

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
 Data File : BF123998.D
 Acq On : 19 May 2021 19:30
 Operator : JU/CG
 Sample : M2437-05 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-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-BF051421.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF123998.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.152	4	11	16	rVB	92349	119081	20.23%	1.496%
2	5.104	510	513	517	rVB2	11437	11477	1.95%	0.144%
3	5.510	578	582	586	rBV	312095	265058	45.04%	3.329%
4	6.516	749	753	760	rBV	347480	276960	47.06%	3.478%
5	6.904	815	819	822	rBV	413839	305569	51.92%	3.838%
6	7.457	910	913	917	rBV	217109	182467	31.00%	2.292%
7	8.186	1034	1037	1040	rBV	530780	398898	67.78%	5.010%
8	8.210	1040	1041	1044	rVB	26453	13121	2.23%	0.165%
9	9.263	1216	1220	1223	rBV	369877	300272	51.02%	3.771%
10	9.375	1235	1239	1248	rVB3	29771	31605	5.37%	0.397%
11	9.651	1284	1286	1290	rBV	38044	34124	5.80%	0.429%
12	9.804	1309	1312	1315	rBV	25287	20470	3.48%	0.257%
13	9.945	1332	1336	1339	rBV	500809	417589	70.95%	5.244%
14	9.975	1339	1341	1346	rVB	15334	14405	2.45%	0.181%
15	10.492	1425	1429	1432	rBV4	12044	14109	2.40%	0.177%
16	10.727	1465	1469	1474	rBV	203604	176407	29.97%	2.215%
17	10.857	1487	1491	1494	rVB4	10617	15507	2.63%	0.195%
18	11.186	1546	1547	1551	rVV4	9803	10653	1.81%	0.134%
19	11.233	1551	1555	1558	rVB5	11932	14662	2.49%	0.184%
20	11.292	1558	1565	1568	rBV7	7994	18611	3.16%	0.234%
21	11.327	1568	1571	1574	rVV3	13168	13498	2.29%	0.170%
22	11.427	1585	1588	1591	rBV	427861	409760	69.62%	5.146%
23	11.451	1591	1592	1597	rVV	195005	136483	23.19%	1.714%
24	11.504	1597	1601	1605	rVV	45453	43572	7.40%	0.547%
25	11.657	1622	1627	1629	rBV5	12190	14017	2.38%	0.176%
26	11.727	1634	1639	1641	rVV5	15147	19316	3.28%	0.243%
27	11.763	1641	1645	1651	rVB6	16167	20101	3.42%	0.252%
28	11.933	1671	1674	1675	rVV	36645	32121	5.46%	0.403%
29	11.957	1675	1678	1682	rVV2	57443	63630	10.81%	0.799%
30	12.004	1683	1686	1689	rVV2	18244	23252	3.95%	0.292%
31	12.045	1689	1693	1695	rVV	63037	74408	12.64%	0.934%
32	12.063	1695	1696	1701	rVB2	38688	31815	5.41%	0.400%
33	12.163	1711	1713	1716	rVB4	16442	12932	2.20%	0.162%
34	12.198	1716	1719	1721	rBV3	10030	11574	1.97%	0.145%
35	12.227	1721	1724	1726	rVV2	31604	28043	4.76%	0.352%
36	12.245	1726	1727	1731	rVB4	21280	21338	3.63%	0.268%

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Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051421.M
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37	12.380	1747	1750	1753	rVV4	19663	19133	3.25%	0.240%
38	12.410	1753	1755	1757	rVV2	11412	12700	2.16%	0.159%
39	12.480	1764	1767	1770	rVV2	49838	47892	8.14%	0.601%
40	12.516	1770	1773	1776	rVV5	25144	31430	5.34%	0.395%

41	12.563	1779	1781	1787	rVB5	26384	32894	5.59%	0.413%
42	12.645	1792	1795	1800	rBV	276683	242069	41.13%	3.040%
43	12.739	1808	1811	1813	rBV	29034	28240	4.80%	0.355%
44	12.768	1813	1816	1818	rBV3	12967	14707	2.50%	0.185%
45	12.804	1818	1822	1823	rBV3	11054	12850	2.18%	0.161%

46	12.874	1828	1834	1837	rBV	379098	379601	64.50%	4.767%
47	12.927	1840	1843	1847	rVB5	23228	34230	5.82%	0.430%
48	12.986	1847	1853	1855	rBV7	18780	30038	5.10%	0.377%
49	13.010	1855	1857	1864	rVB	289667	303107	51.50%	3.807%
50	13.092	1864	1871	1875	rBV5	43372	78413	13.32%	0.985%

51	13.198	1882	1889	1892	rVB3	53848	101461	17.24%	1.274%
52	13.245	1895	1897	1898	rBV2	14686	12436	2.11%	0.156%
53	13.268	1898	1901	1903	rVB3	21199	18702	3.18%	0.235%
54	13.304	1903	1907	1910	rBV3	73630	61907	10.52%	0.777%
55	13.368	1915	1918	1920	rVV4	10842	16344	2.78%	0.205%

56	13.398	1920	1923	1926	rVV	56868	58241	9.90%	0.731%
57	13.427	1926	1928	1931	rVV2	60002	52173	8.86%	0.655%
58	13.463	1931	1934	1936	rVV4	11727	15473	2.63%	0.194%
59	13.521	1938	1944	1946	rVB7	11913	19785	3.36%	0.248%
60	13.580	1951	1954	1956	rBV4	14357	19778	3.36%	0.248%

61	13.610	1957	1959	1964	rVB6	15497	18059	3.07%	0.227%
62	13.651	1964	1966	1969	rVB4	16774	15571	2.65%	0.196%
63	13.704	1971	1975	1981	rVV4	24998	40650	6.91%	0.511%
64	13.821	1991	1995	1997	rBV3	26578	39181	6.66%	0.492%
65	13.851	1997	2000	2002	rVV	26150	22135	3.76%	0.278%

66	13.868	2002	2003	2007	rVV2	34700	35319	6.00%	0.444%
67	13.915	2007	2011	2015	rVB4	46228	58028	9.86%	0.729%
68	14.068	2030	2037	2040	rBV2	453641	588534	100.00%	7.391%
69	14.098	2040	2042	2048	rVV	175735	166700	28.32%	2.094%
70	14.157	2050	2052	2053	rVV2	18492	15784	2.68%	0.198%

71	14.174	2053	2055	2058	rVV	51965	55231	9.38%	0.694%
72	14.210	2058	2061	2068	rVV2	57146	88510	15.04%	1.112%
73	14.298	2072	2076	2079	rBV5	18501	30934	5.26%	0.388%
74	14.327	2079	2081	2084	rVB4	21195	18102	3.08%	0.227%
75	14.421	2095	2097	2099	rBV2	26480	23025	3.91%	0.289%

76	14.457	2099	2103	2107	rVV3	56953	70233	11.93%	0.882%
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 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 >
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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	14.492	2107	2109	2112	rVB2	42427	34648	5.89%	0.435%
78	14.545	2112	2118	2121	rBV6	36475	66732	11.34%	0.838%
79	14.586	2122	2125	2126	rVV3	28231	25561	4.34%	0.321%
80	14.604	2126	2128	2130	rVB2	29641	26161	4.45%	0.329%
81	14.639	2133	2134	2136	rBV2	13542	11880	2.02%	0.149%
82	14.798	2158	2161	2167	rVB8	18912	25531	4.34%	0.321%
83	14.851	2168	2170	2174	rVV5	9858	12758	2.17%	0.160%
84	14.968	2188	2190	2192	rVB3	15535	13702	2.33%	0.172%
85	15.027	2198	2200	2203	rVB4	15355	13121	2.23%	0.165%
86	15.127	2212	2217	2224	rVB	121136	247034	41.97%	3.102%
87	15.210	2228	2231	2233	rBV4	12354	16589	2.82%	0.208%
88	15.233	2233	2235	2240	rVB2	35017	38464	6.54%	0.483%
89	15.433	2266	2269	2276	rVB	94210	113320	19.25%	1.423%
90	15.498	2276	2280	2287	rVB	138133	168341	28.60%	2.114%
91	15.562	2287	2291	2295	rBV	191375	243249	41.33%	3.055%
92	15.598	2295	2297	2305	rVB4	41031	54976	9.34%	0.690%
93	15.892	2344	2347	2351	rVB6	10183	14904	2.53%	0.187%
94	16.057	2372	2375	2376	rBV3	15644	11439	1.94%	0.144%
95	16.139	2386	2389	2393	rVB6	12400	14580	2.48%	0.183%
96	16.904	2512	2519	2523	rVB9	15250	38664	6.57%	0.486%
97	17.062	2541	2546	2557	rVB3	53287	111745	18.99%	1.403%
98	17.151	2557	2561	2564	rBV5	8567	14657	2.49%	0.184%
99	17.245	2574	2577	2583	rVB8	10266	20119	3.42%	0.253%
100	17.515	2618	2623	2630	rVB3	49622	97899	16.63%	1.229%

