

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062425\
 Data File : BF142831.D
 Acq On : 24 Jun 2025 21:08
 Operator : RC/JU
 Sample : Q2384-01 5X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-01-6202025

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-BF061925.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF142831.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.081	486	491	498	rBV	9540	13102	1.91%	0.169%
2	5.493	556	561	577	rBV	106775	146311	21.34%	1.882%
3	6.504	728	733	747	rBV	106138	146499	21.37%	1.885%
4	6.881	792	797	802	rBV	273717	330466	48.21%	4.252%
5	7.440	887	892	902	rVB	65387	83114	12.12%	1.069%
6	8.163	1010	1015	1026	rBV	326231	407501	59.44%	5.243%
7	8.881	1133	1137	1144	rVB	8594	11705	1.71%	0.151%
8	8.975	1149	1153	1158	rVB	11757	14988	2.19%	0.193%
9	9.239	1192	1198	1202	rBV	146532	182342	26.60%	2.346%
10	9.316	1207	1211	1217	rBV4	4923	7749	1.13%	0.100%
11	9.504	1238	1243	1247	rBV2	16701	24586	3.59%	0.316%
12	9.575	1251	1255	1258	rBV	27064	38281	5.58%	0.493%
13	9.604	1258	1260	1263	rVB	7520	7372	1.08%	0.095%
14	9.698	1271	1276	1285	rBV	14375	26196	3.82%	0.337%
15	9.781	1285	1290	1296	rBV2	42159	57696	8.42%	0.742%
16	9.834	1296	1299	1304	rVB3	5708	8871	1.29%	0.114%
17	9.922	1307	1314	1318	rBV	360997	446490	65.13%	5.745%
18	9.951	1318	1319	1323	rVB	14334	12079	1.76%	0.155%
19	10.028	1328	1332	1336	rBV2	10848	16352	2.39%	0.210%
20	10.122	1343	1348	1352	rBV3	22054	29945	4.37%	0.385%
21	10.163	1352	1355	1359	rVB	21141	24838	3.62%	0.320%
22	10.234	1363	1367	1370	rBV2	22662	32638	4.76%	0.420%
23	10.263	1370	1372	1378	rVB	13120	15720	2.29%	0.202%
24	10.322	1378	1382	1384	rBV3	15594	23943	3.49%	0.308%
25	10.416	1394	1398	1401	rBV3	8524	12766	1.86%	0.164%
26	10.469	1403	1407	1412	rVB3	35030	48784	7.12%	0.628%
27	10.533	1413	1418	1420	rVV3	9951	16868	2.46%	0.217%
28	10.581	1423	1426	1429	rVV3	10685	13020	1.90%	0.168%
29	10.628	1429	1434	1441	rVB3	25301	41077	5.99%	0.529%
30	10.710	1442	1448	1452	rBV	101132	130633	19.06%	1.681%
31	10.745	1452	1454	1459	rVB4	5626	7255	1.06%	0.093%
32	10.833	1463	1469	1475	rVB4	27288	46814	6.83%	0.602%
33	10.910	1475	1482	1484	rBV2	13968	22815	3.33%	0.294%
34	11.016	1495	1500	1503	rBV2	28069	45556	6.65%	0.586%
35	11.045	1503	1505	1508	rVV2	18819	19784	2.89%	0.255%
36	11.098	1512	1514	1517	rVB	14327	14942	2.18%	0.192%

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37	11.145	1518	1522	1524	rBV4	9308	15106	2.20%	0.194%
38	11.216	1531	1534	1538	rBV3	10827	14692	2.14%	0.189%
39	11.263	1538	1542	1545	rVV4	15074	24430	3.56%	0.314%
40	11.304	1545	1549	1555	rVB2	32494	48690	7.10%	0.626%
41	11.410	1562	1567	1570	rBV	390487	571380	83.35%	7.351%
42	11.486	1576	1580	1583	rBV	41098	48712	7.11%	0.627%
43	11.516	1583	1585	1589	rVB3	16793	21074	3.07%	0.271%
44	11.645	1604	1607	1612	rBV3	13519	25824	3.77%	0.332%
45	11.698	1613	1616	1619	rVV2	19030	27583	4.02%	0.355%
46	11.739	1620	1623	1628	rVB	31940	43435	6.34%	0.559%
47	11.828	1634	1638	1642	rVB	22881	30855	4.50%	0.397%
48	11.916	1648	1653	1655	rBV	90663	134430	19.61%	1.730%
49	11.939	1655	1657	1661	rVV	118866	133331	19.45%	1.715%
50	11.986	1662	1665	1668	rVV2	32924	48071	7.01%	0.618%
51	12.028	1668	1672	1680	rVB2	156030	310786	45.34%	3.999%
52	12.210	1699	1703	1705	rBV	39240	54057	7.89%	0.696%
53	12.357	1722	1728	1730	rBV2	33664	66241	9.66%	0.852%
54	12.386	1730	1733	1736	rVV	36186	43860	6.40%	0.564%
55	12.463	1742	1746	1749	rBV	128743	186980	27.28%	2.406%
56	12.498	1749	1752	1758	rVV2	95239	177615	25.91%	2.285%
57	12.551	1758	1761	1770	rVB3	58414	121396	17.71%	1.562%
58	12.627	1770	1774	1778	rBV	348586	427013	62.29%	5.494%
59	12.780	1797	1800	1803	rBV	29623	38830	5.66%	0.500%
60	12.857	1807	1813	1819	rBV	467690	685530	100.00%	8.820%
61	12.910	1819	1822	1826	rVB2	48250	61255	8.94%	0.788%
62	12.992	1833	1836	1844	rVB	184709	280894	40.97%	3.614%
63	13.075	1846	1850	1856	rBV2	74025	129034	18.82%	1.660%
64	13.180	1862	1868	1872	rVB	96693	177060	25.83%	2.278%
65	13.286	1883	1886	1891	rBV	89587	117347	17.12%	1.510%
66	13.380	1899	1902	1905	rBV	66820	84073	12.26%	1.082%
67	13.410	1905	1907	1911	rVB	50068	58461	8.53%	0.752%
68	14.051	2012	2016	2020	rBV2	346070	589963	86.06%	7.591%
69	15.551	2268	2271	2292	rVB2	185838	445237	64.95%	5.728%

