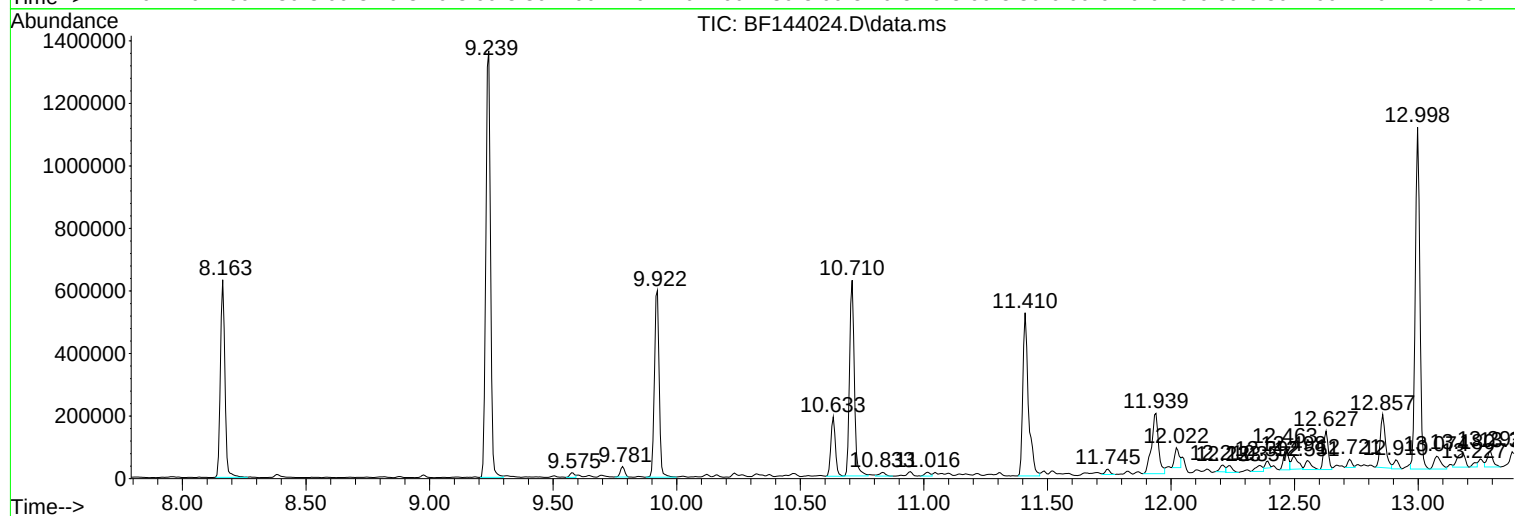
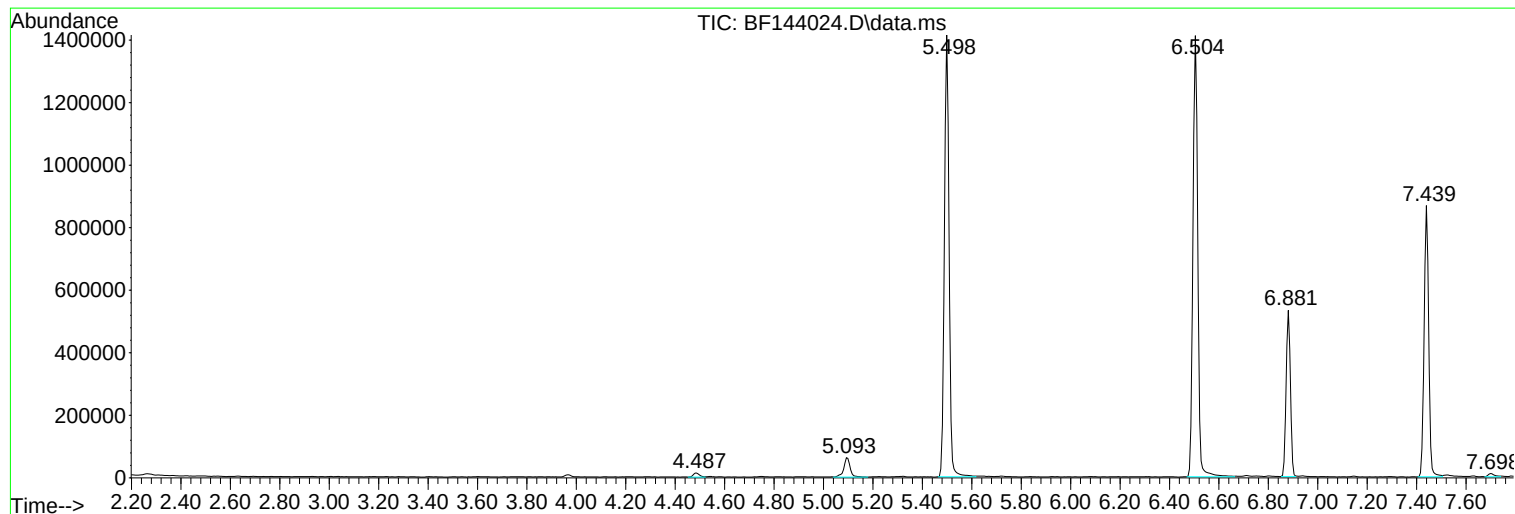
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100625.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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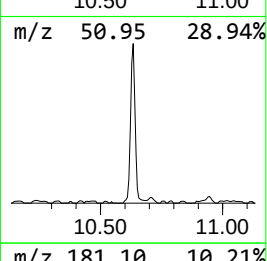
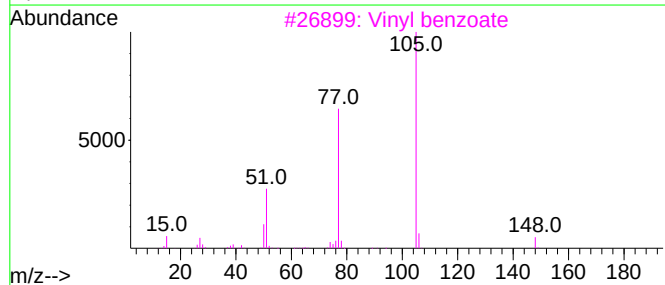
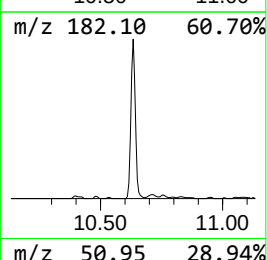
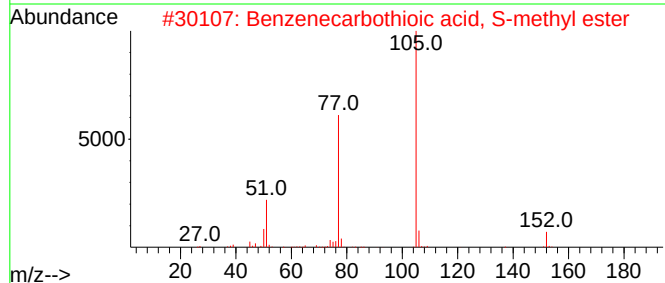
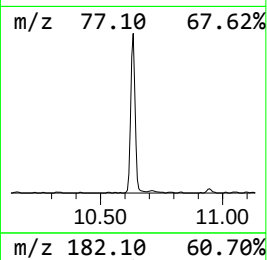
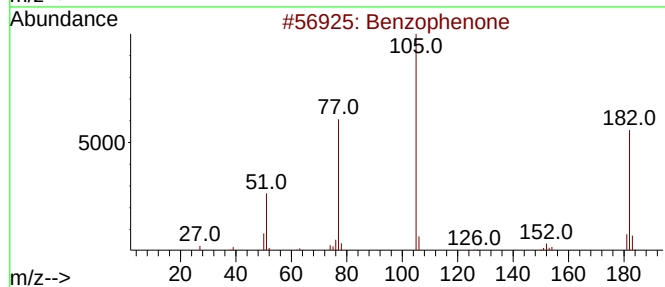
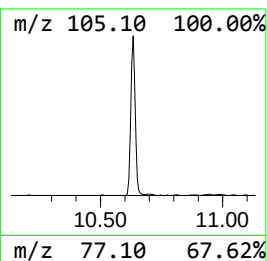
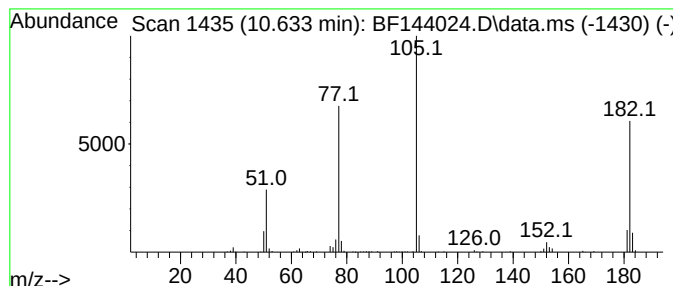
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.633	6.11 ng	241646	Acenaphthene-d10	9.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	49
3			Vinyl benzoate	148	C9H8O2	000769-78-8	47
4			Disulfide, dibenzoyl	274	C14H10O2S2	000644-32-6	47
