

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101325\
 Data File : BF143940.D
 Acq On : 13 Oct 2025 18:58
 Operator : RC/JU
 Sample : Q3323-05
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SED-05-0-2

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: BF143940.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.963	288	301	306	rBV	11477	24594	1.33%	0.105%
2	5.098	485	494	500	rVB2	22452	39883	2.16%	0.170%
3	5.498	553	562	574	rBV	1324192	1789889	97.13%	7.618%
4	5.910	624	632	635	rBV3	9617	19871	1.08%	0.085%
5	6.134	665	670	673	rBV2	21814	32797	1.78%	0.140%
6	6.322	697	702	705	rBV3	13398	23189	1.26%	0.099%
7	6.422	714	719	725	rBV4	20087	47691	2.59%	0.203%
8	6.504	725	733	741	rBV	1304087	1842691	100.00%	7.842%
9	6.634	750	755	758	rBV4	14758	24546	1.33%	0.104%
10	6.716	765	769	771	rBV3	13167	21258	1.15%	0.090%
11	6.845	787	791	792	rVV3	17368	20933	1.14%	0.089%
12	6.881	792	797	803	rVV	548532	714280	38.76%	3.040%
13	6.939	803	807	810	rVV2	18103	26624	1.44%	0.113%
14	6.992	810	816	819	rVV4	28940	57768	3.13%	0.246%
15	7.028	819	822	825	rVV	33793	45547	2.47%	0.194%
16	7.063	825	828	833	rVB	26431	31630	1.72%	0.135%
17	7.151	839	843	847	rVB3	38857	52720	2.86%	0.224%
18	7.275	858	864	867	rBV5	30917	57795	3.14%	0.246%
19	7.363	875	879	882	rBV3	13646	22516	1.22%	0.096%
20	7.439	882	892	902	rVV	759957	1098450	59.61%	4.675%
21	7.522	902	906	910	rBV3	59724	93063	5.05%	0.396%
22	7.581	912	916	920	rVV5	20817	38554	2.09%	0.164%
23	7.663	926	930	931	rBV	29269	36243	1.97%	0.154%
24	7.686	932	934	942	rVB5	56940	85090	4.62%	0.362%
25	7.757	943	946	948	rBV4	21760	28424	1.54%	0.121%
26	7.804	951	954	957	rVB4	40000	43928	2.38%	0.187%
27	7.839	957	960	963	rBV3	28952	36715	1.99%	0.156%
28	7.922	969	974	977	rBV4	24941	45389	2.46%	0.193%
29	7.981	981	984	989	rVB5	39834	58399	3.17%	0.249%
30	8.039	990	994	997	rBV2	42294	61039	3.31%	0.260%
31	8.163	1009	1015	1022	rBV2	628197	1264318	68.61%	5.381%
32	8.222	1022	1025	1028	rVB	133114	152059	8.25%	0.647%
33	8.316	1038	1041	1043	rBV	33649	44221	2.40%	0.188%
34	8.369	1047	1050	1058	rVB3	73297	118451	6.43%	0.504%
35	8.457	1060	1065	1070	rBV6	62212	110832	6.01%	0.472%
36	8.510	1071	1074	1077	rBV3	36561	49383	2.68%	0.210%

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Integrator: RTE

Smoothing : ON

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Max Peaks: 100

Stop Thrs : 0

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

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

37	8.581	1083	1086	1089	rBV	93785	128757	6.99%	0.548%
38	8.669	1098	1101	1104	rBV2	34565	44473	2.41%	0.189%
39	8.704	1104	1107	1112	rVV3	39609	61967	3.36%	0.264%
40	8.881	1134	1137	1140	rVB	69053	76282	4.14%	0.325%

41	8.975	1150	1153	1161	rVB	80887	128285	6.96%	0.546%
42	9.186	1186	1189	1193	rVB	78309	93320	5.06%	0.397%
43	9.239	1193	1198	1203	rVB	1103519	1358535	73.73%	5.782%
44	9.322	1208	1212	1221	rBV2	56825	144332	7.83%	0.614%
45	9.439	1229	1232	1236	rVB2	32010	38774	2.10%	0.165%

46	9.575	1252	1255	1258	rBV	35387	45163	2.45%	0.192%
47	9.639	1263	1266	1269	rVB2	56897	66874	3.63%	0.285%
48	9.786	1287	1291	1295	rVB2	29573	40005	2.17%	0.170%
49	9.922	1309	1314	1317	rBV	593320	741812	40.26%	3.157%
50	9.957	1317	1320	1324	rVB	193114	227753	12.36%	0.969%

51	10.080	1337	1341	1344	rBV3	25406	37697	2.05%	0.160%
52	10.127	1344	1349	1353	rVV	128883	171426	9.30%	0.730%
53	10.239	1365	1368	1370	rBV3	22201	28385	1.54%	0.121%
54	10.333	1380	1384	1390	rBV4	28341	57646	3.13%	0.245%
55	10.469	1403	1407	1413	rBV	146343	195007	10.58%	0.830%

56	10.563	1420	1423	1428	rVB2	26300	42886	2.33%	0.183%
57	10.633	1431	1435	1440	rVB2	185663	239944	13.02%	1.021%
58	10.710	1443	1448	1453	rBV	601390	787841	42.75%	3.353%
59	10.839	1465	1470	1476	rVB2	49322	94237	5.11%	0.401%
60	11.216	1531	1534	1538	rVB	29726	37271	2.02%	0.159%

61	11.269	1538	1543	1546	rBV3	50367	87653	4.76%	0.373%
62	11.310	1546	1550	1556	rVB	81106	117403	6.37%	0.500%
63	11.416	1563	1568	1569	rBV	484363	615019	33.38%	2.618%
64	11.439	1569	1572	1576	rVV	912499	1148958	62.35%	4.890%
65	11.486	1576	1580	1584	rVB	169490	218746	11.87%	0.931%

66	11.645	1603	1607	1614	rVB	60747	90845	4.93%	0.387%
67	11.939	1649	1657	1662	rBV2	229065	441987	23.99%	1.881%
68	11.992	1663	1666	1669	rVV	52580	79899	4.34%	0.340%
69	12.027	1669	1672	1680	rVB2	146987	265279	14.40%	1.129%
70	12.210	1700	1703	1711	rVB2	62134	108303	5.88%	0.461%