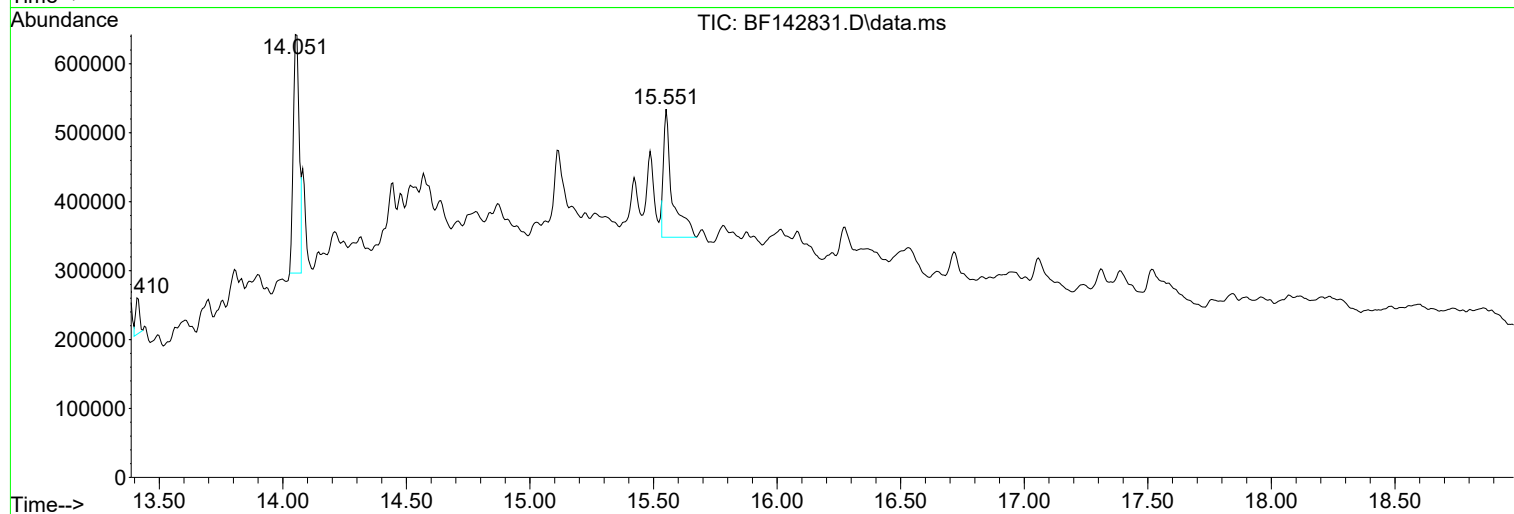
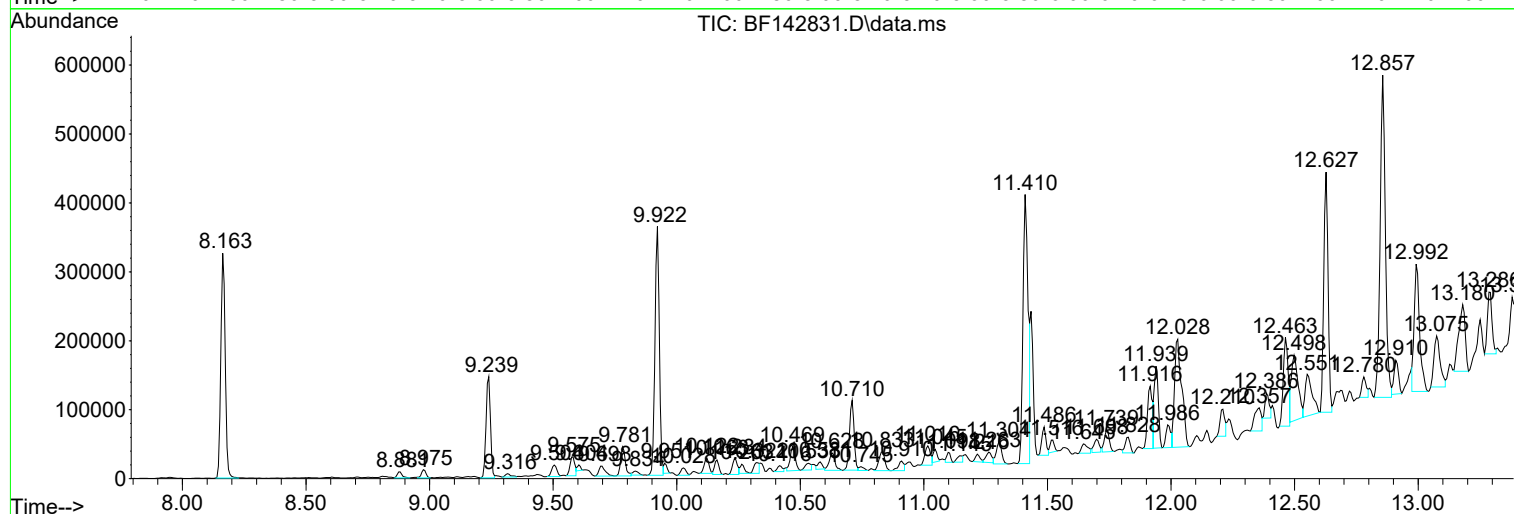
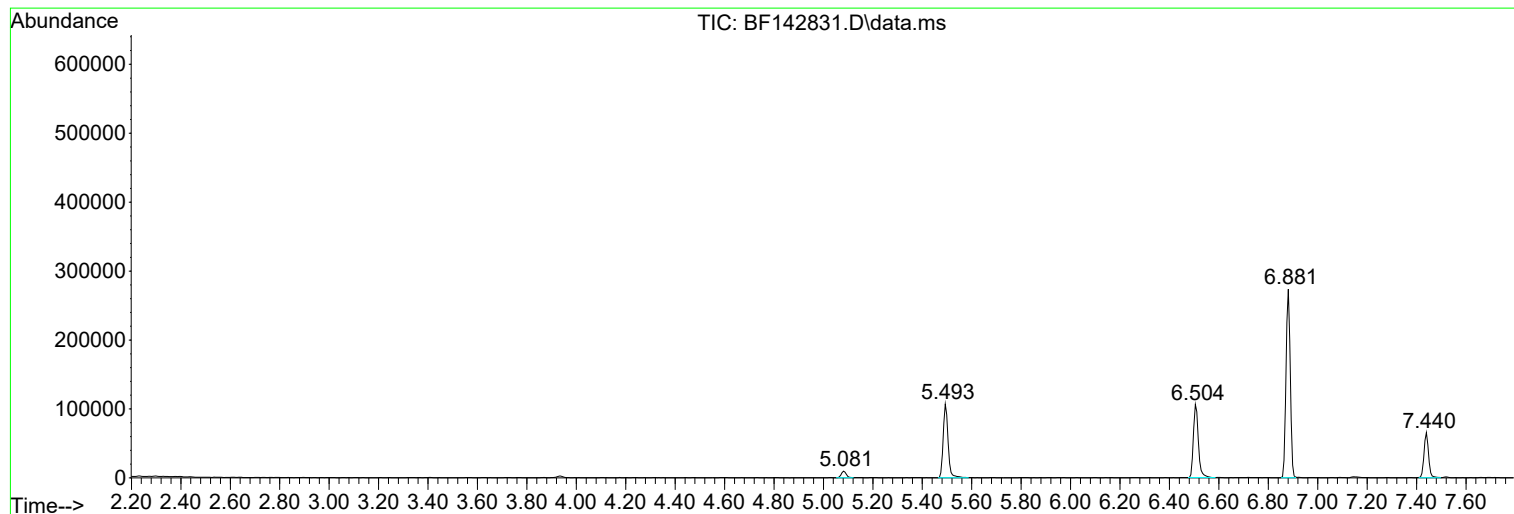
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :
BNA_F
ClientSampleId :
OK-01-6202025

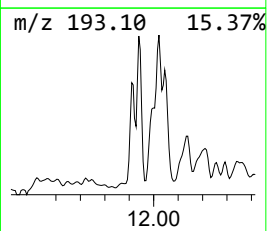
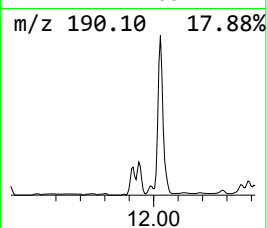
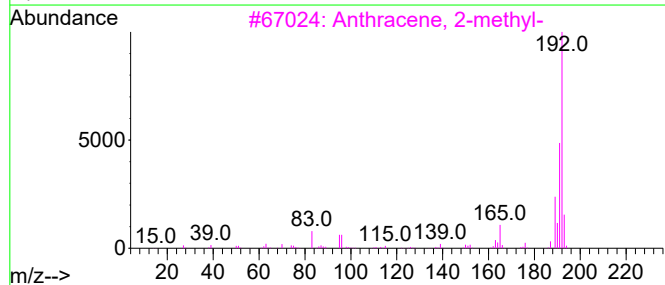
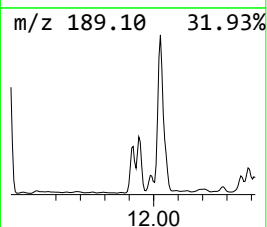
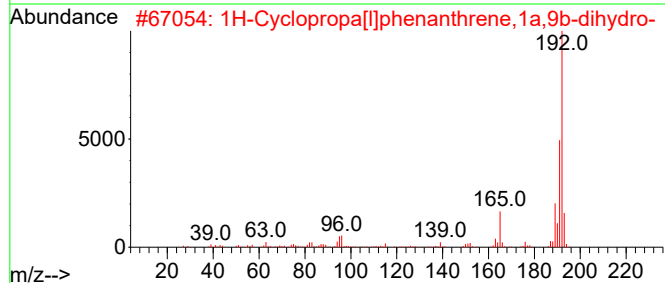
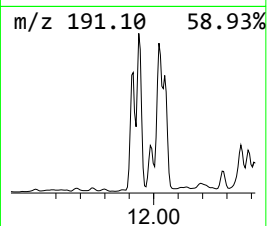
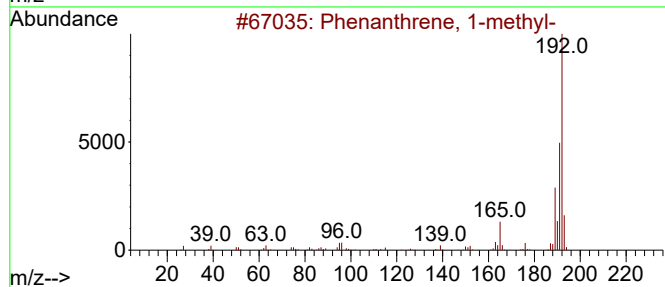
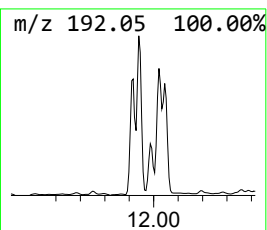
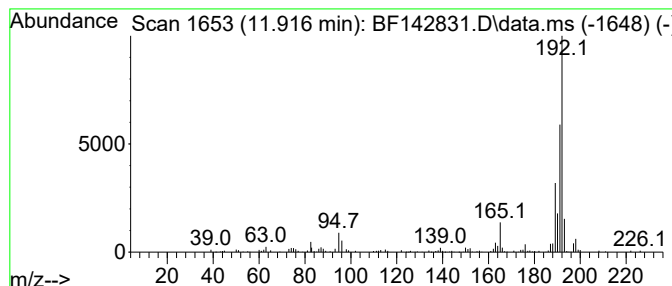
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Phenanthrene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.916	4.71 ng	134430	Phenanthrene-d10	11.410

