

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111425\
 Data File : BF144256.D
 Acq On : 14 Nov 2025 11:46
 Operator : RC/JU
 Sample : Q3624-01 2X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-02-11122025

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: BF144256.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.481	554	559	577	rBV	309567	443958	67.11%	5.914%
2	6.493	726	731	748	rBV	304799	441176	66.69%	5.877%
3	6.863	789	794	807	rVB	236963	298428	45.11%	3.976%
4	7.422	884	889	901	rBV	191811	267616	40.45%	3.565%
5	8.145	1004	1012	1019	rBV	266976	361196	54.60%	4.812%
6	8.734	1108	1112	1116	rBV2	5406	7170	1.08%	0.096%
7	8.863	1130	1134	1141	rVB	14173	18642	2.82%	0.248%
8	8.963	1147	1151	1158	rBV2	14936	20280	3.07%	0.270%
9	9.222	1190	1195	1204	rBV	371123	468763	70.86%	6.245%
10	9.298	1204	1208	1211	rBV	8534	10079	1.52%	0.134%
11	9.486	1235	1240	1246	rBV	14579	23714	3.58%	0.316%
12	9.563	1248	1253	1255	rBV	22399	30723	4.64%	0.409%
13	9.586	1255	1257	1260	rVB	6253	6632	1.00%	0.088%
14	9.681	1269	1273	1279	rBV	10721	16202	2.45%	0.216%
15	9.763	1283	1287	1291	rBV	12115	15497	2.34%	0.206%
16	9.828	1291	1298	1303	rVB2	8913	15296	2.31%	0.204%
17	9.904	1303	1311	1315	rBV	255925	334930	50.63%	4.462%
18	9.934	1315	1316	1325	rVB	31752	31041	4.69%	0.414%
19	10.010	1325	1329	1333	rBV4	6969	10375	1.57%	0.138%
20	10.104	1340	1345	1349	rVV2	23782	34727	5.25%	0.463%
21	10.145	1349	1352	1356	rVB	15950	18614	2.81%	0.248%
22	10.222	1361	1365	1367	rBV2	13947	18376	2.78%	0.245%
23	10.251	1367	1370	1375	rVB	9410	11411	1.72%	0.152%
24	10.328	1375	1383	1387	rBV5	17163	35229	5.33%	0.469%
25	10.404	1391	1396	1398	rBV3	5378	7196	1.09%	0.096%
26	10.451	1398	1404	1410	rBV	36426	50371	7.61%	0.671%
27	10.516	1410	1415	1417	rBV4	6741	11788	1.78%	0.157%
28	10.616	1427	1432	1438	rVB	39965	56615	8.56%	0.754%
29	10.692	1440	1445	1457	rVB	153400	215045	32.51%	2.865%
30	10.810	1461	1465	1475	rVB3	19989	37413	5.66%	0.498%
31	10.892	1475	1479	1481	rBV4	6486	8075	1.22%	0.108%
32	10.928	1481	1485	1489	rVB4	5612	7688	1.16%	0.102%
33	10.998	1492	1497	1500	rBV3	18155	28828	4.36%	0.384%
34	11.028	1500	1502	1506	rVV2	12363	15493	2.34%	0.206%
35	11.086	1509	1512	1515	rVB	8952	10422	1.58%	0.139%
36	11.145	1515	1522	1527	rBV8	7248	19241	2.91%	0.256%

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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-BF110525.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	11.251	1535	1540	1543	rBV2	13535	20500	3.10%	0.273%
38	11.292	1543	1547	1554	rVB	18843	28983	4.38%	0.386%
39	11.392	1559	1564	1567	rBV	219506	334435	50.55%	4.455%
40	11.416	1567	1568	1573	rVB	163907	138230	20.90%	1.841%
41	11.469	1573	1577	1581	rBV	32334	39347	5.95%	0.524%
42	11.504	1581	1583	1589	rVB5	6861	9056	1.37%	0.121%
43	11.633	1600	1605	1609	rBV	12234	21279	3.22%	0.283%
44	11.680	1609	1613	1617	rVB2	10457	14708	2.22%	0.196%
45	11.728	1617	1621	1627	rVB	15753	21555	3.26%	0.287%
46	11.810	1631	1635	1639	rBV	9584	11888	1.80%	0.158%
47	11.898	1645	1650	1652	rBV	45097	64326	9.72%	0.857%
48	11.922	1652	1654	1659	rVB	58260	71550	10.82%	0.953%
49	11.975	1659	1663	1665	rBV2	14610	19351	2.93%	0.258%
50	12.010	1665	1669	1678	rVB	72033	144736	21.88%	1.928%
51	12.086	1678	1682	1686	rBV3	10864	14970	2.26%	0.199%
52	12.192	1696	1700	1703	rBV	18442	24509	3.70%	0.327%
53	12.222	1703	1705	1710	rVB3	12472	16875	2.55%	0.225%
54	12.286	1711	1716	1719	rBV6	3962	7774	1.18%	0.104%
55	12.339	1719	1725	1728	rBV2	14118	24710	3.74%	0.329%
56	12.375	1728	1731	1733	rBV	16755	18016	2.72%	0.240%
57	12.445	1739	1743	1746	rBV	46770	68296	10.32%	0.910%
58	12.480	1746	1749	1755	rVV3	38416	66233	10.01%	0.882%
59	12.539	1755	1759	1767	rVB5	19260	38859	5.87%	0.518%
60	12.610	1767	1771	1776	rBV	214206	267498	40.44%	3.564%
61	12.657	1776	1779	1780	rBV2	5659	7052	1.07%	0.094%
62	12.763	1794	1797	1803	rBV	15476	21830	3.30%	0.291%
63	12.839	1803	1810	1816	rBV	230069	347152	52.48%	4.625%
64	12.892	1816	1819	1823	rVB2	20146	26743	4.04%	0.356%
65	12.980	1826	1834	1842	rVB	297245	424112	64.11%	5.650%
66	13.063	1842	1848	1854	rBV4	31089	66121	9.99%	0.881%
67	13.116	1854	1857	1859	rBV3	9240	10729	1.62%	0.143%
68	13.169	1859	1866	1870	rVB	49917	89883	13.59%	1.197%
69	13.233	1870	1877	1880	rBV2	37906	64593	9.76%	0.861%
70	13.274	1880	1884	1892	rVB2	39846	58288	8.81%	0.777%
71	13.369	1896	1900	1902	rBV	29472	37207	5.62%	0.496%
72	13.398	1902	1905	1908	rVB2	22085	24618	3.72%	0.328%
73	13.551	1928	1931	1933	rVB3	16802	22552	3.41%	0.300%
74	13.792	1968	1972	1974	rBV	24021	34641	5.24%	0.461%
75	14.039	2008	2014	2026	rBV3	321027	661544	100.00%	8.813%
76	14.427	2077	2080	2083	rBV	27486	35707	5.40%	0.476%

