

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081022\
 Data File : BF129707.D
 Acq On : 10 Aug 2022 21:31
 Operator : CG\JU
 Sample : N4119-01 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP1-2-4FT

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-BF072522.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF129707.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.475	539	542	555	rBV	625478	609480	19.98%	1.765%
2	6.492	711	715	729	rBV	611285	633293	20.76%	1.834%
3	6.663	740	744	748	rBV2	24628	33081	1.08%	0.096%
4	6.834	769	773	777	rBV	1337707	1110844	36.42%	3.217%
5	7.398	865	869	875	rBV	458916	366441	12.01%	1.061%
6	8.116	987	991	994	rBV	1533905	1390857	45.60%	4.027%
7	8.139	994	995	998	rVB	123543	67607	2.22%	0.196%
8	8.828	1107	1112	1116	rBV2	56985	58829	1.93%	0.170%
9	8.928	1126	1129	1133	rBV	40973	35499	1.16%	0.103%
10	9.192	1170	1174	1177	rVB	743786	642457	21.06%	1.860%
11	9.533	1228	1232	1234	rBV	37419	38256	1.25%	0.111%
12	9.733	1263	1266	1271	rVB	203364	176016	5.77%	0.510%
13	9.875	1286	1290	1293	rVV	1623305	1239787	40.65%	3.590%
14	9.904	1293	1295	1299	rVB	77851	63202	2.07%	0.183%
15	10.075	1321	1324	1327	rBV2	55384	54752	1.80%	0.159%
16	10.186	1339	1343	1346	rBV2	34013	40814	1.34%	0.118%
17	10.292	1354	1361	1364	rBV4	25998	50684	1.66%	0.147%
18	10.422	1380	1383	1389	rVB2	78536	87361	2.86%	0.253%
19	10.580	1405	1410	1413	rVB	36211	35540	1.17%	0.103%
20	10.669	1417	1425	1430	rBV2	325240	335586	11.00%	0.972%
21	10.774	1438	1443	1450	rVB5	37820	83943	2.75%	0.243%
22	10.969	1472	1476	1479	rBV4	35285	43417	1.42%	0.126%
23	11.098	1493	1498	1500	rBV2	45796	54129	1.77%	0.157%
24	11.122	1500	1502	1508	rVV3	58310	70322	2.31%	0.204%
25	11.174	1508	1511	1514	rVV	61853	72210	2.37%	0.209%
26	11.210	1514	1517	1520	rVV2	67029	75285	2.47%	0.218%
27	11.263	1523	1526	1530	rVB	56747	60677	1.99%	0.176%
28	11.363	1540	1543	1545	rBV	1150921	1006507	33.00%	2.914%
29	11.386	1545	1547	1551	rVV	984197	837579	27.46%	2.425%
30	11.439	1551	1556	1559	rVV	430351	357409	11.72%	1.035%
31	11.474	1559	1562	1567	rVB2	59111	71803	2.35%	0.208%
32	11.598	1578	1583	1587	rBV2	109467	157488	5.16%	0.456%
33	11.657	1591	1593	1597	rVV2	48025	52457	1.72%	0.152%
34	11.698	1597	1600	1604	rVV3	38612	53127	1.74%	0.154%
35	11.745	1604	1608	1611	rVV5	27462	46484	1.52%	0.135%
36	11.869	1625	1629	1631	rBV	168103	166301	5.45%	0.482%

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Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	11.898	1631	1634	1637	rVB	263800	235194	7.71%	0.681%
38	11.939	1637	1641	1644	rBV	149881	152795	5.01%	0.442%
39	11.980	1644	1648	1650	rVV2	293334	326565	10.71%	0.946%
40	11.998	1650	1651	1656	rVB	144567	120948	3.97%	0.350%
41	12.045	1656	1659	1662	rBV2	60812	54387	1.78%	0.157%
42	12.163	1675	1679	1682	rVV	146807	128471	4.21%	0.372%
43	12.192	1682	1684	1687	rVB	65071	51258	1.68%	0.148%
44	12.310	1699	1704	1707	rBV2	61400	82434	2.70%	0.239%
45	12.339	1707	1709	1711	rVV	60465	49156	1.61%	0.142%
46	12.369	1711	1714	1716	rVB	51009	45360	1.49%	0.131%
47	12.416	1716	1722	1724	rBV	193867	208757	6.84%	0.604%
48	12.451	1724	1728	1731	rVV3	107216	182340	5.98%	0.528%
49	12.474	1731	1732	1734	rVB	77518	40111	1.32%	0.116%
50	12.510	1734	1738	1742	rBV2	149894	181400	5.95%	0.525%
51	12.586	1746	1751	1754	rBV	2220975	1979529	64.90%	5.732%
52	12.645	1759	1761	1763	rVB	66111	49748	1.63%	0.144%
53	12.674	1763	1766	1769	rBV	213940	192431	6.31%	0.557%
54	12.733	1769	1776	1778	rBV4	75636	118212	3.88%	0.342%
55	12.757	1778	1780	1783	rVB	82326	71313	2.34%	0.206%
56	12.816	1783	1790	1793	rBV2	2262416	2308792	75.70%	6.685%
57	12.863	1795	1798	1802	rVB2	138408	158277	5.19%	0.458%
58	12.945	1806	1812	1815	rBV2	519360	591782	19.40%	1.714%
59	12.974	1815	1817	1820	rVB	194292	162677	5.33%	0.471%
60	13.027	1821	1826	1830	rBV3	208432	355635	11.66%	1.030%
61	13.116	1837	1841	1842	rBV3	186717	205435	6.74%	0.595%
62	13.139	1842	1845	1848	rVB2	396407	396621	13.00%	1.148%
63	13.168	1848	1850	1854	rBV4	43591	81158	2.66%	0.235%
64	13.204	1854	1856	1859	rBV	213446	180062	5.90%	0.521%
65	13.245	1859	1863	1866	rBV2	305772	343532	11.26%	0.995%
66	13.298	1870	1872	1875	rBV2	68230	79079	2.59%	0.229%
67	13.339	1876	1879	1881	rBV	196888	156114	5.12%	0.452%
68	13.368	1881	1884	1887	rBV2	198806	180435	5.92%	0.522%
69	13.451	1896	1898	1901	rVB	58587	51510	1.69%	0.149%
70	13.515	1906	1909	1910	rBV3	37696	45756	1.50%	0.132%
71	13.621	1924	1927	1929	rBV	71231	72086	2.36%	0.209%
72	13.651	1929	1932	1936	rVB	262956	242400	7.95%	0.702%
73	13.763	1948	1951	1953	rBV2	190319	170929	5.60%	0.495%
74	13.786	1953	1955	1958	rVV	260258	210463	6.90%	0.609%
75	13.815	1958	1960	1965	rVB2	210273	268387	8.80%	0.777%
76	13.863	1965	1968	1971	rVB	234619	205096	6.72%	0.594%

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77	13.933	1975	1980	1985	rBV2	65576	136842	4.49%	0.396%
78	14.010	1985	1993	1996	rBV2	2030548	3050084	100.00%	8.832%
79	14.045	1996	1999	2005	rVV	1374991	1344662	44.09%	3.894%
80	14.104	2005	2009	2011	rVV	333385	382437	12.54%	1.107%
81	14.121	2011	2012	2014	rVV	199690	114134	3.74%	0.330%
82	14.163	2017	2019	2022	rVV	225706	226957	7.44%	0.657%
83	14.210	2025	2027	2029	rVV	106208	88415	2.90%	0.256%
84	14.245	2031	2033	2036	rVV	173325	151297	4.96%	0.438%
85	14.274	2036	2038	2040	rVB2	82383	64788	2.12%	0.188%
86	14.362	2050	2053	2055	rBV2	109446	101120	3.32%	0.293%
87	14.398	2055	2059	2063	rVV	355096	402463	13.20%	1.165%
88	14.433	2063	2065	2068	rVB	161841	122112	4.00%	0.354%
89	14.527	2078	2081	2083	rBV2	182396	171698	5.63%	0.497%
90	14.545	2083	2084	2087	rVB2	153549	108768	3.57%	0.315%
91	14.715	2111	2113	2115	rBV2	74725	73294	2.40%	0.212%
92	14.898	2141	2144	2148	rVB3	119443	148294	4.86%	0.429%
93	15.062	2167	2172	2175	rBV	1144008	1855447	60.83%	5.373%
94	15.168	2186	2190	2193	rVB	310453	317885	10.42%	0.920%
95	15.362	2219	2223	2229	rVB	629546	829186	27.19%	2.401%
96	15.427	2230	2234	2239	rVV	1035781	1276620	41.86%	3.697%
97	15.492	2240	2245	2248	rVV	574019	801277	26.27%	2.320%
98	15.521	2248	2250	2256	rVB2	299036	372147	12.20%	1.078%
99	16.974	2491	2497	2504	rVB	408459	786003	25.77%	2.276%
100	17.427	2568	2574	2582	rVB	282183	592619	19.43%	1.716%

