

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061024\
 Data File : BF138180.D
 Acq On : 10 Jun 2024 19:07
 Operator : CG\JU
 Sample : P2787-01 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-04-06072024

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-BF060524.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF138180.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.163	7	13	19	rVB	2318190	2786037	66.78%	4.173%
2	5.104	510	513	519	rVB	168891	181584	4.35%	0.272%
3	5.504	577	581	585	rBV	3525176	3244827	77.77%	4.860%
4	6.516	745	753	756	rBV	3556764	3399033	81.47%	5.091%
5	6.722	784	788	793	rBV3	141063	151312	3.63%	0.227%
6	6.898	814	818	821	rBV	2249194	1973553	47.30%	2.956%
7	7.451	909	912	916	rVB	2181103	1800047	43.14%	2.696%
8	8.175	1032	1035	1038	rBV	3102581	2523417	60.48%	3.780%
9	8.198	1038	1039	1042	rVB	241610	127834	3.06%	0.191%
10	8.886	1154	1156	1159	rVB	186730	143630	3.44%	0.215%
11	8.986	1170	1173	1177	rVB	214207	173362	4.16%	0.260%
12	9.251	1214	1218	1221	rBV	5431936	3997755	95.82%	5.988%
13	9.334	1228	1232	1234	rBV	75721	74088	1.78%	0.111%
14	9.445	1247	1251	1256	rVB	44386	60085	1.44%	0.090%
15	9.516	1256	1263	1266	rBV	88080	125482	3.01%	0.188%
16	9.586	1273	1275	1278	rVV	170256	135284	3.24%	0.203%
17	9.616	1278	1280	1282	rVB	107273	79483	1.91%	0.119%
18	9.651	1282	1286	1288	rBV3	69062	65203	1.56%	0.098%
19	9.704	1293	1295	1300	rVB	94770	97654	2.34%	0.146%
20	9.792	1306	1310	1314	rVB	197362	173994	4.17%	0.261%
21	9.863	1319	1322	1326	rVB	85512	74212	1.78%	0.111%
22	9.933	1330	1334	1336	rVV	2827323	2528635	60.61%	3.787%
23	9.963	1336	1339	1342	rVB	555899	416719	9.99%	0.624%
24	10.033	1348	1351	1356	rBV2	46235	56917	1.36%	0.085%
25	10.086	1356	1360	1364	rVV4	34267	53172	1.27%	0.080%
26	10.133	1364	1368	1372	rVV	359163	298829	7.16%	0.448%
27	10.169	1372	1374	1381	rVB3	64481	66523	1.59%	0.100%
28	10.245	1384	1387	1390	rBV2	65781	67476	1.62%	0.101%
29	10.339	1399	1403	1404	rBV2	47629	58428	1.40%	0.088%
30	10.357	1404	1406	1409	rVB2	87081	77505	1.86%	0.116%
31	10.475	1423	1426	1432	rVB	699679	585209	14.03%	0.877%
32	10.563	1439	1441	1443	rBV2	82346	59255	1.42%	0.089%
33	10.633	1451	1453	1457	rVB	123087	100012	2.40%	0.150%
34	10.716	1463	1467	1471	rBV	2108381	1697726	40.69%	2.543%
35	10.839	1484	1488	1495	rVB2	111898	167848	4.02%	0.251%
36	11.022	1514	1519	1522	rBV2	121740	159592	3.83%	0.239%

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	11.051	1522	1524	1527	rVV2	86684	76228	1.83%	0.114%
38	11.110	1531	1534	1537	rVB	72361	65055	1.56%	0.097%
39	11.151	1537	1541	1543	rBV3	44217	59657	1.43%	0.089%
40	11.263	1557	1560	1562	rBV	49669	58325	1.40%	0.087%
41	11.316	1566	1569	1573	rVB	258356	247594	5.93%	0.371%
42	11.416	1582	1586	1588	rBV	2705648	2258317	54.13%	3.382%
43	11.439	1588	1590	1594	rVV	3927316	3322918	79.64%	4.977%
44	11.492	1596	1599	1602	rVV	1196661	913681	21.90%	1.368%
45	11.522	1602	1604	1607	rVV2	76475	74652	1.79%	0.112%
46	11.645	1618	1625	1628	rBV	496232	445171	10.67%	0.667%
47	11.710	1634	1636	1640	rBV3	108841	103228	2.47%	0.155%
48	11.745	1640	1642	1648	rVB2	98610	91969	2.20%	0.138%
49	11.827	1654	1656	1660	rVB	60712	58118	1.39%	0.087%
50	11.916	1668	1671	1673	rBV	425201	362006	8.68%	0.542%
51	11.945	1673	1676	1680	rVV	699958	597602	14.32%	0.895%
52	11.992	1681	1684	1687	rVV	226150	189262	4.54%	0.283%
53	12.027	1687	1690	1693	rVV	813249	884601	21.20%	1.325%
54	12.051	1693	1694	1698	rVB	324703	201080	4.82%	0.301%
55	12.116	1703	1705	1708	rBV	83071	65587	1.57%	0.098%
56	12.216	1719	1722	1724	rBV	261490	247669	5.94%	0.371%
57	12.233	1724	1725	1729	rVB2	115275	81500	1.95%	0.122%
58	12.363	1745	1747	1749	rVB	91594	64199	1.54%	0.096%
59	12.392	1749	1752	1754	rBV	95533	79099	1.90%	0.118%
60	12.416	1754	1756	1758	rVB2	96817	70653	1.69%	0.106%
61	12.469	1761	1765	1767	rBV	252291	271665	6.51%	0.407%
62	12.504	1767	1771	1773	rVV2	211239	248455	5.95%	0.372%
63	12.551	1776	1779	1784	rBV2	115241	146849	3.52%	0.220%
64	12.633	1789	1793	1796	rBV	4237408	3856662	92.44%	5.776%
65	12.722	1806	1808	1811	rBV	129625	105228	2.52%	0.158%
66	12.780	1815	1818	1822	rVV2	140154	139932	3.35%	0.210%
67	12.857	1825	1831	1837	rBV	3655749	4017822	96.30%	6.018%
68	12.904	1837	1839	1844	rVB2	155848	199696	4.79%	0.299%
69	12.974	1848	1851	1852	rBV2	208223	172066	4.12%	0.258%
70	13.004	1852	1856	1863	rVB	3845964	3598540	86.25%	5.390%
71	13.074	1863	1868	1871	rBV3	227380	386698	9.27%	0.579%
72	13.104	1871	1873	1876	rVB	119647	102622	2.46%	0.154%
73	13.163	1880	1883	1884	rBV2	176094	180597	4.33%	0.270%
74	13.186	1884	1887	1890	rVB	675133	583168	13.98%	0.873%
75	13.251	1893	1898	1901	rBV	633976	554131	13.28%	0.830%
76	13.286	1901	1904	1907	rBV2	354633	347308	8.32%	0.520%