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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-BF110525.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	15.092	2188	2193	2200	rBV	72870	159716	24.14%	2.128%
78	15.392	2242	2244	2250	rVB	43385	59504	8.99%	0.793%
79	15.457	2252	2255	2261	rVV	65877	109747	16.59%	1.462%
80	15.521	2261	2266	2278	rVB	190257	348762	52.72%	4.646%

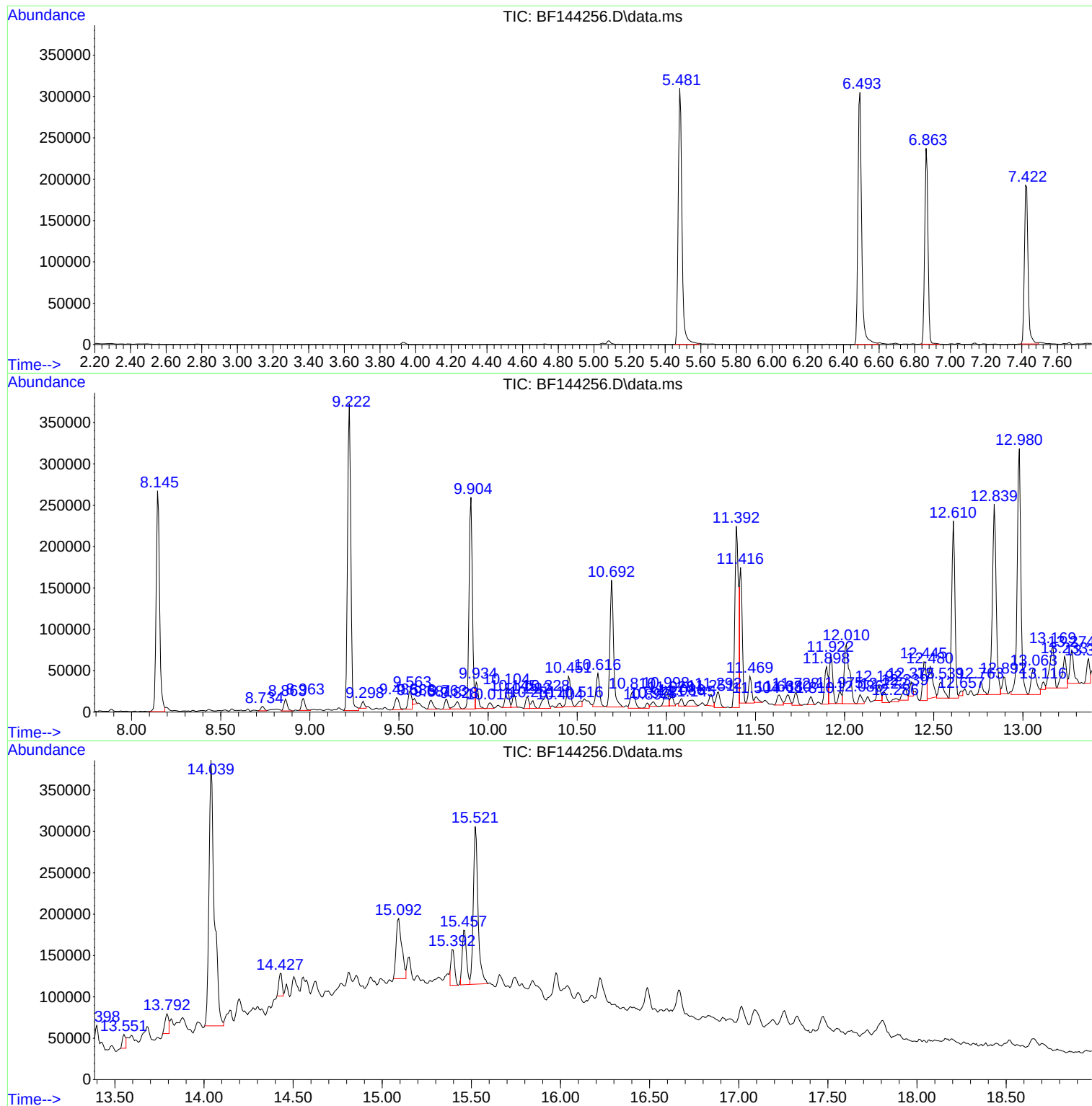
Sum of corrected areas: 7506435

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Data File : BF144256.D
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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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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