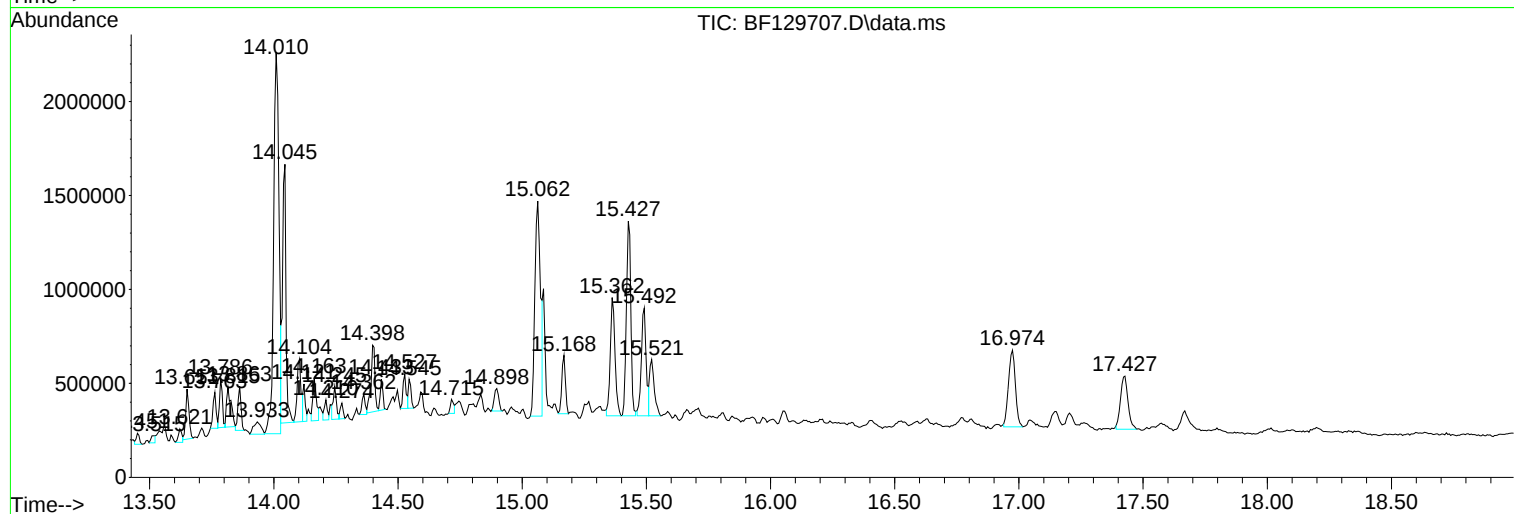
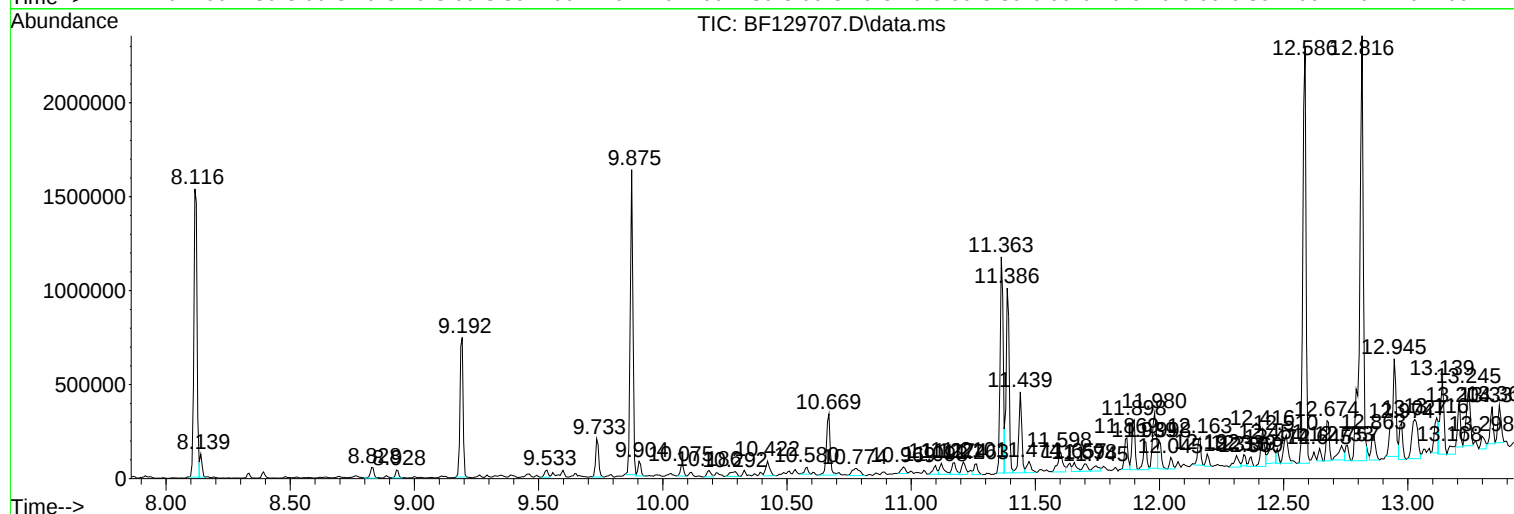
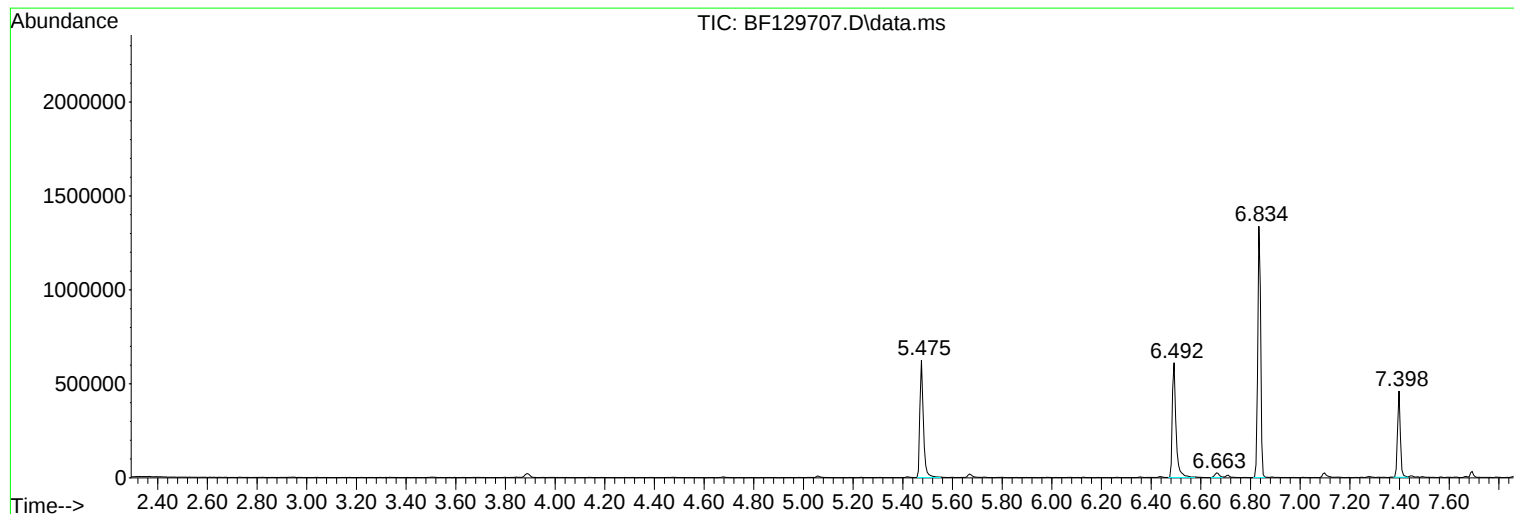
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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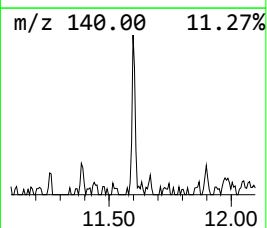
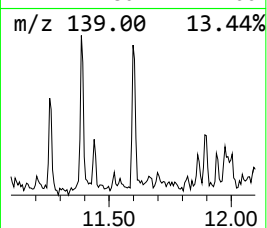
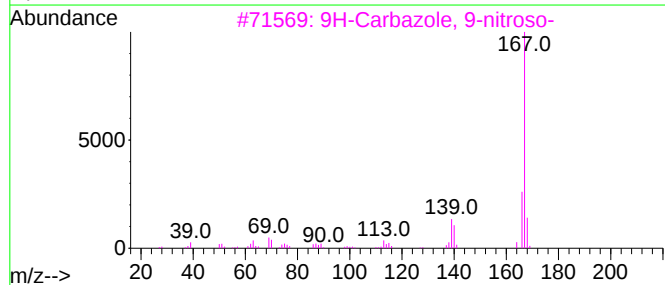
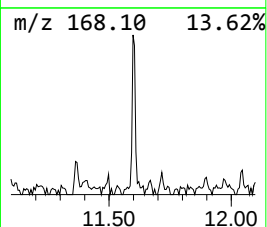
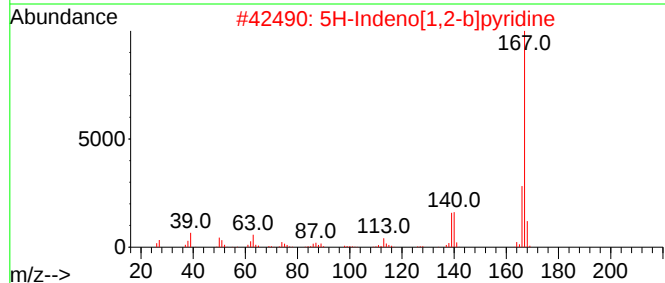
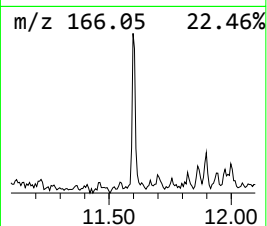
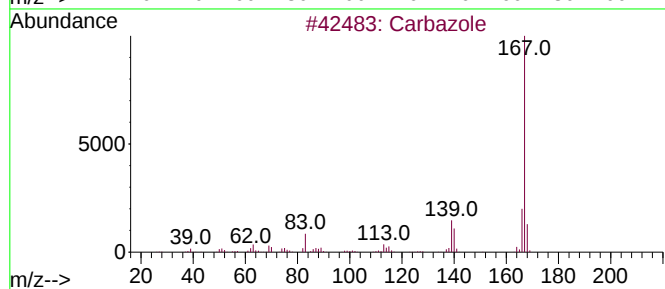
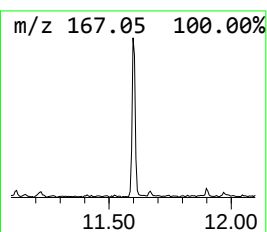
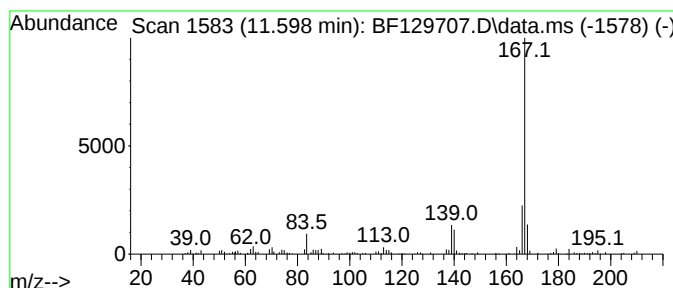
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 5H-Indeno[1,2-b]pyridine Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.598	3.13 ng	157488	Phenanthrene-d10	11.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbazole	167	C12H9N	000086-74-8	96
2			5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	95
3			9H-Carbazole, 9-nitroso-	196	C12H8N2O	002788-23-0	83
4			9H-Carbazole-9-methanol	197	C13H11NO	002409-36-1	80
5			1-Naphthaleneacetonitrile	167	C12H9N	000132-75-2	59



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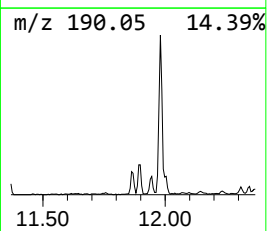
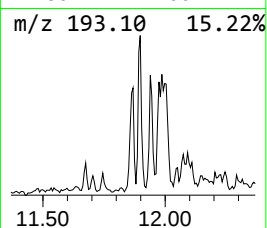
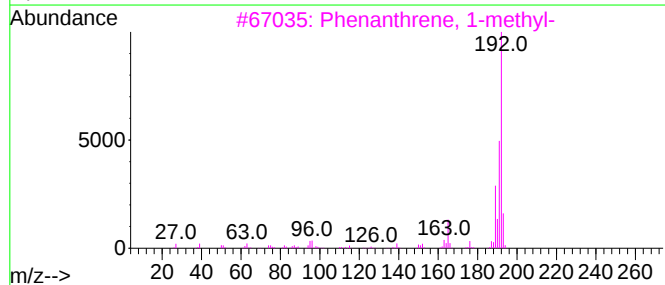
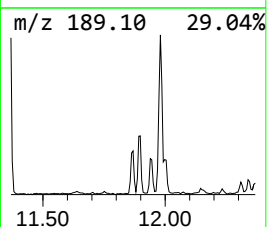
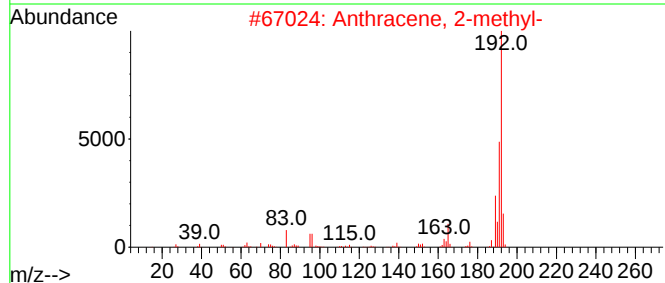
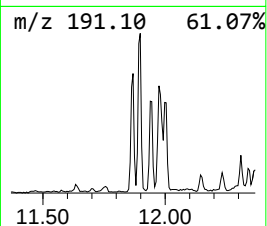
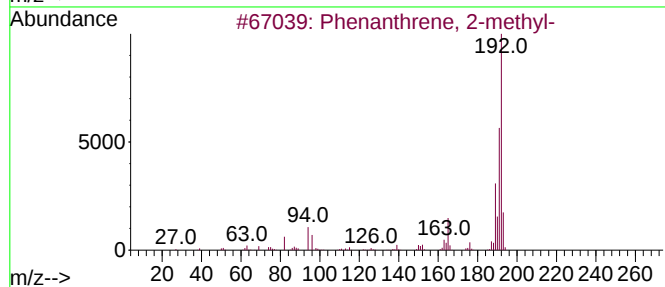
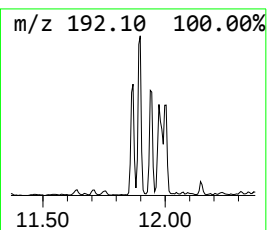
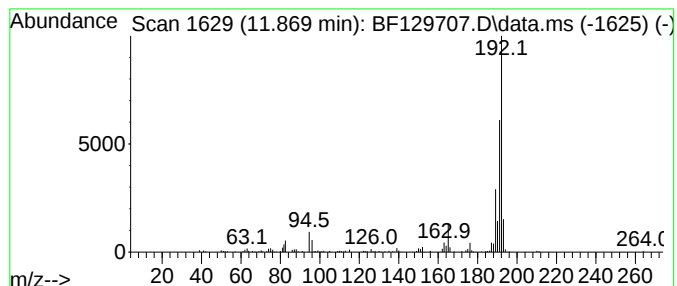
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.868	3.30 ng	166301	Phenanthrene-d10	11.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
4			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	95
5			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	94



