

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112025\
 Data File : BF144296.D
 Acq On : 20 Nov 2025 15:59
 Operator : RC/JU
 Sample : Q3662-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B1-0.0-1.0-20251117

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-BF110525.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF144296.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.069	481	489	503	rVB	91747	131154	19.13%	1.532%
2	5.487	555	560	577	rBV	315970	442076	64.48%	5.165%
3	6.493	726	731	748	rBV	302918	420041	61.26%	4.907%
4	6.863	789	794	800	rBV	183896	224519	32.75%	2.623%
5	7.422	884	889	900	rBV	205429	257977	37.63%	3.014%
6	8.145	1007	1012	1019	rBV	210829	260607	38.01%	3.045%
7	9.216	1189	1194	1199	rBV	339366	417882	60.95%	4.882%
8	9.557	1248	1252	1255	rBV3	5907	8699	1.27%	0.102%
9	9.622	1259	1263	1269	rVB3	5948	9360	1.37%	0.109%
10	9.763	1283	1287	1291	rBV	31668	40042	5.84%	0.468%
11	9.898	1305	1310	1315	rBV	218648	280262	40.88%	3.274%
12	10.098	1340	1344	1349	rBV3	8080	11544	1.68%	0.135%
13	10.145	1349	1352	1356	rVB4	5608	7067	1.03%	0.083%
14	10.216	1360	1364	1367	rBV4	7273	10597	1.55%	0.124%
15	10.304	1374	1379	1385	rBV8	6037	13775	2.01%	0.161%
16	10.451	1401	1404	1409	rVB4	8718	12933	1.89%	0.151%
17	10.610	1427	1431	1438	rVB	49709	70791	10.33%	0.827%
18	10.692	1440	1445	1450	rBV	222065	283118	41.29%	3.308%
19	10.816	1461	1466	1472	rVB4	10347	16364	2.39%	0.191%
20	10.998	1493	1497	1500	rBV4	7971	10552	1.54%	0.123%
21	11.128	1515	1519	1521	rBV5	4738	7007	1.02%	0.082%
22	11.198	1527	1531	1535	rVB	9425	10567	1.54%	0.123%
23	11.286	1542	1546	1551	rVB	15602	22462	3.28%	0.262%
24	11.345	1551	1556	1558	rBV6	4954	8264	1.21%	0.097%
25	11.392	1559	1564	1566	rVV	249038	334681	48.81%	3.910%
26	11.416	1566	1568	1573	rVV	171317	186306	27.17%	2.177%
27	11.469	1573	1577	1580	rVV	26133	35504	5.18%	0.415%
28	11.498	1580	1582	1589	rVB6	9628	14556	2.12%	0.170%
29	11.628	1600	1604	1608	rBV2	10519	17771	2.59%	0.208%
30	11.722	1617	1620	1625	rBV	17676	22515	3.28%	0.263%
31	11.804	1631	1634	1638	rVB	12007	15828	2.31%	0.185%
32	11.845	1638	1641	1644	rBV3	7481	10156	1.48%	0.119%
33	11.916	1644	1653	1660	rVV2	170117	354292	51.68%	4.139%
34	11.969	1660	1662	1664	rVV3	17138	20079	2.93%	0.235%
35	12.004	1664	1668	1677	rVB2	78812	166605	24.30%	1.946%
36	12.080	1677	1681	1686	rBV6	16049	33857	4.94%	0.396%

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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

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37	12.127	1686	1689	1692	rVV2	14719	17722	2.58%	0.207%
38	12.192	1696	1700	1702	rVV	35072	54684	7.98%	0.639%
39	12.222	1702	1705	1711	rVB	34935	51411	7.50%	0.601%
40	12.339	1719	1725	1727	rBV2	20708	37119	5.41%	0.434%
41	12.369	1727	1730	1733	rVV	19615	23521	3.43%	0.275%
42	12.445	1739	1743	1746	rBV	67385	91788	13.39%	1.072%
43	12.480	1746	1749	1755	rVV2	35450	65173	9.51%	0.761%
44	12.533	1755	1758	1766	rVB4	36925	58521	8.54%	0.684%
45	12.610	1766	1771	1775	rBV	387018	494041	72.06%	5.772%
46	12.704	1783	1787	1791	rVB2	40964	51863	7.56%	0.606%
47	12.839	1804	1810	1816	rBV	381412	556202	81.12%	6.498%
48	12.892	1816	1819	1823	rVB2	37996	46719	6.81%	0.546%
49	12.980	1828	1834	1841	rVB	406613	598790	87.34%	6.996%
50	13.057	1842	1847	1854	rBV4	50436	97747	14.26%	1.142%
51	13.163	1858	1865	1870	rVB2	67350	140268	20.46%	1.639%
52	13.233	1874	1877	1880	rVV2	29924	37293	5.44%	0.436%
53	13.269	1880	1883	1891	rVB2	58376	93061	13.57%	1.087%
54	13.363	1896	1899	1902	rVV2	38258	49545	7.23%	0.579%
55	13.398	1902	1905	1908	rVB	36190	42595	6.21%	0.498%
56	13.680	1950	1953	1957	rVB	47643	59785	8.72%	0.698%
57	13.786	1967	1971	1974	rBV3	61066	95384	13.91%	1.114%
58	14.039	2008	2014	2017	rBV2	388816	685615	100.00%	8.010%
59	15.092	2188	2193	2200	rBV	160685	346697	50.57%	4.051%
60	15.398	2241	2245	2250	rVB	91835	145272	21.19%	1.697%
61	15.457	2251	2255	2261	rBV	124000	192859	28.13%	2.253%
62	15.521	2262	2266	2271	rBV	147544	235765	34.39%	2.755%

