

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101317\
 Data File : BF099442.D
 Acq On : 13 Oct 2017 16:30
 Operator : SJ/JU
 Sample : I5780-05 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DUCT-BANK-1-8-COMP

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.063	504	506	519	rBV	25058	31903	2.95%	0.233%
2	5.469	572	575	588	rBV	199313	205694	19.00%	1.501%
3	6.481	744	747	757	rBV	198851	212346	19.62%	1.549%
4	6.622	768	771	777	rBV	289259	227494	21.02%	1.660%
5	6.851	806	810	817	rBV	694879	554859	51.26%	4.048%
6	7.004	833	836	844	rVB	203469	179421	16.58%	1.309%
7	7.416	902	906	913	rBV	124354	125714	11.61%	0.917%
8	8.033	1007	1011	1015	rBB	21163	16946	1.57%	0.124%
9	8.133	1024	1028	1030	rBV	902341	711112	65.70%	5.188%
10	8.151	1030	1031	1035	rVB	110083	86746	8.01%	0.633%
11	8.845	1146	1149	1157	rBV	59899	50978	4.71%	0.372%
12	8.945	1162	1166	1173	rVV	38758	32166	2.97%	0.235%
13	9.204	1207	1210	1218	rBV	268348	201057	18.58%	1.467%
14	9.475	1250	1256	1259	rBV2	12741	16167	1.49%	0.118%
15	9.545	1265	1268	1271	rBV	24495	22489	2.08%	0.164%
16	9.598	1275	1277	1285	rVB	33858	28011	2.59%	0.204%
17	9.751	1299	1303	1307	rBV2	22251	21065	1.95%	0.154%
18	9.886	1320	1326	1330	rBV2	779787	683107	63.11%	4.984%
19	9.922	1330	1332	1335	rVB	82255	64642	5.97%	0.472%
20	10.092	1357	1361	1364	rBV2	43094	40658	3.76%	0.297%
21	10.433	1416	1419	1425	rBV	65320	60592	5.60%	0.442%
22	10.592	1444	1446	1451	rVB2	22837	23218	2.15%	0.169%
23	10.674	1457	1460	1464	rBV	101837	94930	8.77%	0.693%
24	10.792	1475	1480	1488	rBV6	12566	24360	2.25%	0.178%
25	10.986	1506	1513	1515	rBV3	16757	21711	2.01%	0.158%
26	11.110	1529	1534	1536	rBV4	10540	15813	1.46%	0.115%
27	11.133	1536	1538	1542	rVV	14537	14624	1.35%	0.107%
28	11.186	1542	1547	1550	rVV2	22488	24107	2.23%	0.176%
29	11.227	1550	1554	1557	rVV3	18371	26008	2.40%	0.190%
30	11.274	1560	1562	1566	rVB	32428	25942	2.40%	0.189%
31	11.374	1575	1579	1581	rBV	660403	552895	51.08%	4.034%
32	11.398	1581	1583	1587	rVV	640897	511501	47.26%	3.732%
33	11.451	1589	1592	1595	rVV	152727	126401	11.68%	0.922%
34	11.486	1595	1598	1603	rVB3	11261	16726	1.55%	0.122%

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35	11.610	1615	1619	1623	rBV2	52822	57924	5.35%	0.423%
36	11.710	1632	1636	1641	rVB3	15585	20877	1.93%	0.152%
37	11.874	1659	1664	1667	rBV	83988	82960	7.66%	0.605%
38	11.904	1667	1669	1675	rVV	125877	100073	9.25%	0.730%
39	11.951	1675	1677	1680	rVV	41886	34952	3.23%	0.255%
40	11.986	1680	1683	1686	rVV2	122891	141803	13.10%	1.035%
41	12.010	1686	1687	1692	rVB	62110	40486	3.74%	0.295%
42	12.168	1711	1714	1717	rBV	55544	47996	4.43%	0.350%
43	12.204	1717	1720	1724	rVB	46667	41081	3.80%	0.300%
44	12.351	1743	1745	1747	rVB	20847	15021	1.39%	0.110%
45	12.374	1747	1749	1752	rVB	20532	15332	1.42%	0.112%
46	12.421	1754	1757	1760	rBV	59121	57102	5.28%	0.417%
47	12.463	1760	1764	1766	rVV4	29971	41474	3.83%	0.303%
48	12.515	1769	1773	1777	rVB3	42367	46881	4.33%	0.342%
49	12.592	1782	1786	1789	rBV	863836	759244	70.15%	5.539%
50	12.651	1794	1796	1799	rVB3	22893	17924	1.66%	0.131%
51	12.739	1804	1811	1814	rBV4	29933	43613	4.03%	0.318%
52	12.821	1818	1825	1830	rBV	757157	844185	77.99%	6.159%
53	12.868	1830	1833	1837	rVB2	41708	46668	4.31%	0.340%
54	12.951	1841	1847	1850	rBV2	167733	174529	16.12%	1.273%
55	12.980	1850	1852	1856	rVB	41849	36649	3.39%	0.267%
56	13.033	1856	1861	1865	rBV3	48551	85475	7.90%	0.624%
57	13.121	1873	1876	1877	rBV2	46366	48262	4.46%	0.352%
58	13.145	1877	1880	1883	rVB	106672	102899	9.51%	0.751%
59	13.210	1888	1891	1894	rVV	74577	61482	5.68%	0.449%
60	13.251	1894	1898	1900	rVV2	76633	86346	7.98%	0.630%
61	13.268	1900	1901	1905	rVB	20980	19951	1.84%	0.146%
62	13.345	1910	1914	1916	rBV	40492	39212	3.62%	0.286%
63	13.374	1916	1919	1922	rBV	43819	44562	4.12%	0.325%
64	13.627	1959	1962	1964	rBV4	20811	22175	2.05%	0.162%
65	13.657	1964	1967	1971	rVB	75545	66135	6.11%	0.483%
66	13.768	1981	1986	1988	rBV	78263	92536	8.55%	0.675%
67	13.792	1988	1990	1992	rVV	77584	66260	6.12%	0.483%
68	13.821	1992	1995	2000	rVV4	58744	113490	10.49%	0.828%
69	13.862	2000	2002	2006	rVB	71780	68547	6.33%	0.500%
70	13.933	2012	2014	2019	rVB2	18079	22056	2.04%	0.161%
71	14.015	2023	2028	2031	rVV2	777345	1082379	100.00%	7.897%