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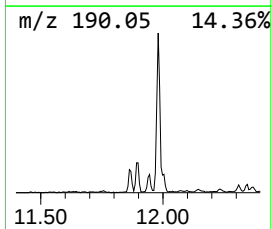
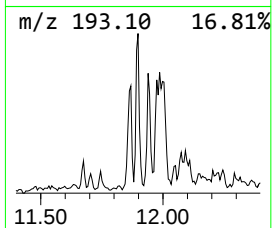
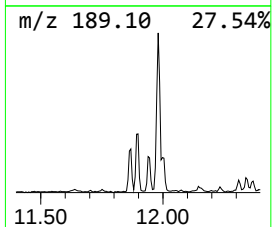
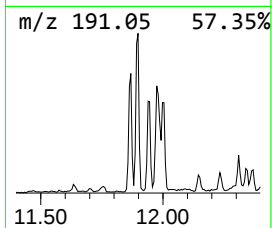
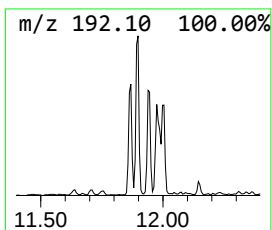
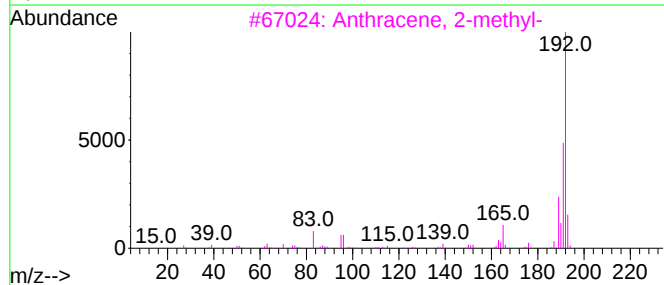
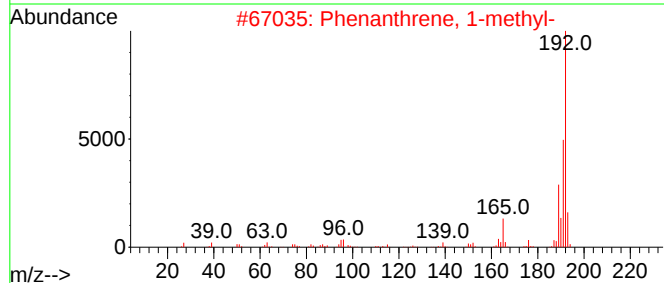
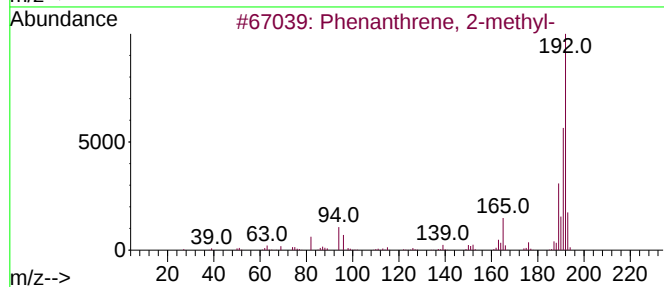
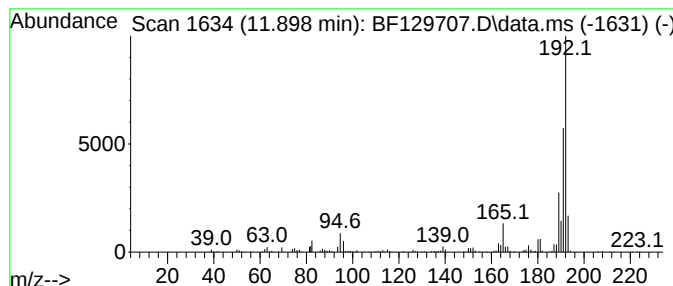
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Phenanthrene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.898	4.67 ng	235194	Phenanthrene-d10	11.363

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081022\
Data File : BF129707.D
Acq On : 10 Aug 2022 21:31
Operator : CG\JU
Sample : N4119-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

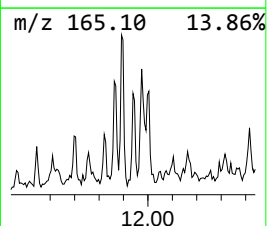
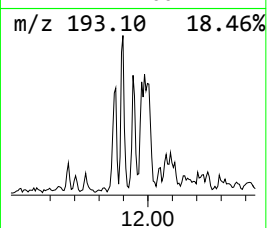
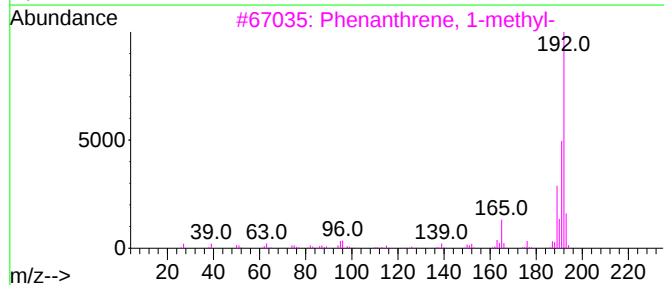
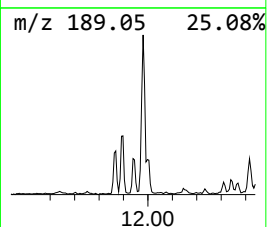
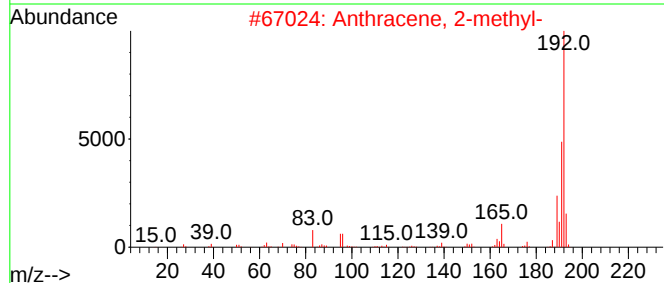
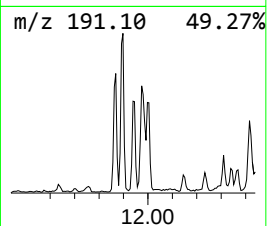
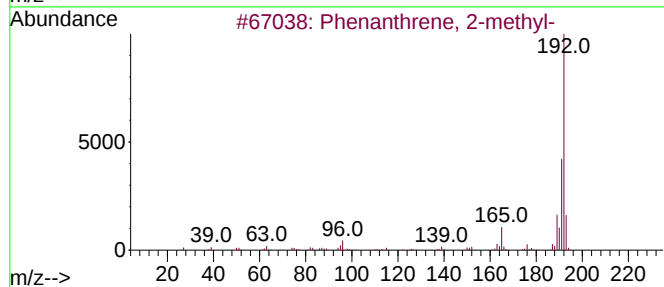
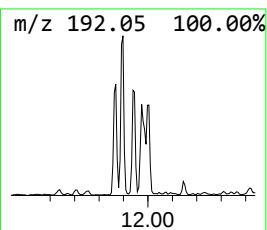
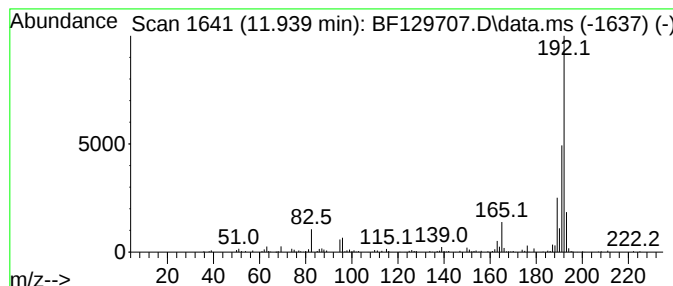
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ClientSampleId :
TP1-2-4FT

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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 Anthracene, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.939	3.04 ng	152795	Phenanthrene-d10	11.363	
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	96
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
4	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081022\
Data File : BF129707.D
Acq On : 10 Aug 2022 21:31
Operator : CG\JU
Sample : N4119-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP1-2-4FT

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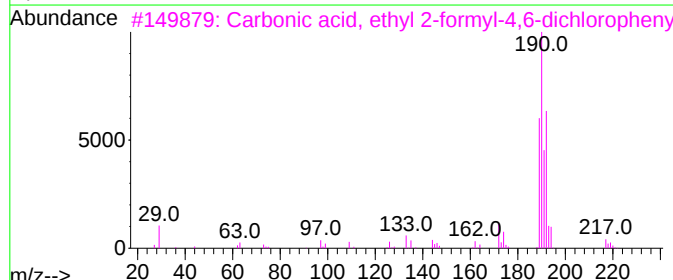
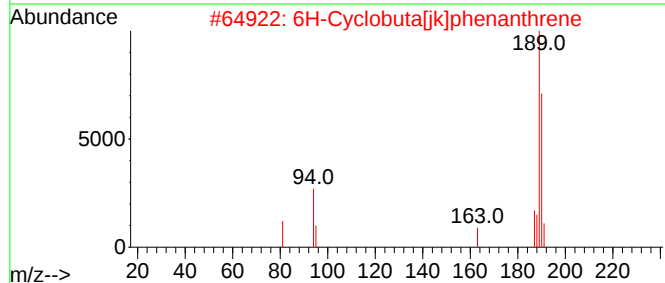
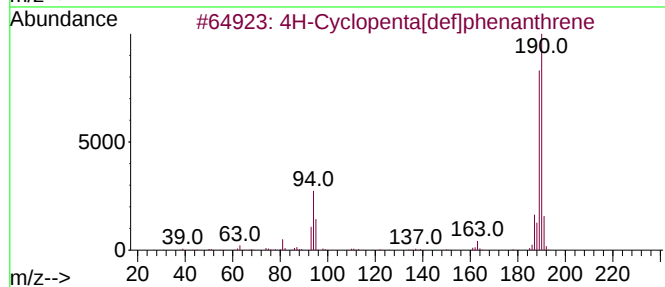
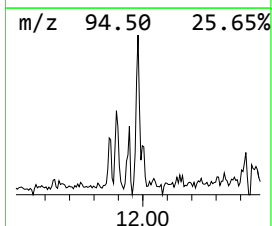
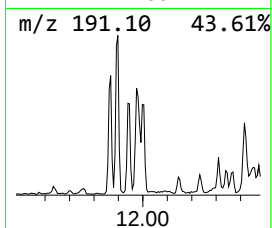
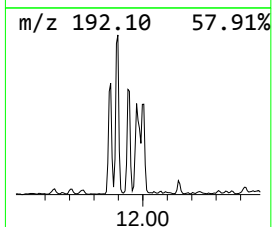
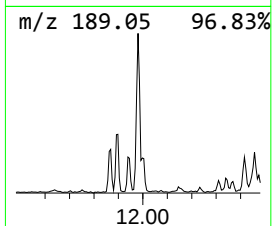
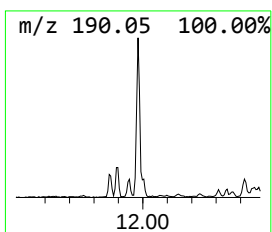
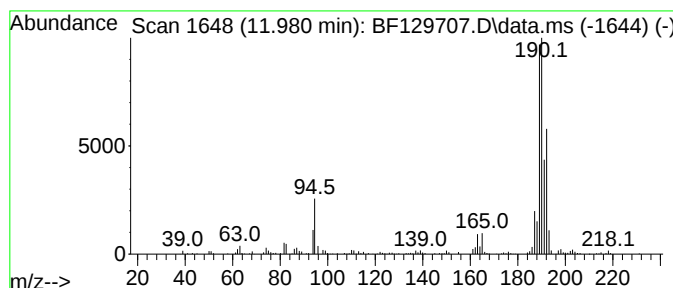
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.980	6.49 ng	326565	Phenanthrene-d10	11.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	58
3			Carbonic acid, ethyl 2-formyl-4,...	262	C10H8Cl2O4	1000331-34-4	50
4			Methyl diselenide	190	C2H6Se2	007101-31-7	41
5			2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081022\
Data File : BF129707.D
Acq On : 10 Aug 2022 21:31
Operator : CG\JU
Sample : N4119-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