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77	13.345	1912	1914	1917	rBV2	68413	76341	1.83%	0.114%
78	13.380	1917	1920	1923	rVB	189126	145311	3.48%	0.218%
79	13.410	1923	1925	1927	rBV	211923	178952	4.29%	0.268%
80	13.804	1990	1992	1995	rVB	204672	167001	4.00%	0.250%
81	13.833	1995	1997	1999	rBV	295776	242280	5.81%	0.363%
82	13.851	1999	2000	2005	rVB2	231616	258811	6.20%	0.388%
83	14.051	2027	2034	2037	rBV2	2887977	4172270	100.00%	6.249%
84	14.080	2037	2039	2042	rVB	1516700	1142832	27.39%	1.712%
85	14.157	2050	2052	2055	rVB	291591	256007	6.14%	0.383%
86	14.192	2056	2058	2061	rBV	139730	147386	3.53%	0.221%
87	14.439	2098	2100	2103	rVB	230360	189629	4.54%	0.284%
88	14.533	2113	2116	2118	rBV2	203551	188150	4.51%	0.282%
89	14.563	2119	2121	2123	rBV	183916	199533	4.78%	0.299%
90	15.098	2208	2212	2215	rBV	953109	1379669	33.07%	2.066%
91	15.404	2261	2264	2271	rVB	549799	706052	16.92%	1.058%
92	15.468	2271	2275	2282	rVV	840499	1095660	26.26%	1.641%
93	15.533	2282	2286	2289	rVV	1170197	1403255	33.63%	2.102%
94	15.563	2289	2291	2298	rVB2	288652	437277	10.48%	0.655%
95	17.015	2534	2538	2548	rVB2	316151	629823	15.10%	0.943%
96	17.468	2611	2615	2625	rVB	240986	437930	10.50%	0.656%

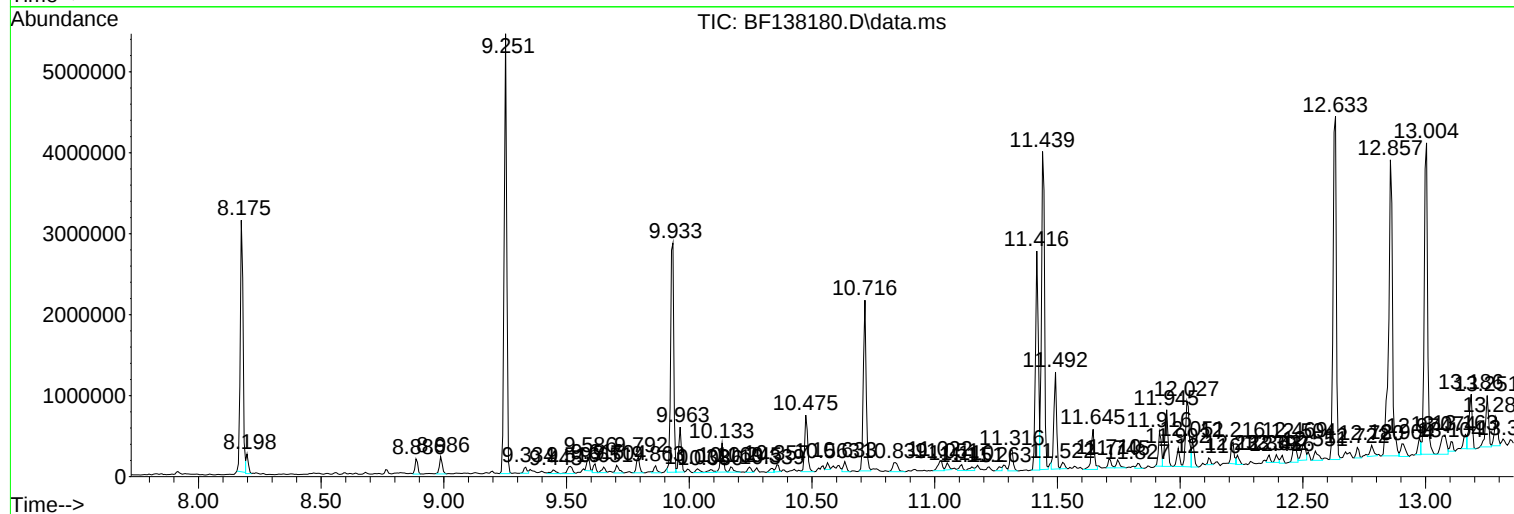
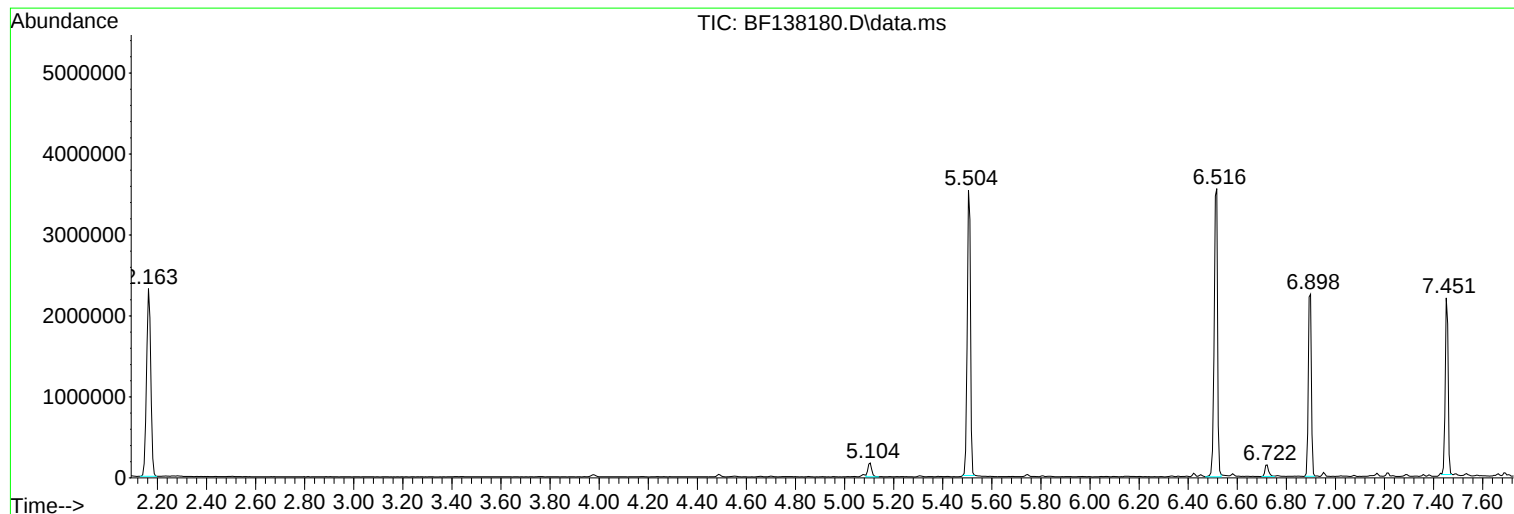
Sum of corrected areas: 66765501