Sum of corrected areas: 7962579

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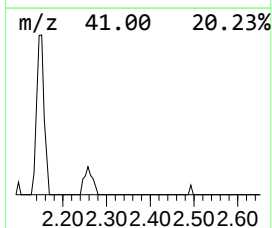
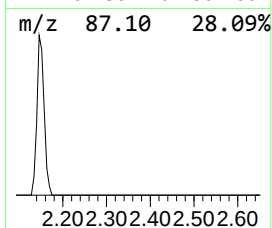
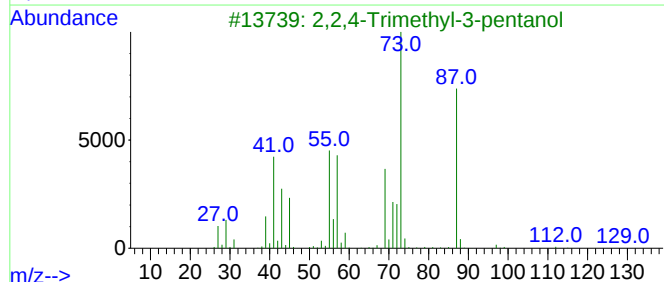
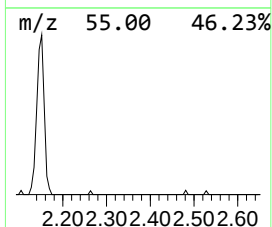
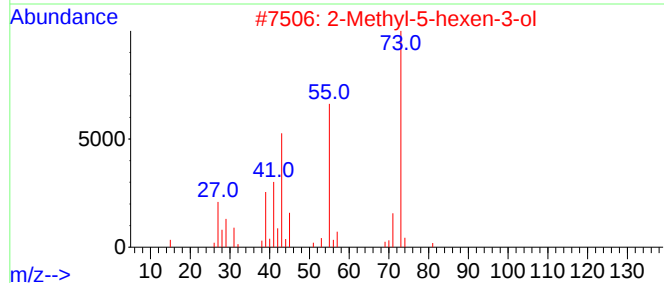
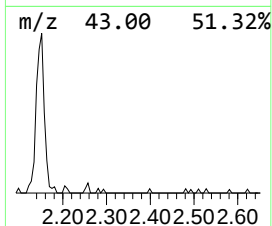
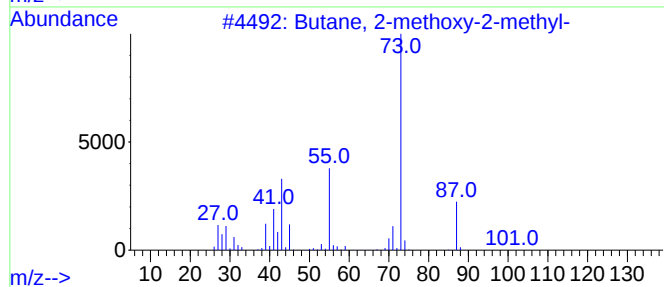
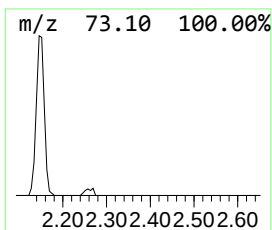
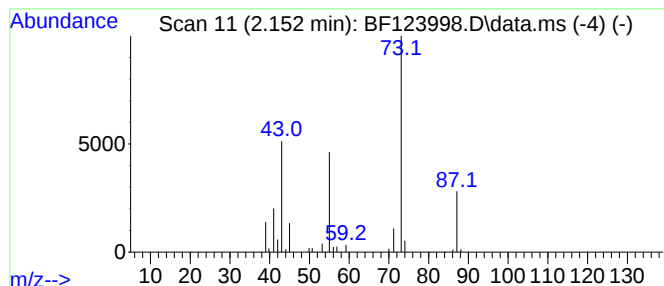
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.152	7.79 ng	119081	1,4-Dichlorobenzene-d4	6.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	40
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	16
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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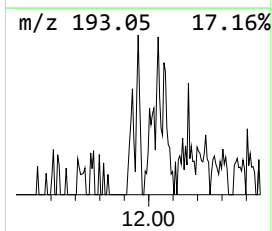
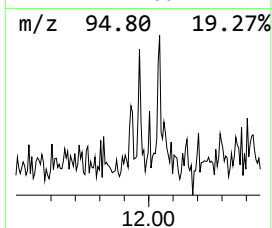
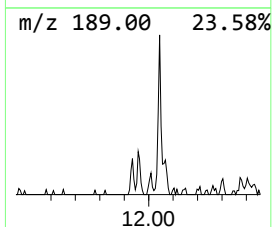
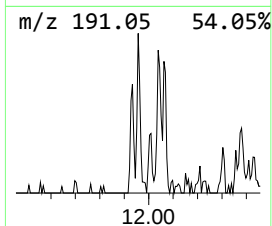
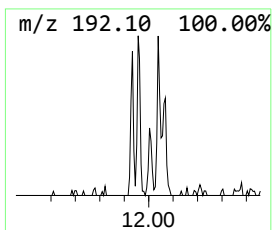
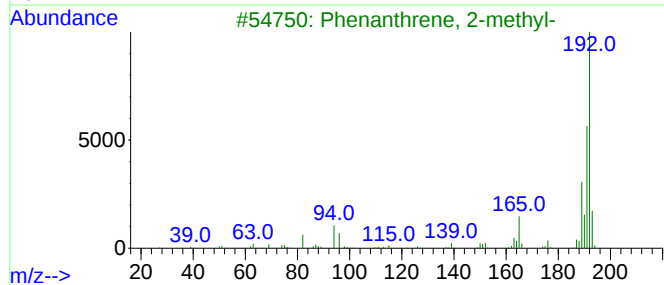
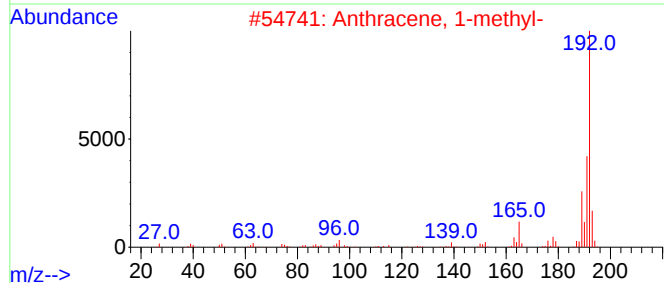
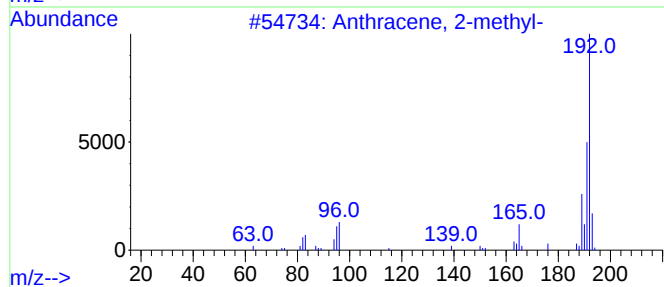
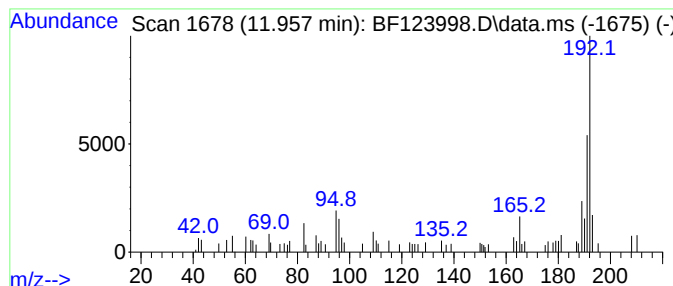
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Anthracene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.957	3.11 ng	63630	Phenanthrene-d10	11.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
5			Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	93