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111425\
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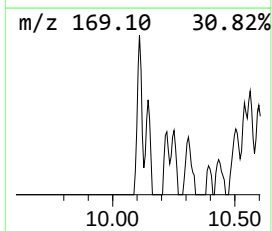
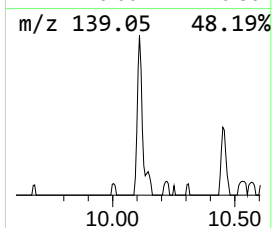
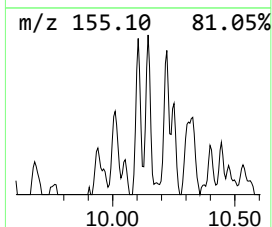
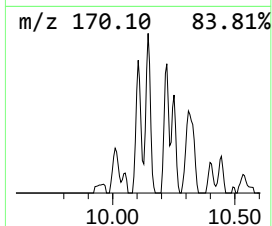
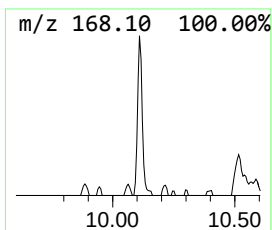
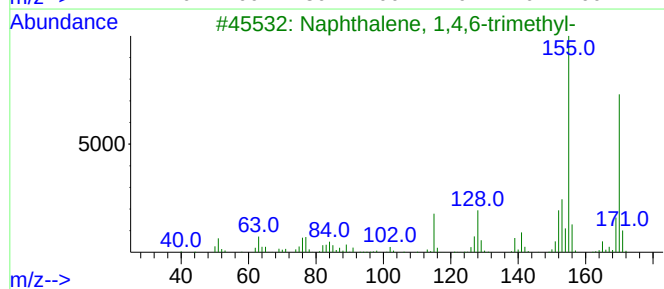
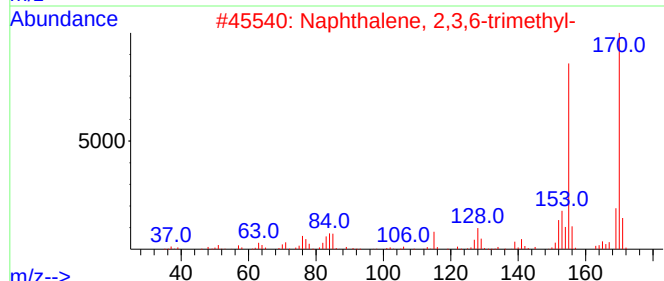
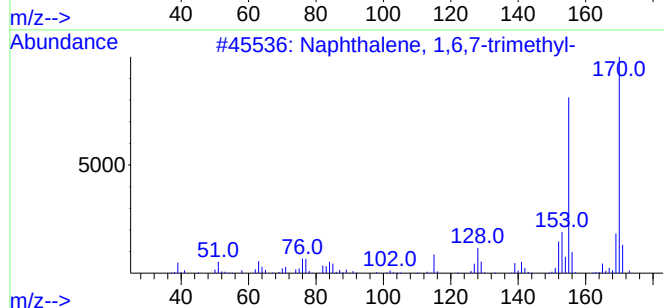
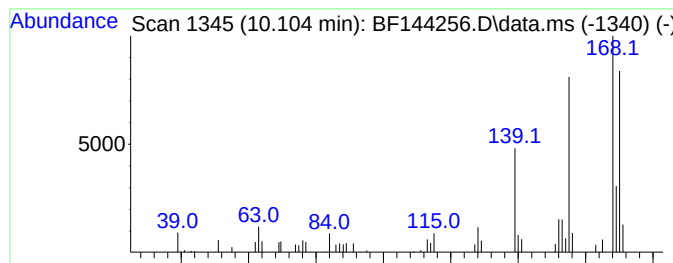
Instrument :
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ClientSampleId :
OK-02-11122025

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 1 Naphthalene, 1,6,7-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.104	2.07 ng	34727	Acenaphthene-d10	9.904	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	93
2	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	91
3	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	78
4	1-(2,2-Dimethylcyclopropyl)-2-ph...	170	C13H14	1000294-91-1	60
5	4,6,8-Trimethylazulene	170	C13H14	000941-81-1	53