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060524.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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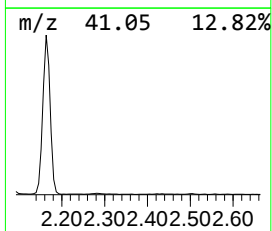
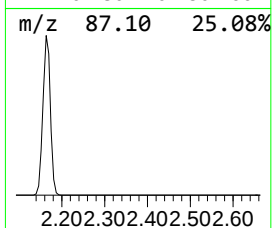
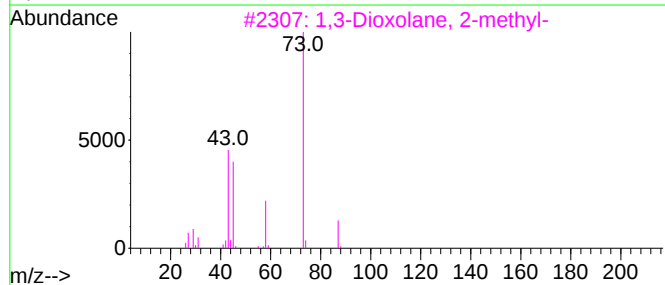
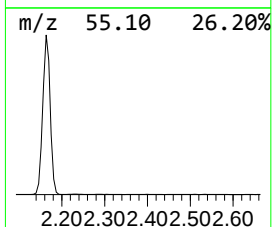
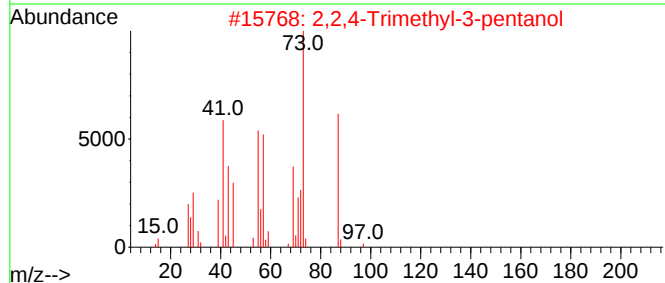
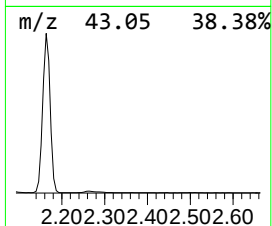
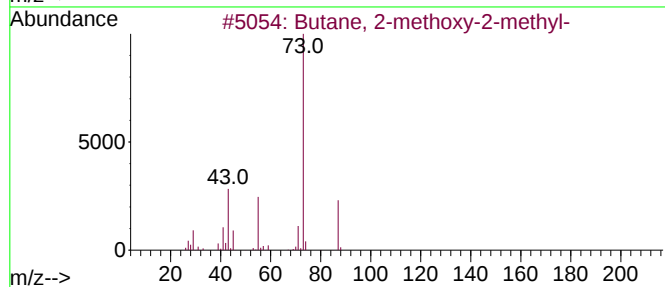
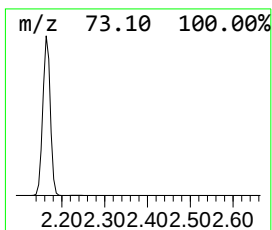
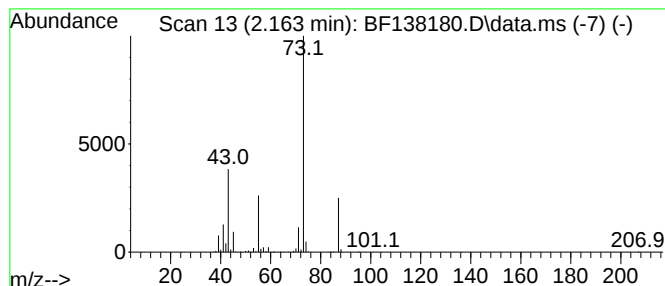
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060524.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.163	28.23 ng	2786040	1,4-Dichlorobenzene-d4	6.898

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	10



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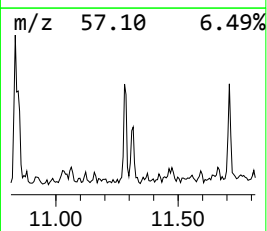
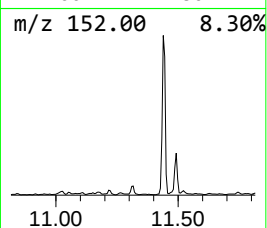
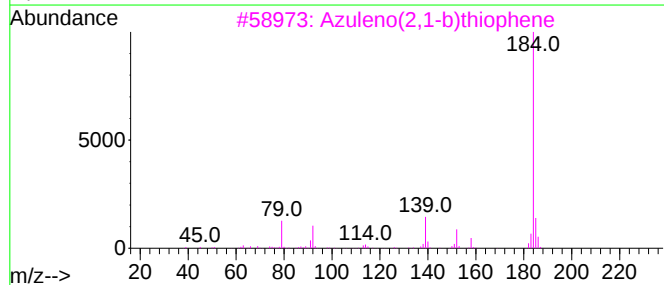
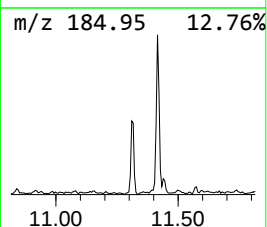
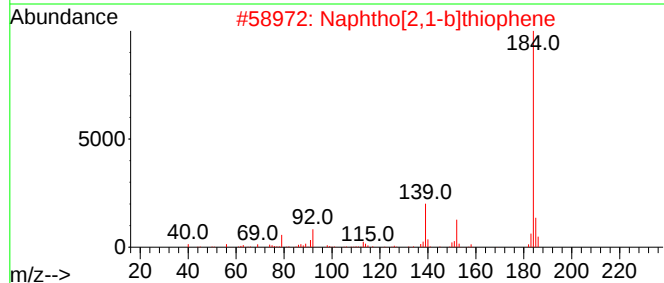
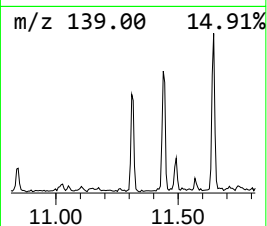
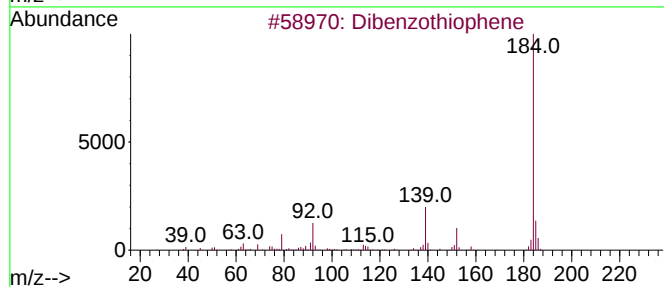
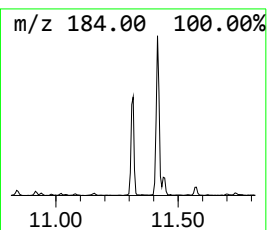
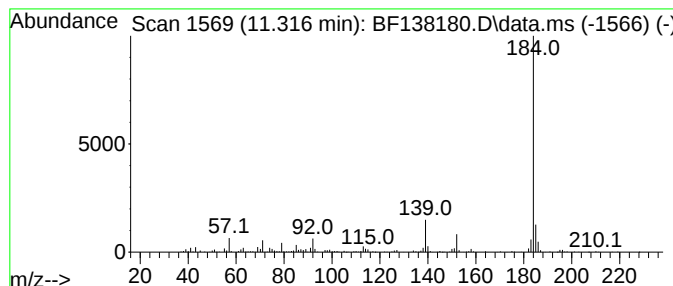
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Dibenzothiophene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.316	2.19 ng	247594	Phenanthrene-d10	11.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzothiophene	184	C12H8S	000132-65-0	96
2			Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	93
3			Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	91
4			Dibenzothiophene, 5-oxide	200	C12H8OS	001013-23-6	78
5			1,1'-Biphenyl, 2-methoxy-	184	C13H12O	000086-26-0	72