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
2		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	93
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
4		Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
5		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	91



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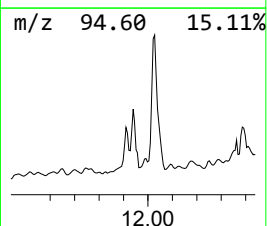
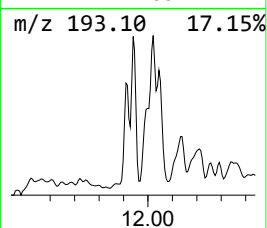
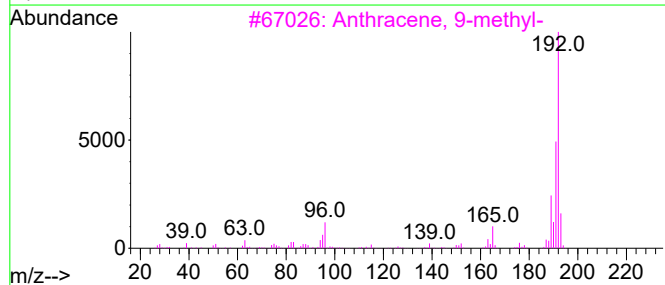
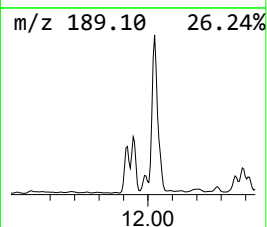
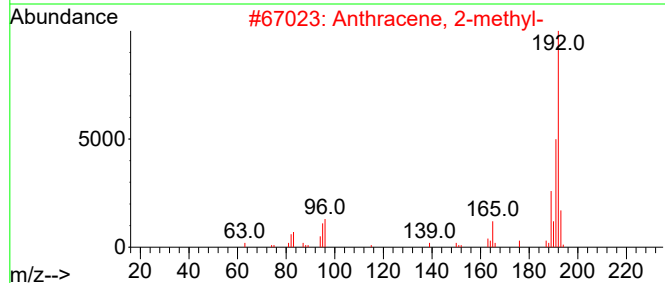
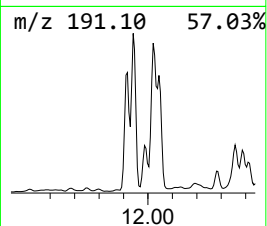
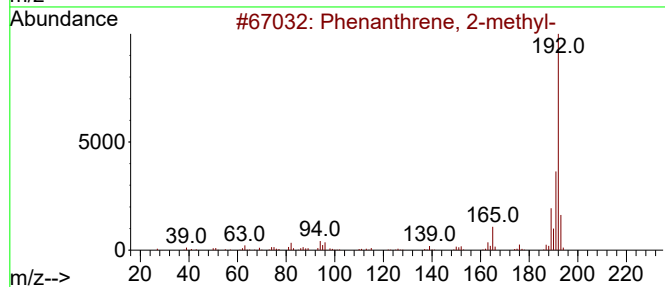
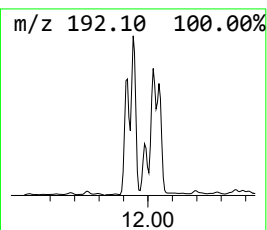
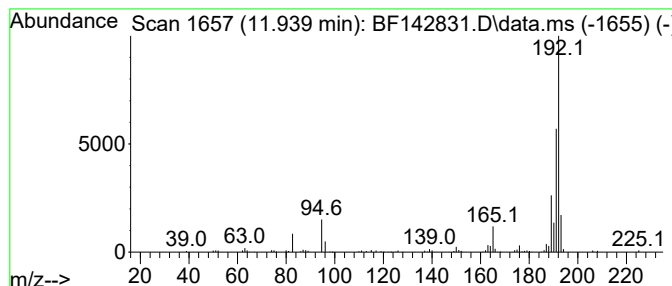
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.M
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 Phenanthrene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.939	4.67 ng	133331	Phenanthrene-d10	11.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
4			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	91
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91



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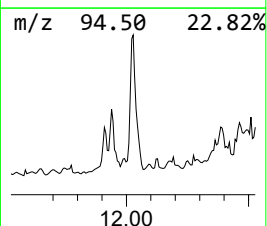
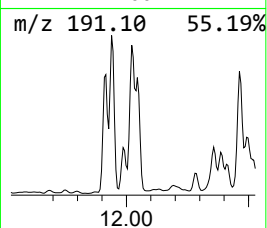
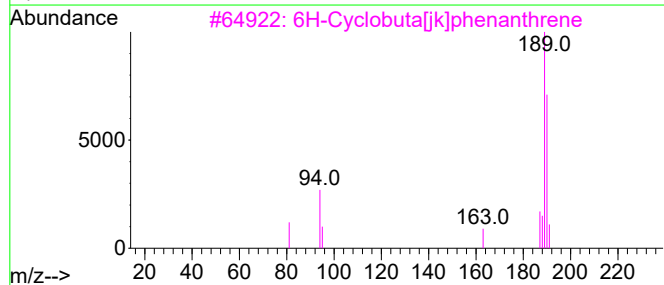
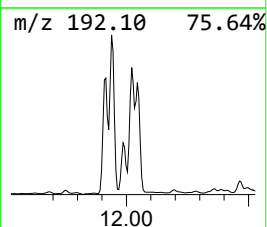
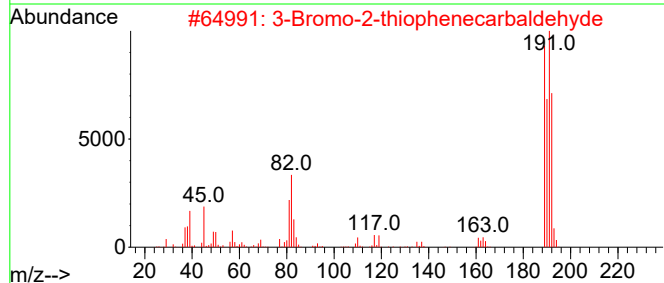
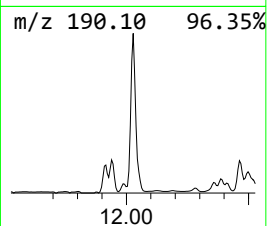
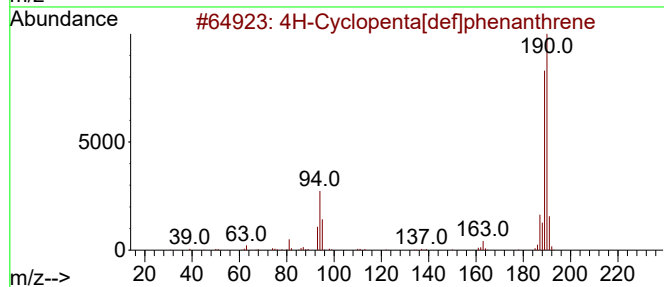
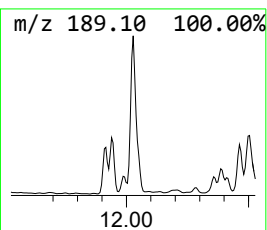
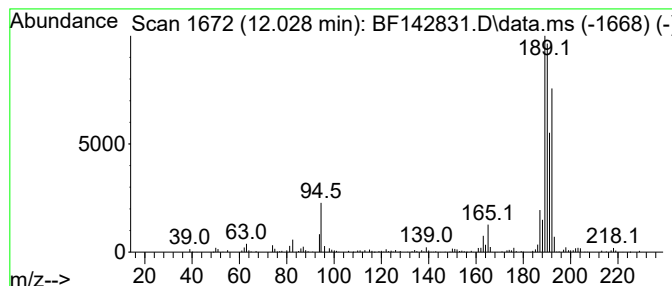
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Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.028	10.88 ng	310786	Phenanthrene-d10	11.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	91
2			3-Bromo-2-thiophenecarbaldehyde	190	C5H3BrOS	000930-96-1	53
3			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	47
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	38
5			Anthracene, 2-methyl-	192	C15H12	000613-12-7	38