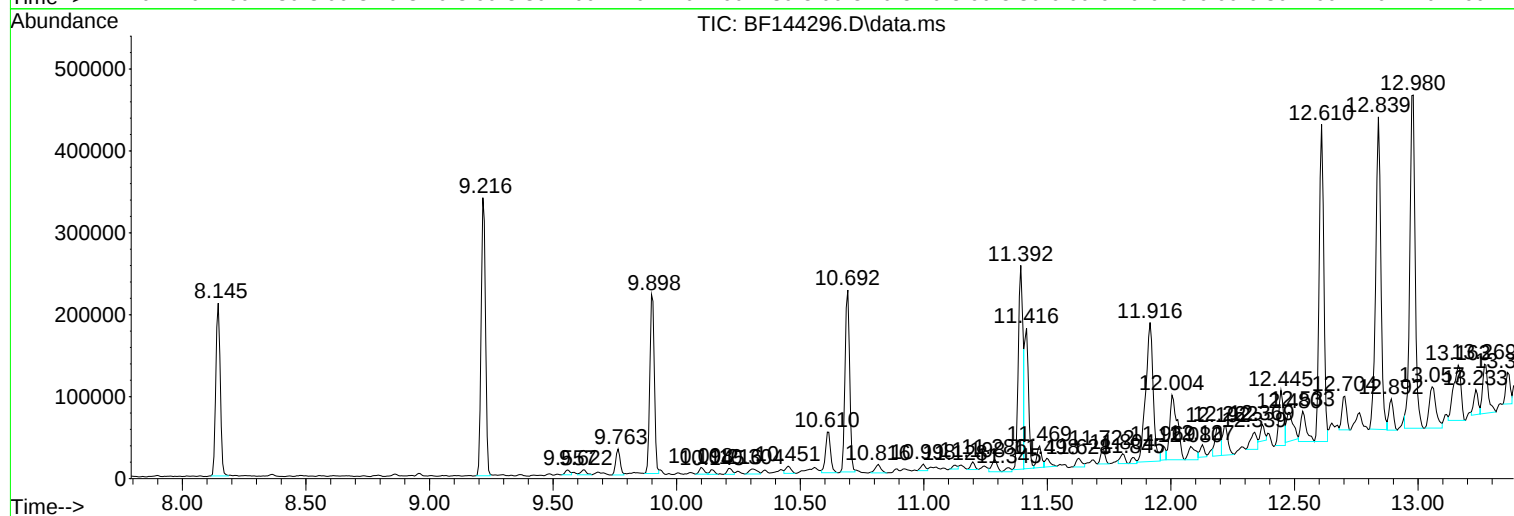
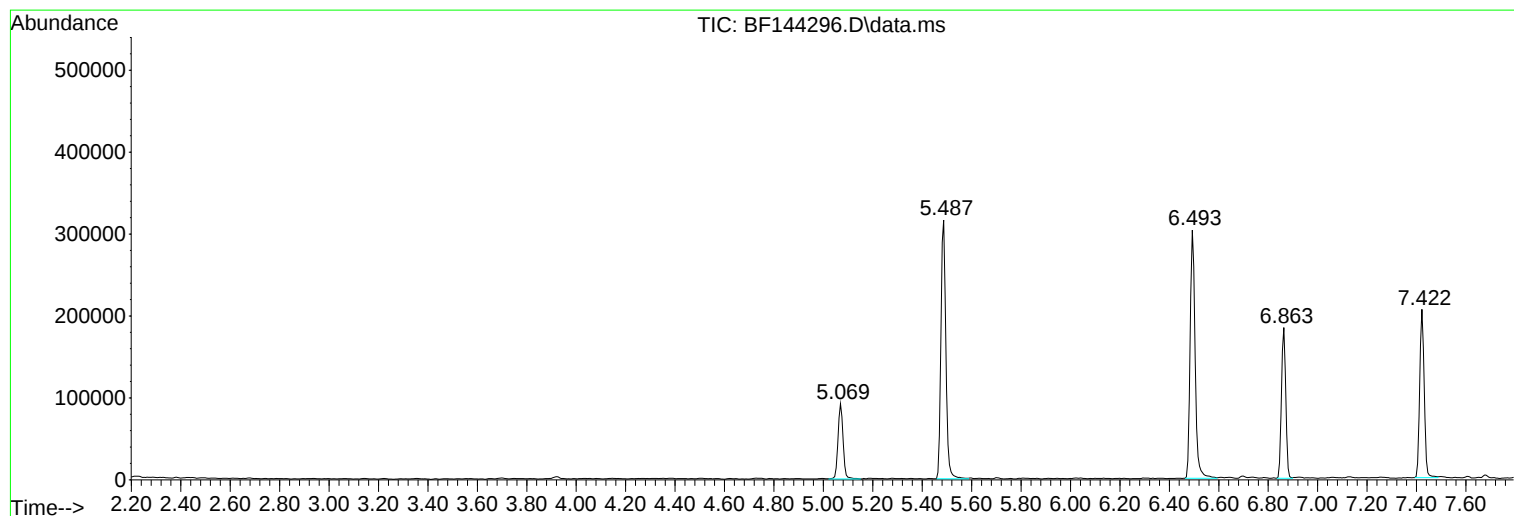
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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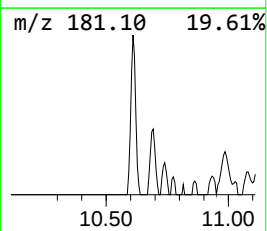
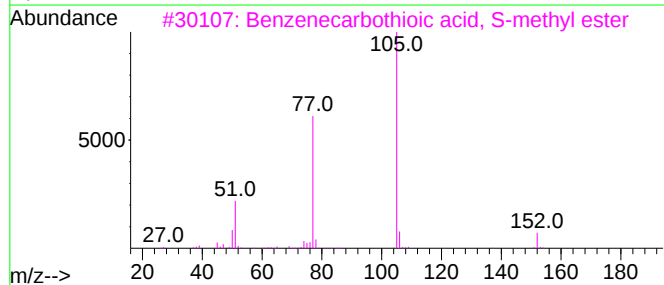
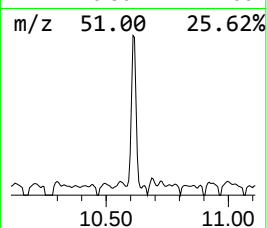
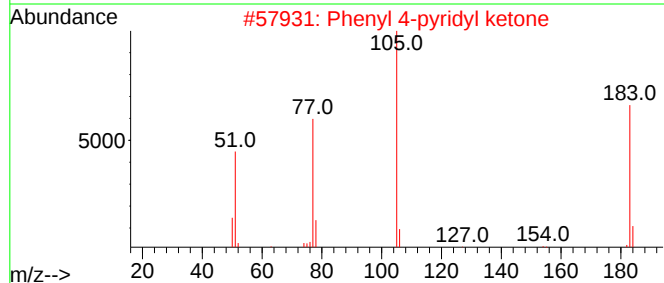
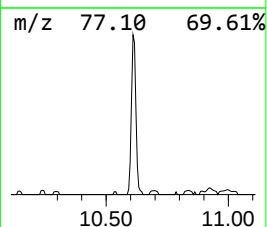
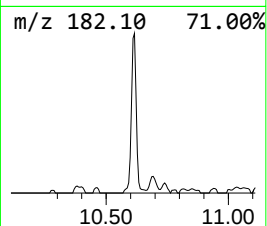
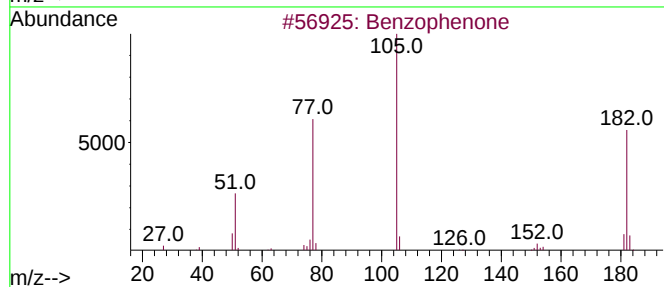
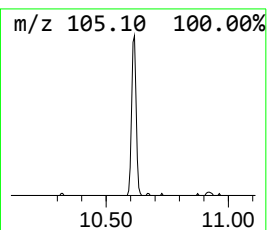
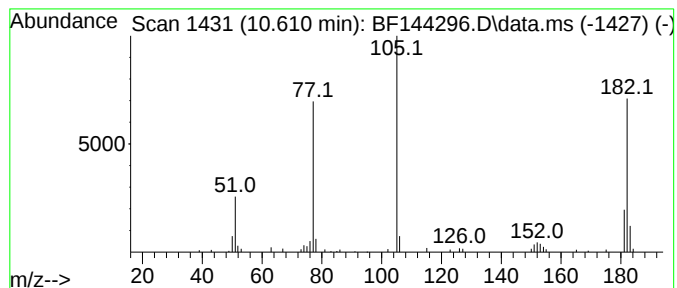
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.610	5.05 ng	70791	Acenaphthene-d10	9.898