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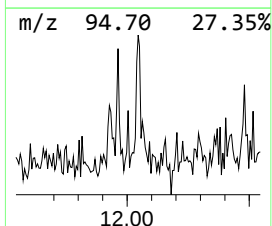
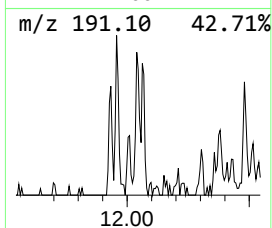
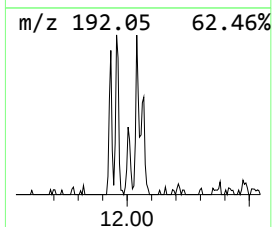
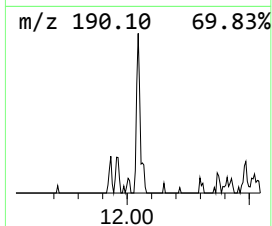
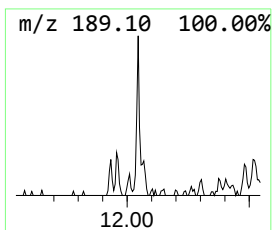
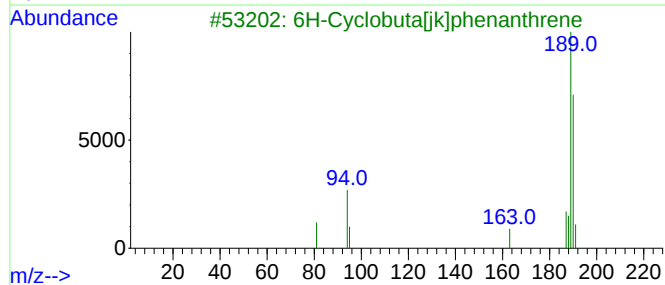
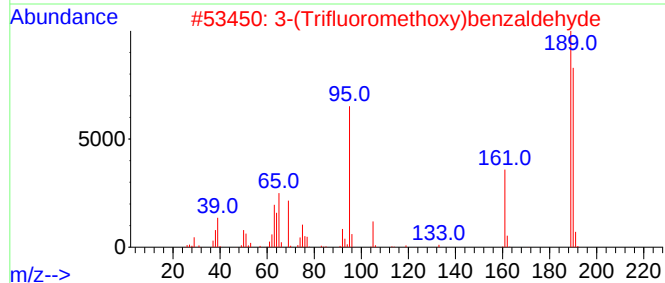
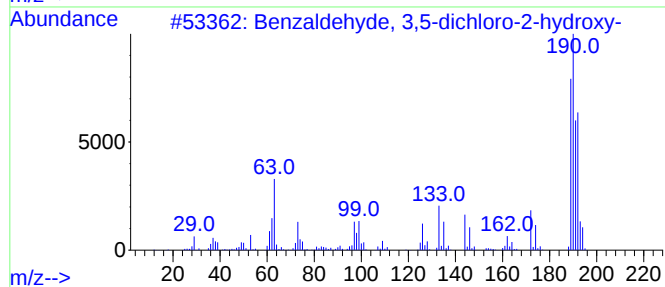
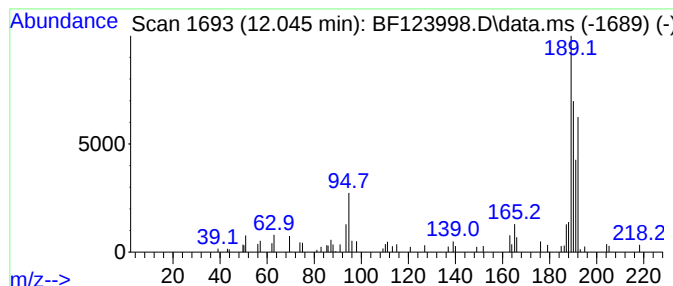
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown12.045 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.045	3.63 ng	74408	Phenanthrene-d10	11.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	49
2			3-(Trifluoromethoxy)benzaldehyde	190	C8H5F3O2	052771-21-8	43
3			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	43
4			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	38
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	30



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ALS Vial : 16 Sample Multiplier: 1

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ClientSampleId :
TP-2

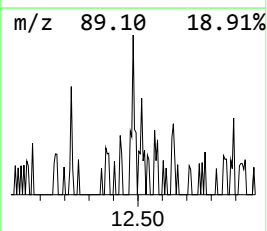
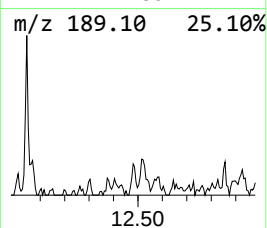
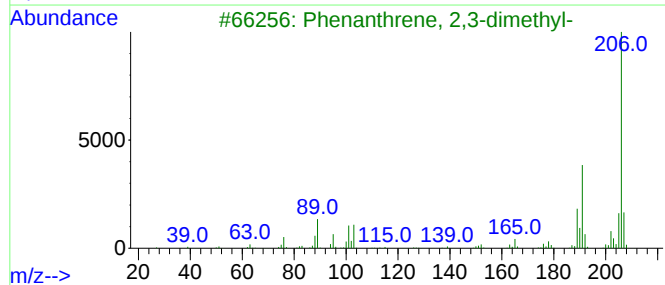
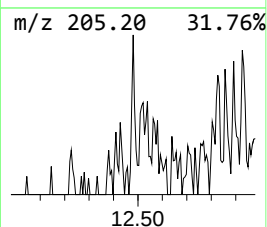
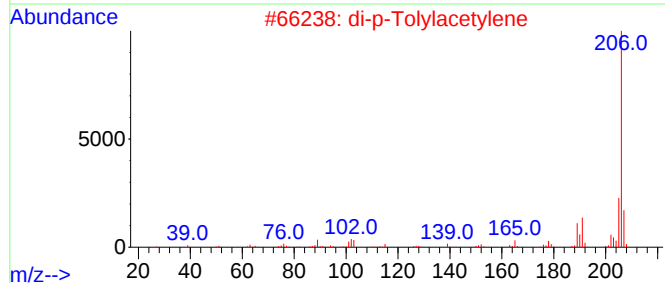
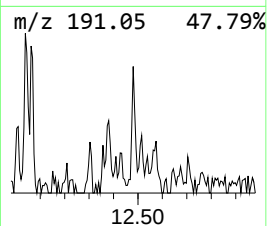
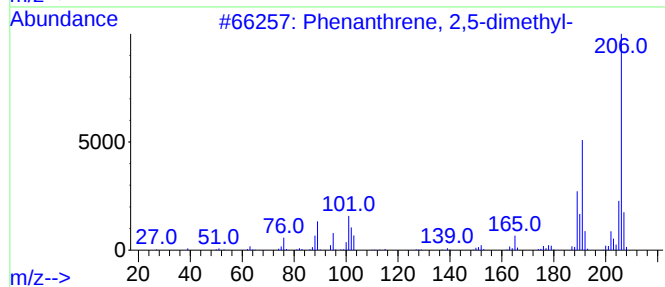
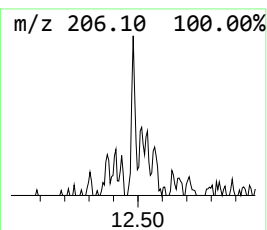
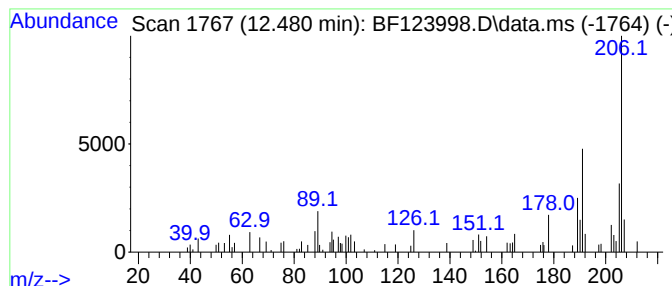
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Phenanthrene, 2,5-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.480	2.34 ng	47892	Phenanthrene-d10	11.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	91
2			di-p-Tolylacetylene	206	C16H14	002789-88-0	87
3			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	83
4			[14]Annulene, 1,6:8,13-bis(metha...	206	C16H14	085385-68-8	80
5			3-Methyl-1-phenyl-1H-indene	206	C16H14	022360-63-0	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
Data File : BF123998.D
Acq On : 19 May 2021 19:30
Operator : JU/CG
Sample : M2437-05 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