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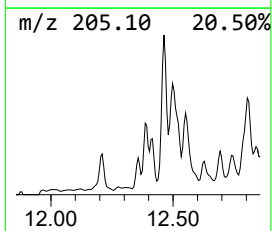
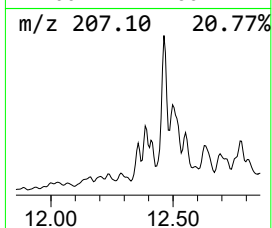
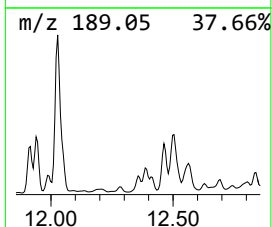
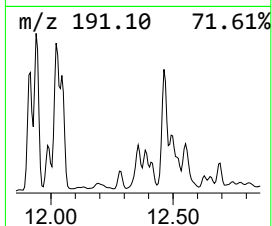
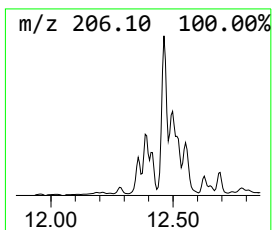
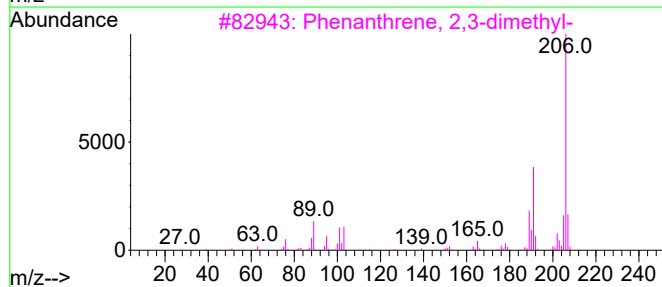
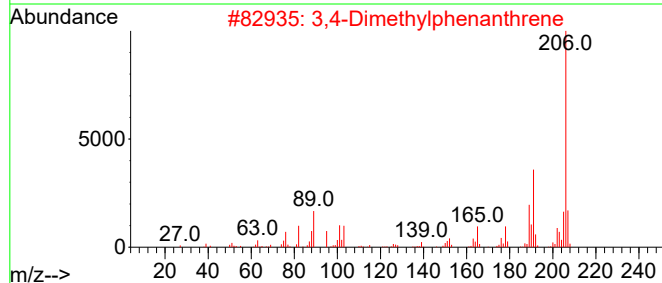
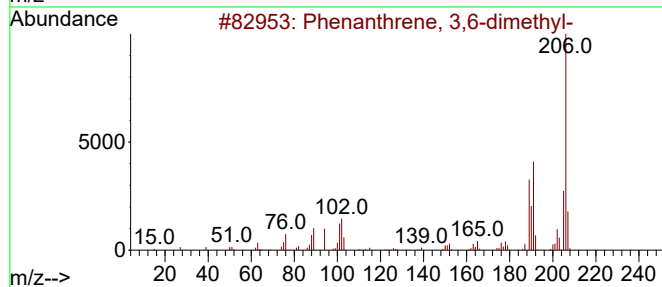
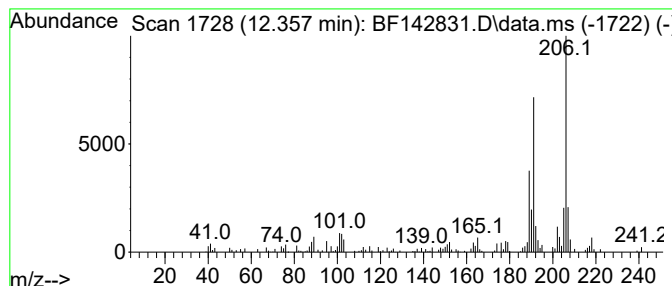
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Phenanthrene, 3,6-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.357	2.32 ng	66241	Phenanthrene-d10	11.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	93
2			3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	91
3			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	90
4			Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	87
5			Anthracene, 2-ethyl-	206	C16H14	052251-71-5	87



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Data File : BF142831.D
Acq On : 24 Jun 2025 21:08
Operator : RC/JU
Sample : Q2384-01 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OK-01-6202025

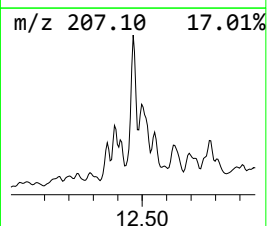
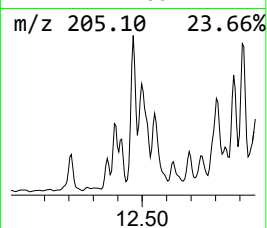
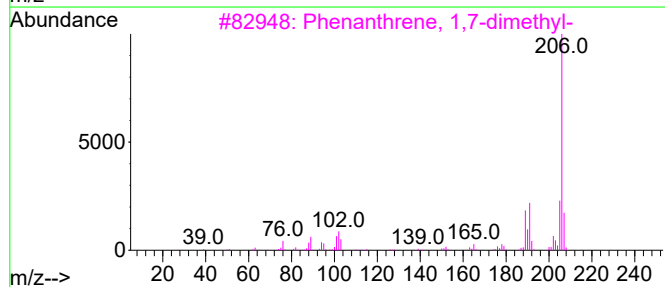
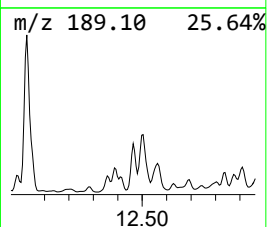
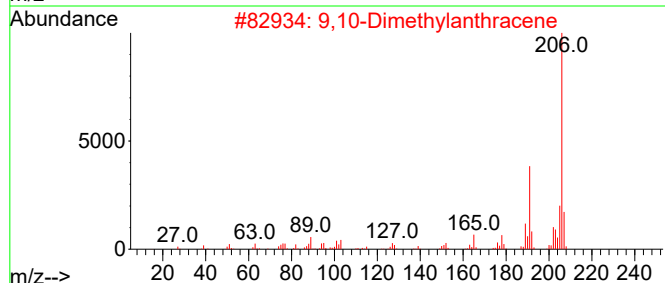
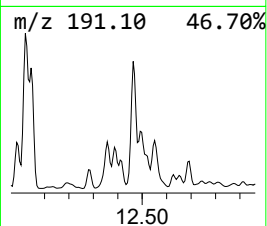
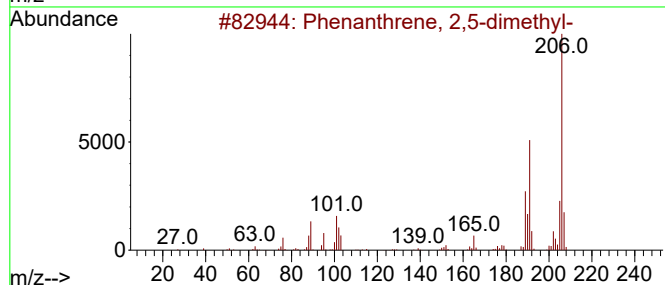
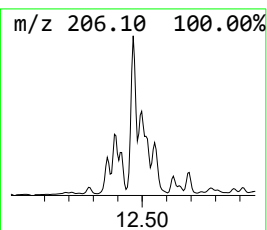
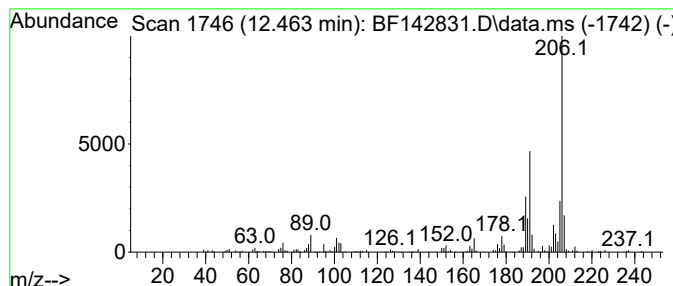
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.463	6.54 ng	186980	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	96
3			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	90
4			di-p-Tolylacetylene	206	C16H14	002789-88-0	90
5			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062425\
Data File : BF142831.D
Acq On : 24 Jun 2025 21:08
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Instrument :
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ClientSampleId :
OK-01-6202025