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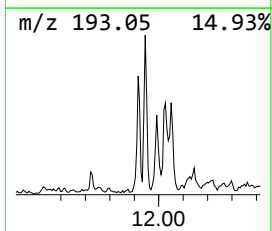
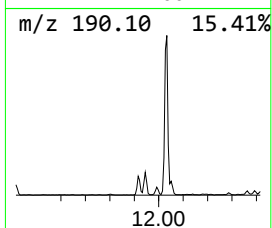
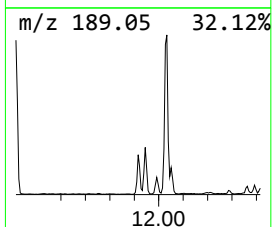
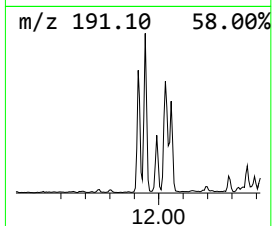
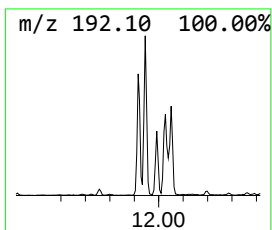
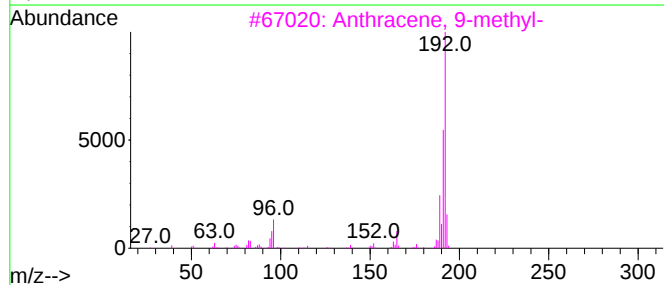
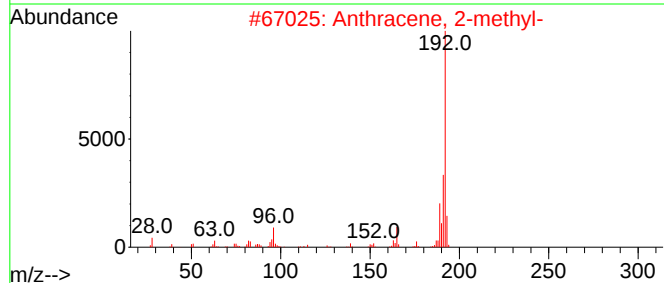
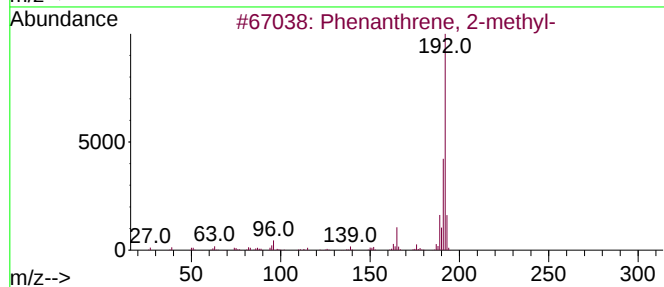
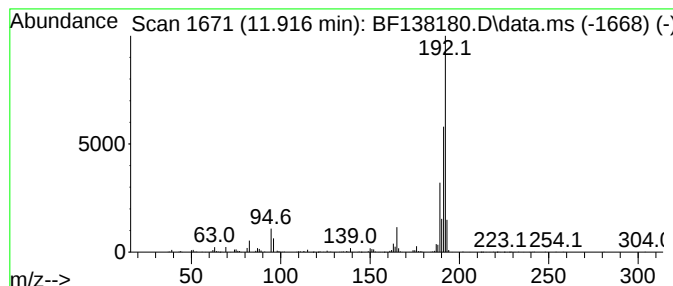
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Phenanthrene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.916	3.21 ng	362006	Phenanthrene-d10	11.416

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		Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
5		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	90