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ClientSampleId :
TP1-2-4FT

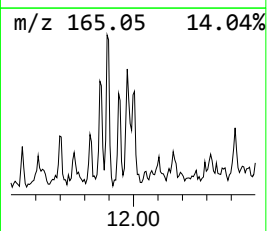
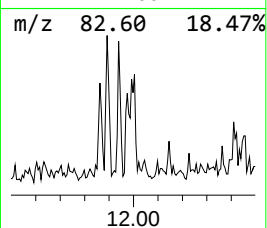
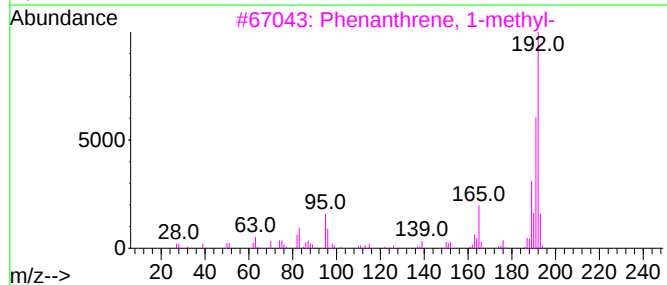
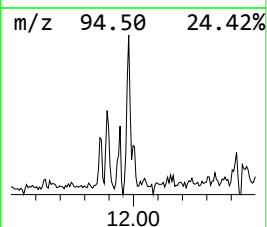
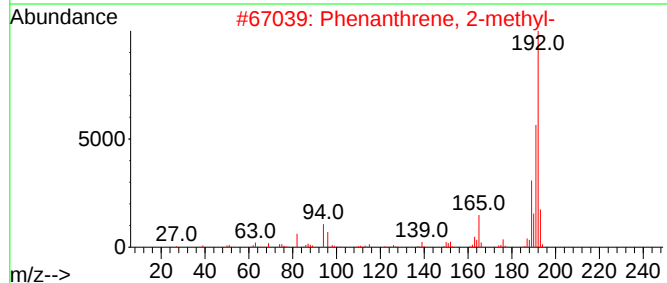
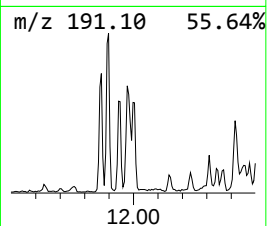
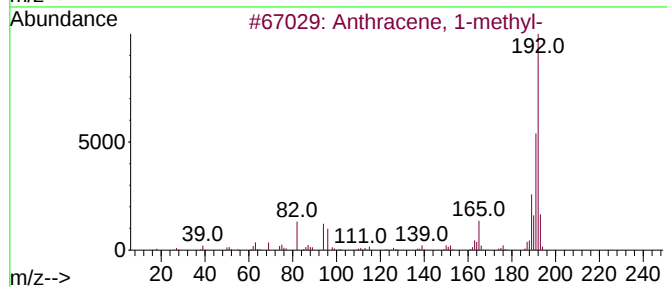
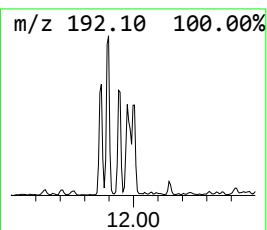
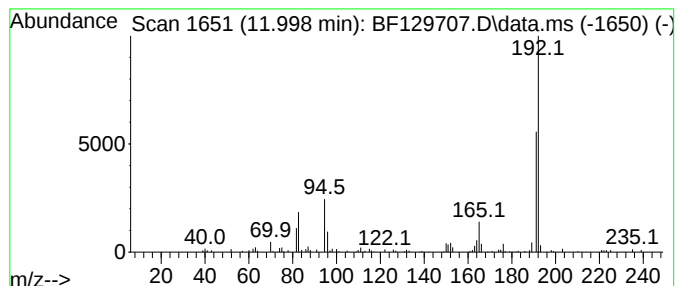
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Anthracene, 1-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.998	2.40 ng	120948	Phenanthrene-d10	11.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 1-methyl-	192	C15H12	000610-48-0	80
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	74
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	68
4			Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	64
5			Anthracene, 2-methyl-	192	C15H12	000613-12-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081022\
Data File : BF129707.D
Acq On : 10 Aug 2022 21:31
Operator : CG\JU
Sample : N4119-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

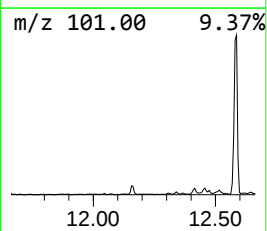
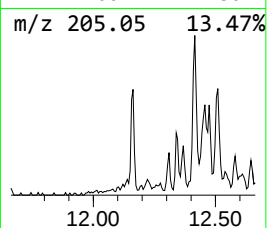
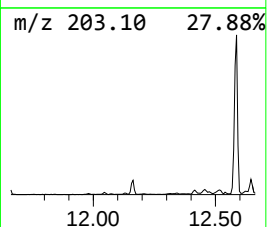
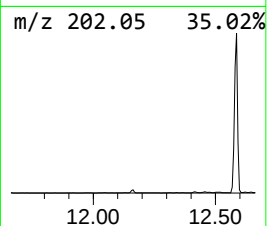
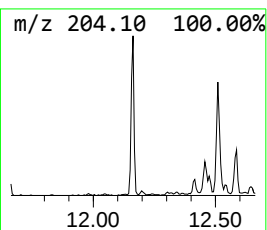
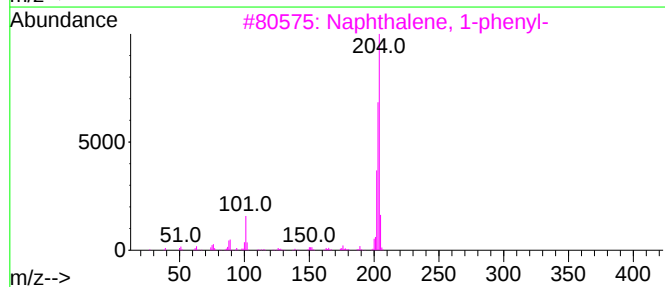
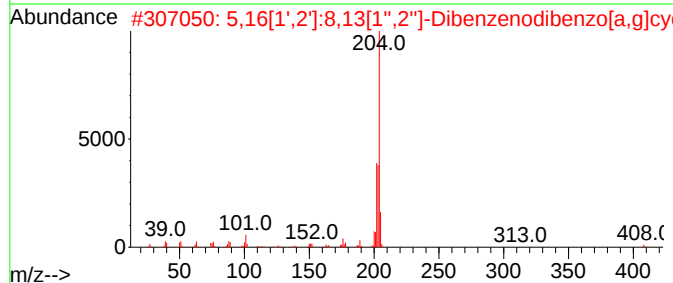
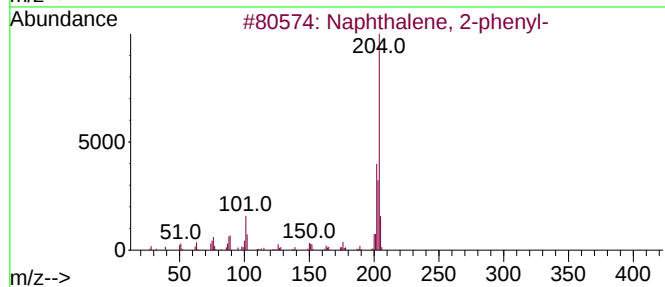
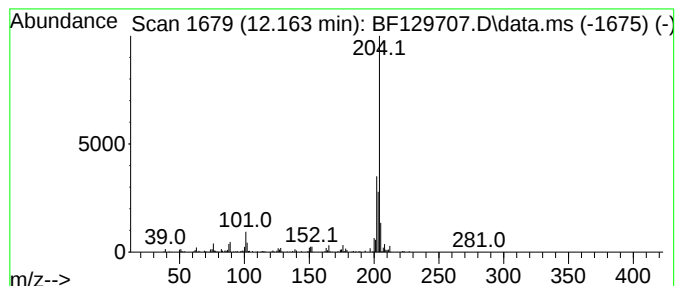
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ClientSampleId :
TP1-2-4FT

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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 Naphthalene, 2-phenyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.163	2.55 ng	128471	Phenanthrene-d10	11.363	
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	Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	64
4	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	59
5	9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081022\
Data File : BF129707.D
Acq On : 10 Aug 2022 21:31
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Sample : N4119-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

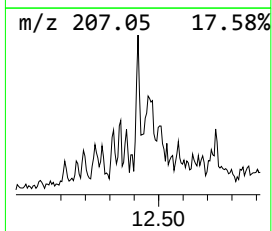
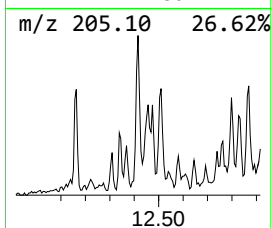
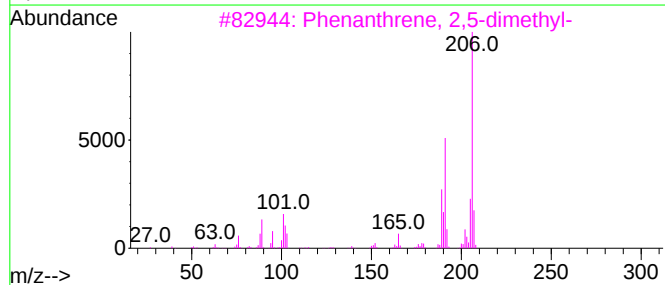
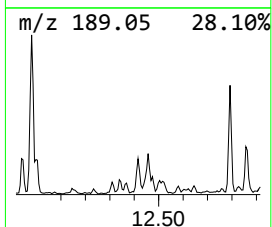
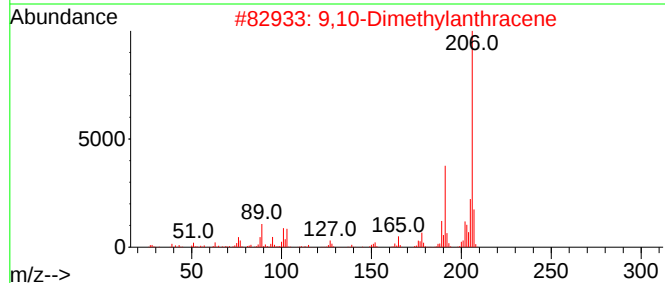
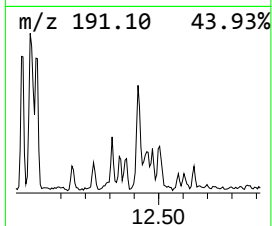
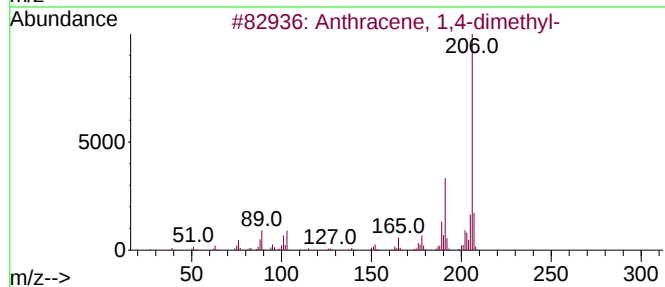
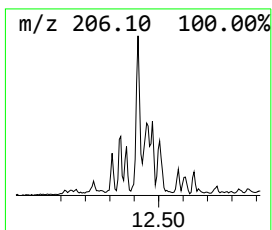
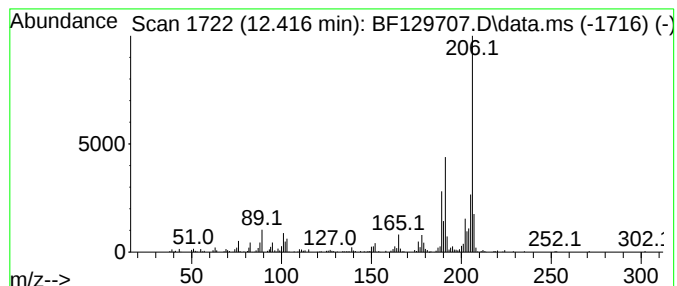
Instrument :
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ClientSampleId :
TP1-2-4FT