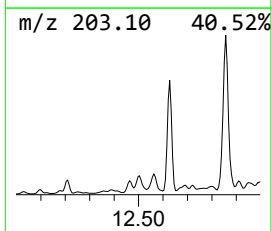
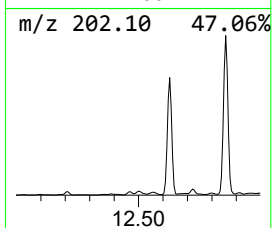
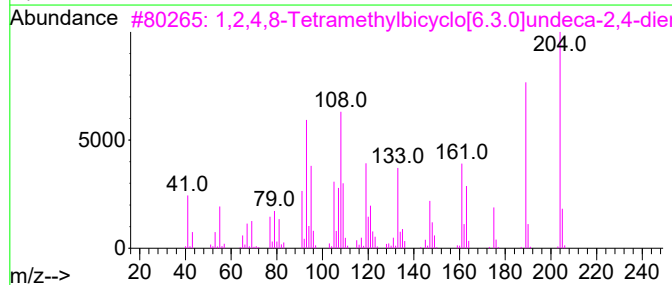
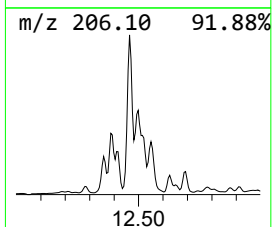
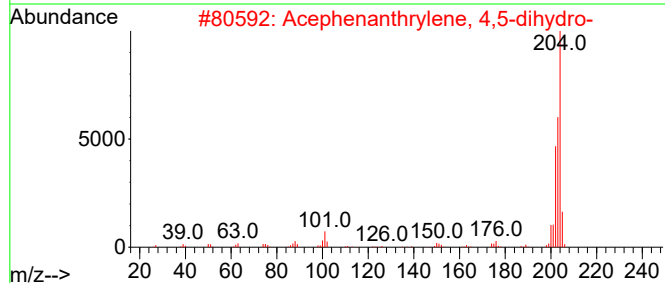
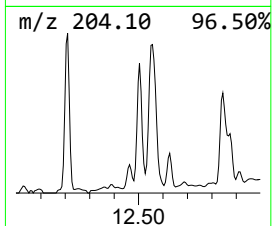
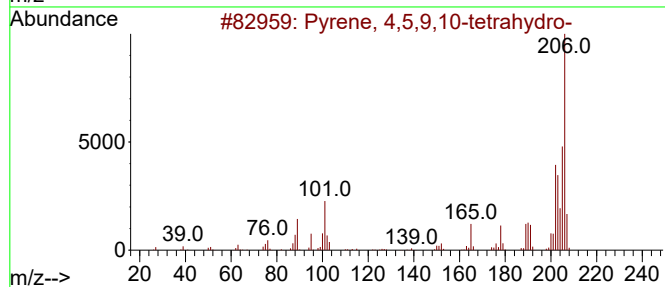
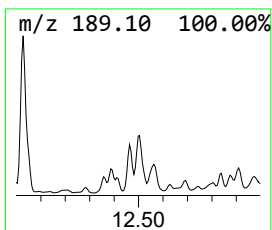
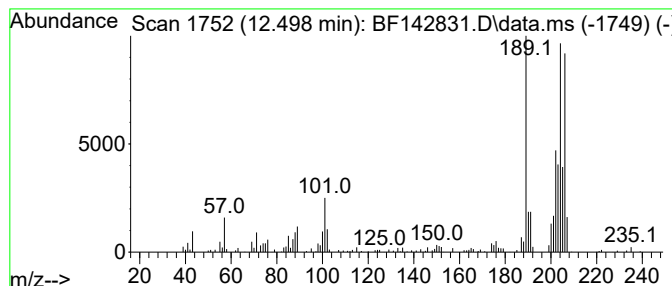
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown12.498 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.498	6.22 ng	177615	Phenanthrene-d10	11.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 4,5,9,10-tetrahydro-	206	C16H14	000781-17-9	42
2			Acephenanthrylene, 4,5-dihydro-	204	C16H12	006232-48-0	38
3			1,2,4,8-Tetramethylbicyclo[6.3.0...]	204	C15H24	137235-51-9	35
4			Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	35
5			5,16[1',2']: ⁸ ,13[1',2'']-Dibenz...	408	C32H24	005672-97-9	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062425\
Data File : BF142831.D
Acq On : 24 Jun 2025 21:08
Operator : RC/JU
Sample : Q2384-01 5X
Misc :
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Instrument :
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ClientSampleId :
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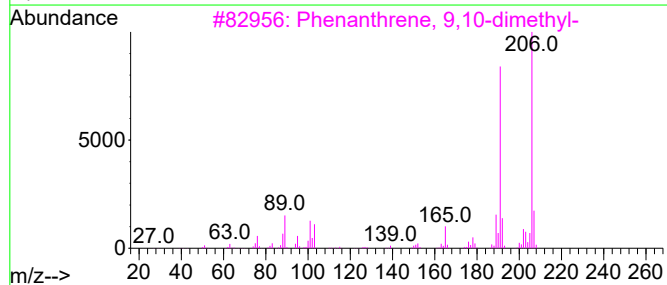
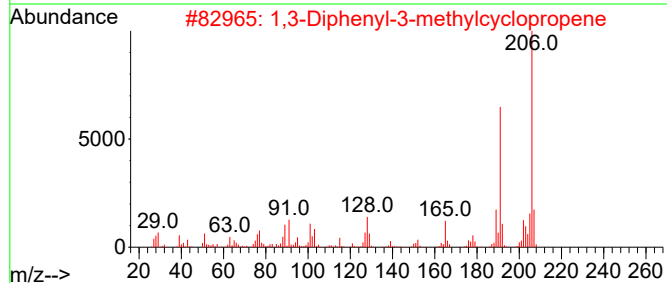
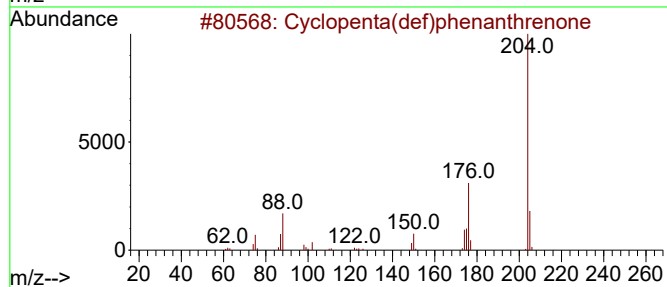
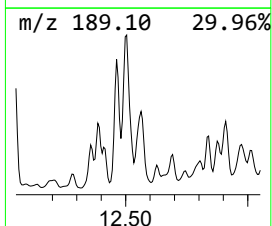
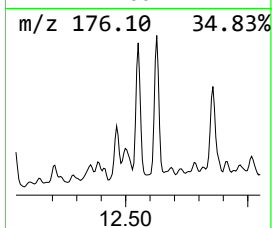
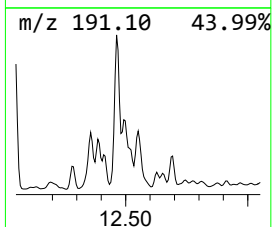
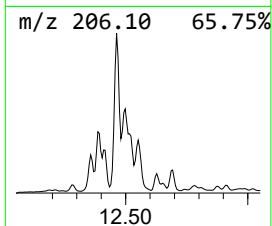
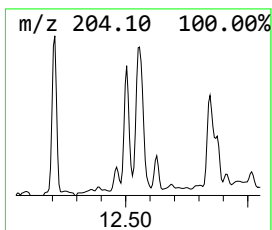
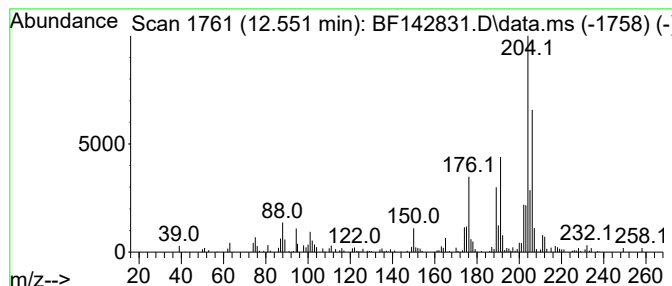
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Cyclopenta(def)phenanthrenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.551	4.25 ng	121396	Phenanthrene-d10	11.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	83
2			1,3-Diphenyl-3-methylcyclopropene	206	C16H14	086544-79-8	80
3			Phenanthrene, 9,10-dimethyl-	206	C16H14	000604-83-1	51
4			9,10-Dimethylanthracene	206	C16H14	000781-43-1	46
5			[14]Annulene, 1,6:8,13-bis(metha...	206	C16H14	085385-68-8	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062425\
Data File : BF142831.D
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Sample : Q2384-01 5X
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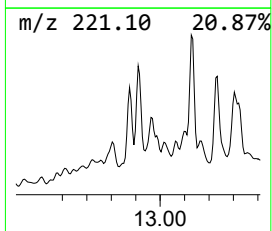
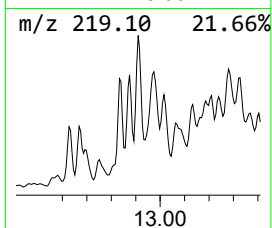
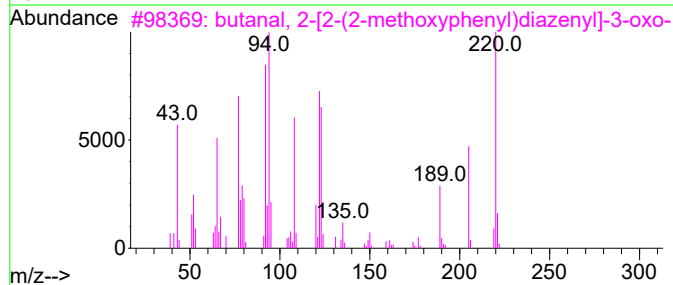
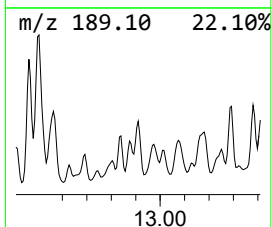
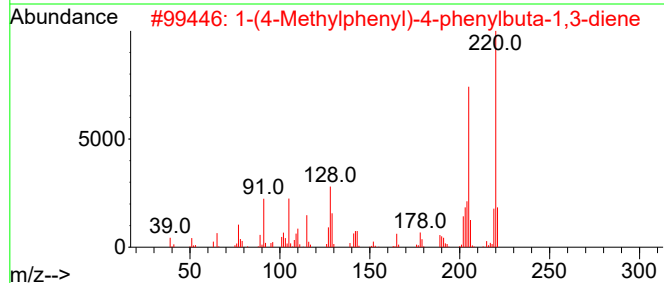
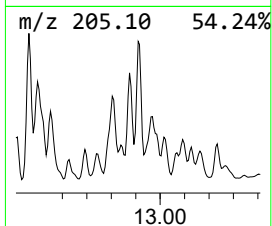
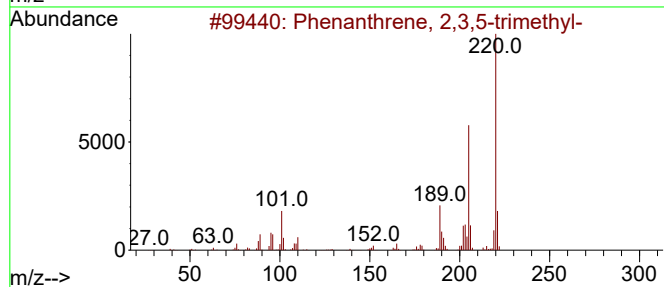
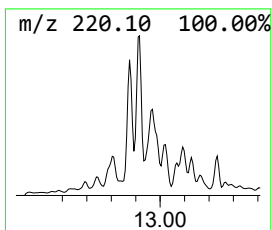
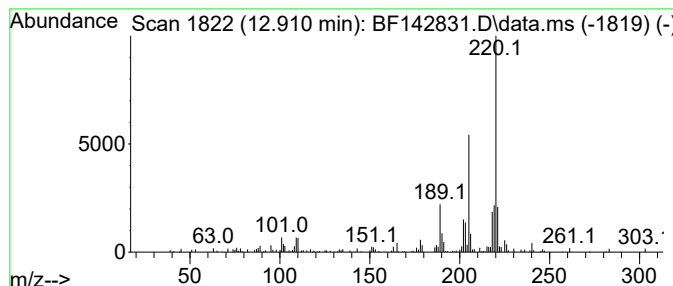
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phenanthrene, 2,3,5-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.910	2.08 ng	61255	Chrysene-d12	14.057	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,3,5-trimethyl-	220	C17H16	003674-73-5	76
2	1-(4-Methylphenyl)-4-phenylbuta-...	220	C17H16	037985-11-8	68
3	butanal, 2-[2-(2-methoxyphenyl)d...	220	C11H12N2O3	1000396-24-5	64
4	9-Ethyl-10-methylanthracene	220	C17H16	019713-49-6	58
5	4-Phenyl-2-naphthol	220	C16H12O	036159-74-7	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062425\
Data File : BF142831.D
Acq On : 24 Jun 2025 21:08
Operator : RC/JU
Sample : Q2384-01 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