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Acq On : 10 Jun 2024 19:07
Operator : CG\JU
Sample : P2787-01 2X
Misc :
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ClientSampleId :
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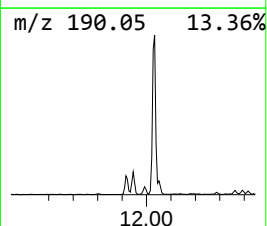
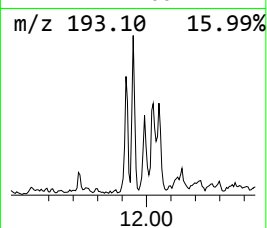
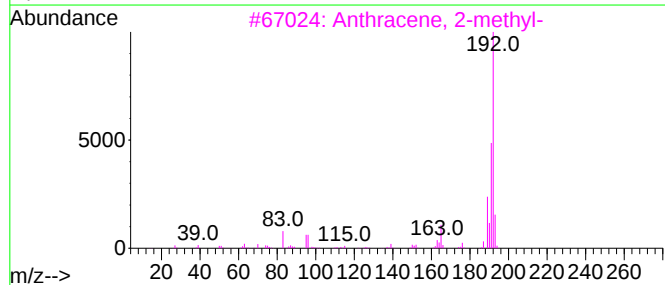
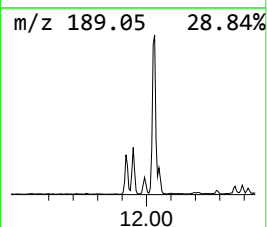
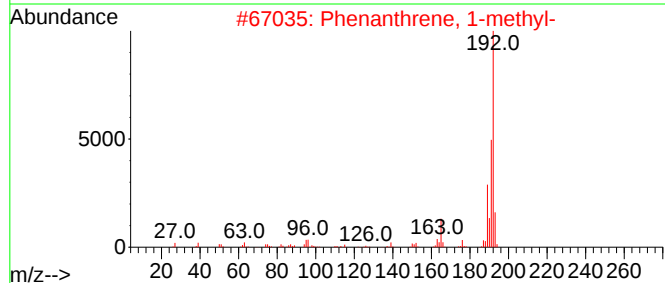
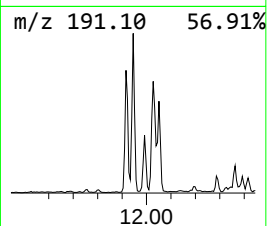
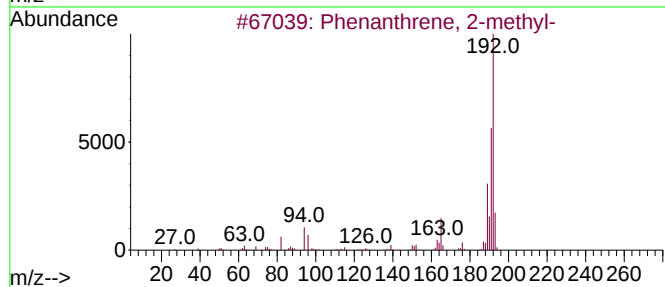
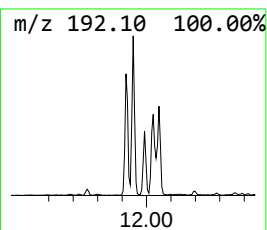
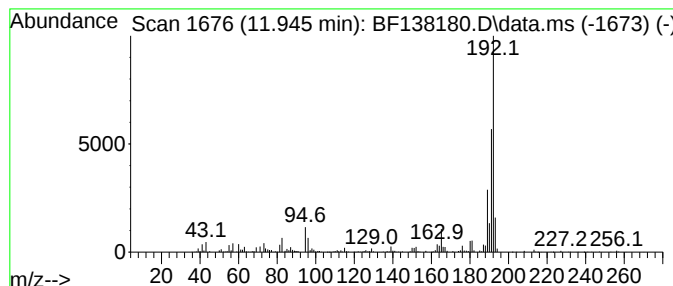
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Phenanthrene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.945	5.29 ng	597602	Phenanthrene-d10	11.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4			Anthracene, 9-methyl-	192	C15H12	000779-02-2	94
5			Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061024\
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Sample : P2787-01 2X
Misc :
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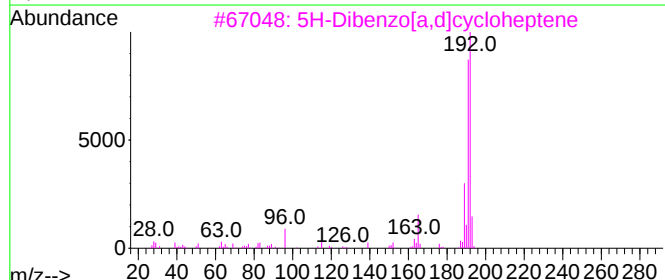
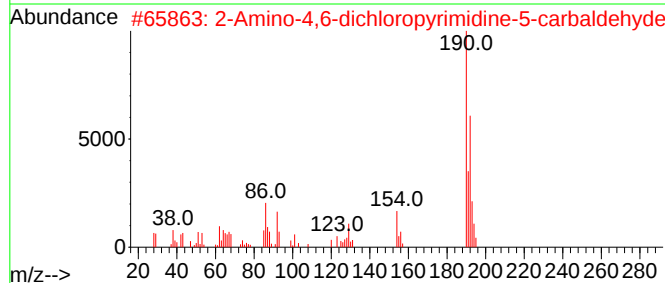
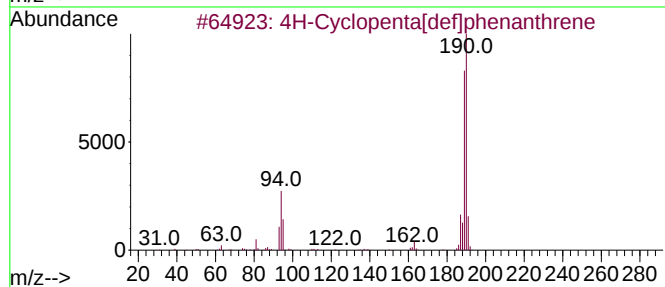
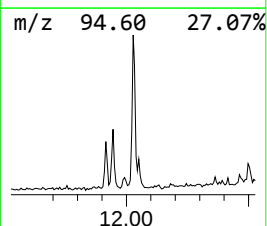
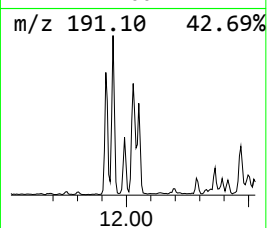
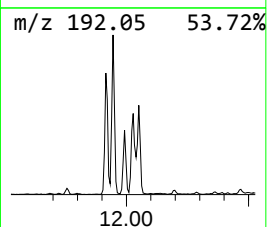
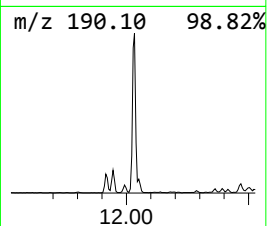
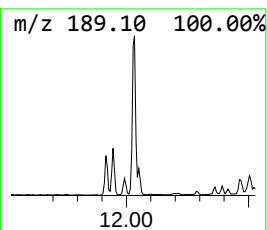
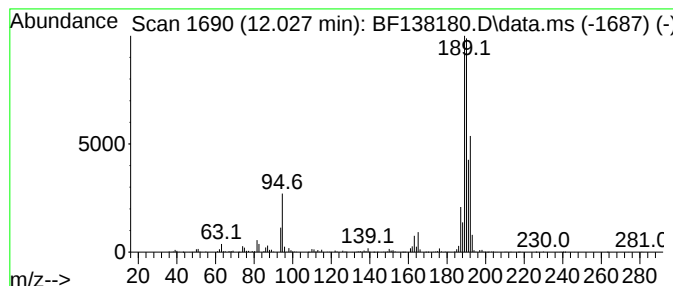
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.028	7.83 ng	884601	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	2-Amino-4,6-dichloropyrimidine-5...		191	C5H3Cl2N3O	005604-46-6	38
3	5H-Dibenzo[a,d]cycloheptene		192	C15H12	000256-81-5	18
4	1H-Indene, 2-phenyl-		192	C15H12	004505-48-0	18
5	Benzene, 1,1'-(1,2-propadienyld...		192	C15H12	014251-57-1	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061024\
Data File : BF138180.D
Acq On : 10 Jun 2024 19:07
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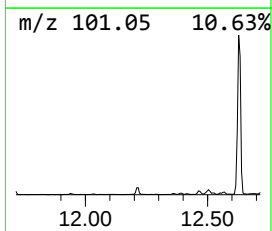
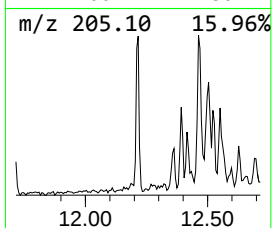
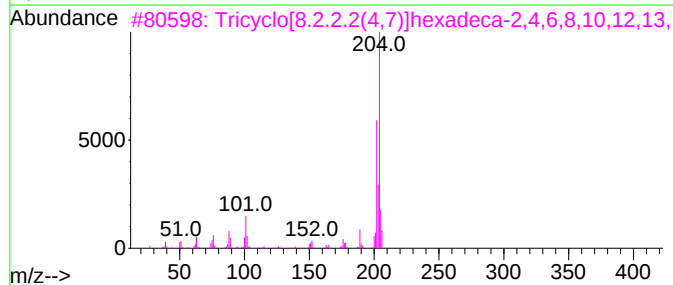
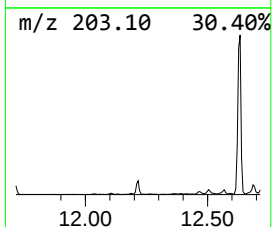
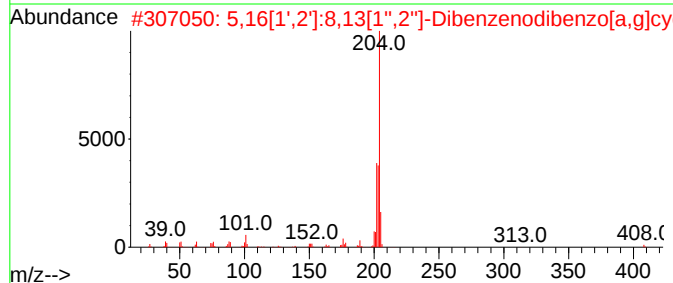
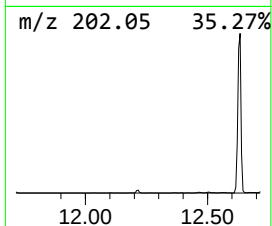
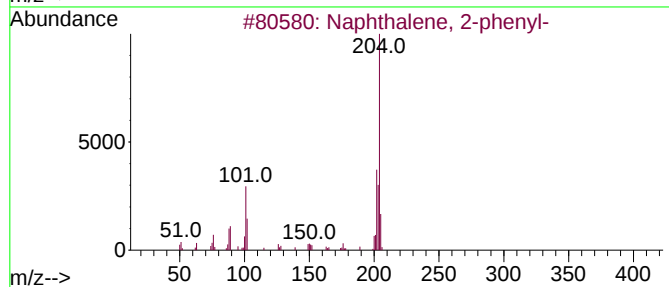
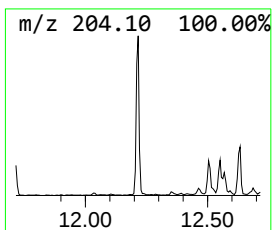
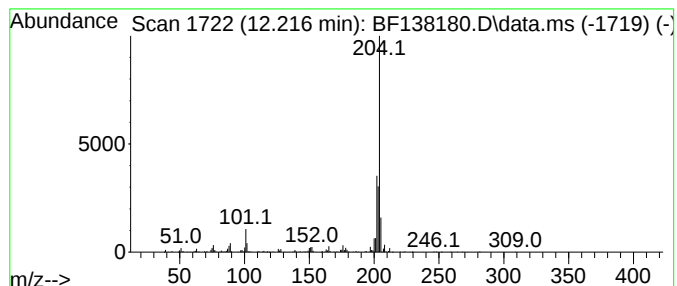
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Naphthalene, 2-phenyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.216	2.19 ng	247669	Phenanthrene-d10	11.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	90
2			5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	87
3			Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	83
4			Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	64
5			9-Acridinecarbonitrile	204	C14H8N2	005326-19-2	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061024\
Data File : BF138180.D
Acq On : 10 Jun 2024 19:07
Operator : CG\JU
Sample : P2787-01 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