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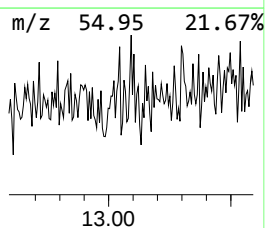
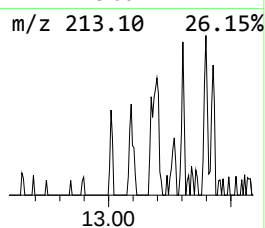
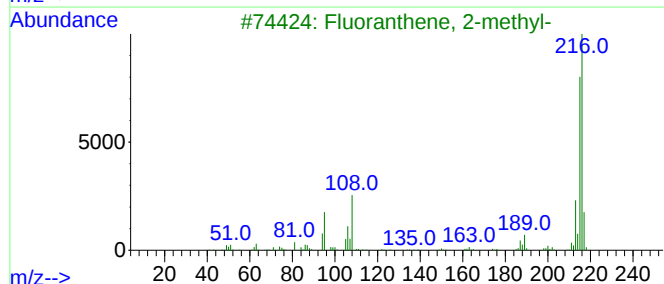
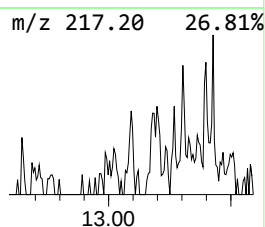
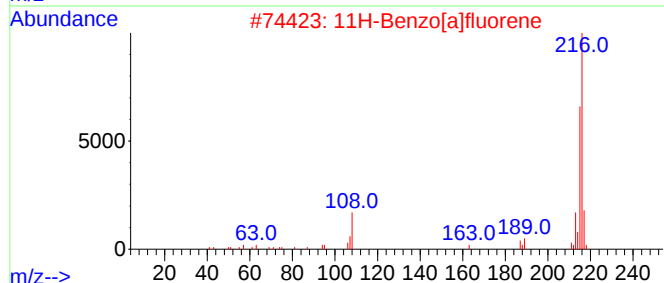
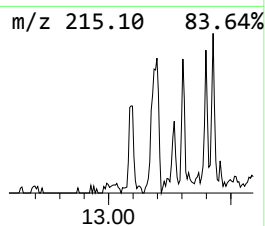
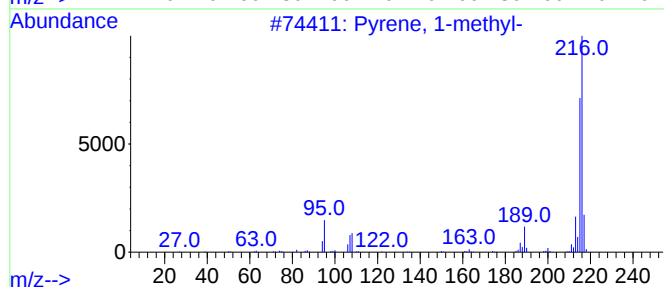
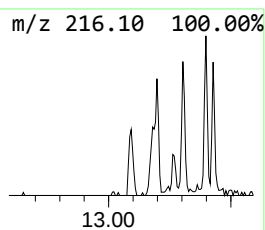
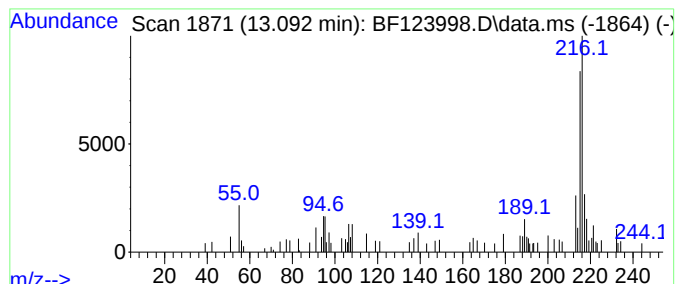
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Pyrene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.092	2.66 ng	78413	Chrysene-d12	14.074

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	68
2			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	64
3			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	53
4			3-Pyridinamine, 6-chloro-N-(phen...	216	C12H9ClN2	005325-71-3	53
5			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
Data File : BF123998.D
Acq On : 19 May 2021 19:30
Operator : JU/CG
Sample : M2437-05 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
TP-2

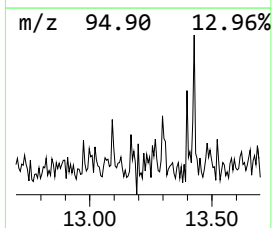
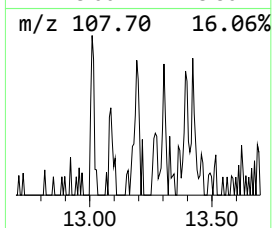
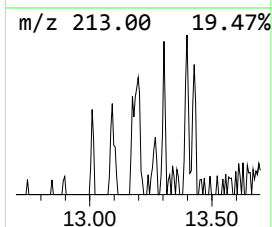
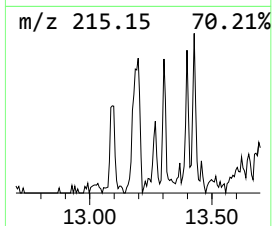
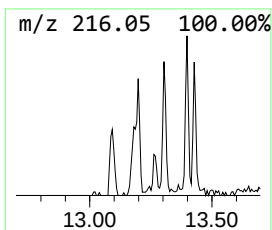
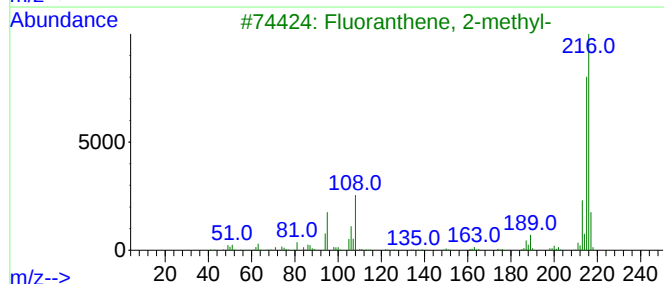
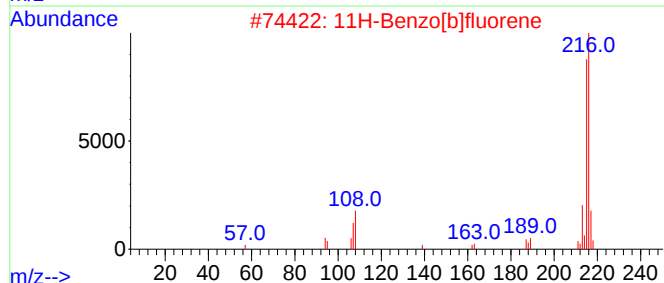
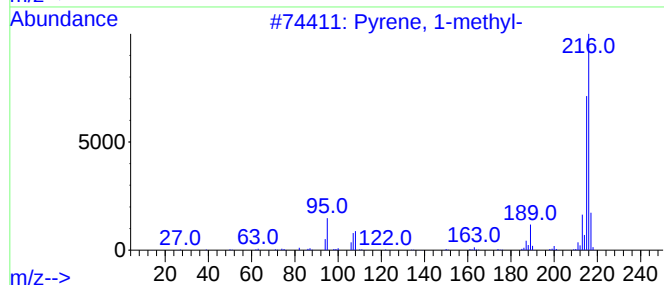
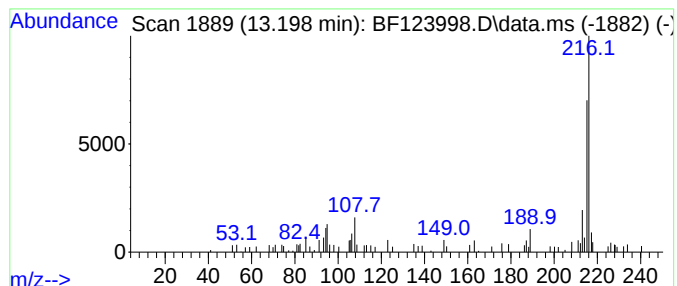
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 11H-Benzo[b]fluorene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.198	3.45 ng	101461	Chrysene-d12	14.074