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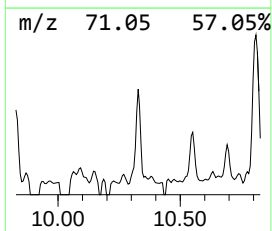
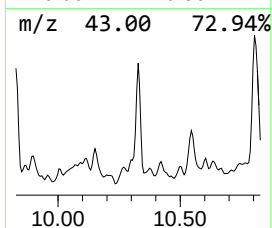
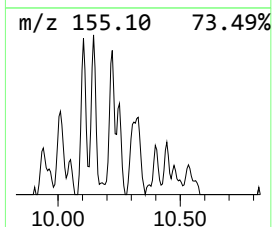
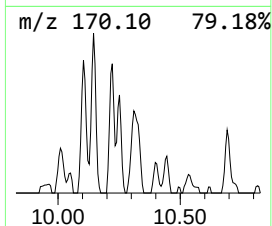
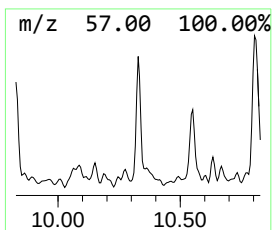
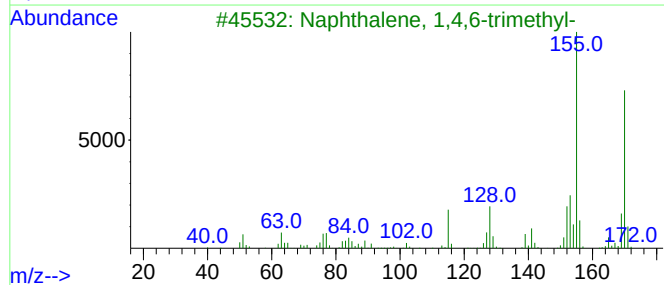
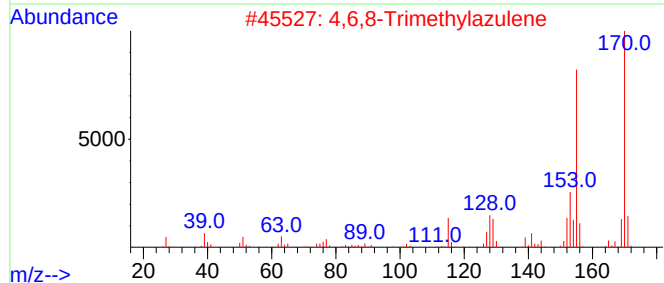
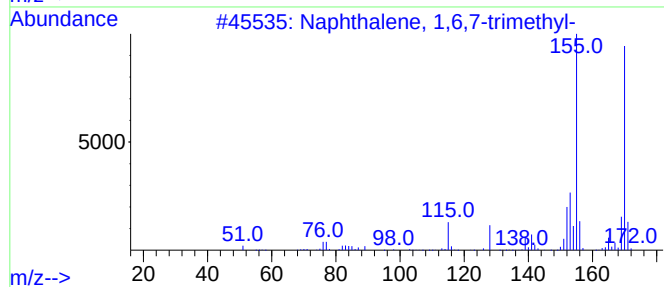
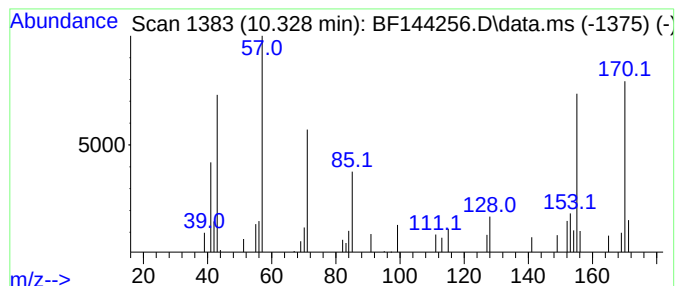
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 4,6,8-Trimethylazulene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.328	2.10 ng	35229	Acenaphthene-d10	9.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	93
2			4,6,8-Trimethylazulene	170	C13H14	000941-81-1	93
3			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	76
4			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	76
5			Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	76



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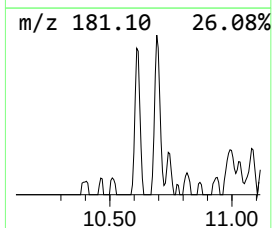
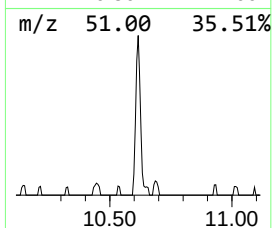
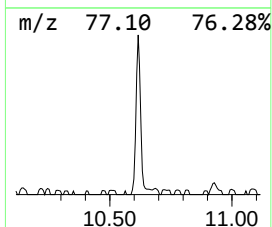
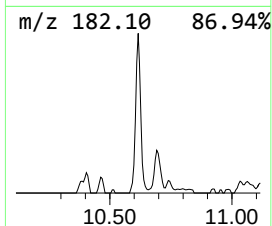
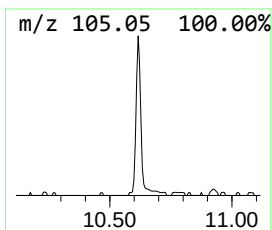
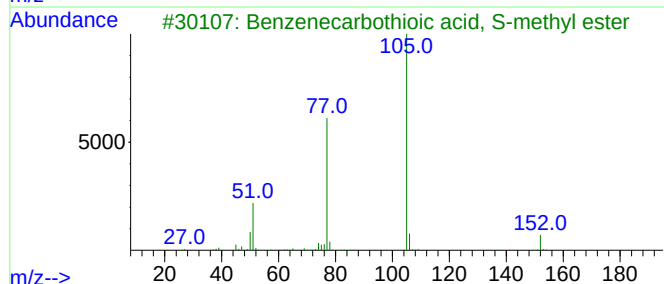
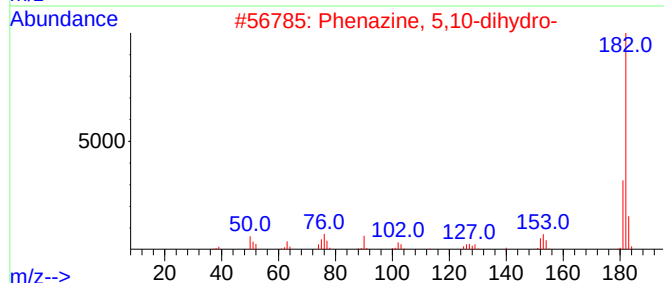
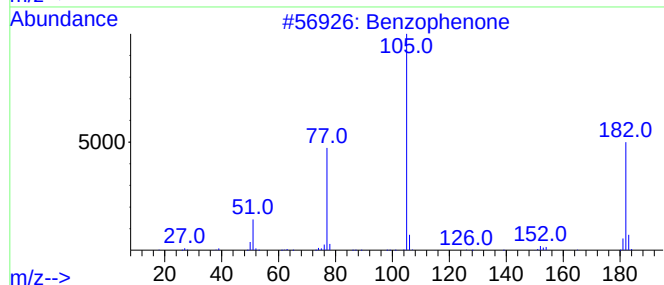
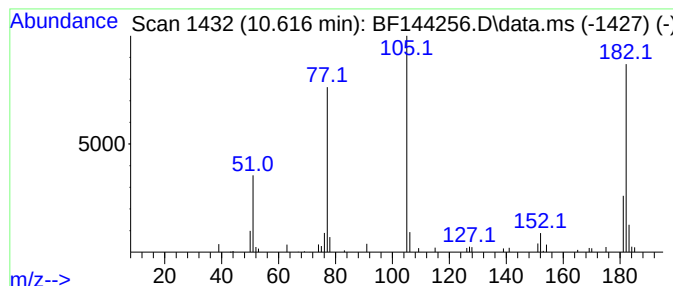
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Benzophenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.616	3.38 ng	56615	Acenaphthene-d10	9.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	91
2			Phenazine, 5,10-dihydro-	182	C12H10N2	000613-32-1	46
3			Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	43
4			S-Benzoyl(thiohydroxylamine)	153	C7H7NOS	025740-80-1	43
5			Benzoylformic acid	150	C8H6O3	000611-73-4	38