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	93
2			Phenyl 4-pyridyl ketone	183	C12H9NO	014548-46-0	43
3			Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	43
4			3,3'-Dimethylbiphenyl	182	C14H14	000612-75-9	38
5			Benzoylformic acid	150	C8H6O3	000611-73-4	38



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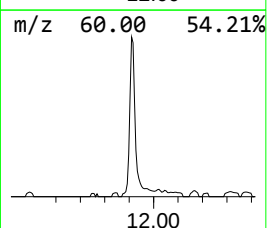
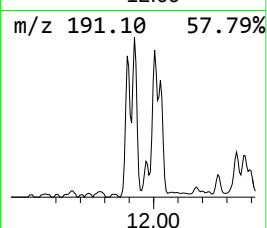
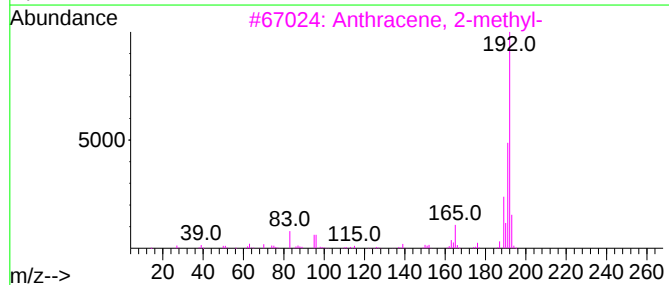
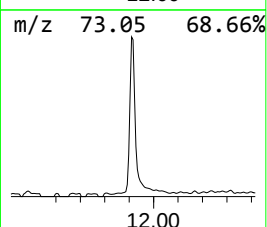
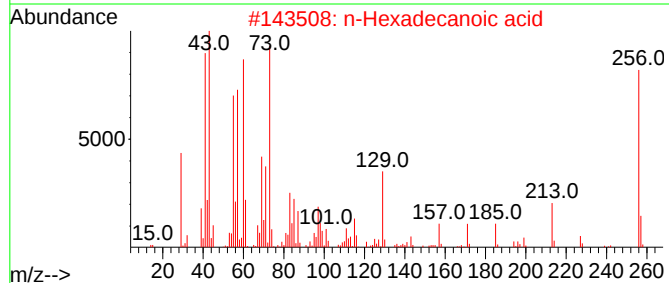
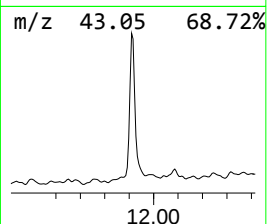
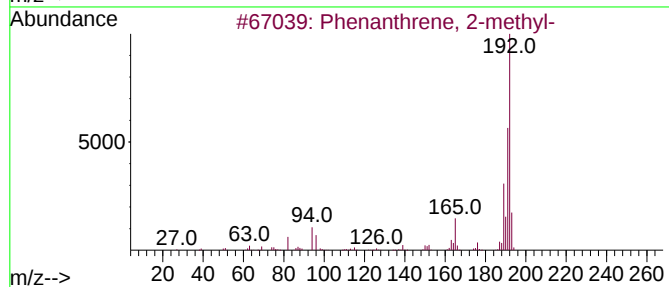
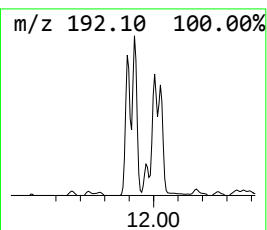
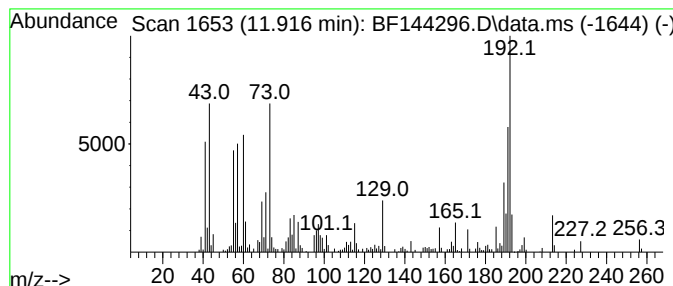
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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Peak Number 3 Phenanthrene, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.916	21.17 ng	354292	Phenanthrene-d10	11.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	92
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	91
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
5			Tridecanoic acid	214	C13H26O2	000638-53-9	90