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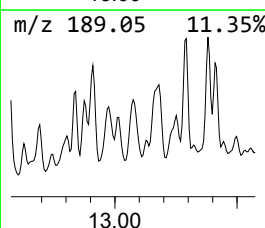
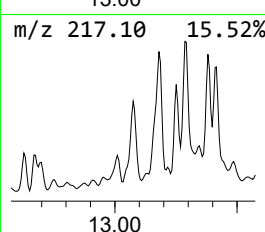
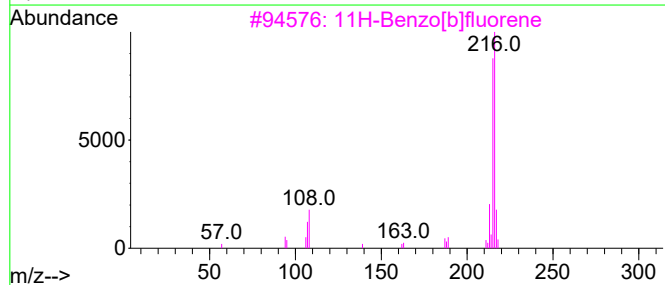
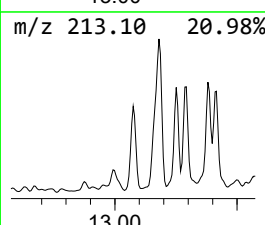
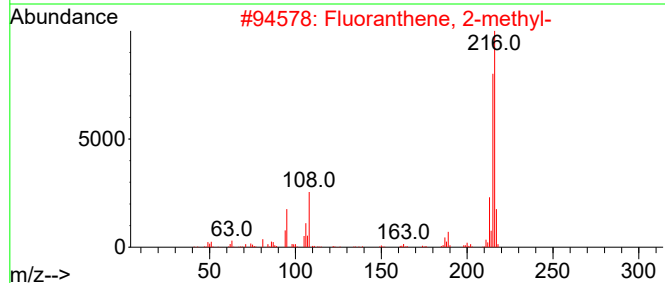
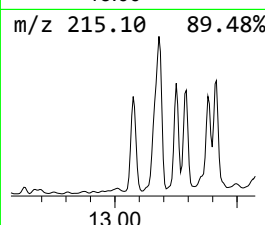
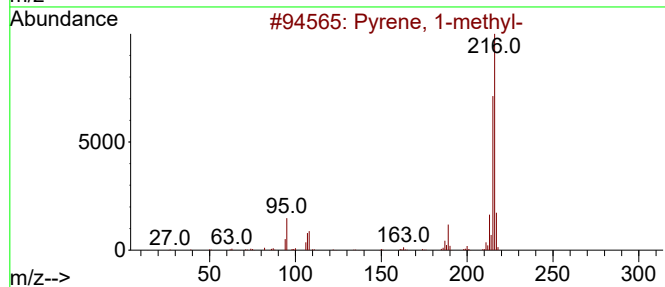
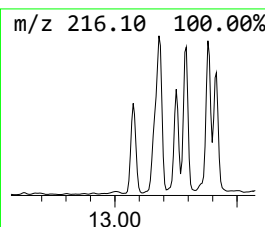
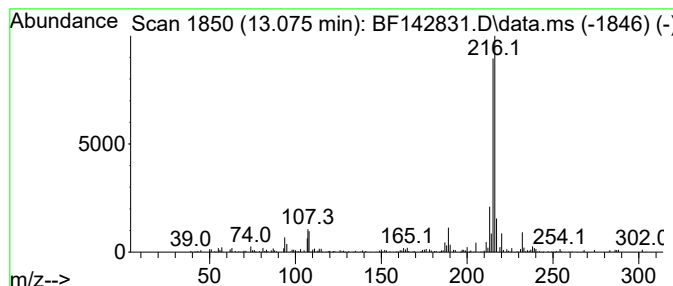
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Pyrene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.075	4.37 ng	129034	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	90
3			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
4			Pyrene, 4-methyl-	216	C17H12	003353-12-6	87
5			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062425\
Data File : BF142831.D
Acq On : 24 Jun 2025 21:08
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Sample : Q2384-01 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