5			1,2-Propanedione, 1-phenyl-	148	C9H8O2	000579-07-7	47



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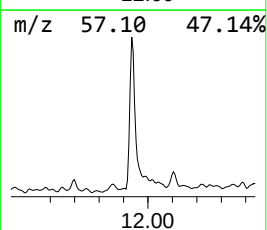
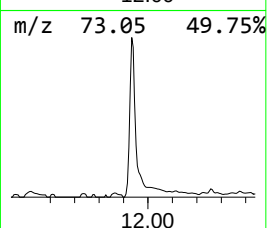
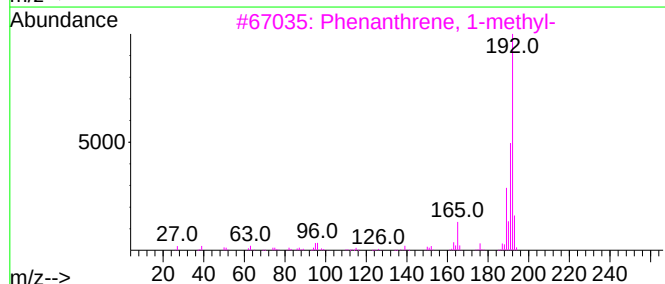
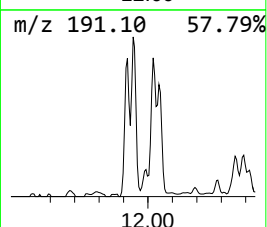
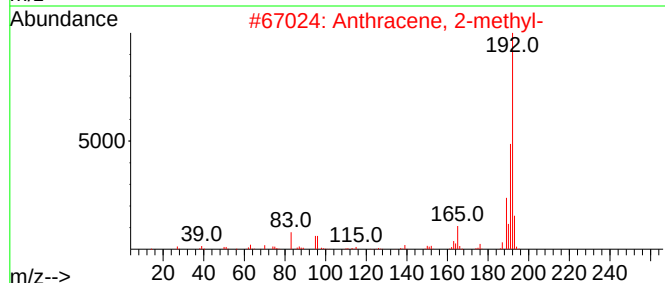
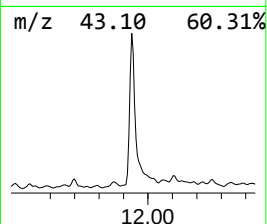
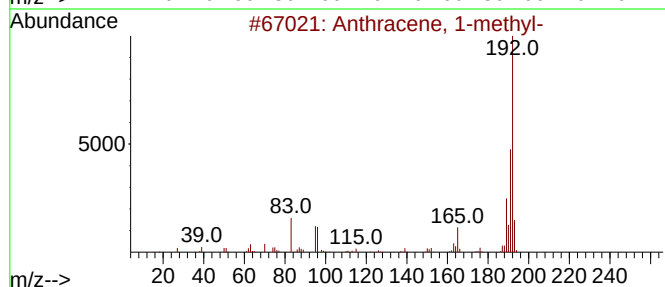
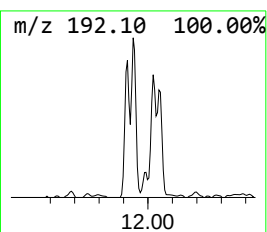
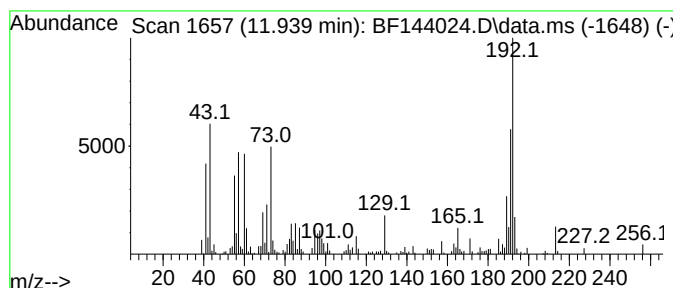
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100625.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Anthracene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.939	9.43 ng	382232	Phenanthrene-d10	11.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 1-methyl-	192	C15H12	000610-48-0	96