71	12.469	1744	1747	1750	rBV	35201	46147	2.50%	0.196%
72	12.633	1770	1775	1779	rBV	737024	956425	51.90%	4.071%
73	12.863	1807	1814	1819	rBV	657351	1037122	56.28%	4.414%
74	12.998	1832	1837	1845	rVB	726171	1039335	56.40%	4.423%
75	13.186	1866	1869	1873	rVB	95755	125053	6.79%	0.532%

76	13.257	1878	1881	1883	rVV2	48034	54683	2.97%	0.233%
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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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	13.292	1884	1887	1891	rVV2	50362	64780	3.52%	0.276%
78	14.057	2011	2017	2021	rBV2	692995	1243209	67.47%	5.291%
79	15.110	2191	2196	2205	rBV	293969	669647	36.34%	2.850%
80	15.415	2244	2248	2254	rVB	161230	258910	14.05%	1.102%
81	15.480	2254	2259	2265	rVV	245105	380409	20.64%	1.619%
82	15.545	2265	2270	2287	rVB	483748	933724	50.67%	3.974%
83	17.039	2519	2524	2533	rVB3	91829	207519	11.26%	0.883%
84	17.492	2596	2601	2611	rVB	68004	155798	8.45%	0.663%

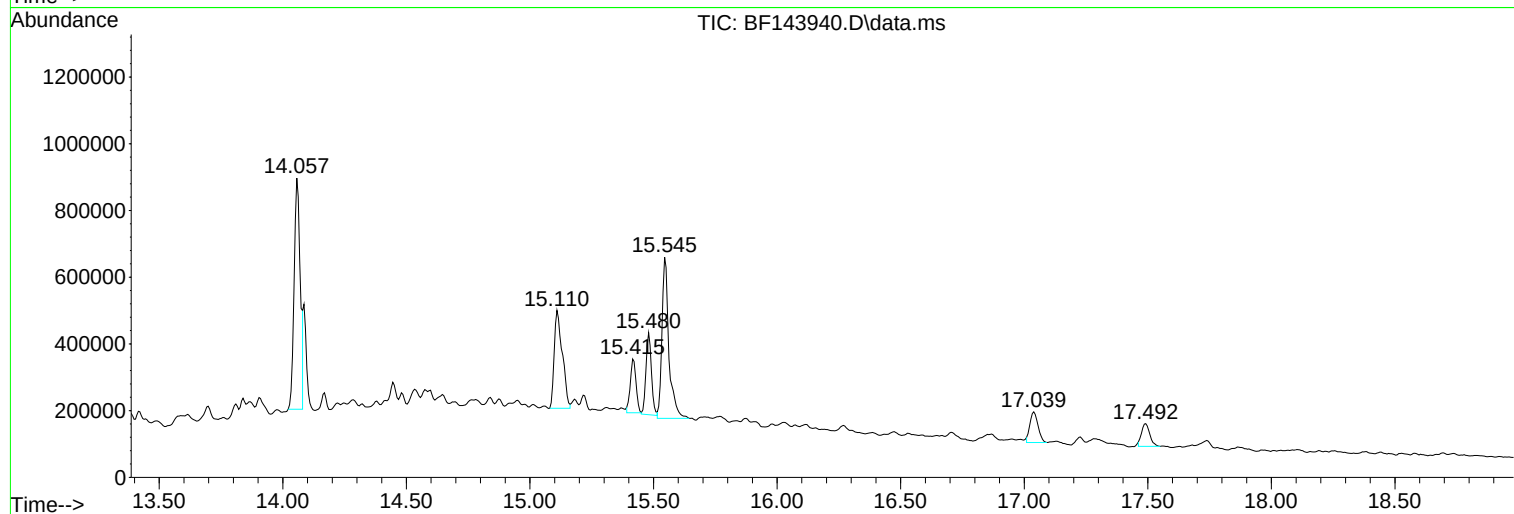
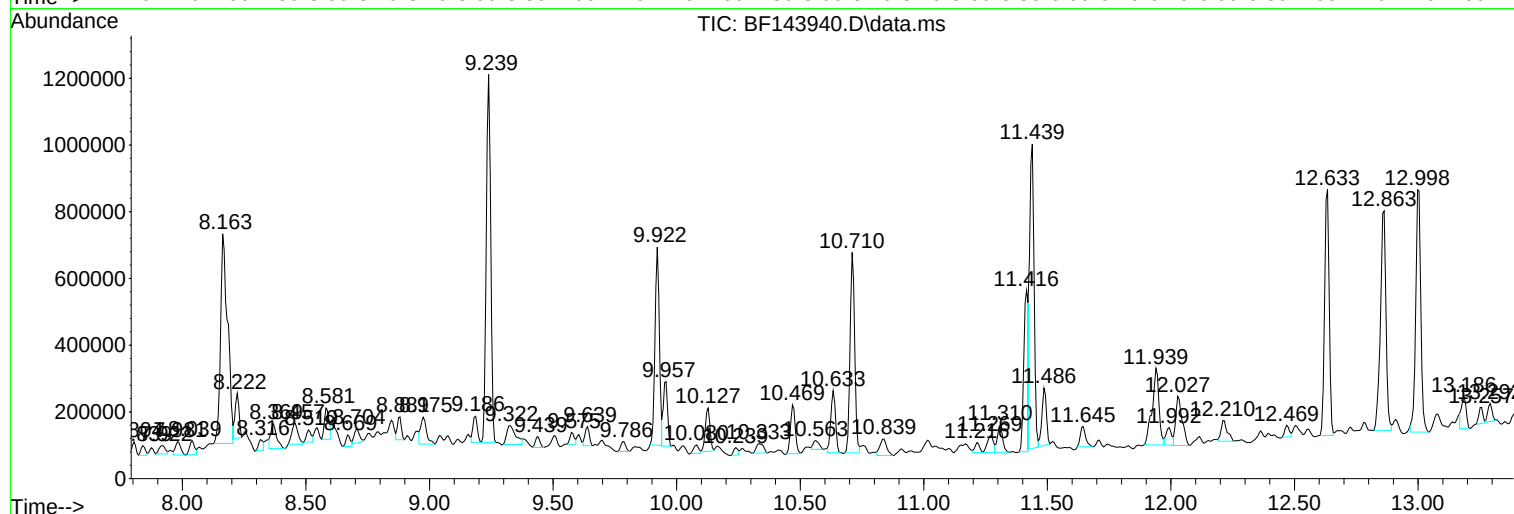
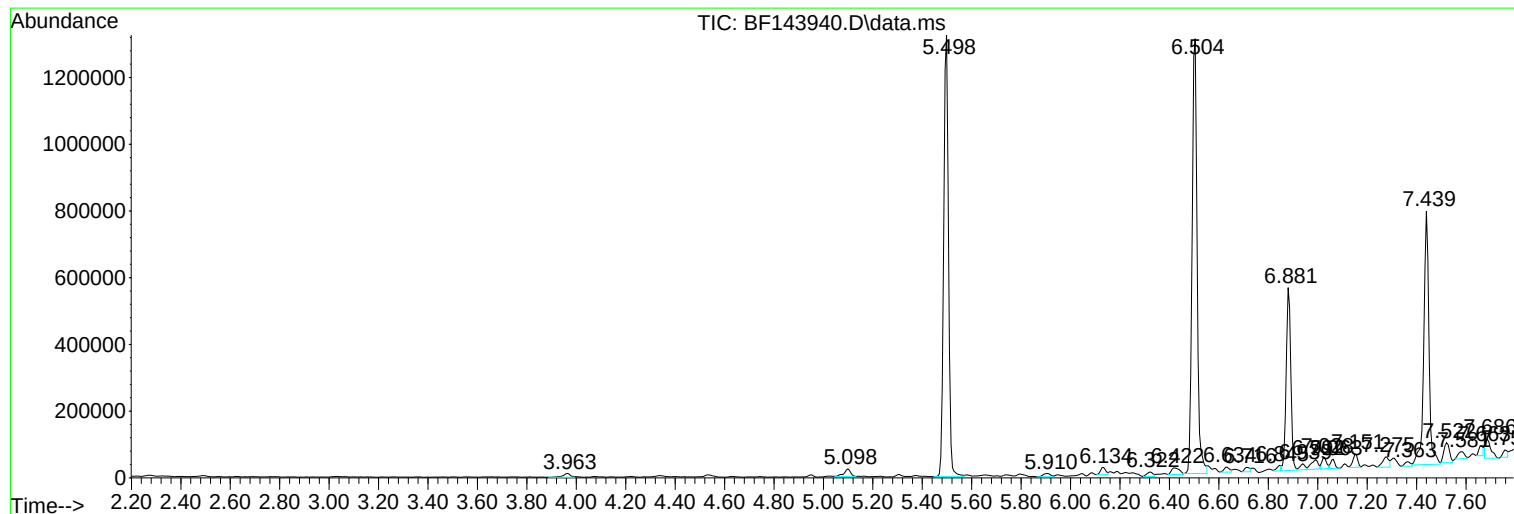
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ALS Vial : 11 Sample Multiplier: 1