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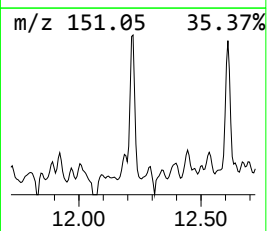
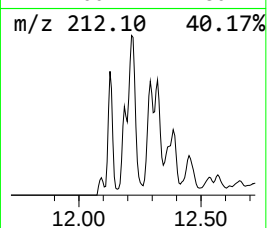
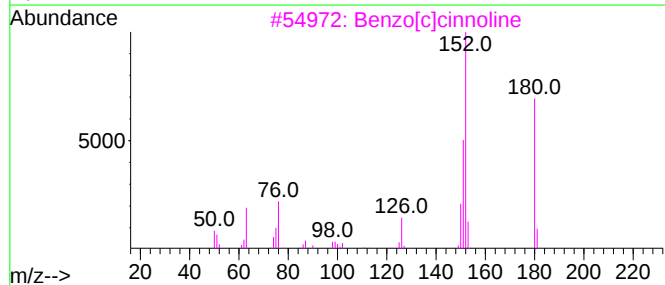
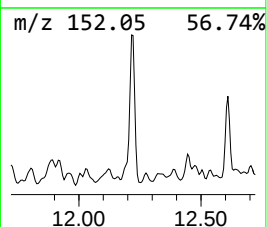
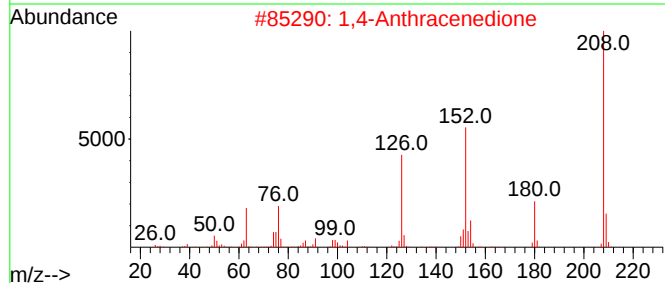
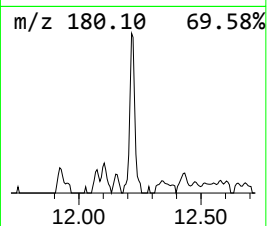
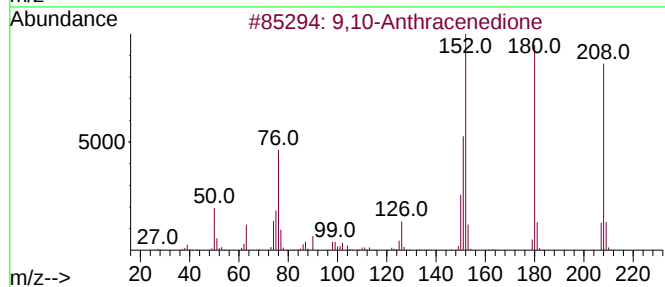
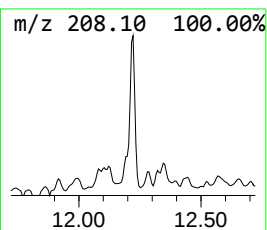
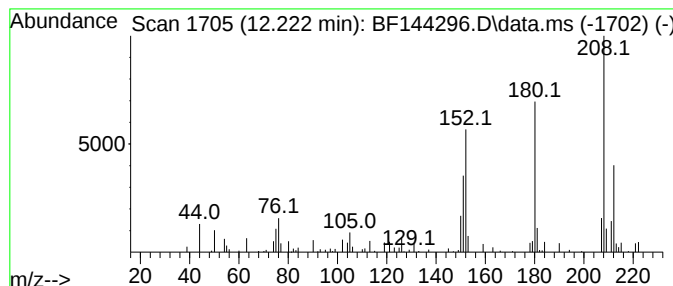
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 9,10-Anthracenedione Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.222	3.07 ng	51411	Phenanthrene-d10	11.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
2			1,4-Anthracenedione	208	C14H8O2	000635-12-1	43
3			Benzo[c]cinnoline	180	C12H8N2	000230-17-1	38
4			5,6,7,8-Tetrahydro-1-phenylapht...	208	C16H16	067064-63-5	38
5			9H-Fluoren-9-one	180	C13H8O	000486-25-9	38