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
5			Anthracene, 9-methyl-	192	C15H12	000779-02-2	92



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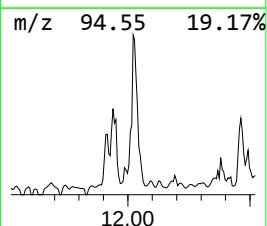
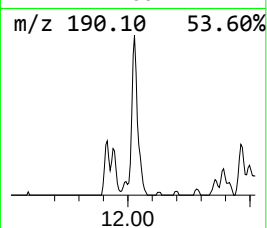
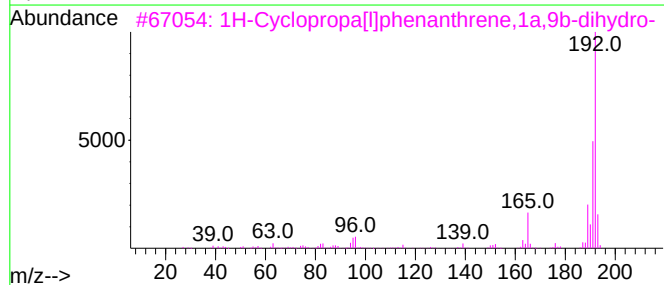
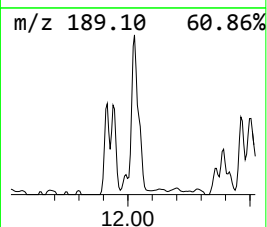
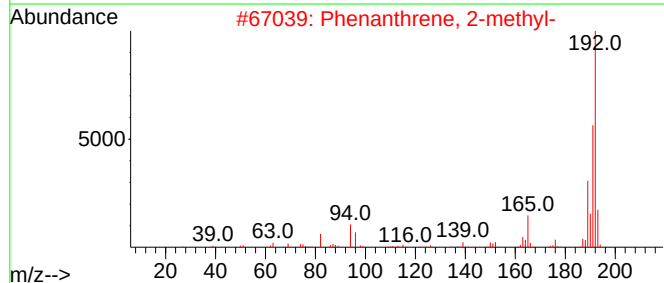
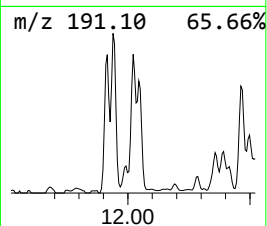
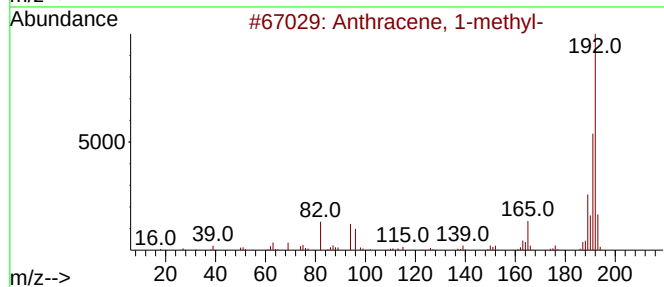
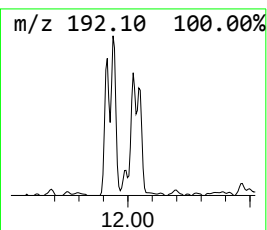
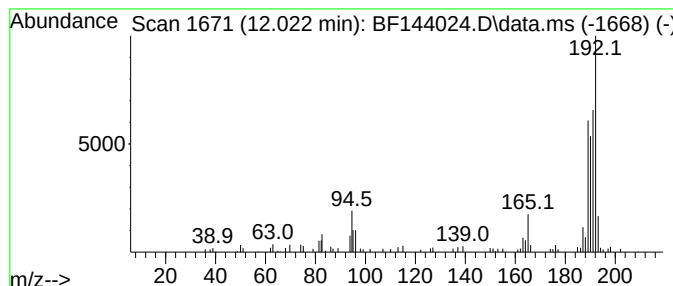
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100625.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.022	2.11 ng	85478	Phenanthrene-d10	11.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 1-methyl-	192	C15H12	000610-48-0	96
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