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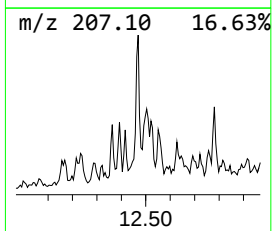
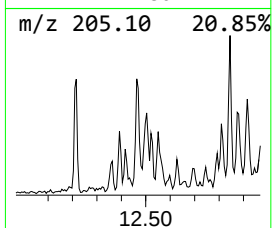
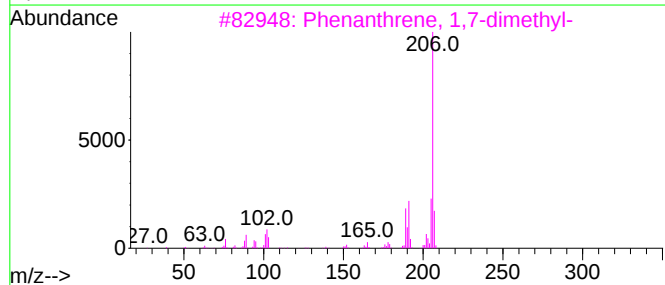
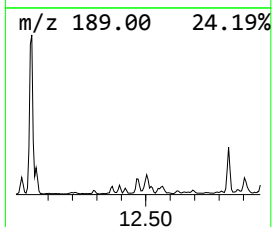
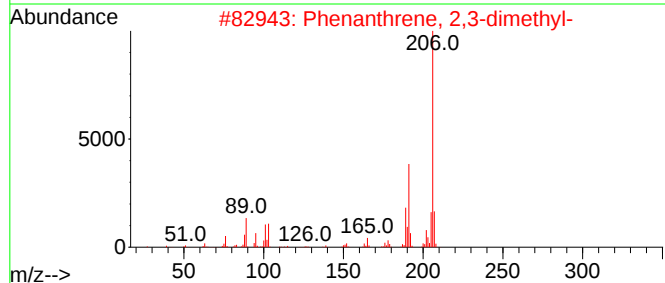
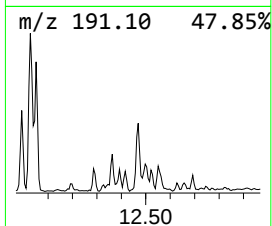
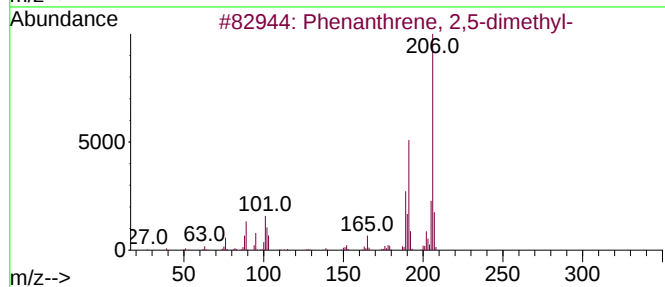
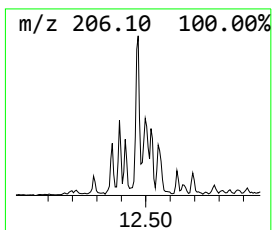
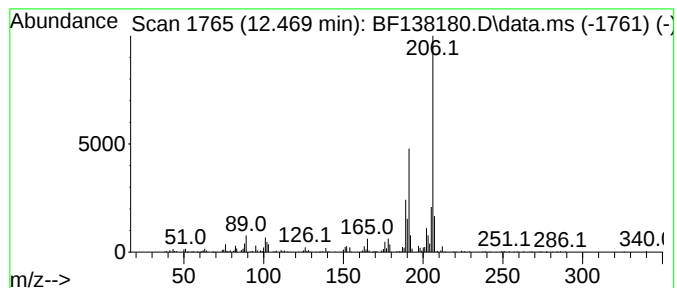
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Phenanthrene, 2,5-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.469	2.41 ng	271665	Phenanthrene-d10	11.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	90
3			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	89
4			9,10-Dimethylanthracene	206	C16H14	000781-43-1	89
5			3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	86