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72	14.045	2031	2033	2040	rVV	385880	351064	32.43%	2.561%
73	14.121	2044	2046	2049	rVV	74693	61310	5.66%	0.447%
74	14.168	2051	2054	2056	rVV4	27954	38689	3.57%	0.282%
75	14.209	2059	2061	2063	rVV	34200	33483	3.09%	0.244%
76	14.245	2063	2067	2070	rVV3	48629	67178	6.21%	0.490%
77	14.274	2070	2072	2075	rVB4	17414	16483	1.52%	0.120%
78	14.333	2080	2082	2085	rBV3	18032	20304	1.88%	0.148%
79	14.362	2085	2087	2089	rBV	35858	33941	3.14%	0.248%
80	14.404	2089	2094	2097	rVB	81806	85117	7.86%	0.621%
81	14.433	2097	2099	2103	rBV	38785	40615	3.75%	0.296%
82	14.527	2113	2115	2117	rBV2	42786	35595	3.29%	0.260%
83	14.545	2117	2118	2122	rVB	36279	32607	3.01%	0.238%
84	14.721	2145	2148	2151	rBV4	33691	47705	4.41%	0.348%
85	14.898	2175	2178	2182	rVB2	41564	45288	4.18%	0.330%
86	15.004	2194	2196	2199	rVB	19024	15672	1.45%	0.114%
87	15.056	2201	2205	2209	rBV	354232	587254	54.26%	4.285%
88	15.168	2221	2224	2228	rBV	74820	79870	7.38%	0.583%
89	15.256	2236	2239	2240	rBV2	27367	28413	2.63%	0.207%
90	15.362	2253	2257	2264	rVB	227570	301699	27.87%	2.201%
91	15.427	2264	2268	2273	rBV	334470	407433	37.64%	2.973%
92	15.492	2273	2279	2282	rBV	497798	639661	59.10%	4.667%
93	15.521	2282	2284	2290	rVB2	93057	117362	10.84%	0.856%
94	15.909	2348	2350	2353	rBV3	17740	17083	1.58%	0.125%
95	16.051	2372	2374	2380	rVB4	19982	23873	2.21%	0.174%
96	16.403	2432	2434	2443	rVB6	27624	50355	4.65%	0.367%
97	16.974	2524	2531	2541	rBV2	162263	355228	32.82%	2.592%
98	17.145	2556	2560	2565	rBV	32397	58775	5.43%	0.429%
99	17.421	2601	2607	2619	rVB	118697	259345	23.96%	1.892%
100	17.674	2646	2650	2662	rVB6	41201	114975	10.62%	0.839%