3			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	90
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	78
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	78



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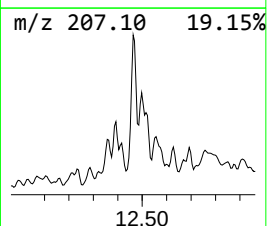
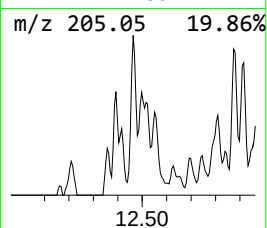
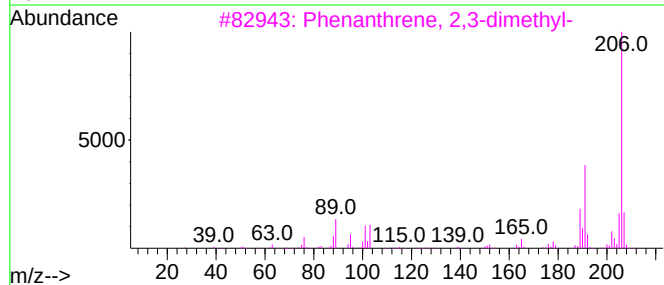
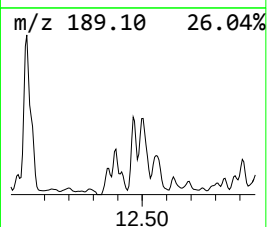
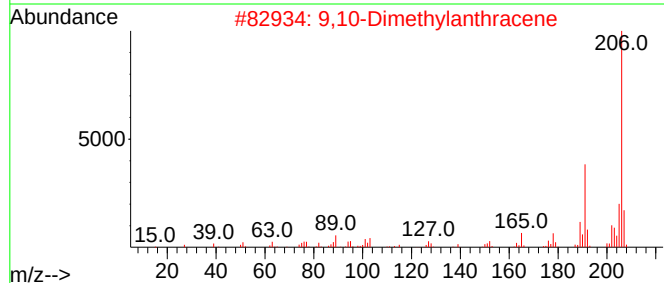
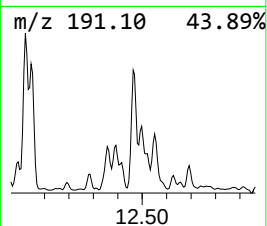
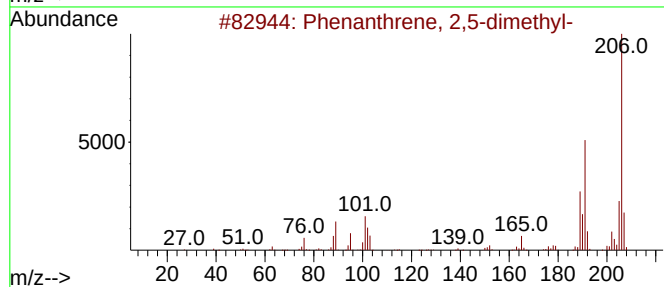
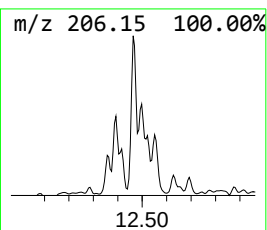
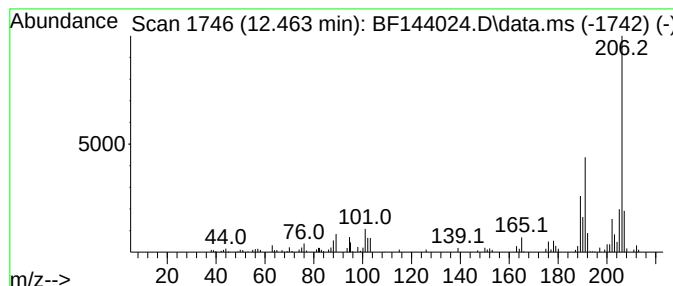
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Phenanthrene, 2,5-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.463	2.40 ng	97066	Phenanthrene-d10	11.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	97