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	89
3			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	89
4			Pyrene, 2-methyl-	216	C17H12	003442-78-2	83
5			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
Data File : BF123998.D
Acq On : 19 May 2021 19:30
Operator : JU/CG
Sample : M2437-05 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
TP-2

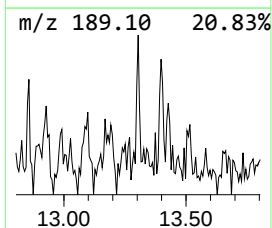
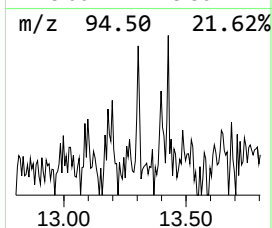
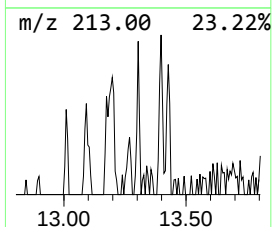
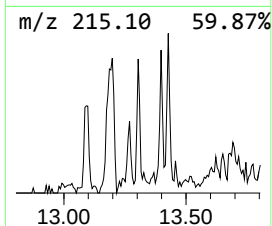
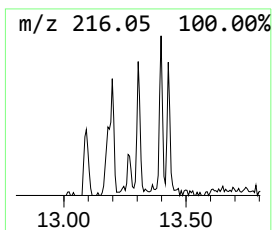
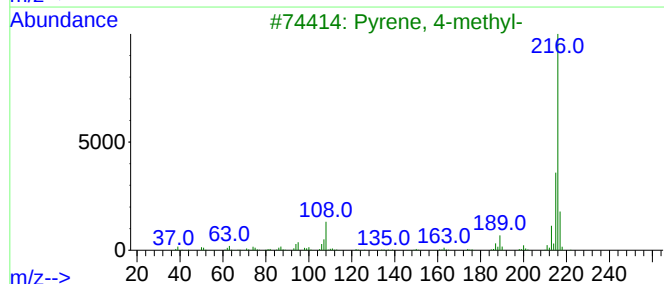
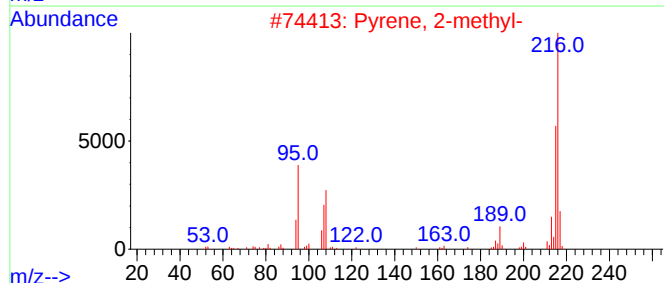
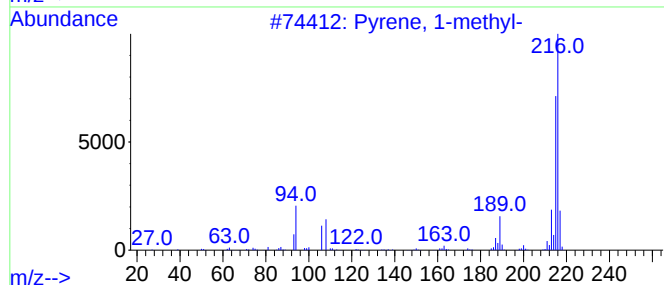
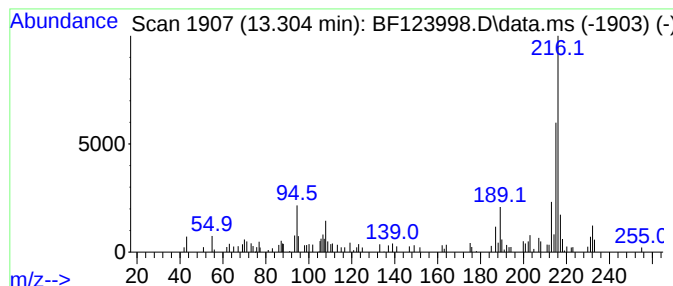
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Pyrene, 2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.304	2.10 ng	61907	Chrysene-d12	14.074

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
2			Pyrene, 2-methyl-	216	C17H12	003442-78-2	81
3			Pyrene, 4-methyl-	216	C17H12	003353-12-6	64
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	64
5			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
Data File : BF123998.D
Acq On : 19 May 2021 19:30
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Sample : M2437-05 5X
Misc :
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Instrument :
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ClientSampled :
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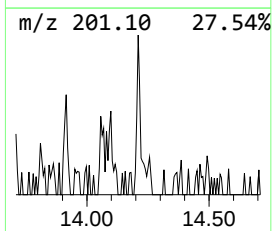
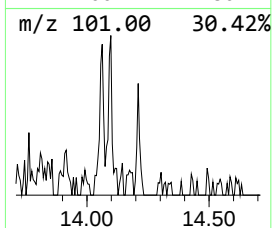
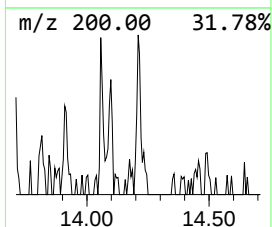
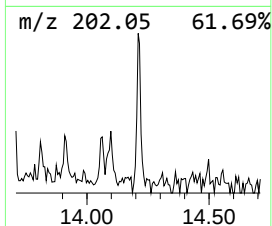
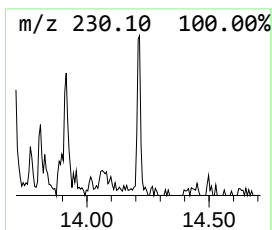
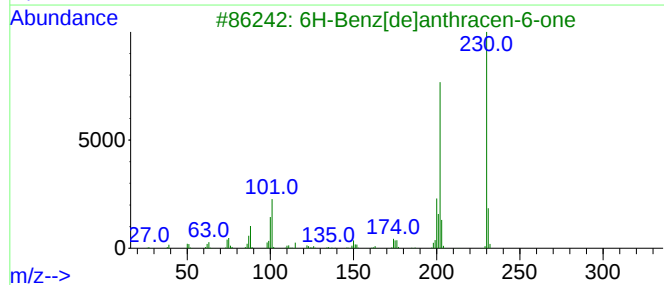
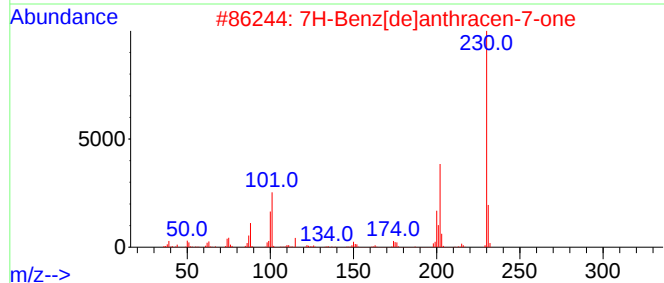
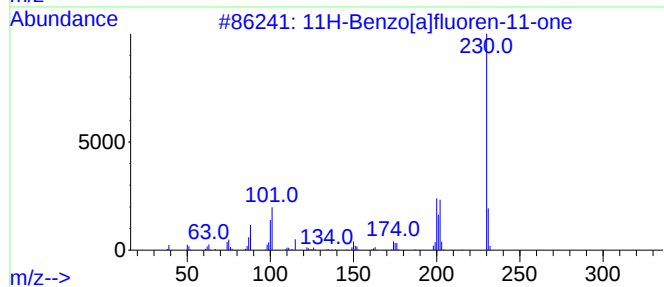
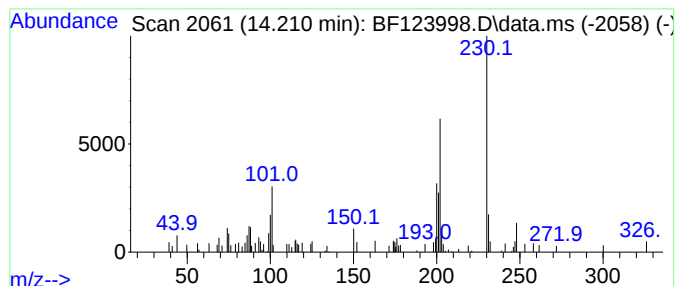
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 11H-Benzo[a]fluoren-11-one Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.210	3.01 ng	88510	Chrysene-d12	14.074