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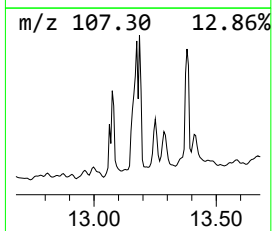
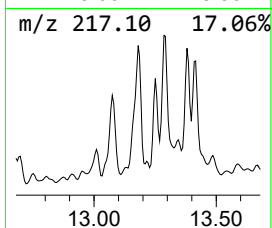
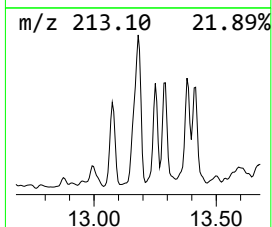
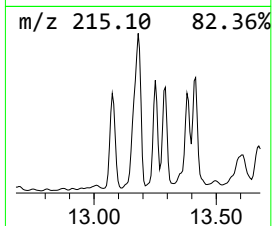
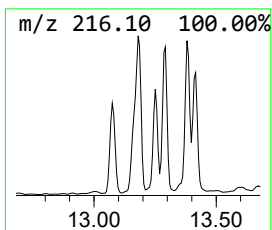
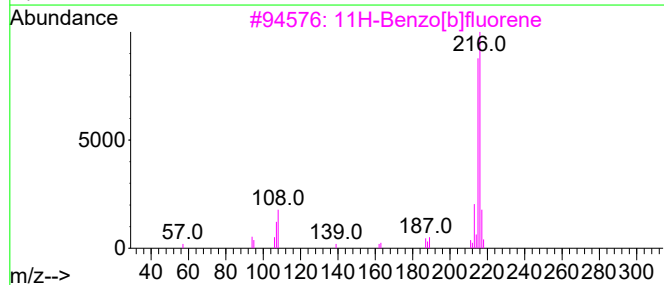
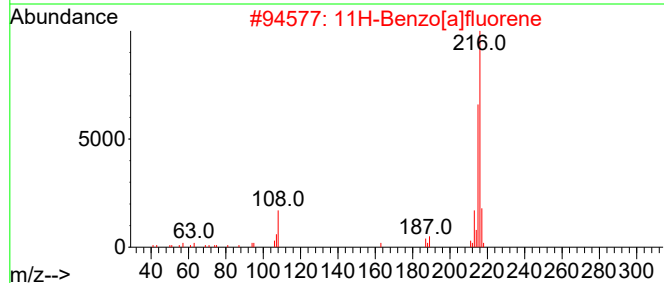
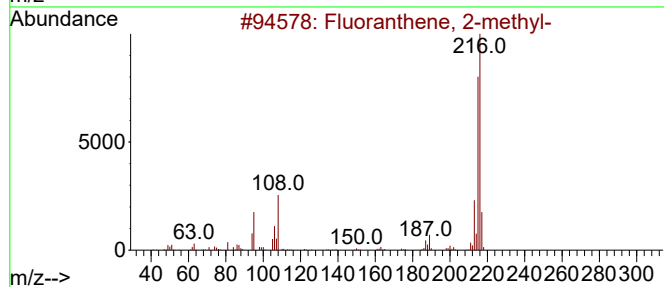
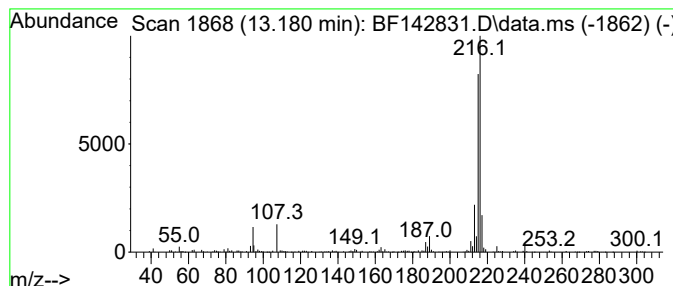
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Fluoranthene, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.180	6.00 ng	177060	Chrysene-d12	14.057

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	93
2			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
3			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
4			Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062425\
Data File : BF142831.D
Acq On : 24 Jun 2025 21:08
Operator : RC/JU
Sample : Q2384-01 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