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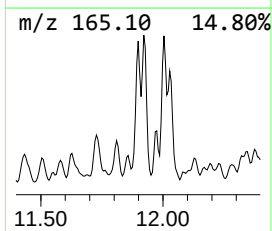
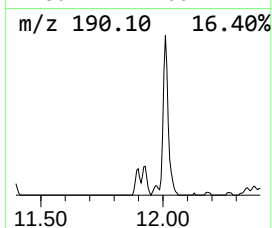
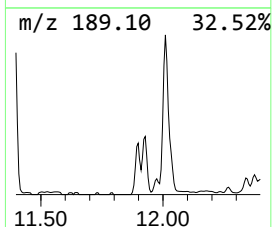
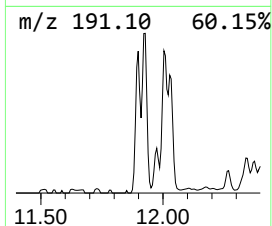
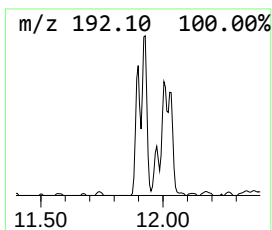
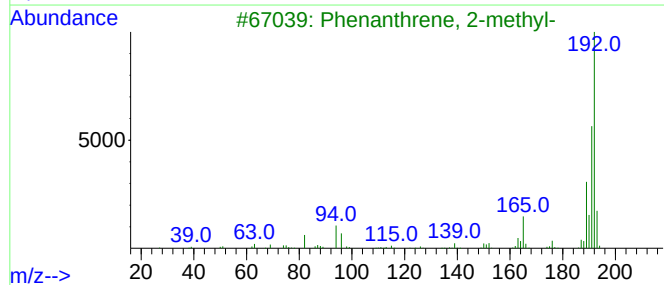
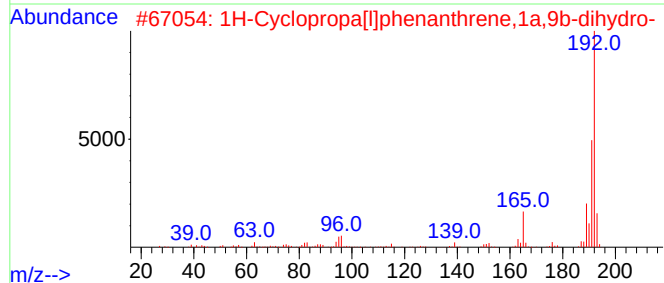
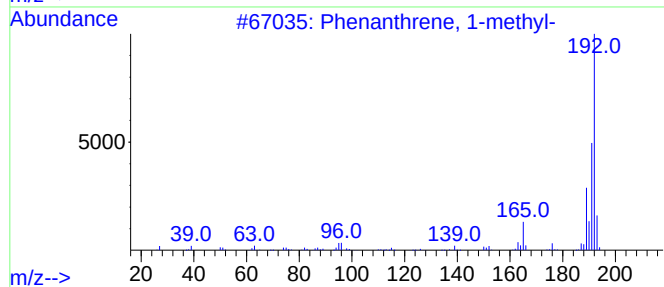
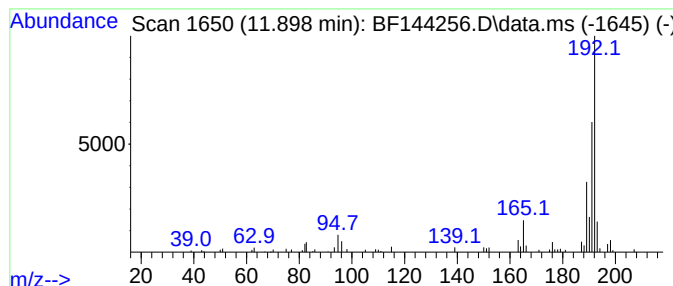
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.898	3.85 ng	64326	Phenanthrene-d10	11.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4			Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
5			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111425\
Data File : BF144256.D
Acq On : 14 Nov 2025 11:46
Operator : RC/JU
Sample : Q3624-01 2X
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OK-02-11122025