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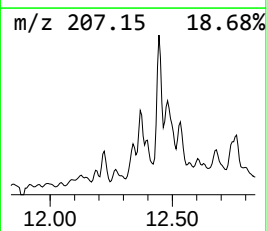
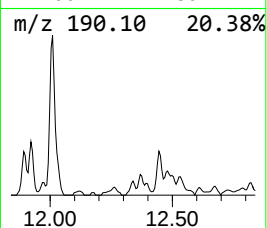
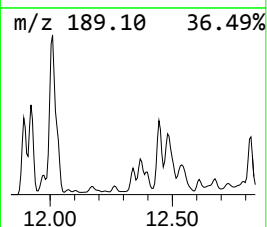
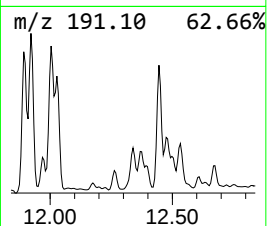
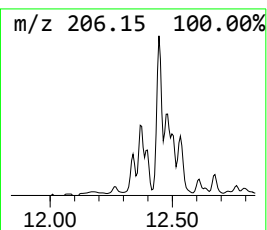
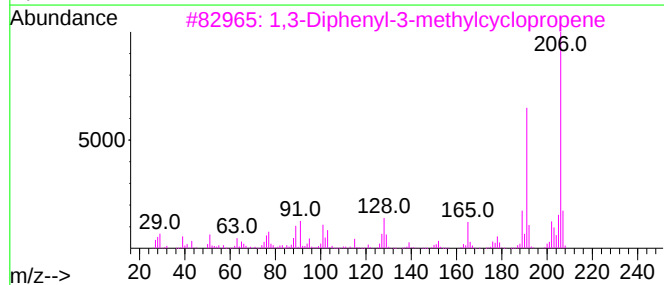
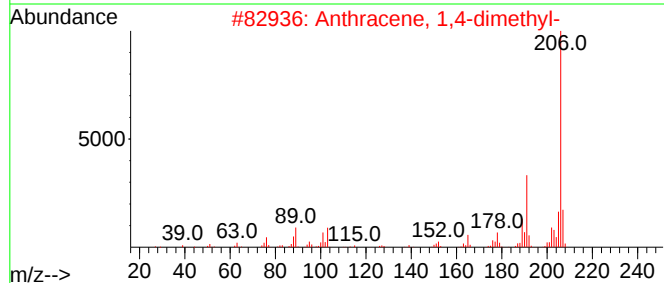
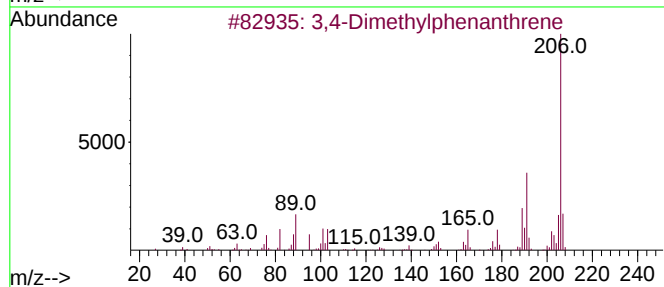
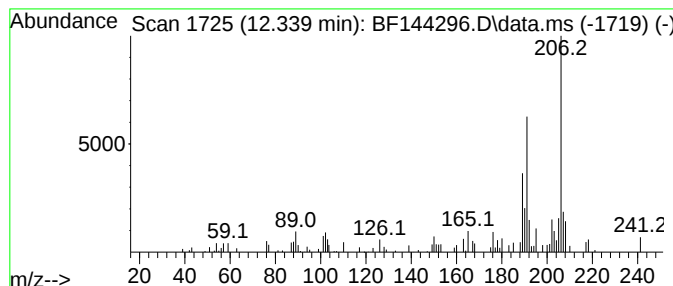
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TIC Library : C:\Database\NIST20.L
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Peak Number 8 3,4-Dimethylphenanthrene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.339	2.22 ng	37119	Phenanthrene-d10	11.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	93
2			Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	92
3			1,3-Diphenyl-3-methylcyclopropene	206	C16H14	086544-79-8	90
4			3-Methyl-1-phenyl-1H-indene	206	C16H14	022360-63-0	90
5			Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	87



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ClientSampleId :
B1-0.0-1.0-20251117

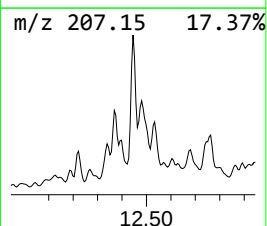
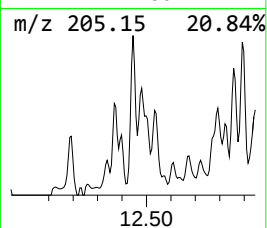
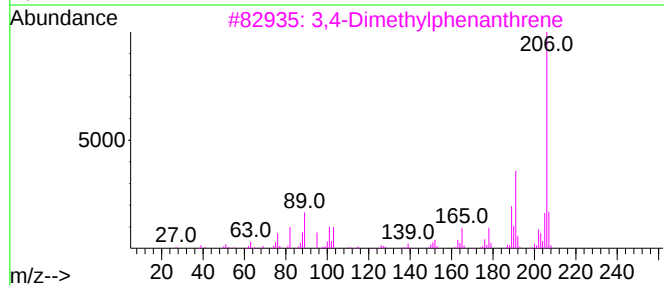
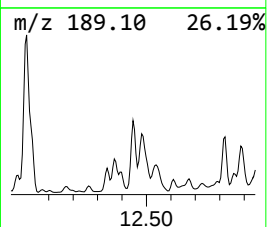
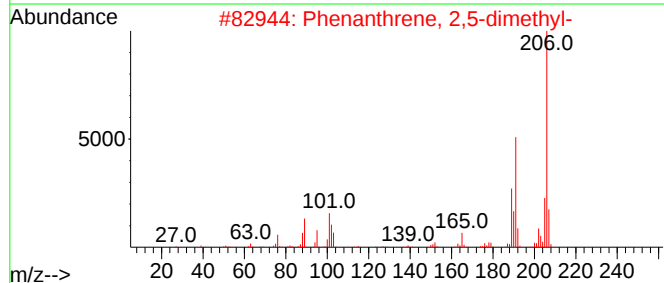
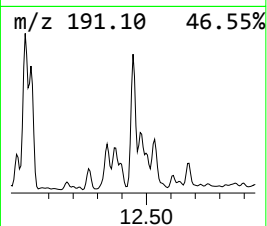
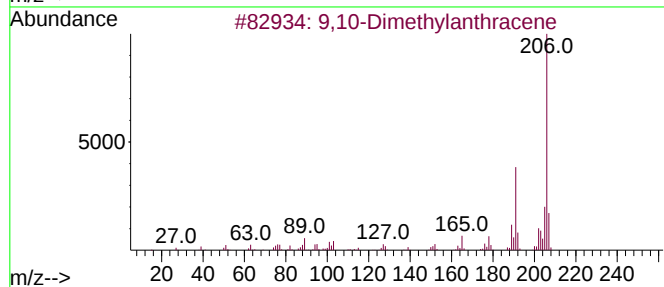
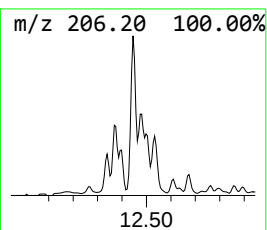
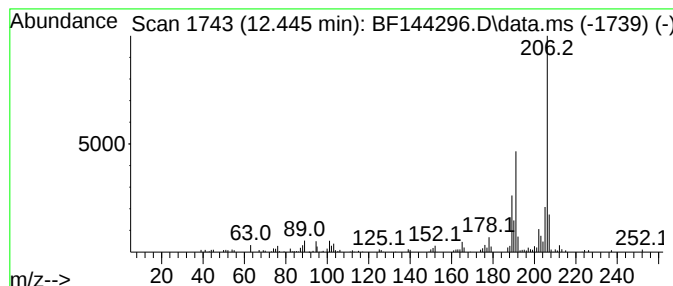
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 9,10-Dimethylantracene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.445	5.49 ng	91788	Phenanthrene-d10	11.392