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Anthracene, 1,4-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.416	4.15 ng	208757	Phenanthrene-d10		11.363	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Anthracene, 1,4-dimethyl-		206	C16H14	000781-92-0	93
2	9,10-Dimethylanthracene		206	C16H14	000781-43-1	93
3	Phenanthrene, 2,5-dimethyl-		206	C16H14	003674-66-6	93
4	Phenanthrene, 2,7-dimethyl-		206	C16H14	001576-69-8	93
5	Phenanthrene, 3,6-dimethyl-		206	C16H14	001576-67-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081022\
Data File : BF129707.D
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Instrument :
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ClientSampleId :
TP1-2-4FT

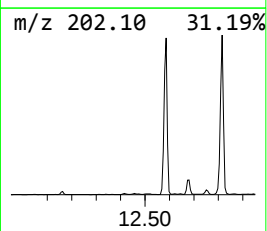
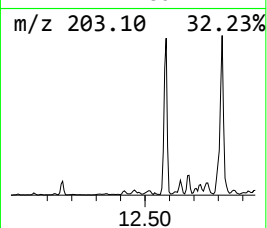
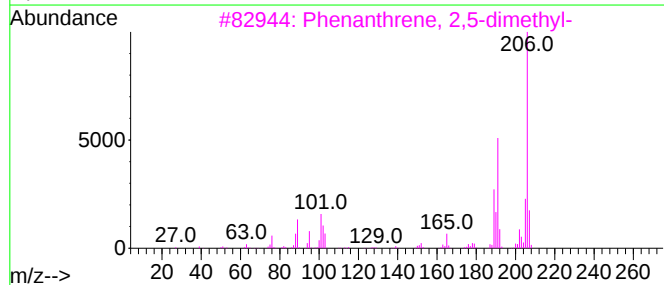
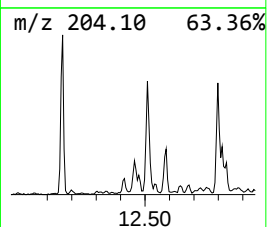
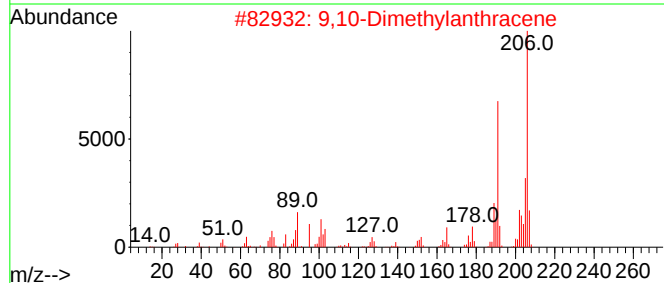
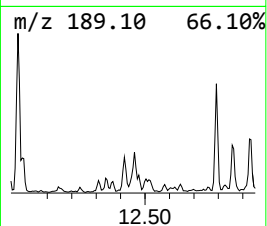
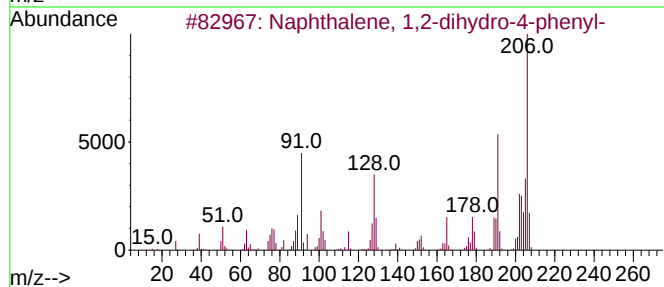
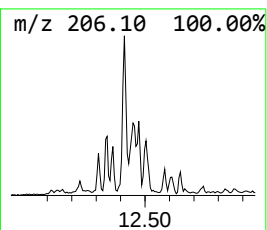
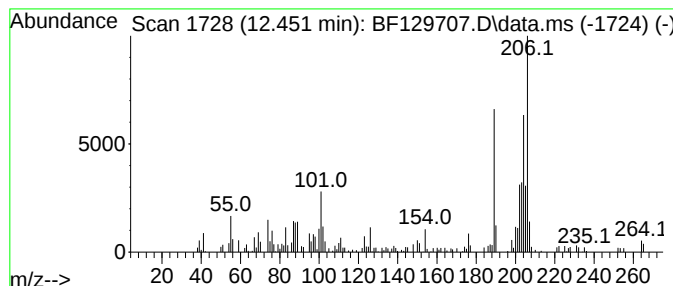
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown12.451 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.451	3.62 ng	182340	Phenanthrene-d10	11.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2-dihydro-4-phenyl-	206	C16H14	007469-40-1	45
2			9,10-Dimethylantracene	206	C16H14	000781-43-1	45
3			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	43
4			1H-Indene, 2-methyl-3-phenyl-	206	C16H14	035099-60-6	43
5			Naphthalene, 1,2-dihydro-1-phenyl-	206	C16H14	016606-46-5	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081022\
Data File : BF129707.D
Acq On : 10 Aug 2022 21:31
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ALS Vial : 19 Sample Multiplier: 1

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ClientSampleId :
TP1-2-4FT

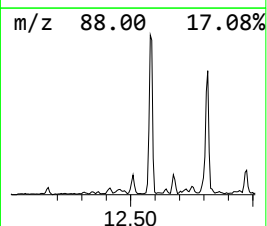
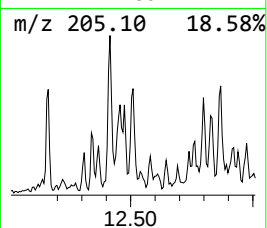
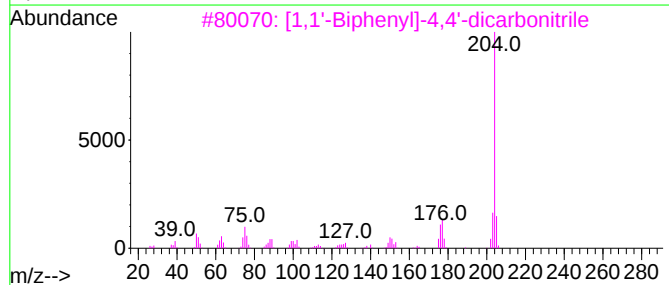
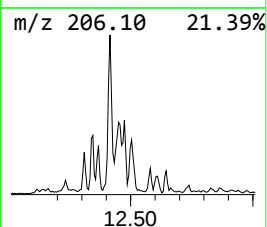
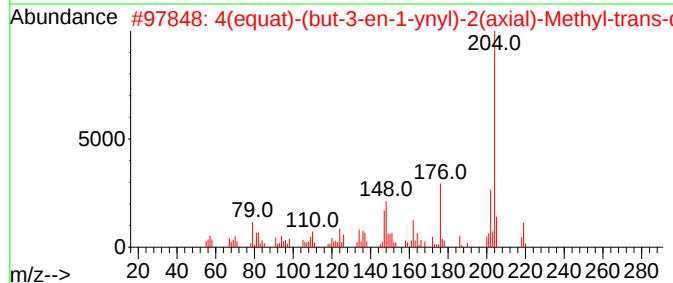
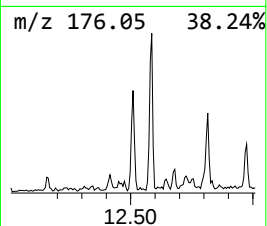
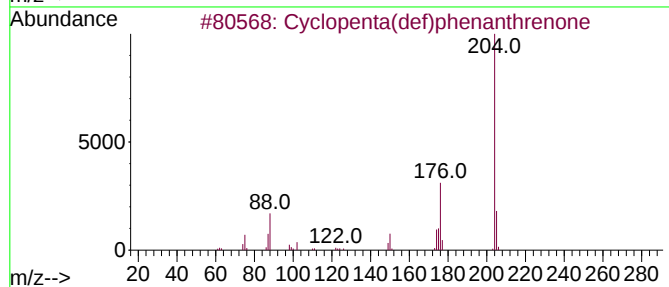
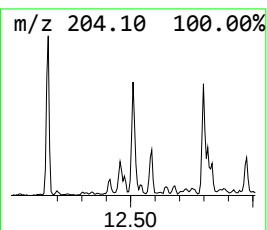
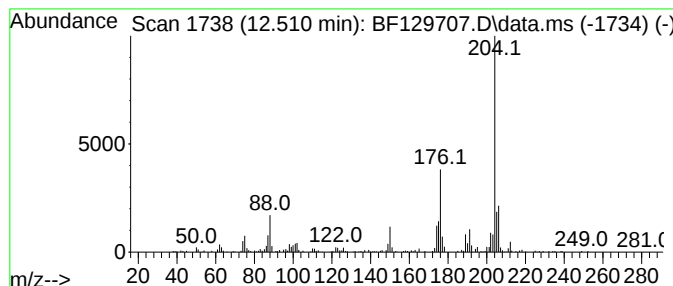
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Cyclopenta(def)phenanthrenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.510	3.60 ng	181400	Phenanthrene-d10	11.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	96
2			4(equat)-(but-3-en-1-ynyl)-2(axi...	219	C14H21NO	038423-22-2	50
3			[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	49
4			1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	41
5			Acetamide, 2-(3,4-dihydro-1-isoq...	332	C17H14F2N2OS	1000270-76-1	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081022\
Data File : BF129707.D
Acq On : 10 Aug 2022 21:31
Operator : CG\JU
Sample : N4119-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP1-2-4FT