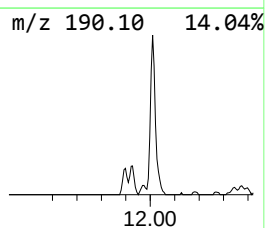
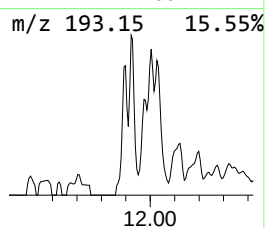
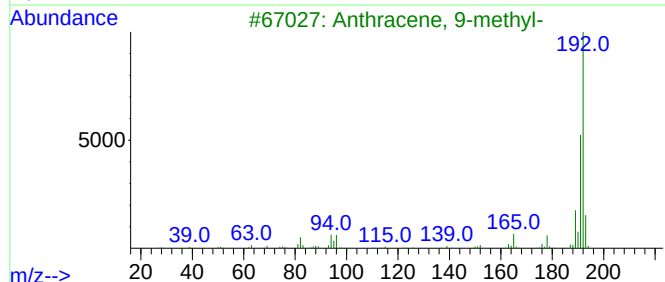
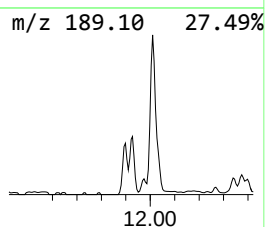
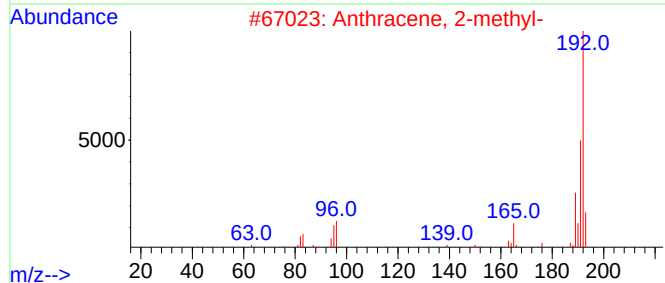
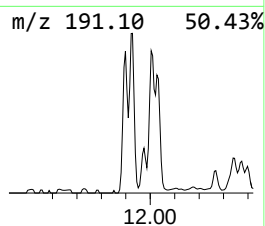
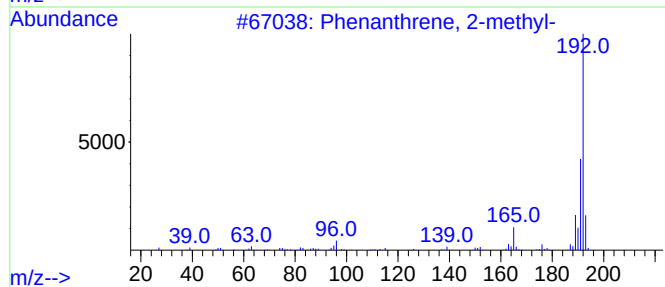
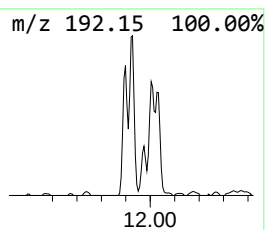
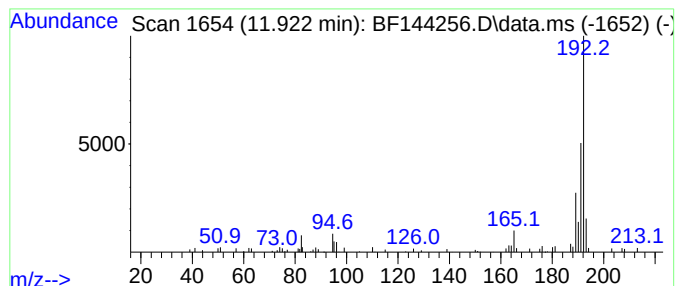
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.922	4.28 ng	71550	Phenanthrene-d10	11.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
4			1H-Indene, 1-phenyl-	192	C15H12	001961-96-2	93
5			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111425\
Data File : BF144256.D
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Operator : RC/JU
Sample : Q3624-01 2X
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Instrument :
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ClientSampleId :
OK-02-11122025