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9,10-Dimethylantracene	206	C16H14	000781-43-1	95
2		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	94
3		3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	90
4		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	90
5		Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112025\
Data File : BF144296.D
Acq On : 20 Nov 2025 15:59
Operator : RC/JU
Sample : Q3662-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

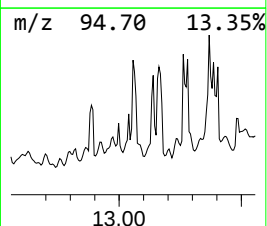
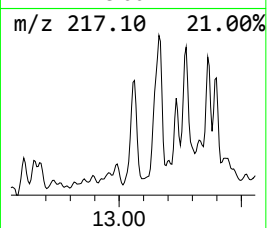
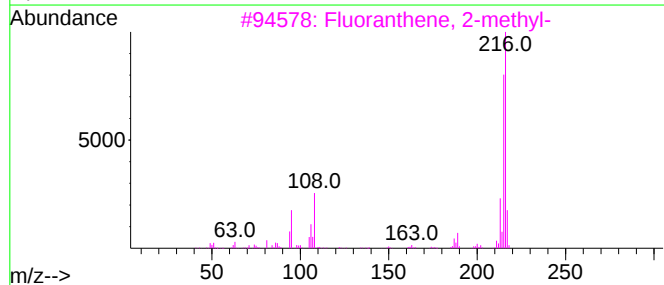
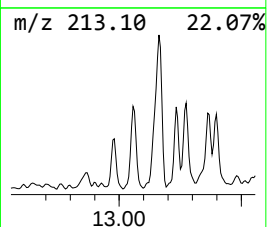
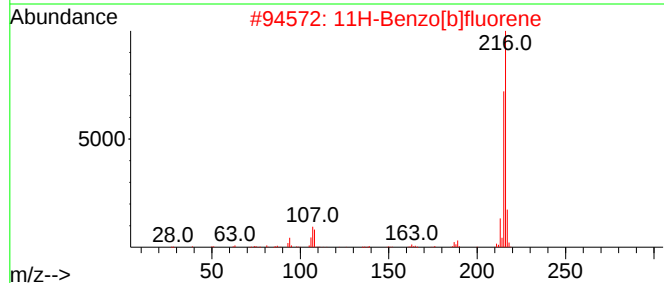
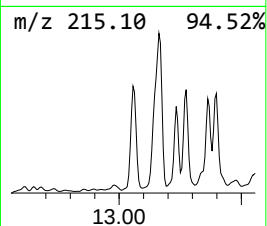
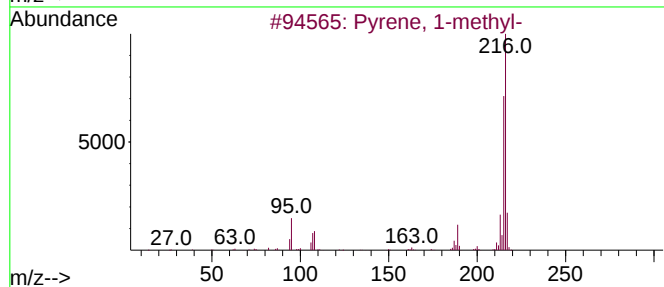
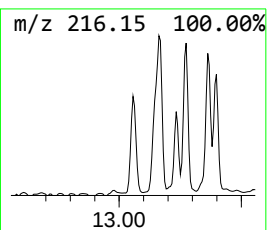
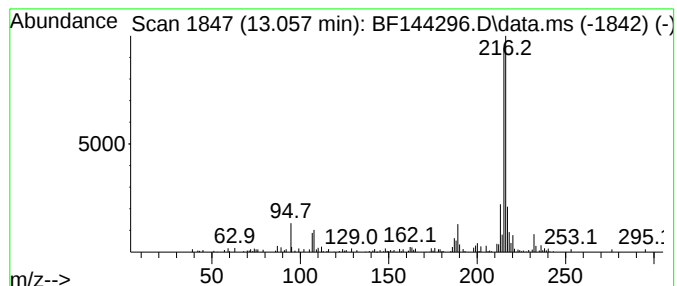
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.057	2.85 ng	97747	Chrysene-d12	14.039	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	96
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112025\
Data File : BF144296.D
Acq On : 20 Nov 2025 15:59
Operator : RC/JU
Sample : Q3662-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B1-0.0-1.0-20251117