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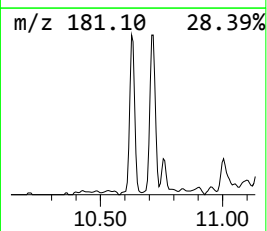
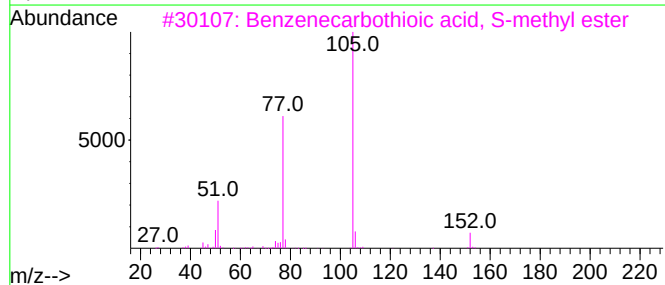
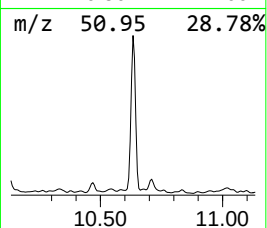
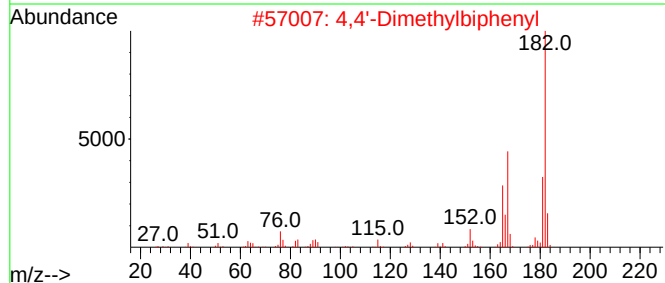
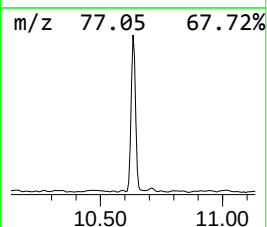
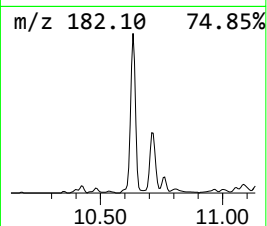
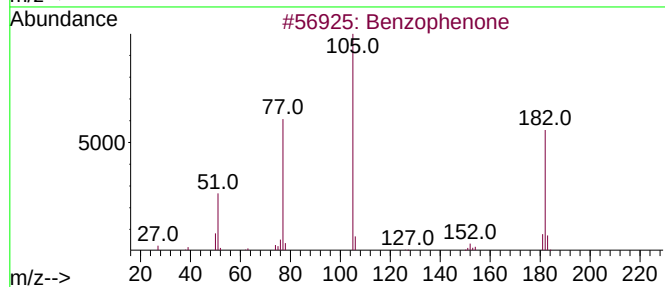
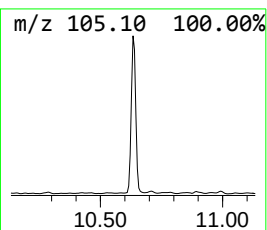
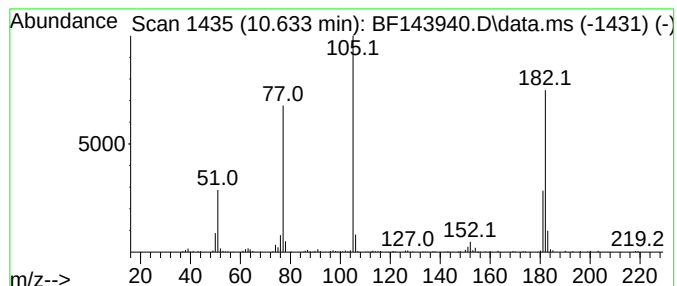
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100625.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.633	6.47 ng	239944	Acenaphthene-d10	9.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	94
2			4,4'-Dimethylbiphenyl	182	C14H14	000613-33-2	43
3			Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	41
4			Benzoic acid, 2,4-dimethoxy-	182	C9H10O4	000091-52-1	35
5			Methyl benzilate	242	C15H14O3	000076-89-1	35