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	95
2			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	95
3			6H-Benz[de]anthracen-6-one	230	C17H10O	080252-14-8	70
4			4-Pyrenecarbaldehyde	230	C17H10O	022245-51-8	68
5			1-Pyrenecarboxaldehyde	230	C17H10O	003029-19-4	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
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Acq On : 19 May 2021 19:30
Operator : JU/CG
Sample : M2437-05 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

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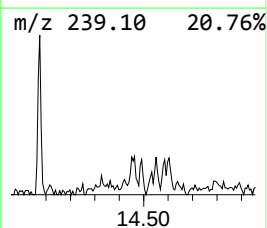
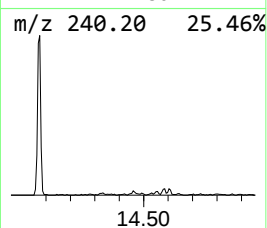
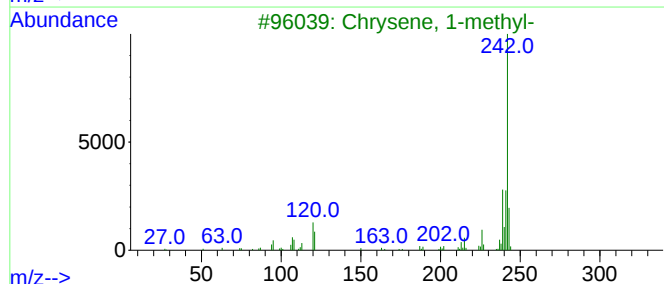
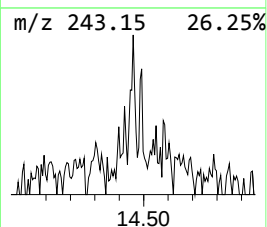
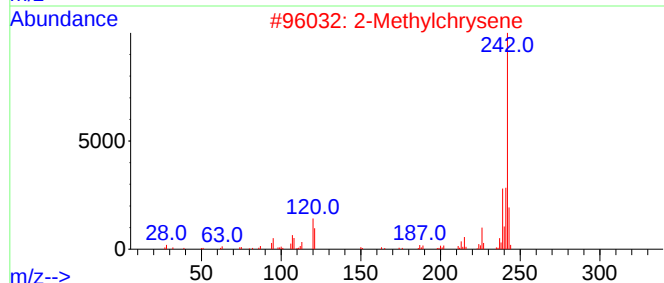
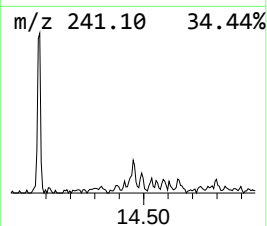
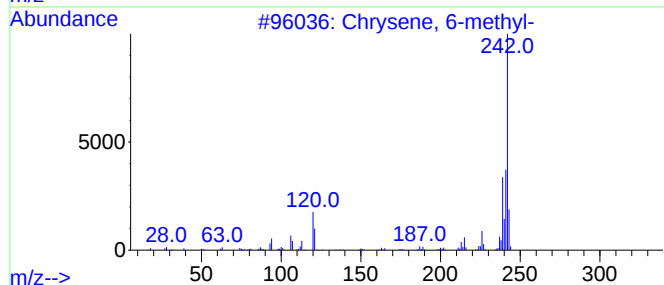
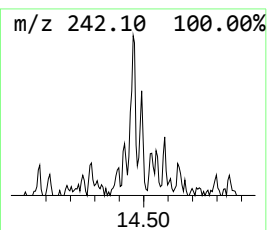
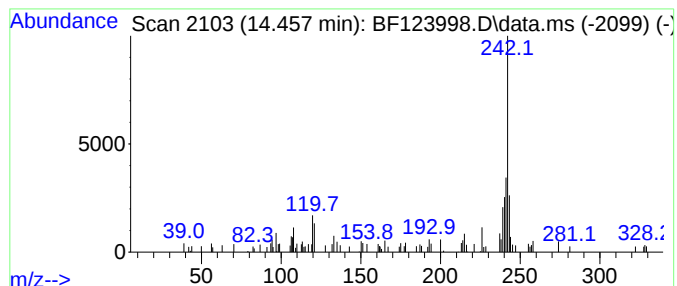
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Chrysene, 6-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.457	2.39 ng	70233	Chrysene-d12	14.074

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Chrysene, 6-methyl-	242	C19H14	001705-85-7	90
2			2-Methylchrysene	242	C19H14	003351-32-4	76
3			Chrysene, 1-methyl-	242	C19H14	003351-28-8	68
4			Benz[a]anthracene, 8-methyl-	242	C19H14	002381-31-9	64
5			Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
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Acq On : 19 May 2021 19:30
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Misc :
ALS Vial : 16 Sample Multiplier: 1

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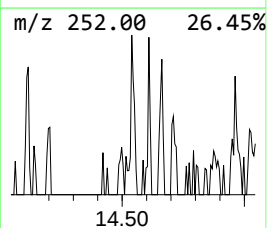
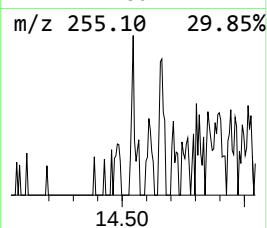
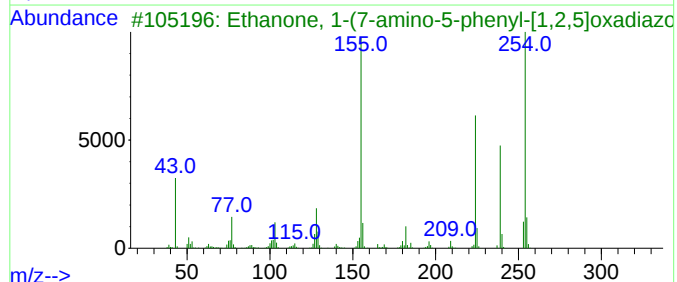
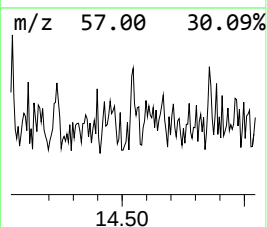
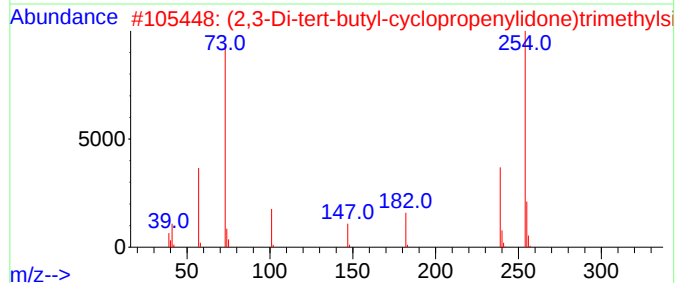
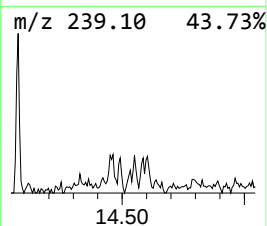
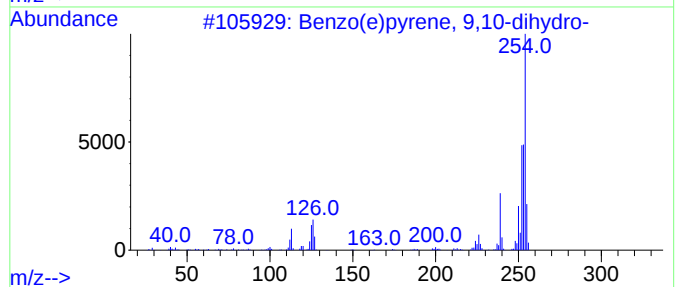
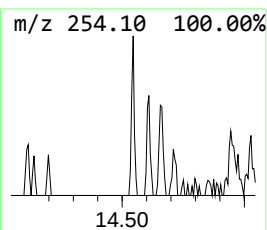
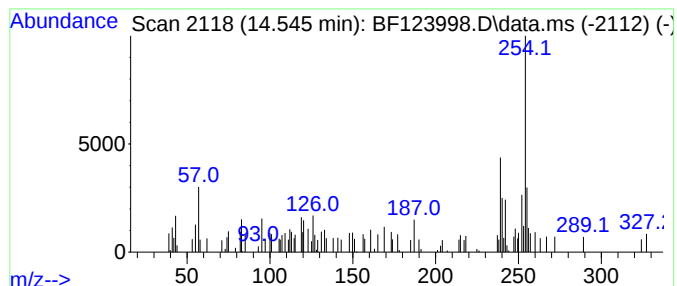
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Benzo(e)pyrene, 9,10-dihydro- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.545	2.27 ng	66732	Chrysene-d12	14.074