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Data File : BF138180.D
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ClientSampleId :
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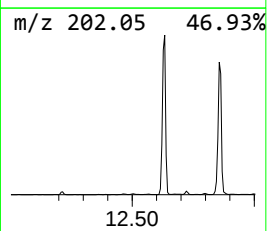
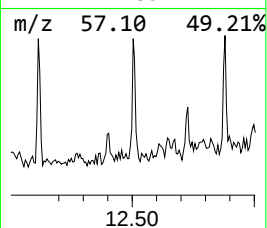
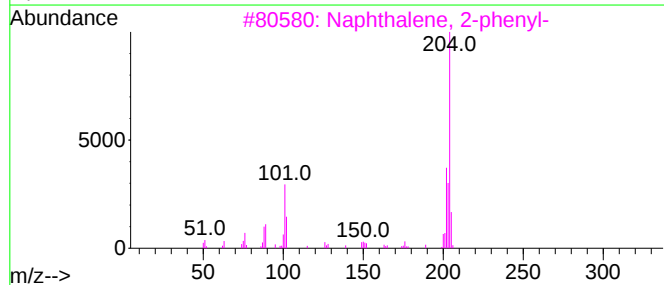
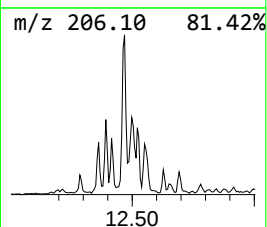
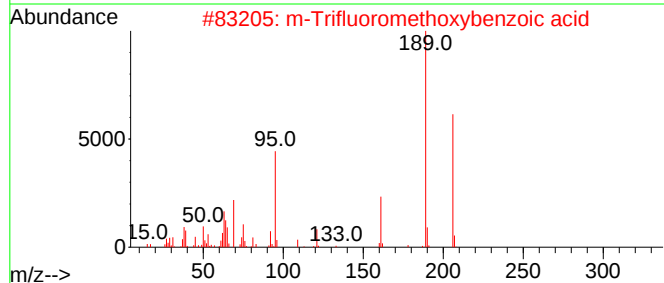
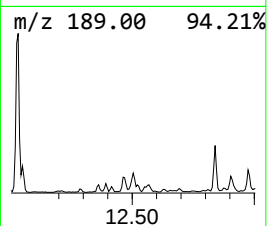
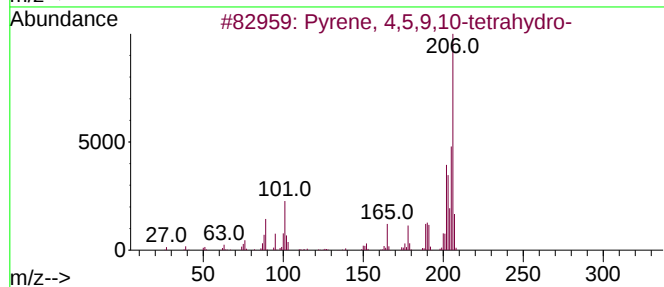
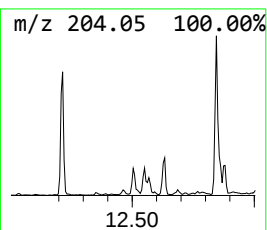
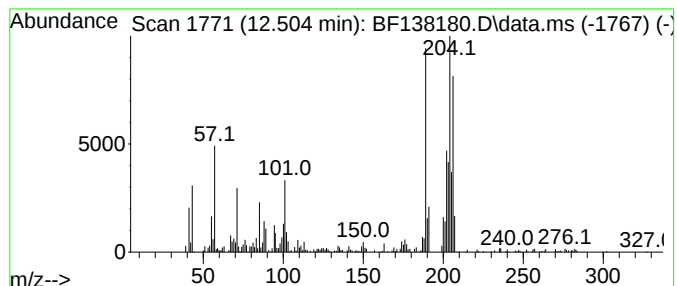
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 unknown12.504 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.504	2.20 ng	248455	Phenanthrene-d10	11.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 4,5,9,10-tetrahydro-	206	C16H14	000781-17-9	38
2			m-Trifluoromethoxybenzoic acid	206	C8H5F3O3	001014-81-9	30
3			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	30
4			Benzene, 1,3,5-tris(1-methylethyl)-	204	C15H24	000717-74-8	30
5			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061024\
Data File : BF138180.D
Acq On : 10 Jun 2024 19:07
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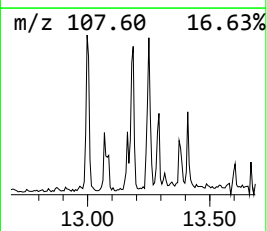
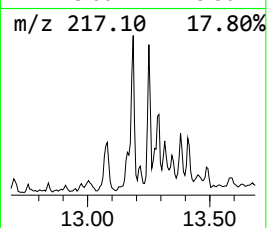
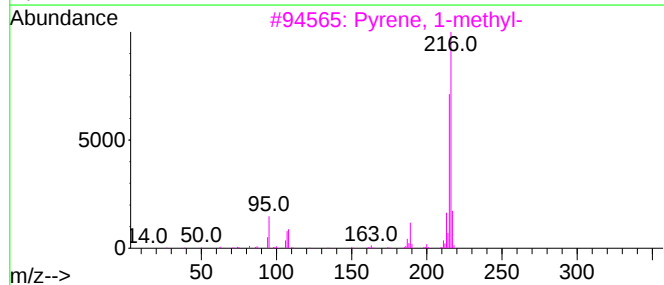
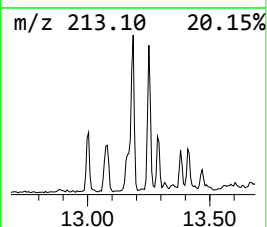
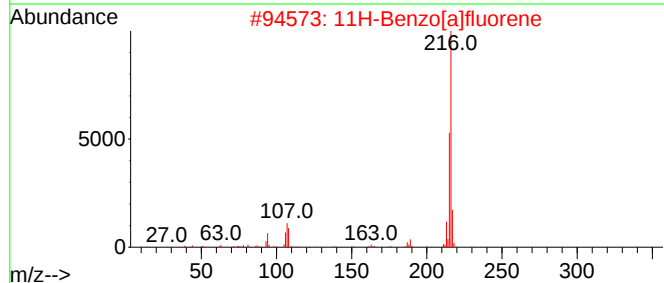
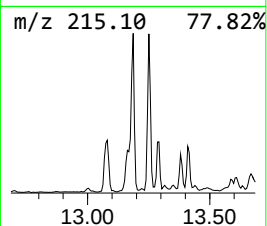
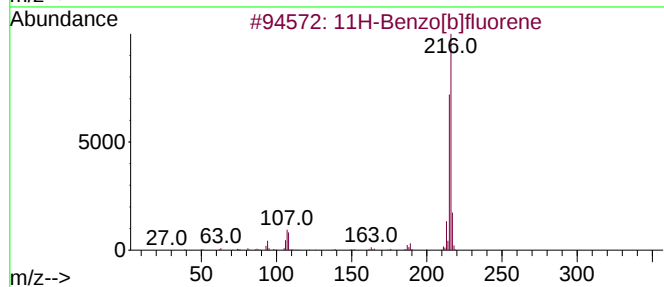
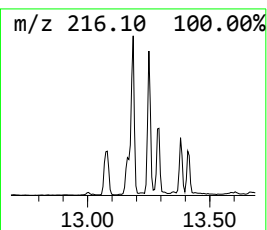
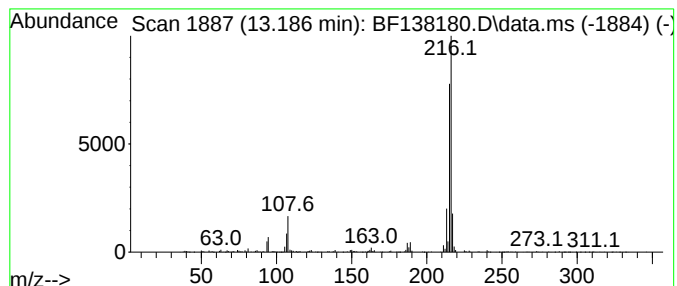
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 11H-Benzo[b]fluorene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.186	2.80 ng	583168	Chrysene-d12	14.057