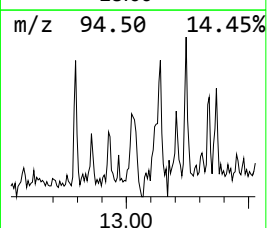
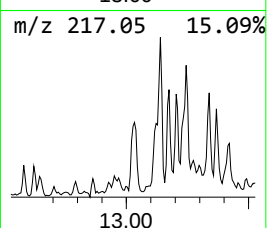
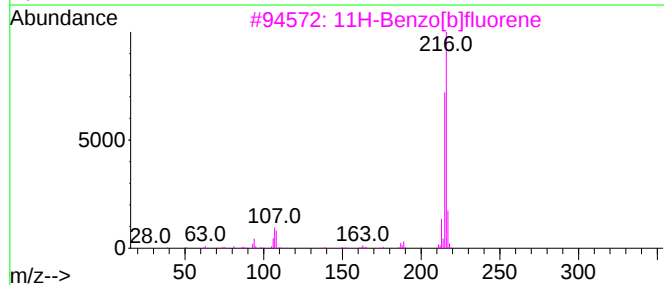
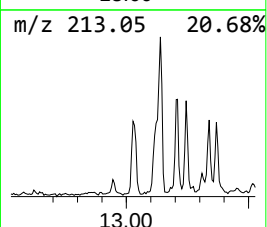
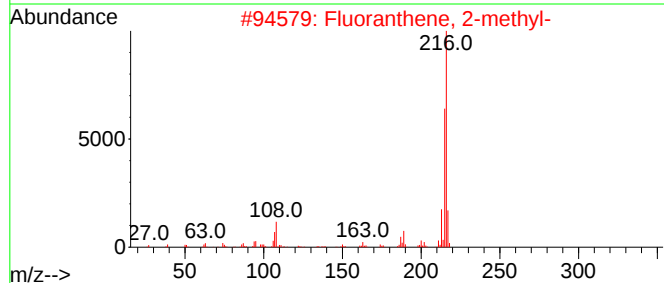
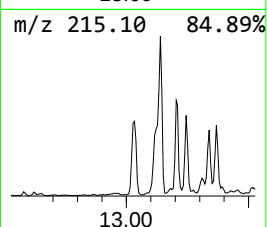
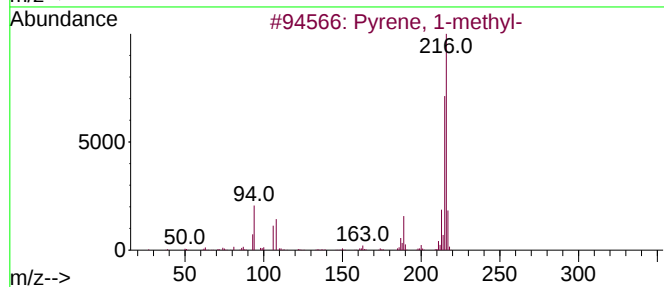
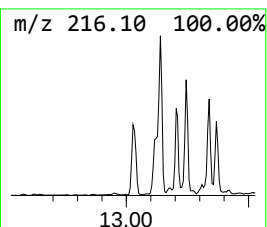
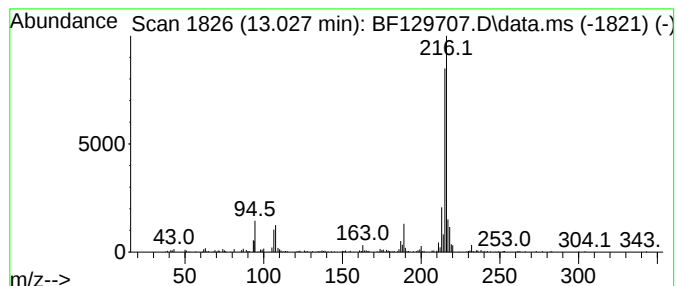
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Pyrene, 1-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.027	2.33 ng	355635	Chrysene-d12	14.015

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	97
2			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	93
3			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
4			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	81
5			Pyrene, 2-methyl-	216	C17H12	003442-78-2	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081022\
Data File : BF129707.D
Acq On : 10 Aug 2022 21:31
Operator : CG\JU
Sample : N4119-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

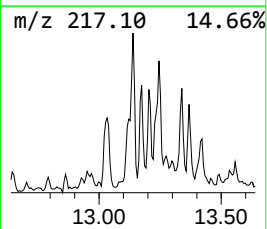
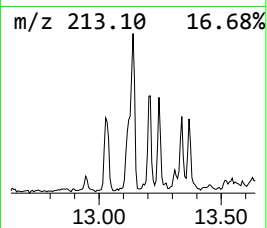
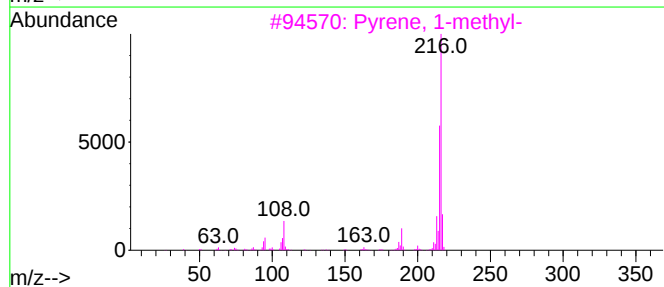
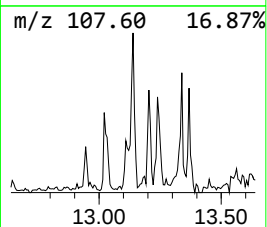
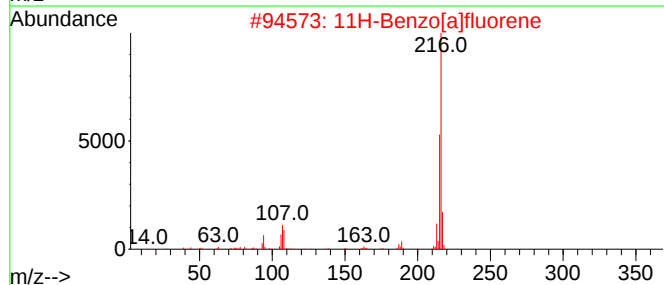
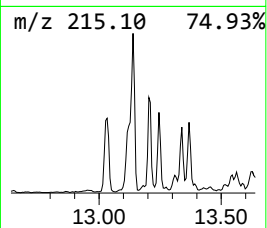
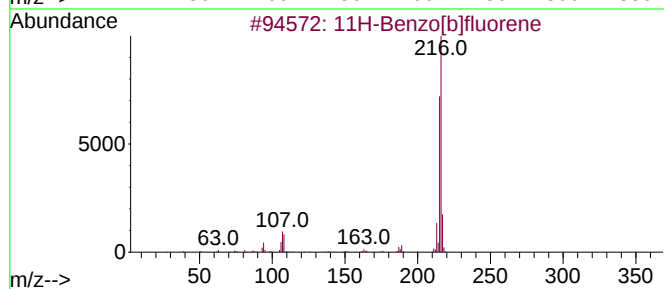
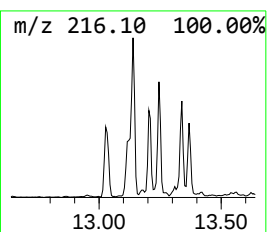
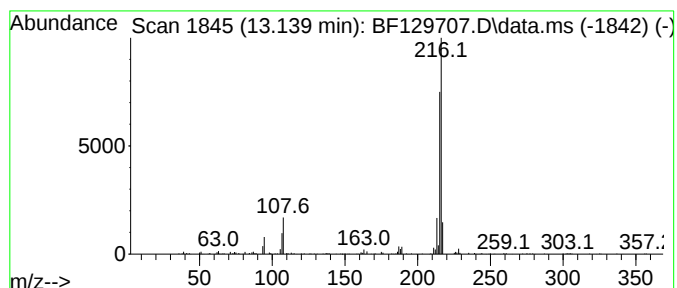
Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 11H-Benzo[b]fluorene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.139	2.60 ng	396621	Chrysene-d12	14.015	
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	90
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	83
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	83
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081022\
Data File : BF129707.D
Acq On : 10 Aug 2022 21:31
Operator : CG\JU
Sample : N4119-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

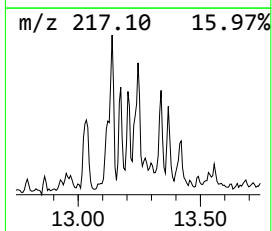
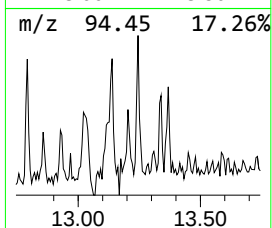
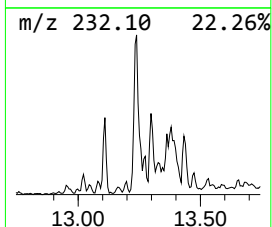
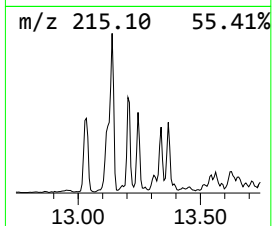
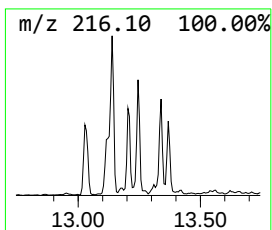
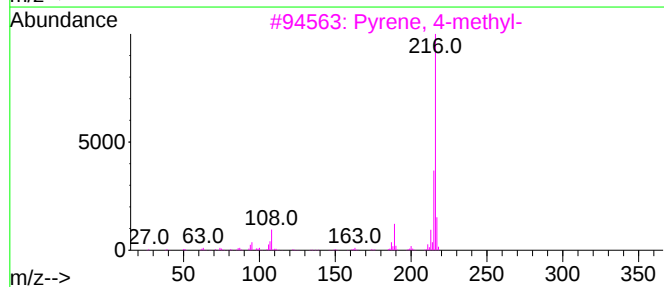
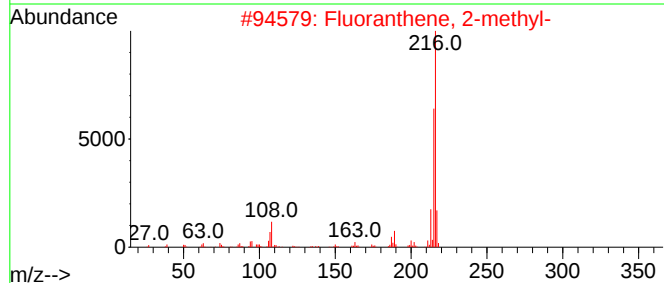
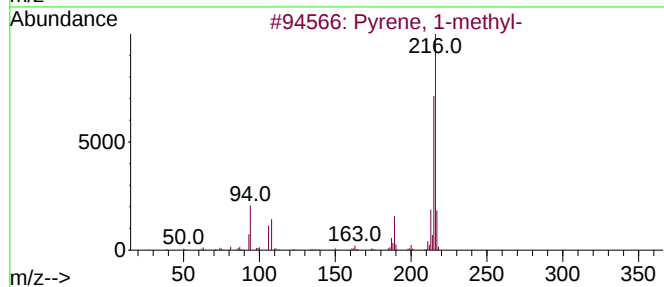
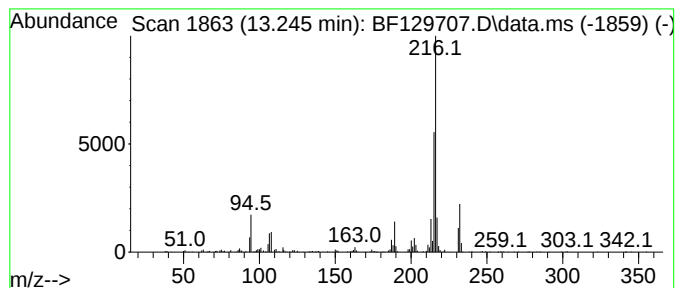
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Fluoranthene, 2-methyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.245	2.25 ng	343532	Chrysene-d12	14.015	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	96
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	95
3	Pyrene, 4-methyl-	216	C17H12	003353-12-6	93
4	Pyrene, 2-methyl-	216	C17H12	003442-78-2	93
5	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081022\
Data File : BF129707.D
Acq On : 10 Aug 2022 21:31
Operator : CG\JU
Sample : N4119-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TP1-2-4FT