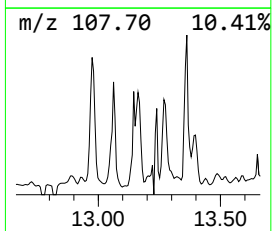
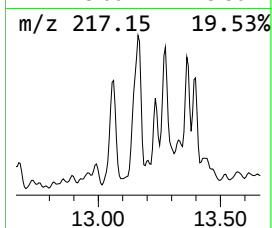
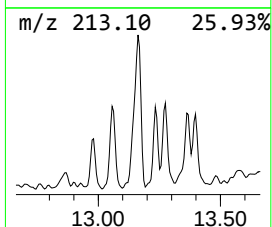
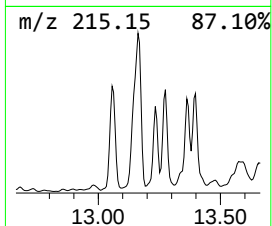
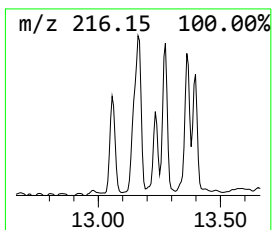
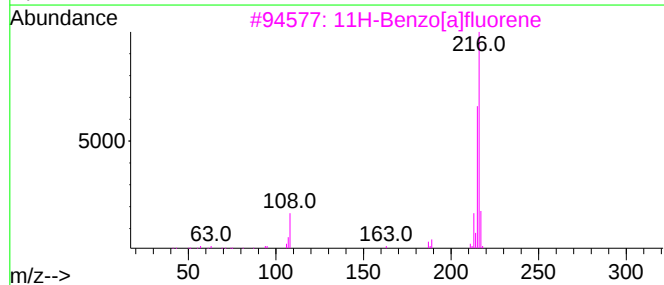
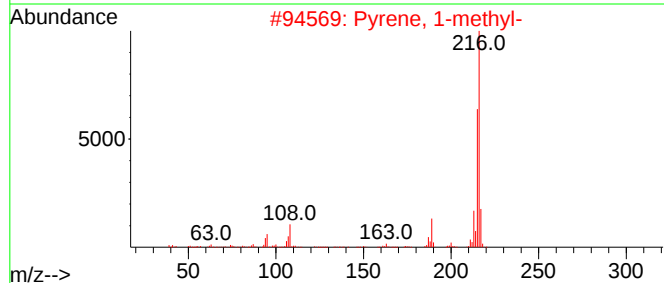
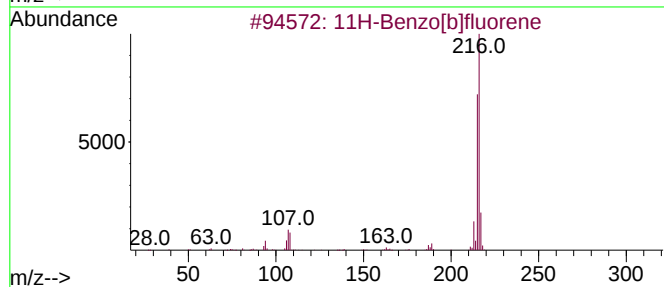
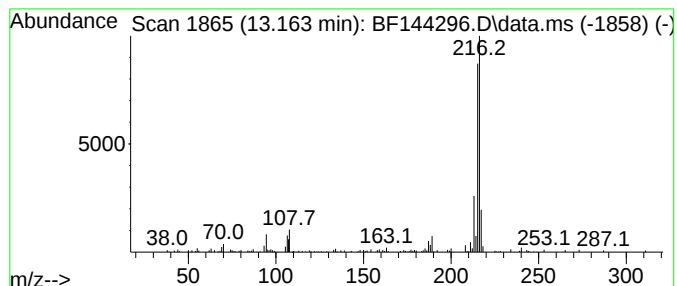
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 11H-Benzo[b]fluorene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.163	4.09 ng	140268	Chrysene-d12	14.039

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112025\
Data File : BF144296.D
Acq On : 20 Nov 2025 15:59
Operator : RC/JU
Sample : Q3662-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

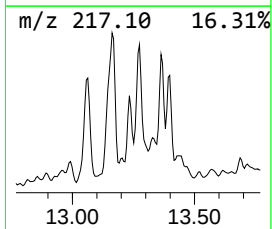
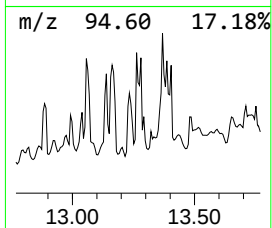
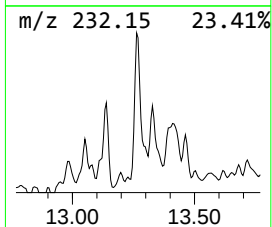
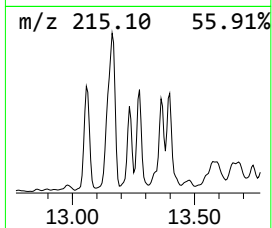
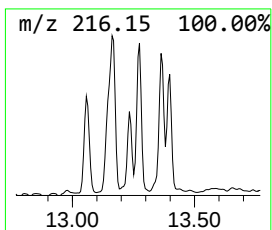
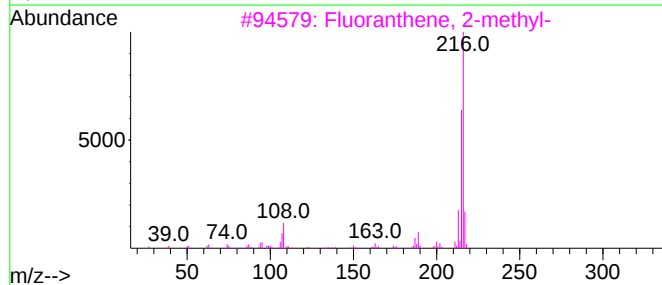
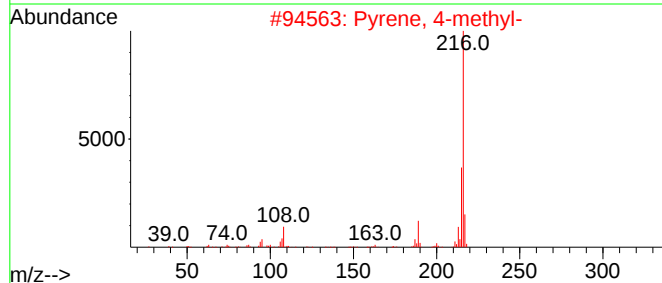
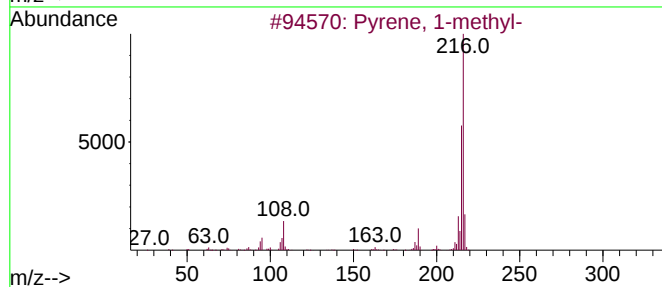
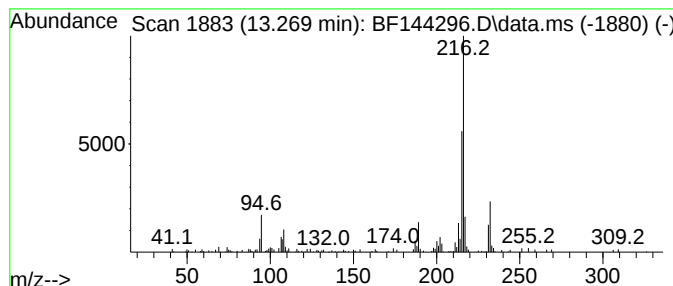
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Pyrene, 4-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.269	2.71 ng	93061	Chrysene-d12	14.039	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	96
2	Pyrene, 4-methyl-	216	C17H12	003353-12-6	93
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	92
4	Pyrene, 2-methyl-	216	C17H12	003442-78-2	89
5	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112025\
Data File : BF144296.D
Acq On : 20 Nov 2025 15:59
Operator : RC/JU
Sample : Q3662-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
B1-0.0-1.0-20251117