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo(e)pyrene, 9,10-dihydro-	254	C20H14	066788-01-0	53
2			(2,3-Di-tert-butyl-cyclopropenyl...	254	C14H27PSi	1000306-16-6	47
3			Ethanone, 1-(7-amino-5-phenyl-[1...	254	C13H10N4O2	1000316-41-1	43
4			Fluorene, 9-(2-pyridinyl)methylene-	255	C19H13N	002871-27-4	38
5			Benzo(a)pyrene, 7,8-dihydro-	254	C20H14	017573-23-8	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
Data File : BF123998.D
Acq On : 19 May 2021 19:30
Operator : JU/CG
Sample : M2437-05 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-2

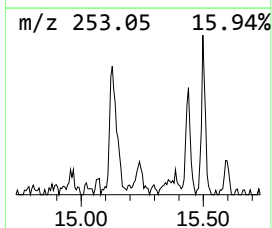
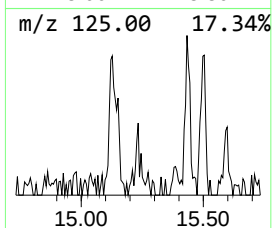
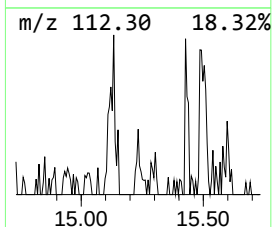
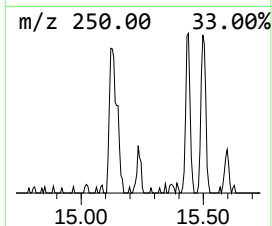
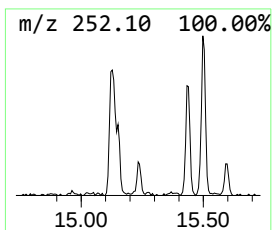
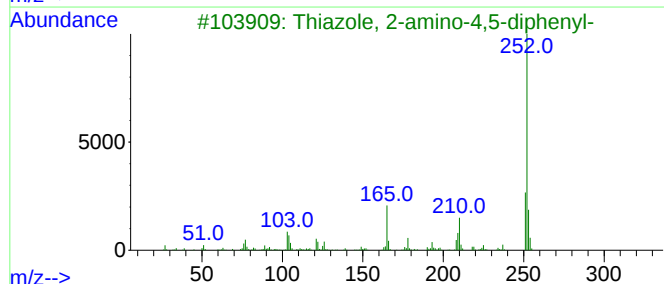
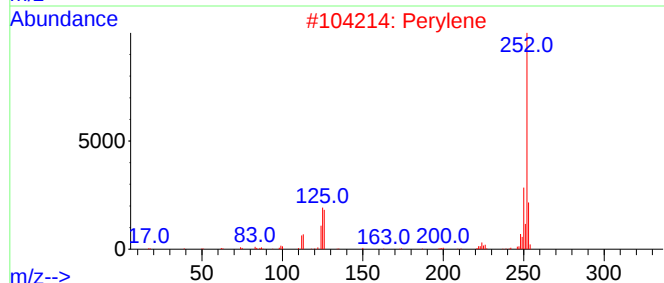
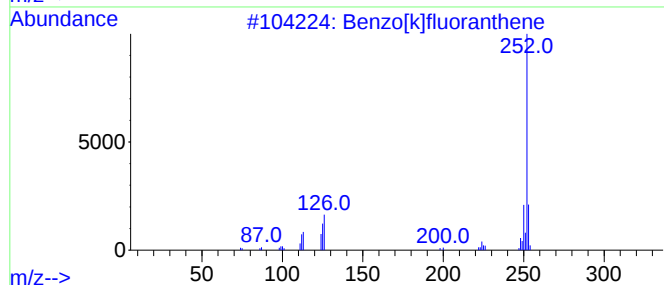
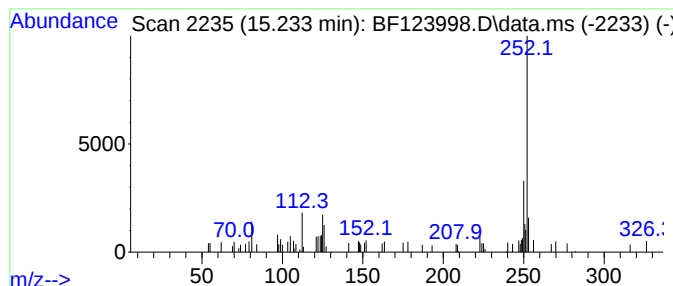
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Perylene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.233	3.16 ng	38464	Perylene-d12	15.562

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[k]fluoranthene	252	C20H12	000207-08-9	53
2		Perylene	252	C20H12	000198-55-0	53
3		Thiazole, 2-amino-4,5-diphenyl-	252	C15H12N2S	006318-74-7	52
4		3H-Pyrrolo[3,2-f]quinoline, 1,2,...	252	C17H20N2	1000338-03-5	52
5		1H-1,3-Benzimidazole, 2-tricyclo...	252	C17H20N2	1000319-80-9	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
Data File : BF123998.D
Acq On : 19 May 2021 19:30
Operator : JU/CG
Sample : M2437-05 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-2