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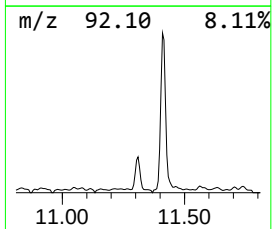
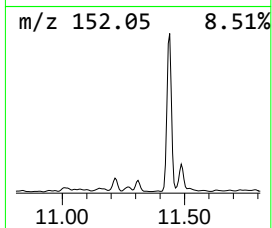
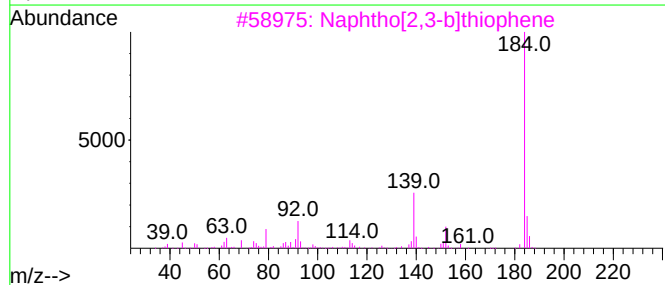
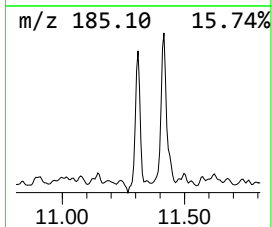
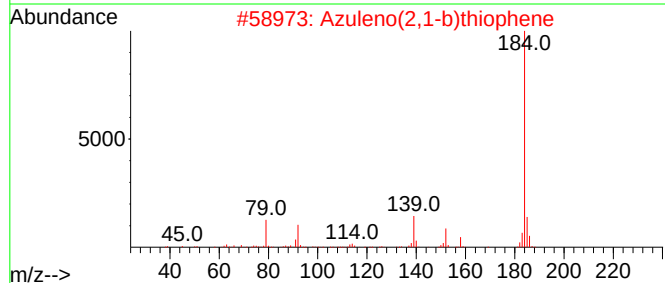
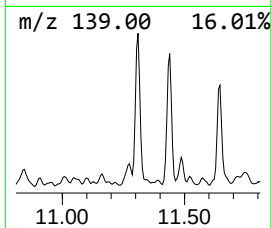
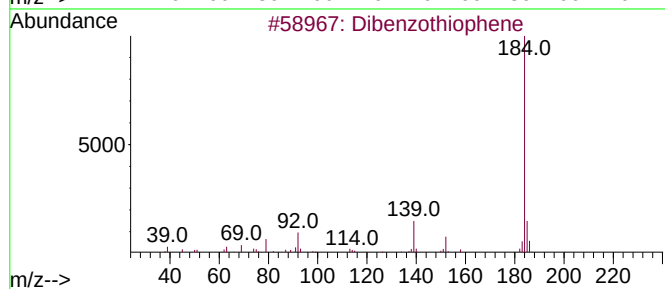
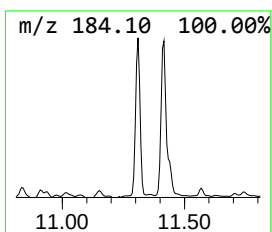
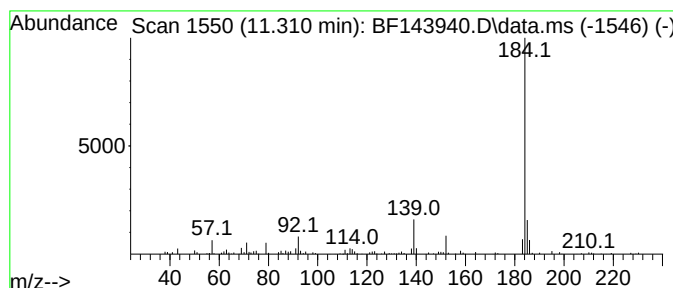
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100625.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Dibenzothiophene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.310	3.82 ng	117403	Phenanthrene-d10	11.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzothiophene	184	C12H8S	000132-65-0	97
2			Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	91
3			Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	90
4			Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	87
5			Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	87