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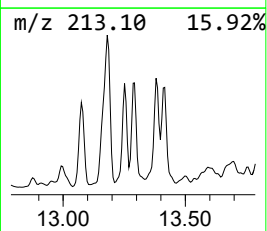
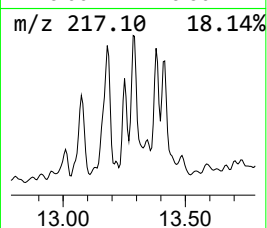
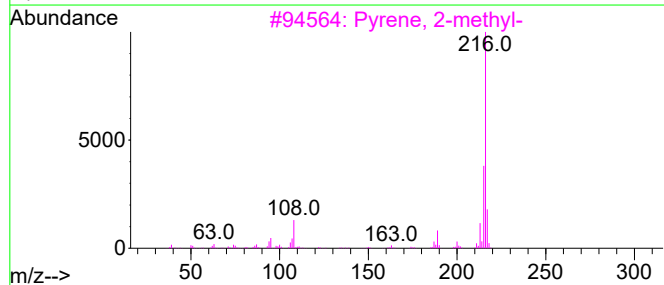
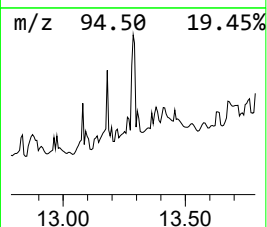
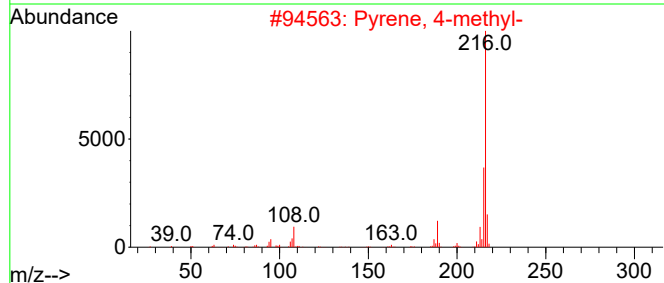
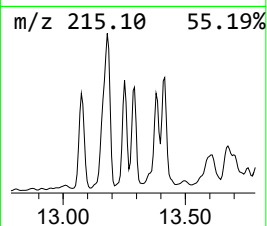
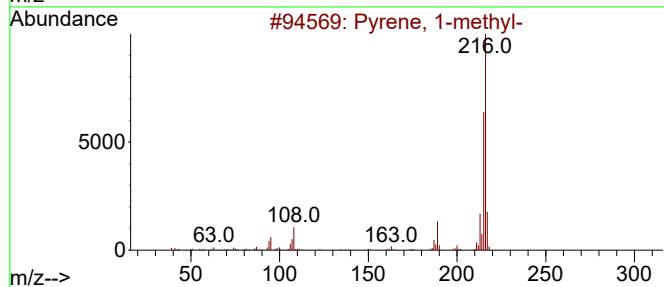
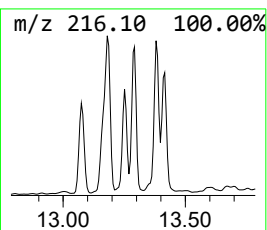
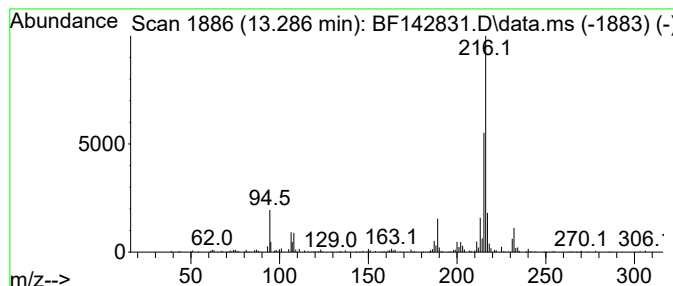
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Pyrene, 4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.286	3.98 ng	117347	Chrysene-d12	14.057

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	96
3			Pyrene, 2-methyl-	216	C17H12	003442-78-2	93
4			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	81
5			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062425\
Data File : BF142831.D
Acq On : 24 Jun 2025 21:08
Operator : RC/JU
Sample : Q2384-01 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OK-01-6202025

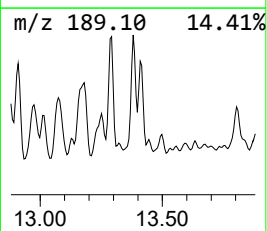
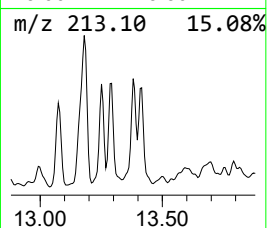
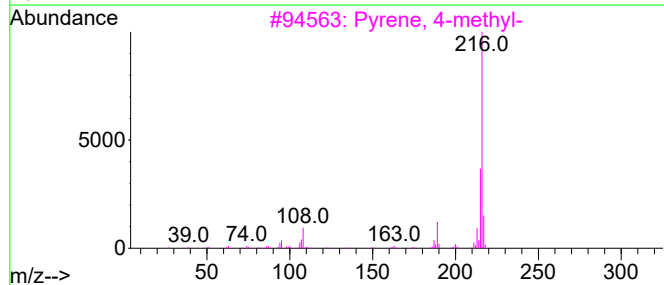
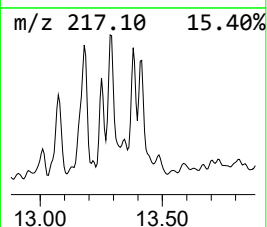
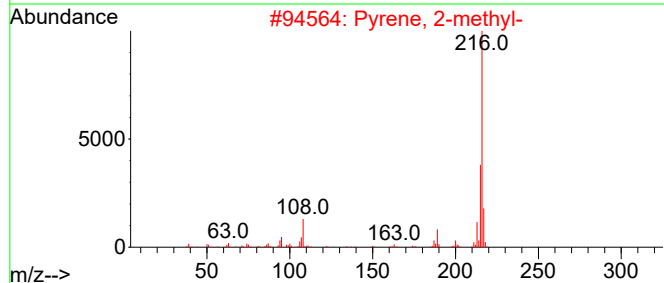
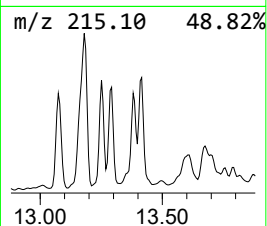
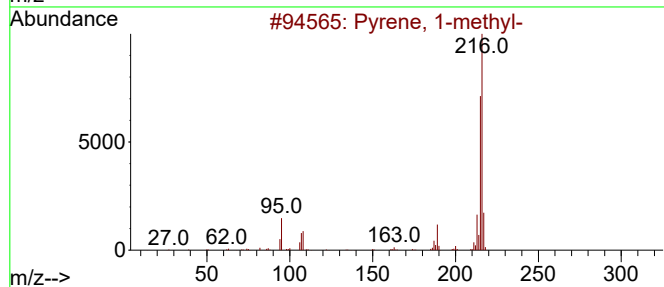
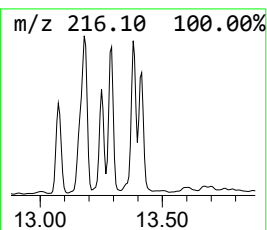
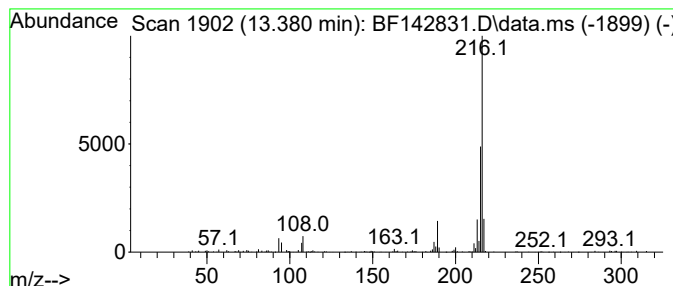
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Pyrene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.380	2.85 ng	84073	Chrysene-d12	14.057

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	96
2			Pyrene, 2-methyl-	216	C17H12	003442-78-2	94
3			Pyrene, 4-methyl-	216	C17H12	003353-12-6	94
4			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
5			1,3-Dihydro-5,6-dimethyl-1-(2-pr...	216	C12H12N2S	080276-22-8	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062425\
Data File : BF142831.D
Acq On : 24 Jun 2025 21:08
Operator : RC/JU
Sample : Q2384-01 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OK-01-6202025

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061925.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
Phenanthrene, 1...	11.916	4.7	ng	134430	4	11.410	571380	20.0
Phenanthrene, 2...	11.939	4.7	ng	133331	4	11.410	571380	20.0
4H-Cyclopenta[d]...	12.028	10.9	ng	310786	4	11.410	571380	20.0
Phenanthrene, 3...	12.357	2.3	ng	66241	4	11.410	571380	20.0
Phenanthrene, 2...	12.463	6.5	ng	186980	4	11.410	571380	20.0
unknown12.498	12.498	6.2	ng	177615	4	11.410	571380	20.0
Cyclopenta(def)...	12.551	4.3	ng	121396	4	11.410	571380	20.0
Phenanthrene, 2...	12.910	2.1	ng	61255	5	14.057	589963	20.0
Pyrene, 1-methyl-	13.075	4.4	ng	129034	5	14.057	589963	20.0
Fluoranthene, 2...	13.180	6.0	ng	177060	5	14.057	589963	20.0
Pyrene, 4-methyl-	13.286	4.0	ng	117347	5	14.057	589963	20.0
Pyrene, 2-methyl-	13.380	2.9	ng	84073	5	14.057	589963	20.0