2			9,10-Dimethylanthracene	206	C16H14	000781-43-1	94
3			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	94
4			3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	94
5			di-p-Tolylacetylene	206	C16H14	002789-88-0	93



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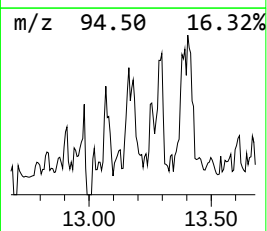
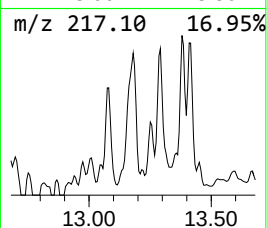
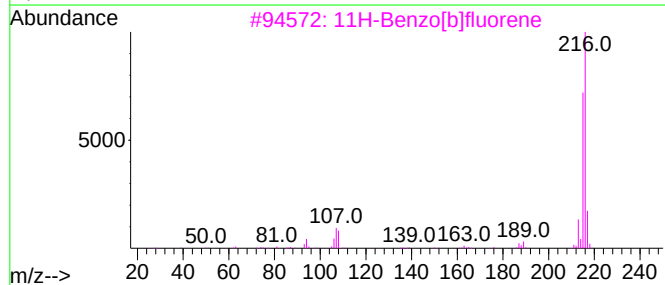
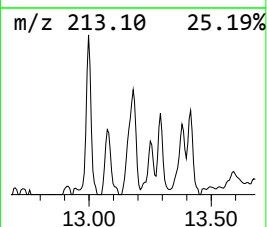
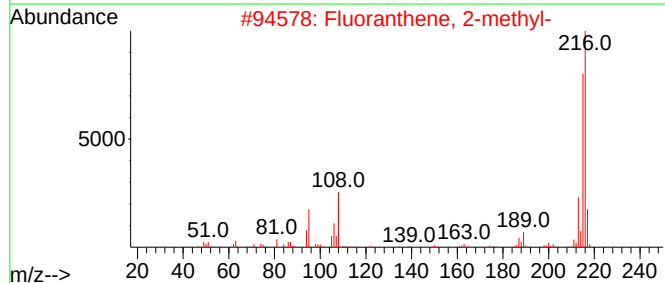
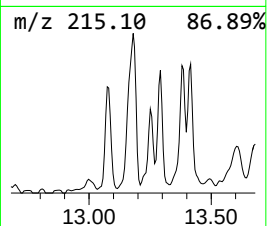
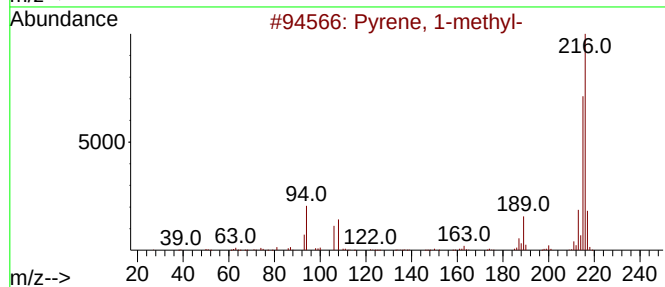
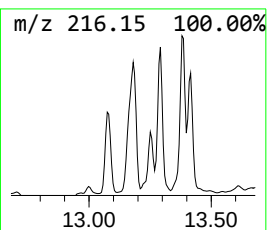
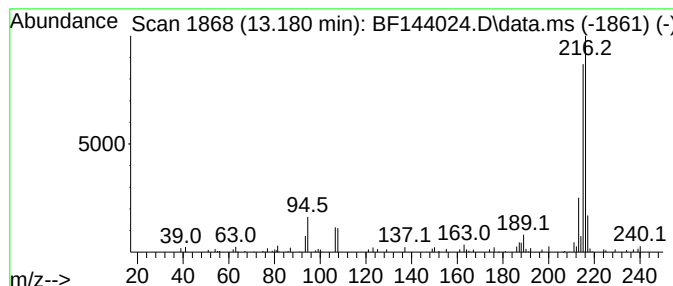
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Pyrene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.180	2.05 ng	98622	Chrysene-d12	14.057

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
2			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
3			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
4			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	81
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102125\
Data File : BF144024.D
Acq On : 21 Oct 2025 20:56
Operator : RC/JU
Sample : Q3403-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-18

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100625.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	10.633	6.1	ng	241646	3	9.922	791250	20.0
Anthracene, 1-m...	11.939	9.4	ng	382232	4	11.410	810403	20.0
Phenanthrene, 2...	12.022	2.1	ng	85478	4	11.410	810403	20.0
Phenanthrene, 2...	12.463	2.4	ng	97066	4	11.410	810403	20.0
Pyrene, 1-methyl-	13.180	2.0	ng	98622	5	14.057	963314	20.0