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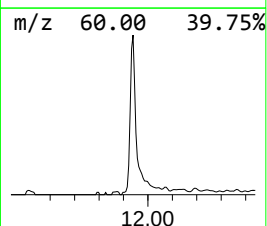
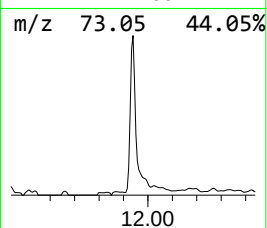
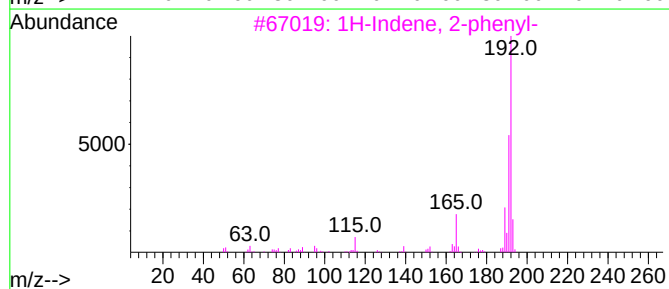
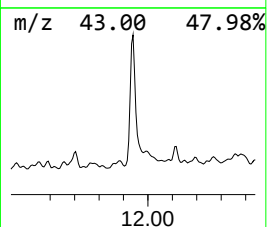
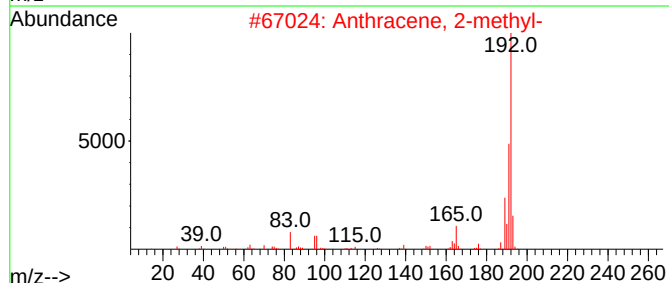
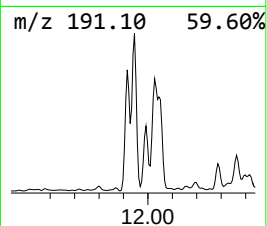
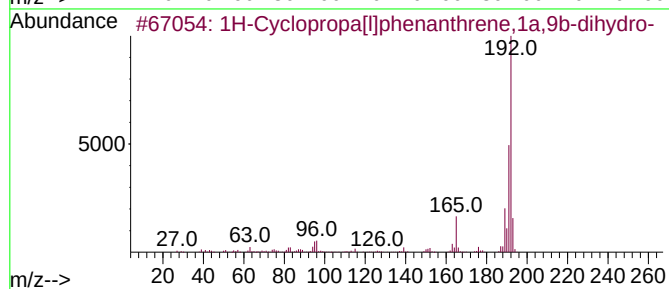
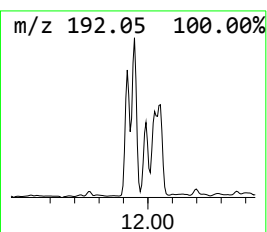
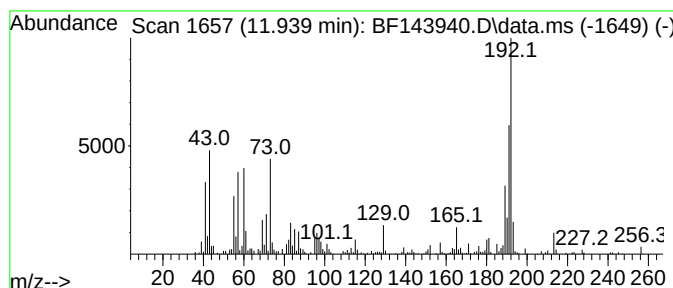
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100625.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1H-Cyclopropa[l]phenanthren... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.939	14.37 ng	441987	Phenanthrene-d10		11.416	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1H-Cyclopropa[1]phenanthrene,1a,...		192	C15H12	000949-41-7	90
2	Anthracene, 2-methyl-		192	C15H12	000613-12-7	90
3	1H-Indene, 2-phenyl-		192	C15H12	004505-48-0	87
4	Phenanthrene, 1-methyl-		192	C15H12	000832-69-9	83
5	Anthracene, 1-methyl-		192	C15H12	000610-48-0	78



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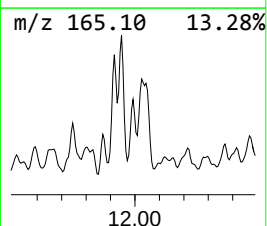
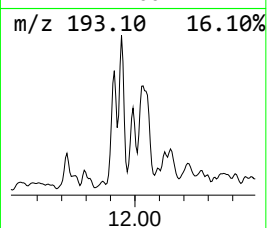
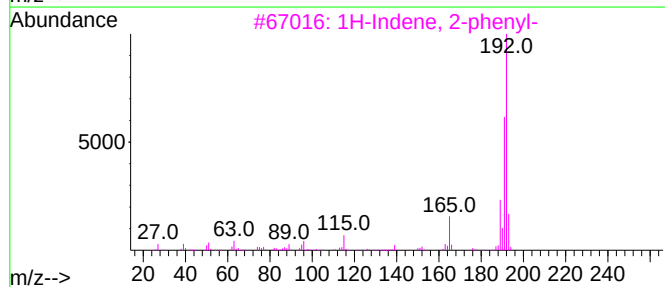
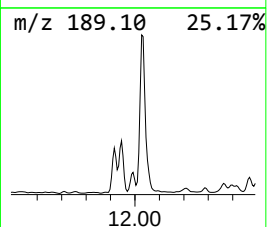
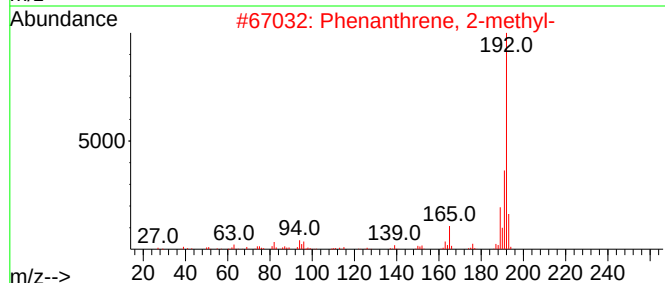
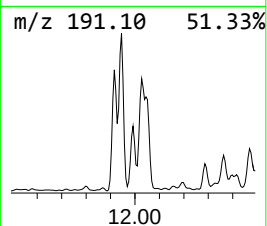
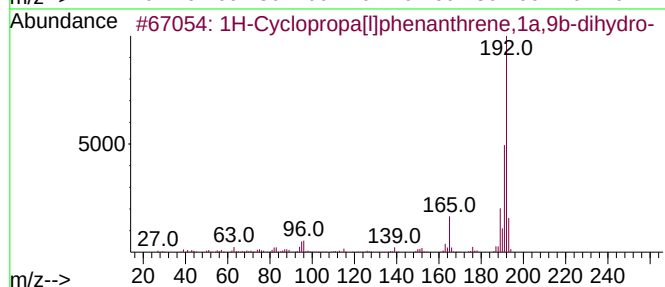
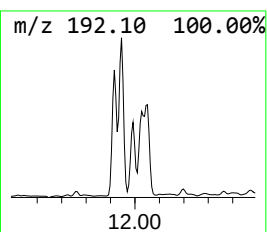
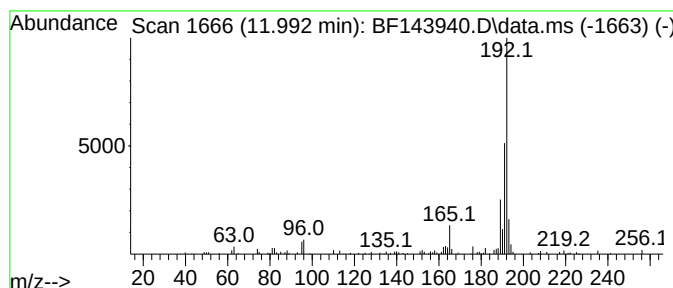
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenanthrene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.992	2.60 ng	79899	Phenanthrene-d10	11.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	96
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	94
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
5			Anthracene, 9-methyl-	192	C15H12	000779-02-2	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101325\
Data File : BF143940.D
Acq On : 13 Oct 2025 18:58
Operator : RC/JU
Sample : Q3323-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