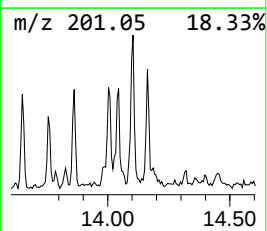
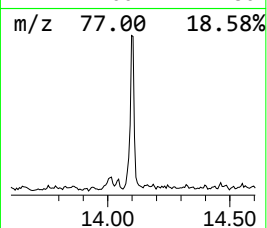
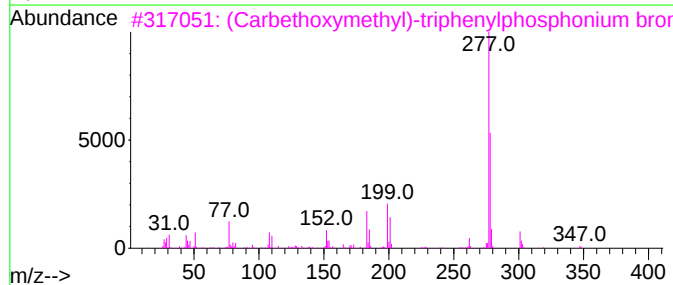
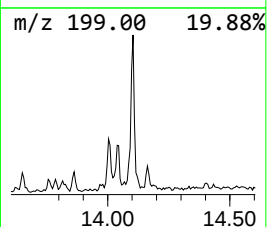
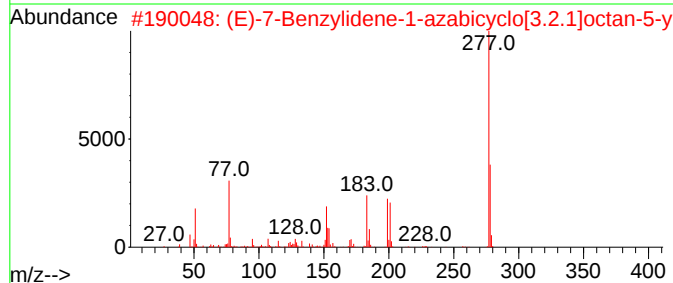
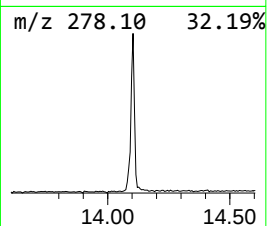
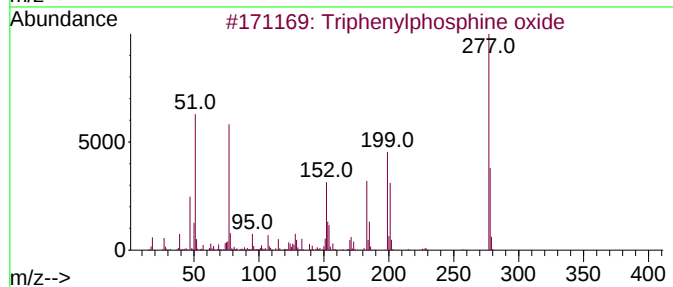
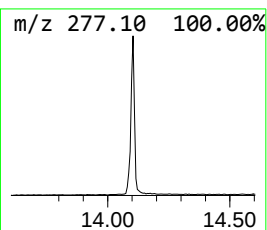
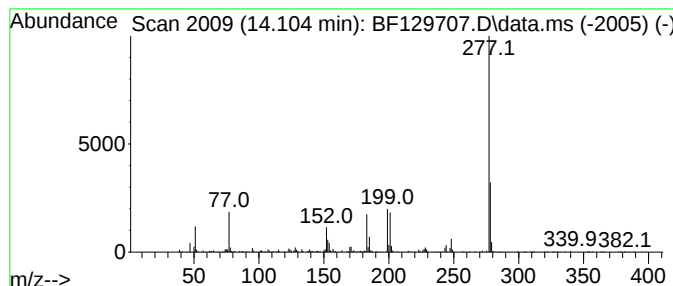
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Triphenylphosphine oxide Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.104	2.51 ng	382437	Chrysene-d12	14.015

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	96
2			(E)-7-Benzylidene-1-azabicyclo[3...]	293	C15H19N03S	1000483-83-4	86
3			(Carbethoxymethyl)-triphenylphos...	428	C22H22BrO2P	001530-45-6	47
4			2-Phenyl-6-(trifluoromethyl)-1H...	277	C15H10F3NO	1000387-59-3	40
5			5-quinolinol, 8-[2-(3,5-dimethyl...	277	C17H15N3O	1000397-10-2	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081022\
Data File : BF129707.D
Acq On : 10 Aug 2022 21:31
Operator : CG\JU
Sample : N4119-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP1-2-4FT

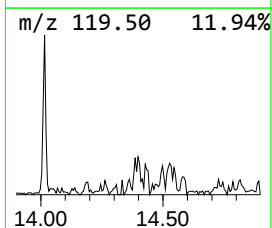
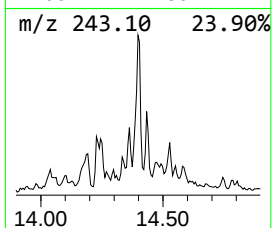
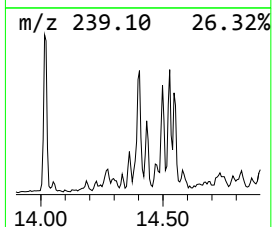
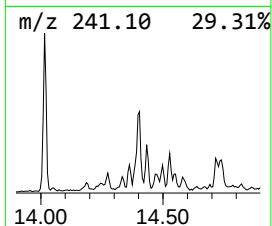
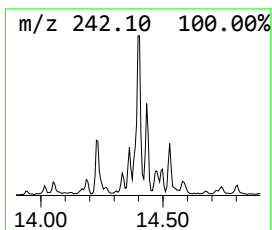
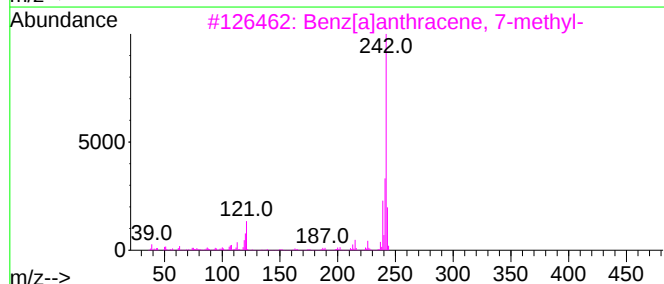
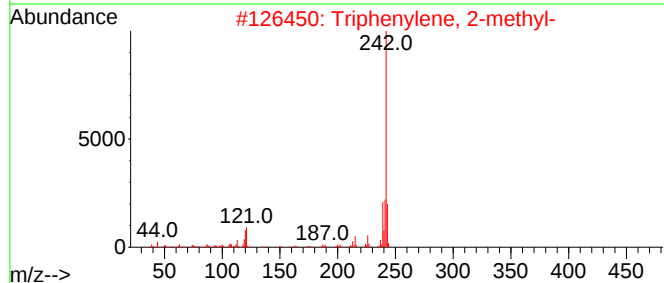
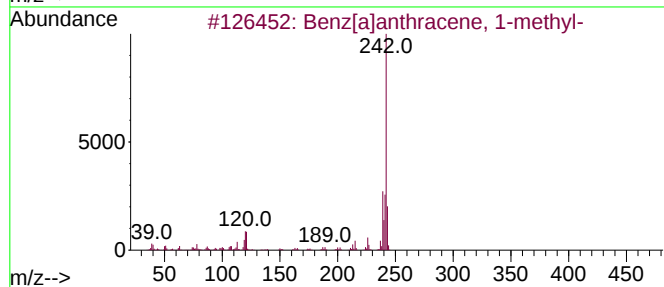
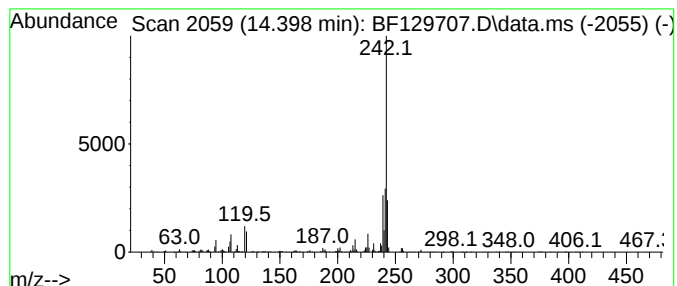
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Benz[a]anthracene, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.398	2.64 ng	402463	Chrysene-d12	14.015

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benz[a]anthracene, 1-methyl-	242	C19H14	002498-77-3	97
2			Triphenylene, 2-methyl-	242	C19H14	001705-84-6	97
3			Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	95
4			Benz[a]anthracene, 2-methyl-	242	C19H14	002498-76-2	95
5			Chrysene, 3-methyl-	242	C19H14	003351-31-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081022\
Data File : BF129707.D
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Misc :
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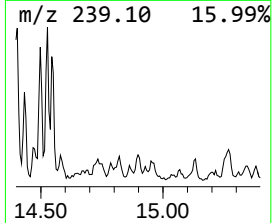
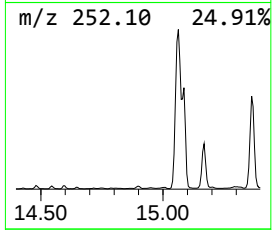
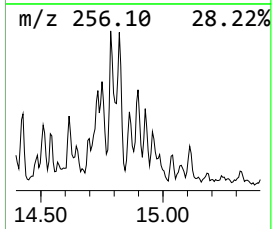
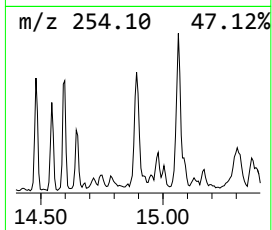
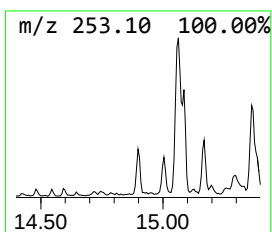
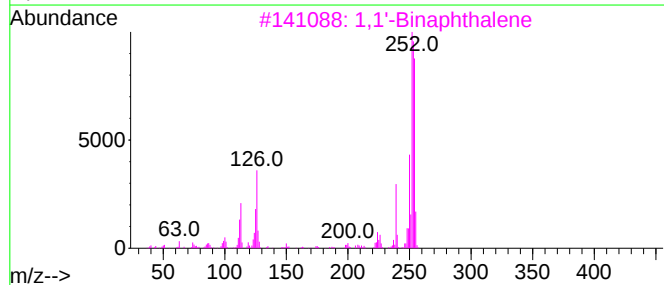
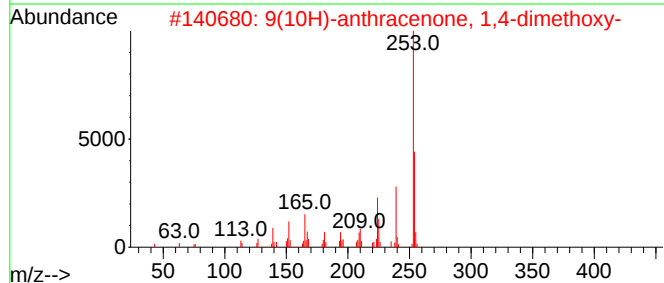
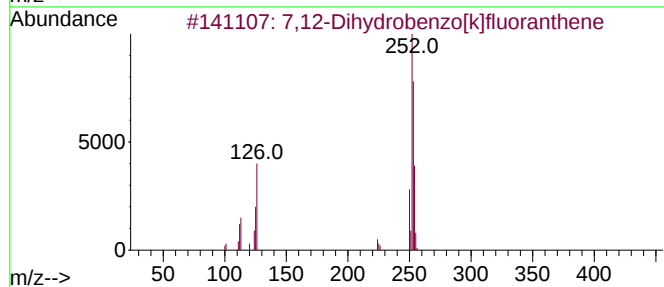
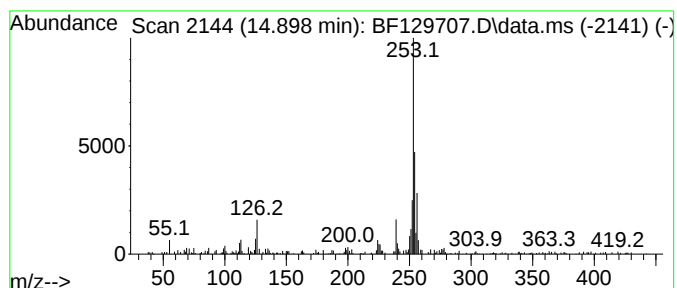
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 7,12-Dihydrobenzo[k]fluoran... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.898	3.70 ng	148294	Perylene-d12	15.492	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,12-Dihydrobenzo[k]fluoranthene	254	C20H14	1000080-17-7	70
2	9(10H)-anthracenone, 1,4-dimethoxy-	254	C16H14O3	1010396-70-2	53
3	1,1'-Binaphthalene	254	C20H14	000604-53-5	50
4	Benzofuran-5,6-diol-3-one, 2-ben...	254	C15H10O4	1000128-65-2	50
5	N-(4-Phenyl-1,3-thiazol-2-yl)-2-...	253	C14H11N3S	1000485-04-5	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081022\
Data File : BF129707.D
Acq On : 10 Aug 2022 21:31
Operator : CG\JU
Sample : N4119-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP1-2-4FT