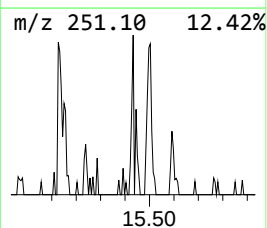
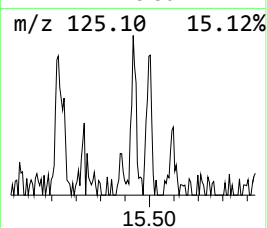
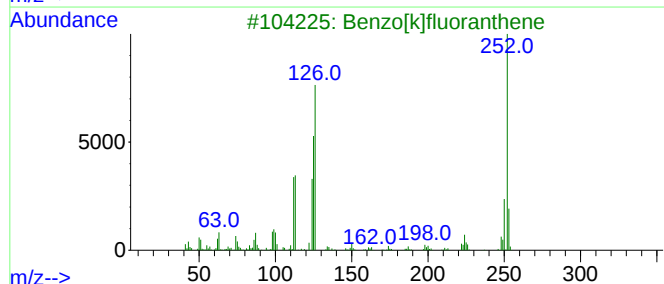
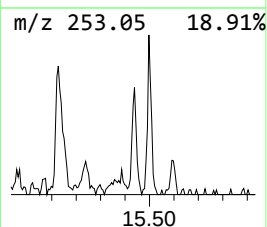
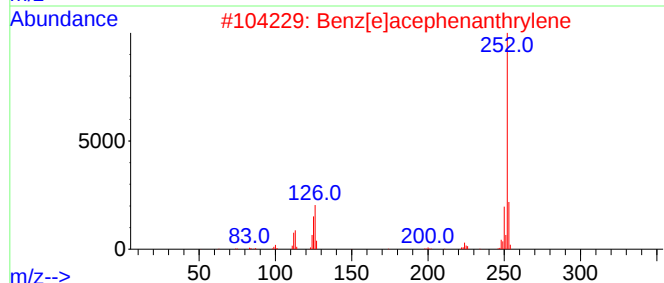
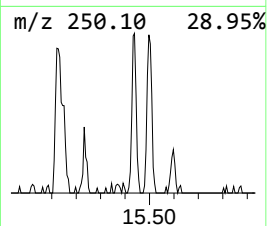
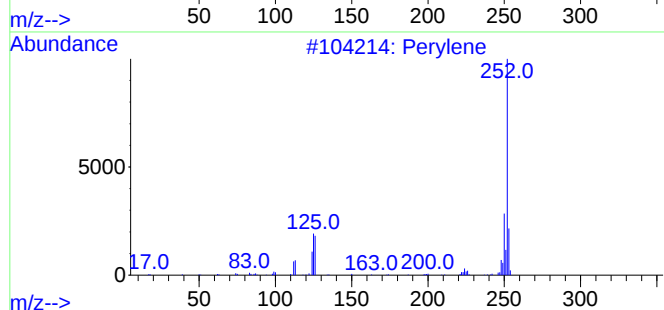
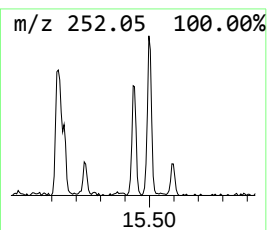
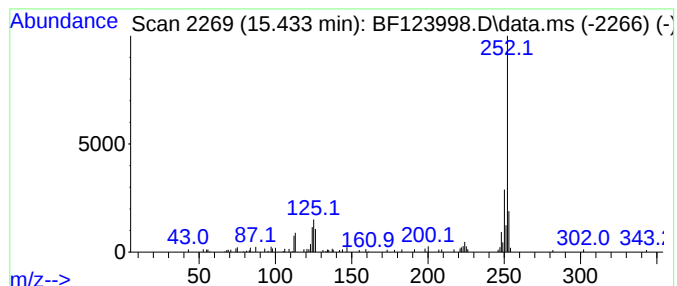
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Benzo[j]fluoranthene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.433	9.32 ng	113320	Perylene-d12	15.562

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Perylene	252	C20H12	000198-55-0	90
2			Benz[e]acephenanthrylene	252	C20H12	000205-99-2	90
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	86
4			Benzo[j]fluoranthene	252	C20H12	000205-82-3	86
5			Benzo[e]pyrene	252	C20H12	000192-97-2	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
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Sample : M2437-05 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-2

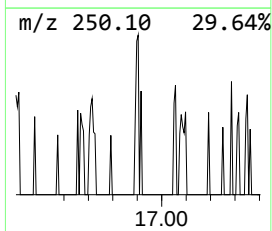
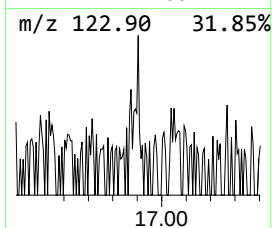
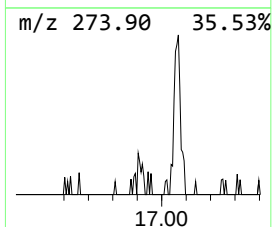
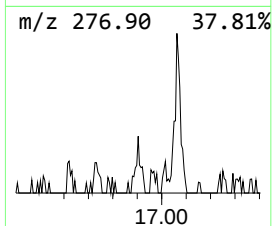
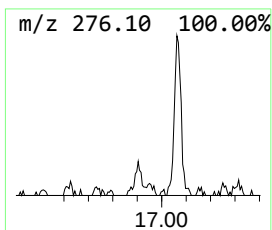
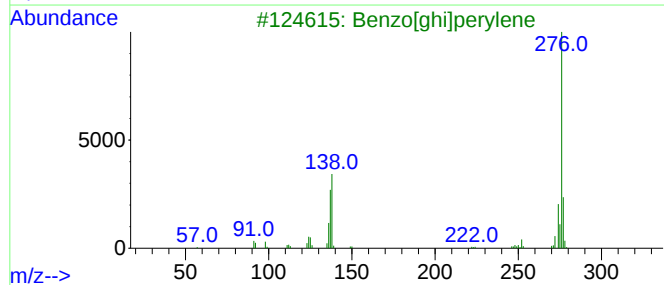
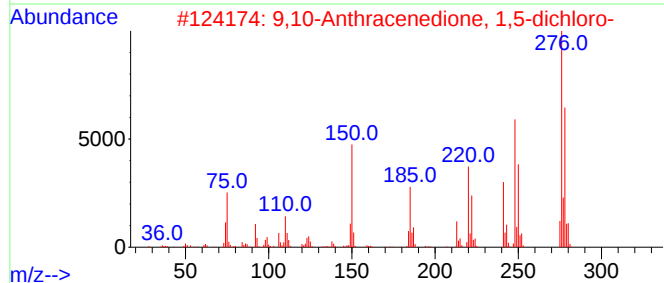
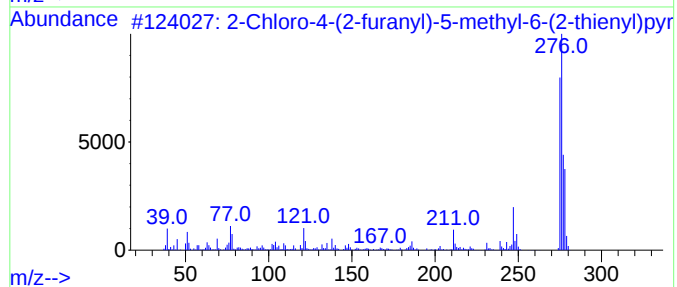
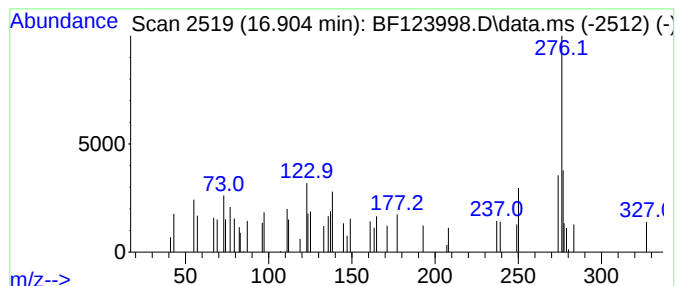
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 unknown16.904 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.904	3.18 ng	38664	Perylene-d12	15.562

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Chloro-4-(2-furanyl)-5-methyl-...	276	C13H9ClN2O5	131022-80-5	43
2		9,10-Anthracenedione, 1,5-dichloro-	276	C14H6Cl2O2	000082-46-2	32
3		Benzo[ghi]perylene	276	C22H12	000191-24-2	30
4		Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	27
5		Indeno[1,2,3-cd]fluoranthene	276	C22H12	000193-43-1	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051921\
Data File : BF123998.D
Acq On : 19 May 2021 19:30
Operator : JU/CG
Sample : M2437-05 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.152	7.8	ng	119081	1	6.904	305569	20.0
Anthracene, 2-m...	11.957	3.1	ng	63630	4	11.427	409760	20.0
unknown12.045	12.045	3.6	ng	74408	4	11.427	409760	20.0
Phenanthrene, 2...	12.480	2.3	ng	47892	4	11.427	409760	20.0
Pyrene, 1-methyl-	13.092	2.7	ng	78413	5	14.074	588534	20.0
11H-Benzo[b]flu...	13.198	3.5	ng	101461	5	14.074	588534	20.0
Pyrene, 2-methyl-	13.304	2.1	ng	61907	5	14.074	588534	20.0
11H-Benzo[a]flu...	14.210	3.0	ng	88510	5	14.074	588534	20.0
Chrysene, 6-met...	14.457	2.4	ng	70233	5	14.074	588534	20.0
Benzo(e)pyrene,...	14.545	2.3	ng	66732	5	14.074	588534	20.0
Perylene	15.233	3.2	ng	38464	6	15.562	243249	20.0
Benzo[j]fluoran...	15.433	9.3	ng	113320	6	15.562	243249	20.0
unknown16.904	16.904	3.2	ng	38664	6	15.562	243249	20.0