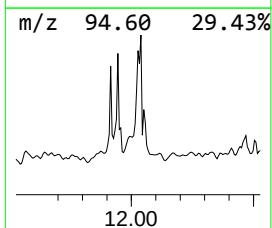
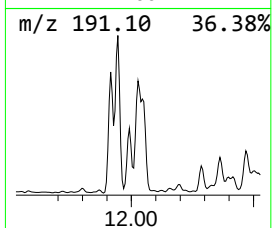
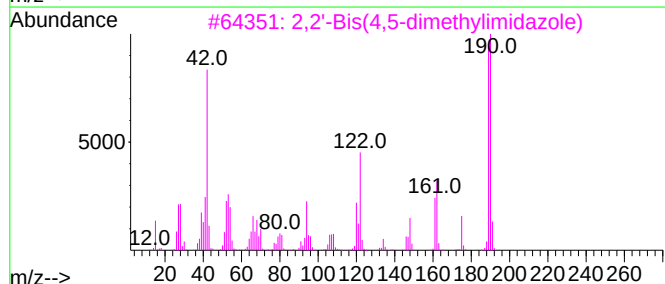
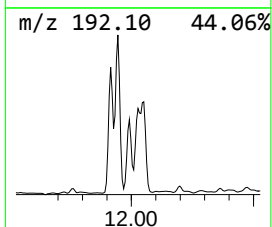
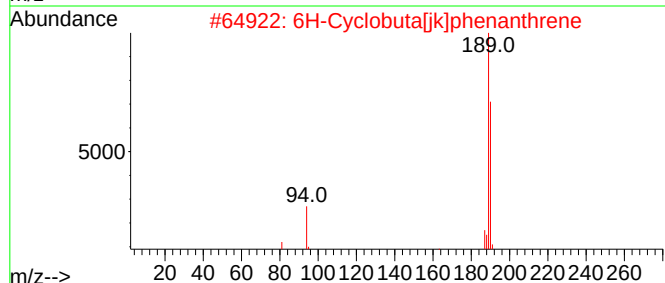
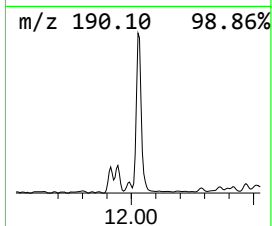
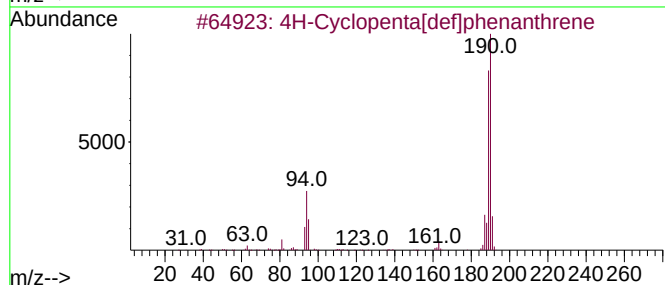
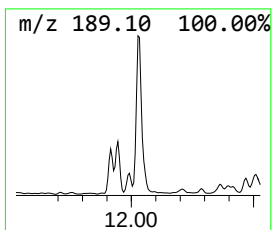
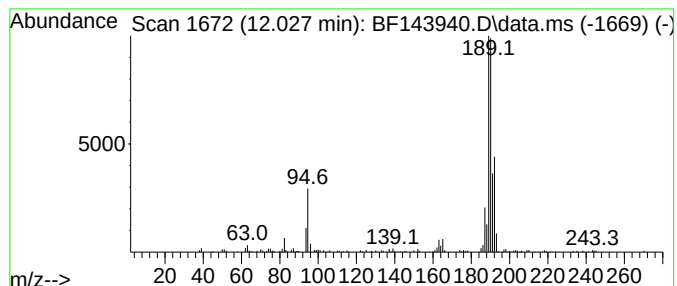
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.027	8.63 ng	265279	Phenanthrene-d10		11.416	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene		190	C15H10	000203-64-5	93
2	6H-Cyclobuta[jk]phenanthrene		190	C15H10	083469-43-6	45
3	2,2'-Bis(4,5-dimethylimidazole)		190	C10H14N4	069286-06-2	42
4	Carbonic acid, ethyl 2-formyl-4,...		262	C10H8Cl2O4	1000331-34-4	33
5	2-Amino-4,6-dichloropyrimidine-5...		191	C5H3Cl2N3O	005604-46-6	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101325\
Data File : BF143940.D
Acq On : 13 Oct 2025 18:58
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Sample : Q3323-05
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Instrument :
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ClientSampleId :
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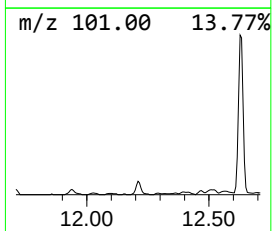
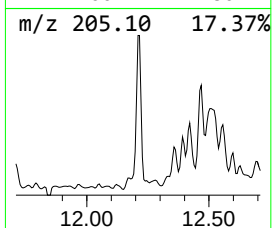
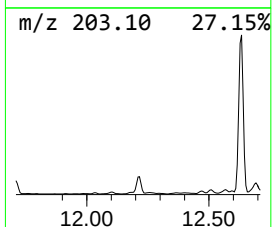
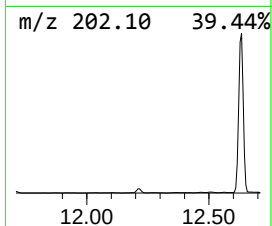
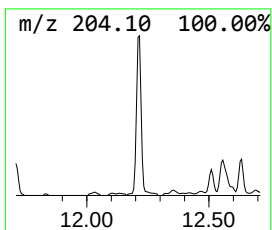
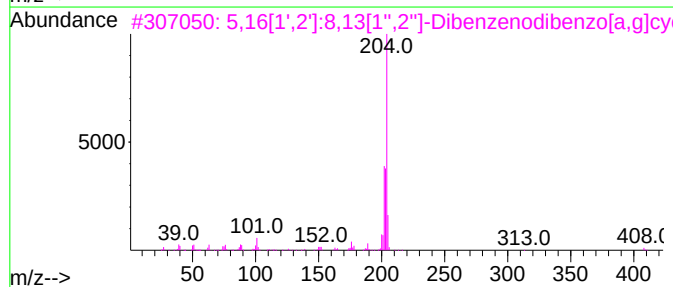
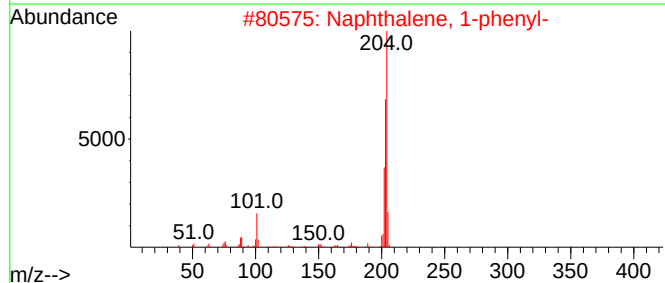
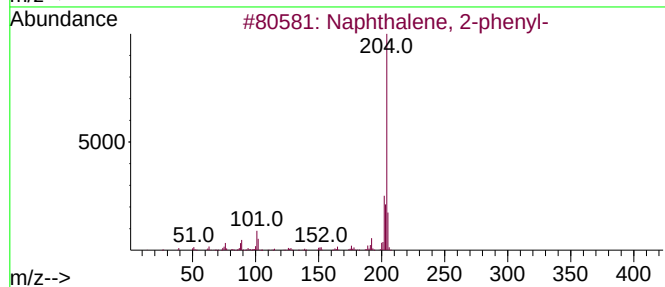
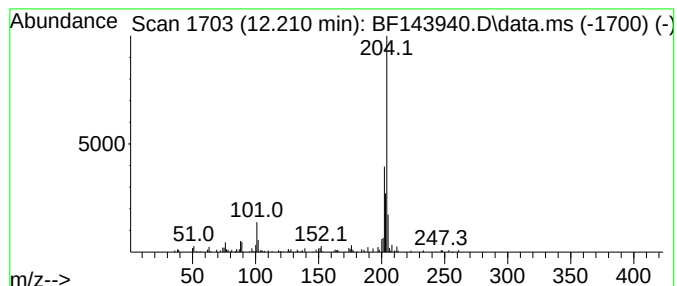
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Naphthalene, 2-phenyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.210	3.52 ng	108303	Phenanthrene-d10	11.416