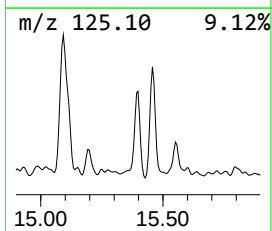
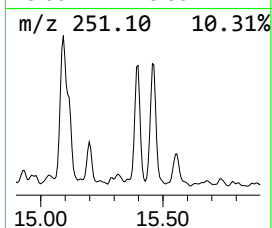
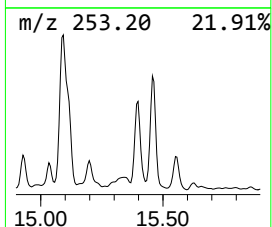
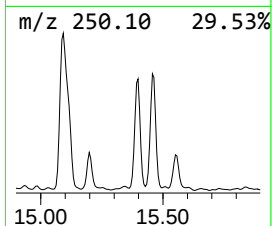
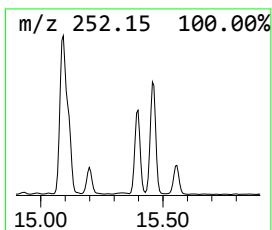
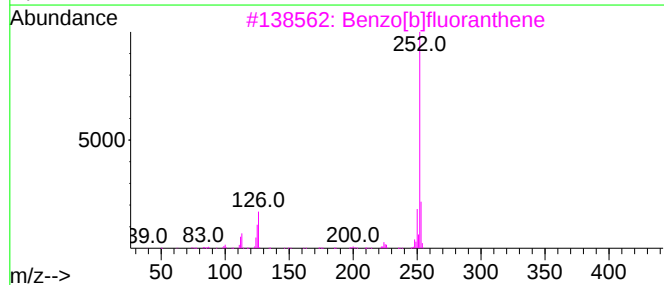
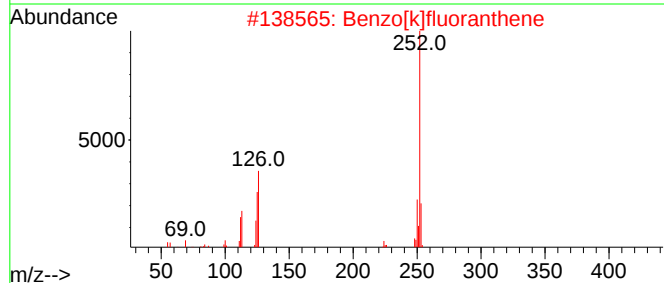
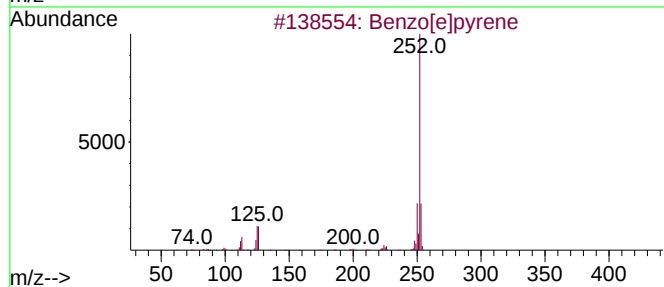
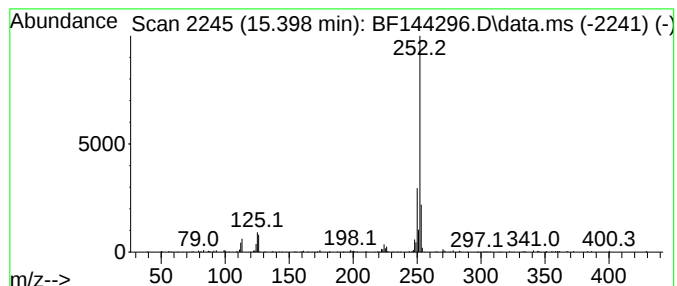
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.398	12.32 ng	145272	Perylene-d12	15.521

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[e]pyrene	252	C20H12	000192-97-2	96
2		Benzo[k]fluoranthene	252	C20H12	000207-08-9	94
3		Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
4		Benzo[a]pyrene	252	C20H12	000050-32-8	91
5		Benzo[j]fluoranthene	252	C20H12	000205-82-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112025\
Data File : BF144296.D
Acq On : 20 Nov 2025 15:59
Operator : RC/JU
Sample : Q3662-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B1-0.0-1.0-20251117

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.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
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Phenanthrene, 2...	11.916	21.2	ng	354292	4	11.392	334681	20.0
9,10-Anthracene...	12.222	3.1	ng	51411	4	11.392	334681	20.0
3,4-Dimethylphe...	12.339	2.2	ng	37119	4	11.392	334681	20.0
9,10-Dimethylan...	12.445	5.5	ng	91788	4	11.392	334681	20.0
Pyrene, 1-methyl-	13.057	2.9	ng	97747	5	14.039	685615	20.0
11H-Benzo[b]flu...	13.163	4.1	ng	140268	5	14.039	685615	20.0
Pyrene, 4-methyl-	13.269	2.7	ng	93061	5	14.039	685615	20.0
Benzo[e]pyrene	15.398	12.3	ng	145272	6	15.521	235765	20.0