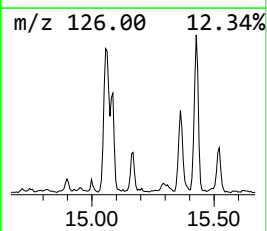
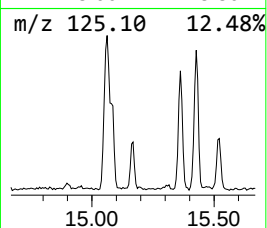
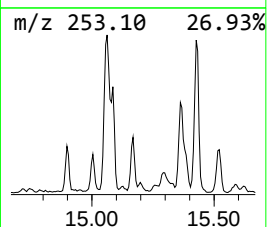
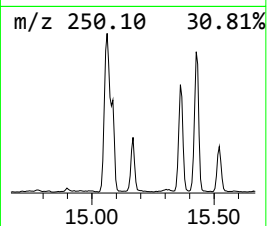
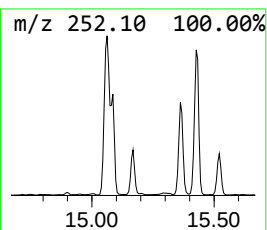
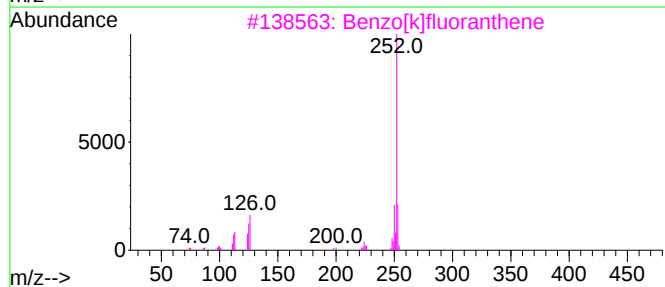
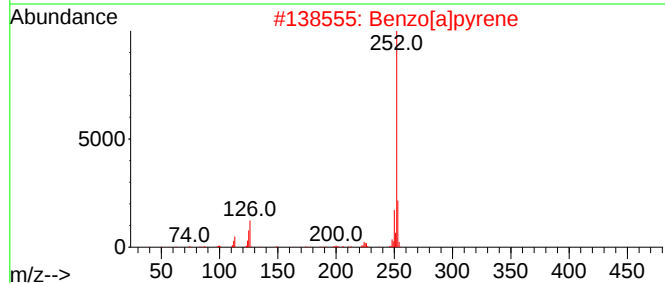
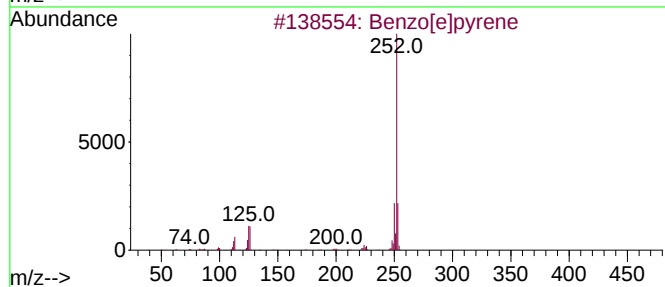
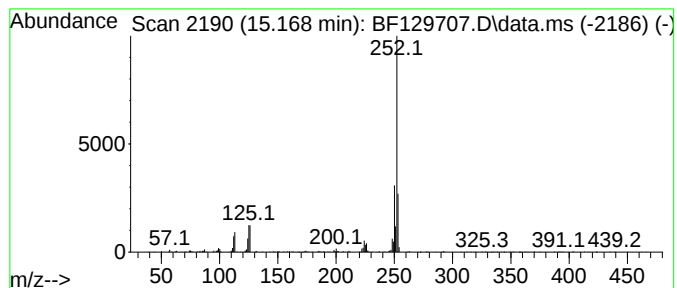
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.168	7.93 ng	317885	Perylene-d12	15.492

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[a]pyrene	252	C20H12	000050-32-8	95
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	94
4			Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
5			Perylene	252	C20H12	000198-55-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081022\
Data File : BF129707.D
Acq On : 10 Aug 2022 21:31
Operator : CG\JU
Sample : N4119-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP1-2-4FT

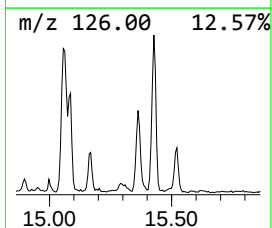
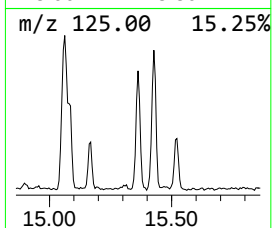
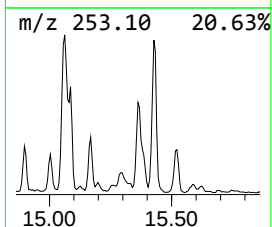
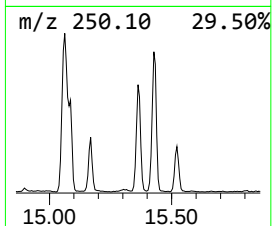
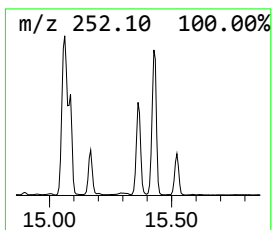
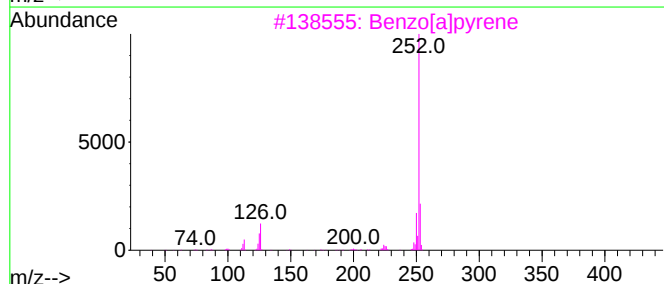
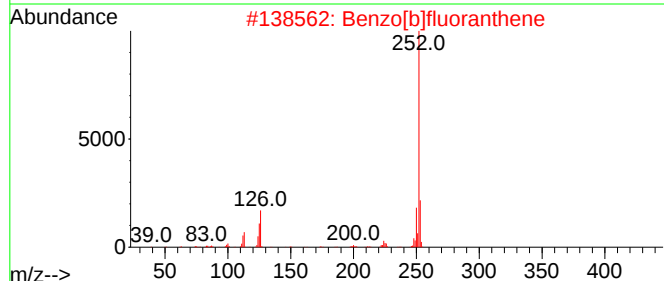
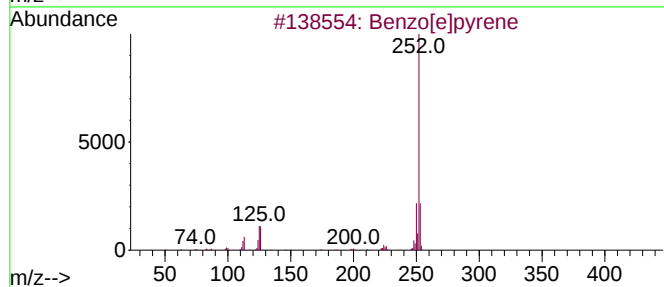
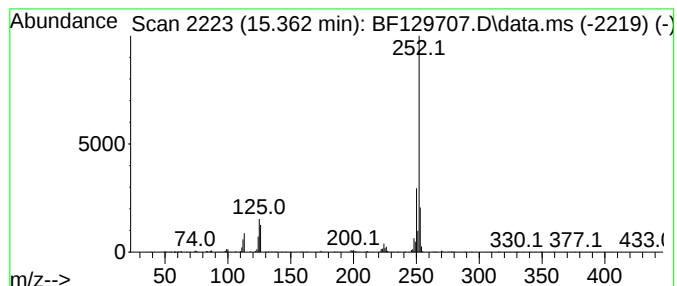
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Perylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.362	20.70 ng	829186	Perylene-d12	15.492

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[e]pyrene	252	C20H12	000192-97-2	99
2		Benzo[b]fluoranthene	252	C20H12	000205-99-2	98
3		Benzo[a]pyrene	252	C20H12	000050-32-8	97
4		Perylene	252	C20H12	000198-55-0	96
5		Benzo[k]fluoranthene	252	C20H12	000207-08-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081022\
 Data File : BF129707.D
 Acq On : 10 Aug 2022 21:31
 Operator : CG\JU
 Sample : N4119-01 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP1-2-4FT

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.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
5H-Indeno[1,2-b...	11.598	3.1	ng	157488	4	11.363	1006510	20.0
Phenanthrene, 2...	11.868	3.3	ng	166301	4	11.363	1006510	20.0
Phenanthrene, 1...	11.898	4.7	ng	235194	4	11.363	1006510	20.0
Anthracene, 2-m...	11.939	3.0	ng	152795	4	11.363	1006510	20.0
4H-Cyclopenta[d...	11.980	6.5	ng	326565	4	11.363	1006510	20.0
Anthracene, 1-m...	11.998	2.4	ng	120948	4	11.363	1006510	20.0
Naphthalene, 2-...	12.163	2.5	ng	128471	4	11.363	1006510	20.0
Anthracene, 1,4...	12.416	4.2	ng	208757	4	11.363	1006510	20.0
unknown12.451	12.451	3.6	ng	182340	4	11.363	1006510	20.0
Cyclopenta(def)...	12.510	3.6	ng	181400	4	11.363	1006510	20.0
Pyrene, 1-methyl-	13.027	2.3	ng	355635	5	14.015	3050080	20.0
11H-Benzo[b]flu...	13.139	2.6	ng	396621	5	14.015	3050080	20.0
Fluoranthene, 2...	13.245	2.3	ng	343532	5	14.015	3050080	20.0
Triphenylphosph...	14.104	2.5	ng	382437	5	14.015	3050080	20.0
Benz[a]anthrace...	14.398	2.6	ng	402463	5	14.015	3050080	20.0
7,12-Dihydroben...	14.898	3.7	ng	148294	6	15.492	801277	20.0
Benzo[e]pyrene	15.168	7.9	ng	317885	6	15.492	801277	20.0
Perylene	15.362	20.7	ng	829186	6	15.492	801277	20.0