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	95
2		Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	90
3		5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	90
4		6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	90
5		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101325\
Data File : BF143940.D
Acq On : 13 Oct 2025 18:58
Operator : RC/JU
Sample : Q3323-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

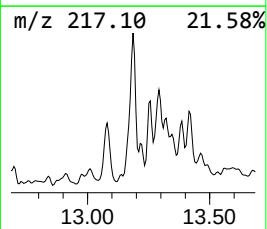
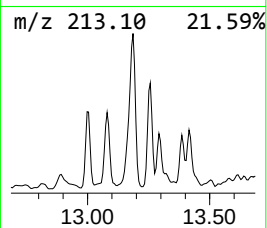
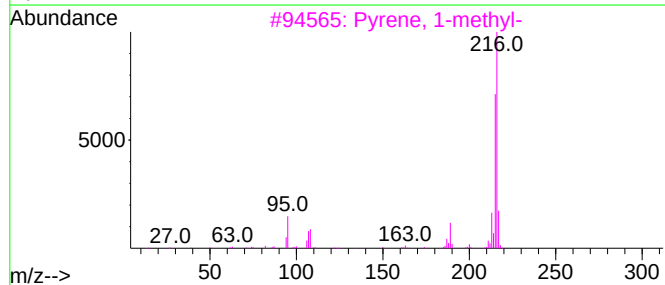
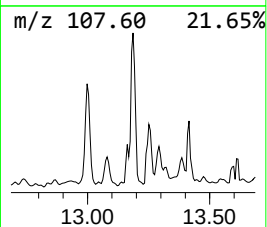
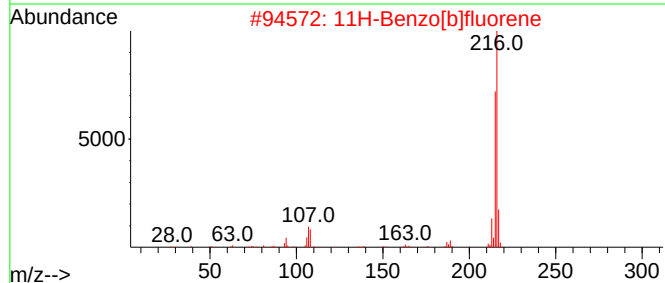
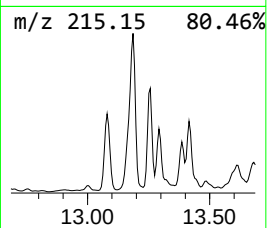
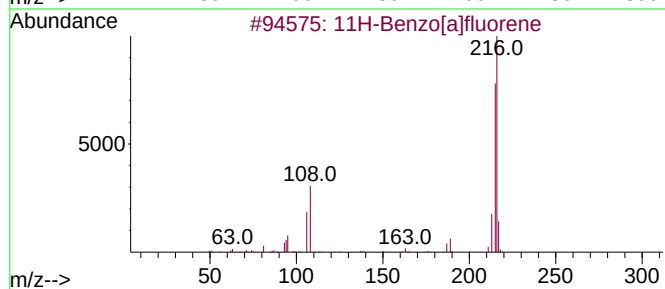
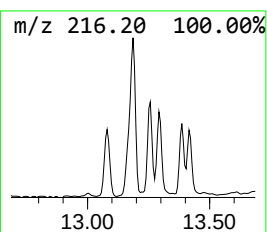
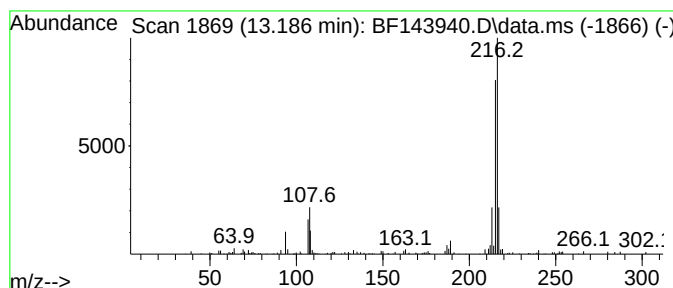
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 11H-Benzo[a]fluorene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.186	2.01 ng	125053	Chrysene-d12	14.063	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	74
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101325\
Data File : BF143940.D
Acq On : 13 Oct 2025 18:58
Operator : RC/JU
Sample : Q3323-05
Misc :
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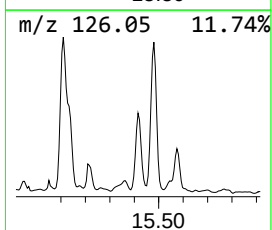
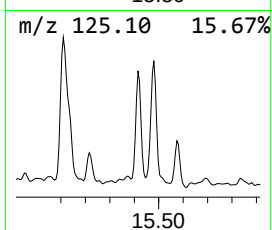
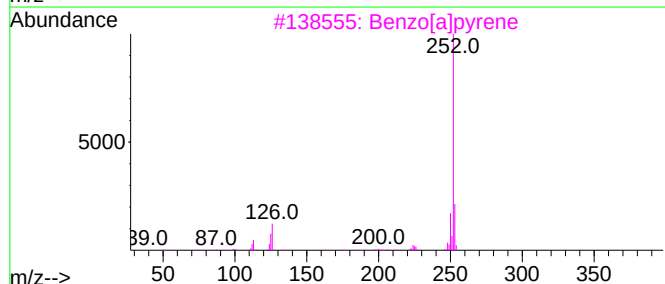