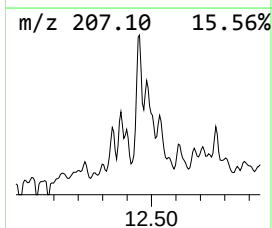
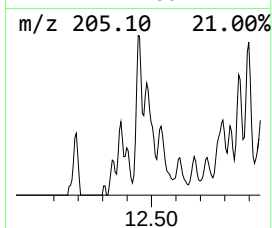
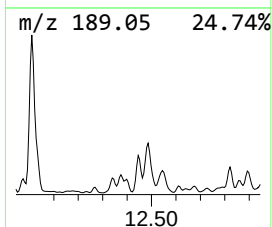
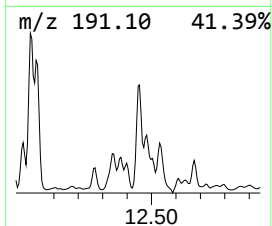
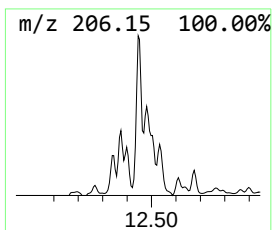
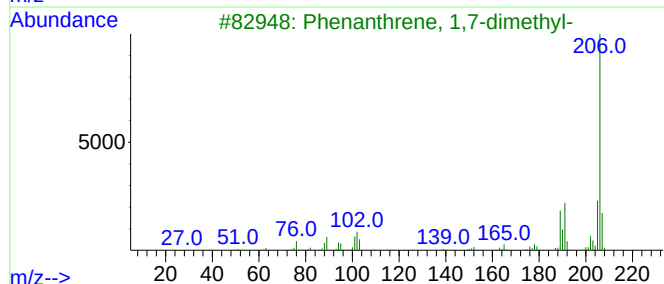
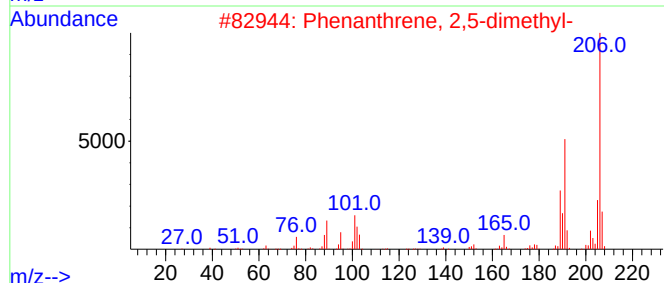
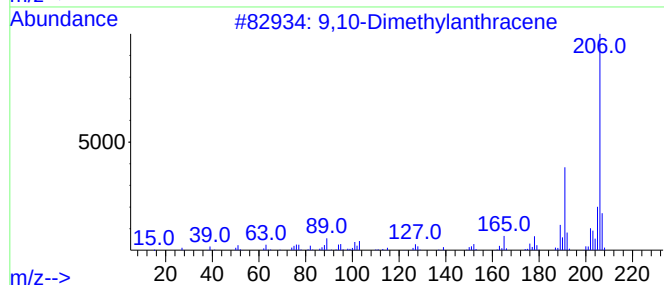
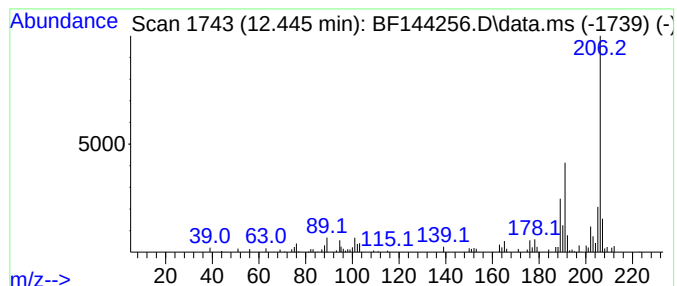
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 9,10-Dimethylantracene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.445	4.08 ng	68296	Phenanthrene-d10	11.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Dimethylantracene	206	C16H14	000781-43-1	96
2			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	95
3			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	93
4			3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	93
5			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111425\
Data File : BF144256.D
Acq On : 14 Nov 2025 11:46
Operator : RC/JU
Sample : Q3624-01 2X
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OK-02-11122025

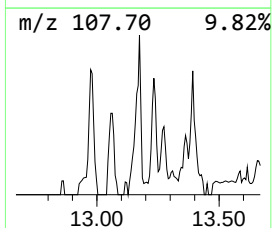
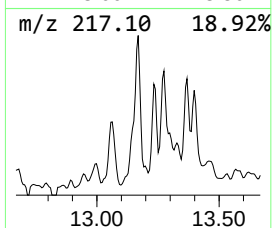
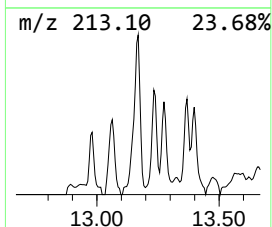
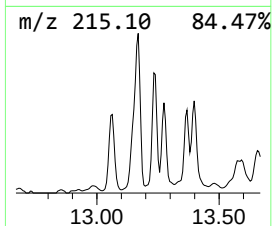
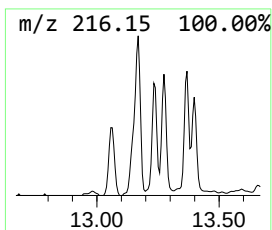
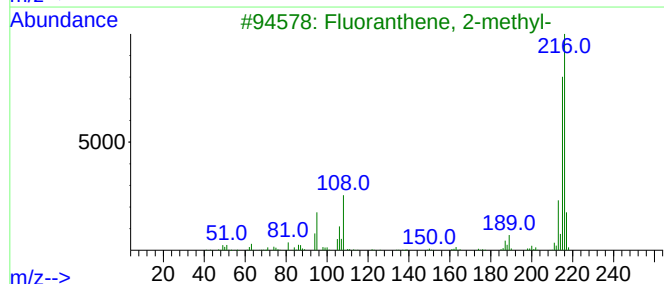
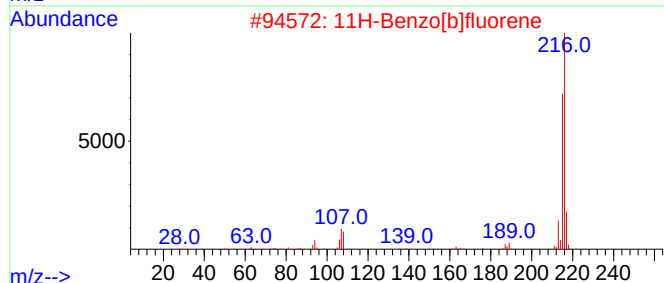
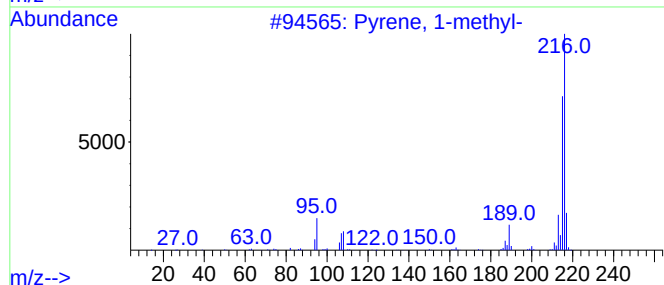
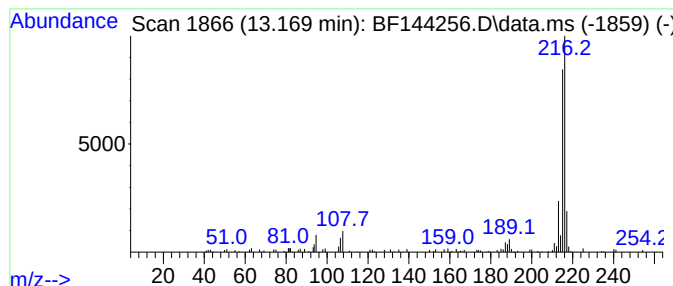
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Pyrene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.169	2.72 ng	89883	Chrysene-d12	14.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
3			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
4			Pyrene, 2-methyl-	216	C17H12	003442-78-2	87
5			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111425\
Data File : BF144256.D
Acq On : 14 Nov 2025 11:46
Operator : RC/JU
Sample : Q3624-01 2X
Misc :
ALS Vial : 5 Sample Multiplier: 1