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
2			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
3			Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	81
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061024\
Data File : BF138180.D
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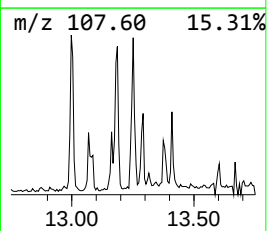
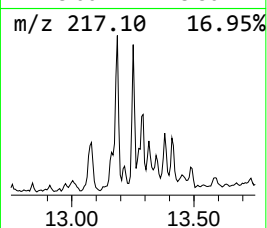
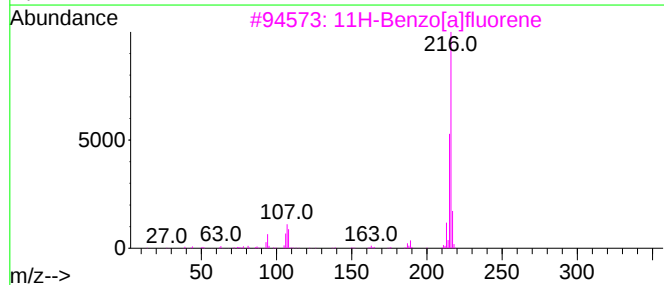
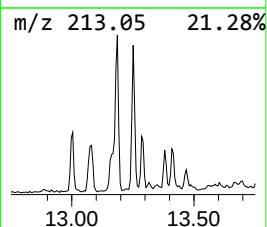
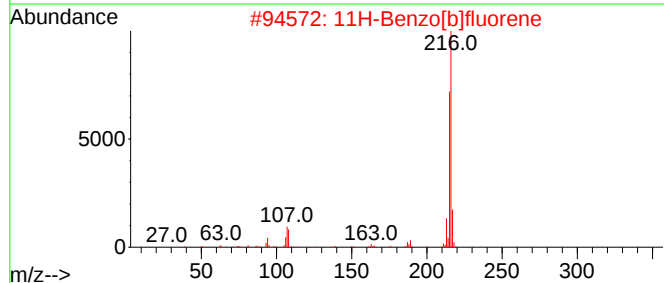
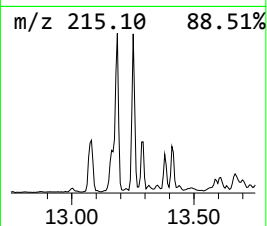
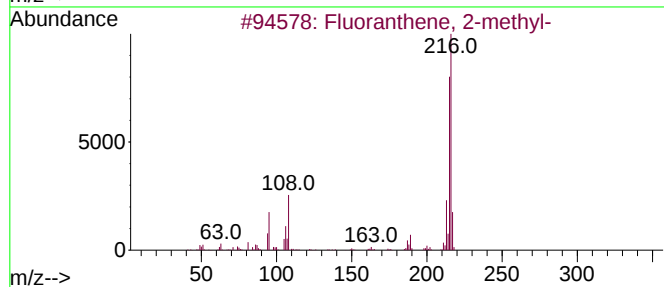
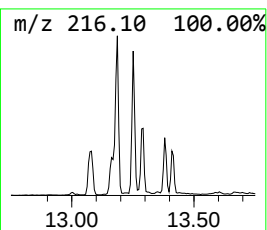
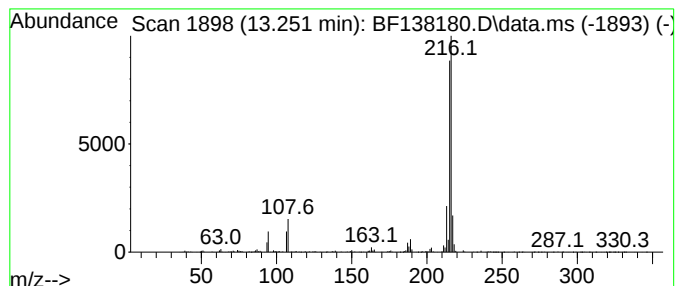
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Fluoranthene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.251	2.66 ng	554131	Chrysene-d12	14.057

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	95
2		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
3		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
4		Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
5		7H-Benzo[c]fluorene	216	C17H12	000205-12-9	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061024\
Data File : BF138180.D
Acq On : 10 Jun 2024 19:07
Operator : CG\JU
Sample : P2787-01 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-06072024

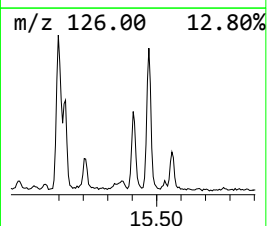
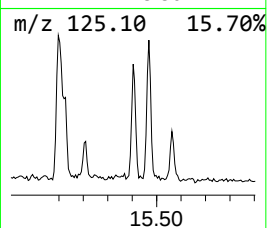
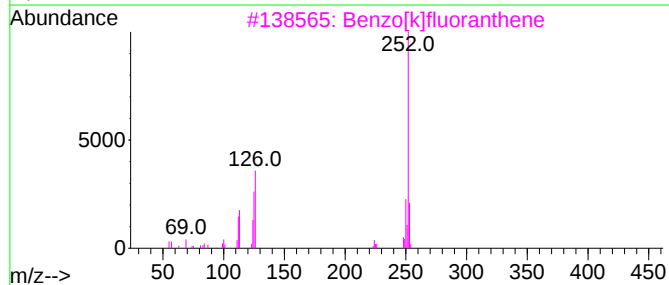
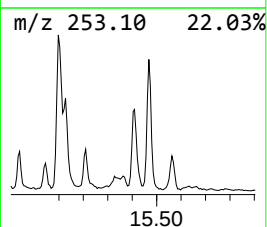
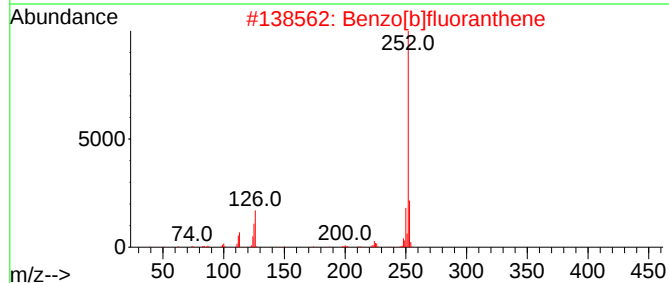
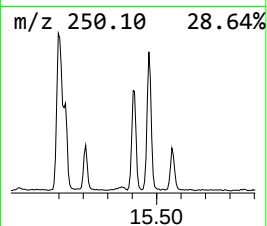
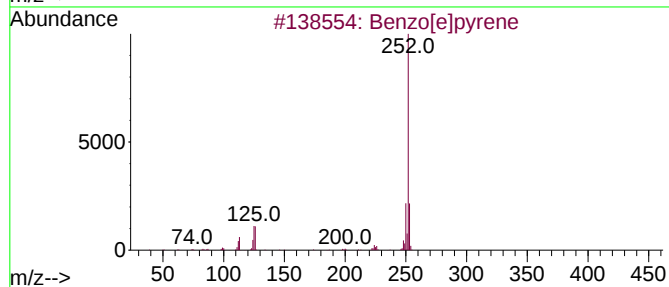
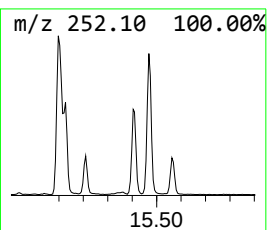
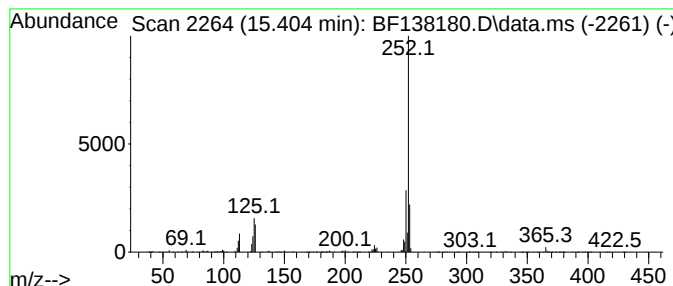
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.404	10.06 ng	706052	Perylene-d12	15.533

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	98
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	94
4			Benzo[j]fluoranthene	252	C20H12	000205-82-3	94
5			Perylene	252	C20H12	000198-55-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061024\
Data File : BF138180.D
Acq On : 10 Jun 2024 19:07
Operator : CG\JU
Sample : P2787-01 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-06072024

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060524.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
Butane, 2-metho...	2.163	28.2	ng	2786040	1	6.898	1973550	20.0
Dibenzothiophene	11.316	2.2	ng	247594	4	11.416	2258320	20.0
Phenanthrene, 2...	11.916	3.2	ng	362006	4	11.416	2258320	20.0
Phenanthrene, 1...	11.945	5.3	ng	597602	4	11.416	2258320	20.0
4H-Cyclopenta[d...	12.028	7.8	ng	884601	4	11.416	2258320	20.0
Naphthalene, 2-...	12.216	2.2	ng	247669	4	11.416	2258320	20.0
Phenanthrene, 2...	12.469	2.4	ng	271665	4	11.416	2258320	20.0
unknown12.504	12.504	2.2	ng	248455	4	11.416	2258320	20.0
11H-Benzo[b]flu...	13.186	2.8	ng	583168	5	14.057	4172270	20.0
Fluoranthene, 2...	13.251	2.7	ng	554131	5	14.057	4172270	20.0
Benzo[e]pyrene	15.404	10.1	ng	706052	6	15.533	1403260	20.0