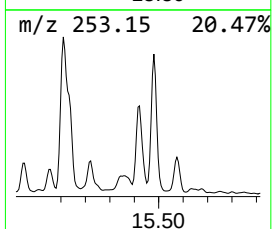
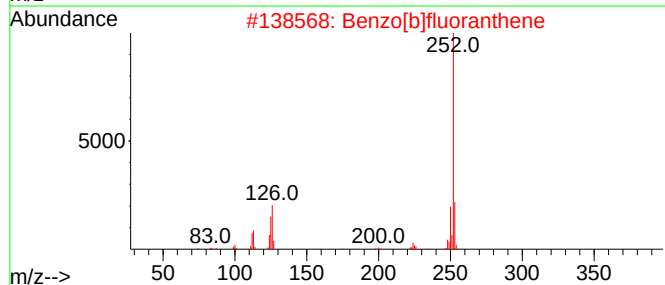
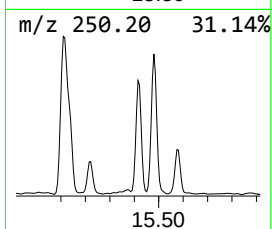
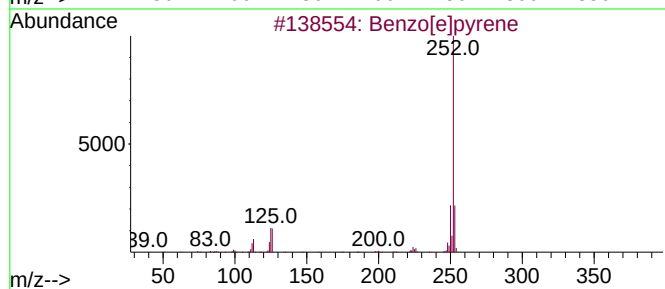
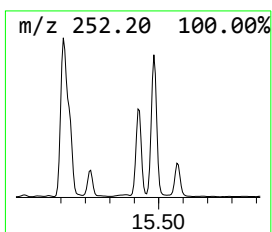
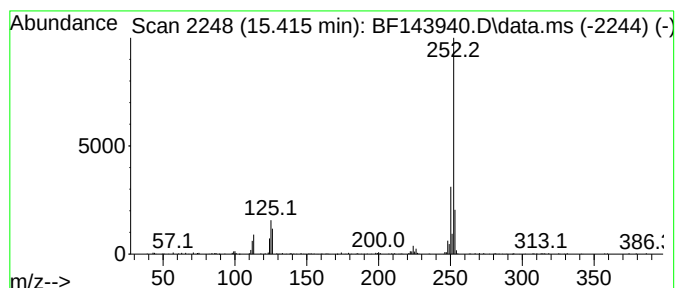
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100625.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.415	5.55 ng	258910	Perylene-d12	15.545	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2	Benzo[b]fluoranthene	252	C20H12	000205-99-2	95
3	Benzo[a]pyrene	252	C20H12	000050-32-8	93
4	Benzo[k]fluoranthene	252	C20H12	000207-08-9	91
5	Benzo[j]fluoranthene	252	C20H12	000205-82-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101325\
Data File : BF143940.D
Acq On : 13 Oct 2025 18:58
Operator : RC/JU
Sample : Q3323-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SED-05-0-2

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
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Dibenzothiophene	11.310	3.8	ng	117403	4	11.416	615019	20.0
1H-Cyclopropa[1...	11.939	14.4	ng	441987	4	11.416	615019	20.0
Phenanthrene, 2...	11.992	2.6	ng	79899	4	11.416	615019	20.0
4H-Cyclopenta[d...	12.027	8.6	ng	265279	4	11.416	615019	20.0
Naphthalene, 2-...	12.210	3.5	ng	108303	4	11.416	615019	20.0
11H-Benzo[a]flu...	13.186	2.0	ng	125053	5	14.063	1243210	20.0
Benzo[e]pyrene	15.415	5.5	ng	258910	6	15.545	933724	20.0