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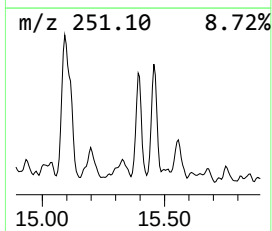
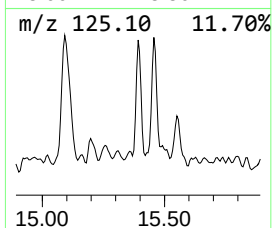
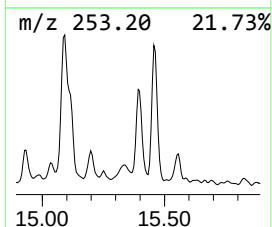
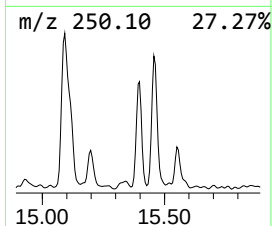
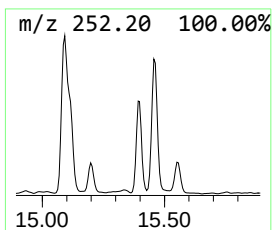
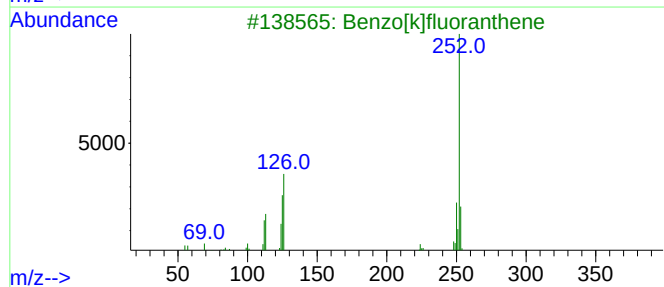
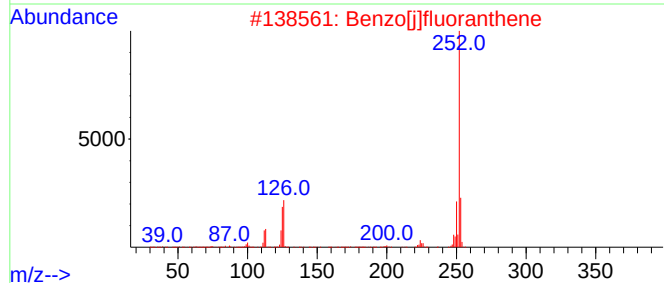
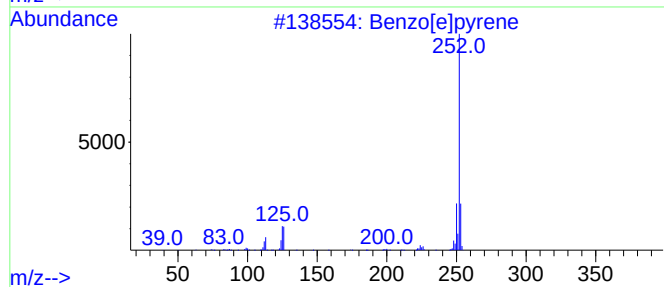
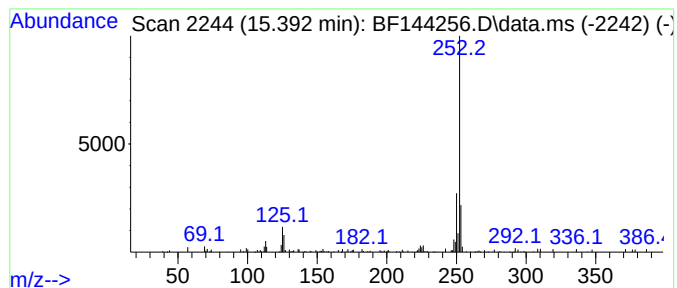
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzo[e]pyrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.392	3.41 ng	59504	Perylene-d12	15.521

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	95
2			Benzo[j]fluoranthene	252	C20H12	000205-82-3	94
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
4			Benzo[a]pyrene	252	C20H12	000050-32-8	93
5			Benzo[b]fluoranthene	252	C20H12	000205-99-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111425\
Data File : BF144256.D
Acq On : 14 Nov 2025 11:46
Operator : RC/JU
Sample : Q3624-01 2X
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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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\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
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4,6,8-Trimethyl...	10.328	2.1	ng	35229	3	9.904	334930	20.0
Benzophenone	10.616	3.4	ng	56615	3	9.904	334930	20.0
Phenanthrene, 1...	11.898	3.9	ng	64326	4	11.392	334435	20.0
Phenanthrene, 2...	11.922	4.3	ng	71550	4	11.392	334435	20.0
9,10-Dimethylan...	12.445	4.1	ng	68296	4	11.392	334435	20.0
Pyrene, 1-methyl-	13.169	2.7	ng	89883	5	14.045	661544	20.0
Benzo[e]pyrene	15.392	3.4	ng	59504	6	15.521	348762	20.0