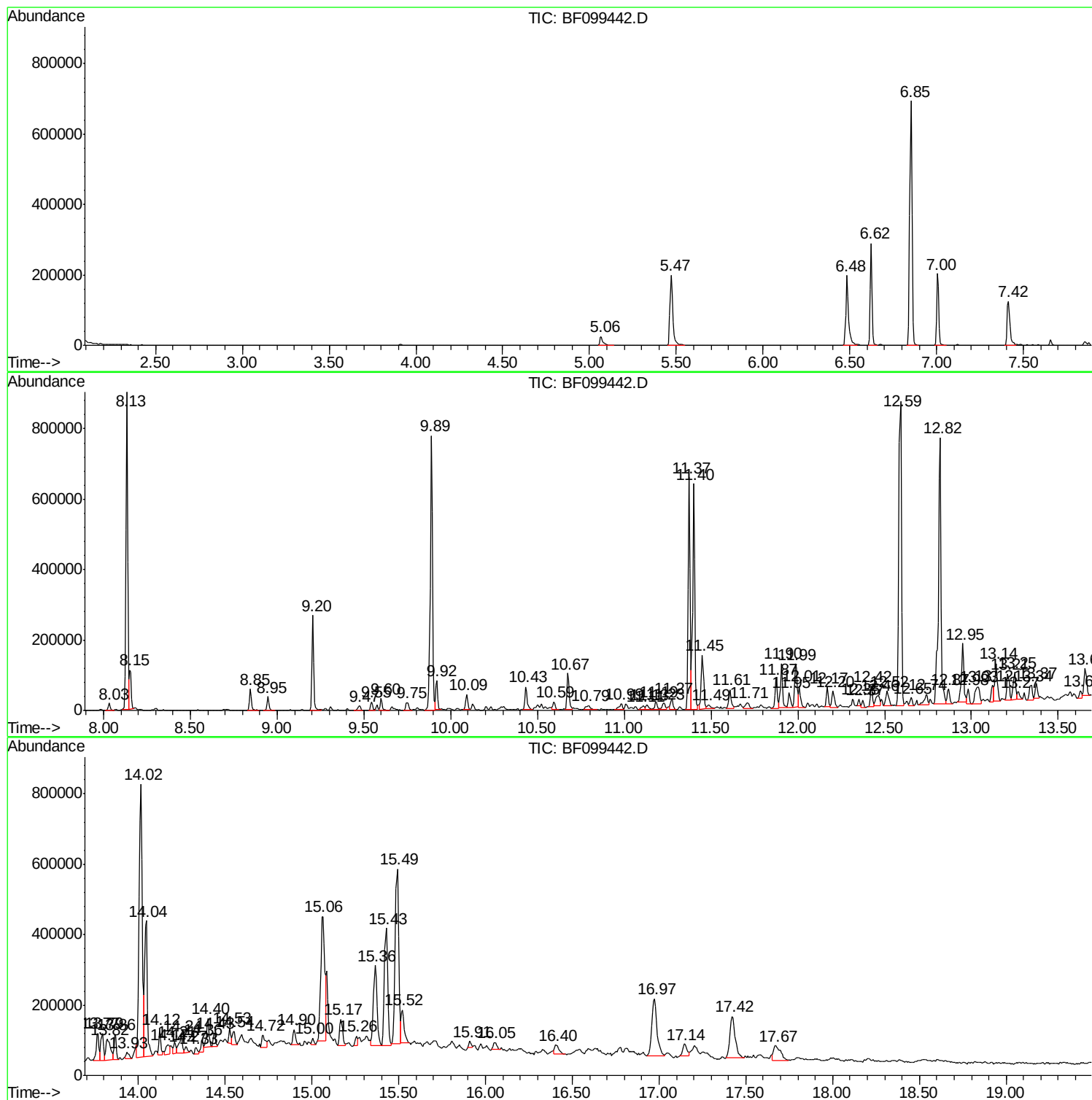
Sum of corrected areas: 13706321

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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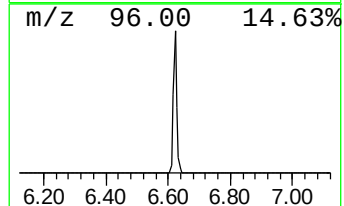
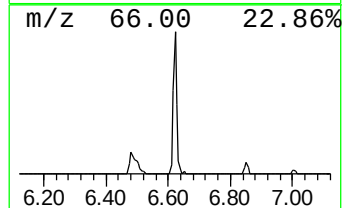
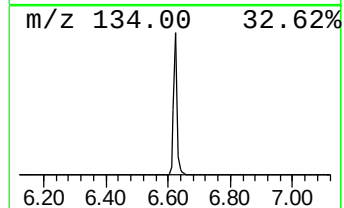
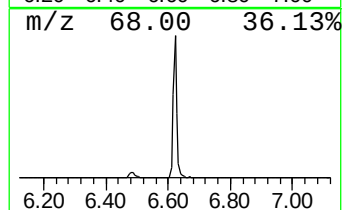
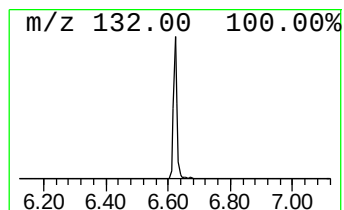
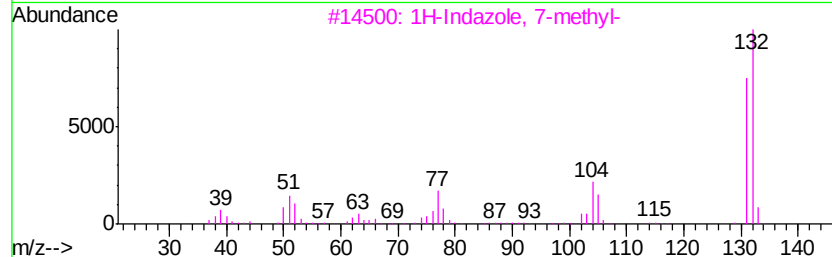
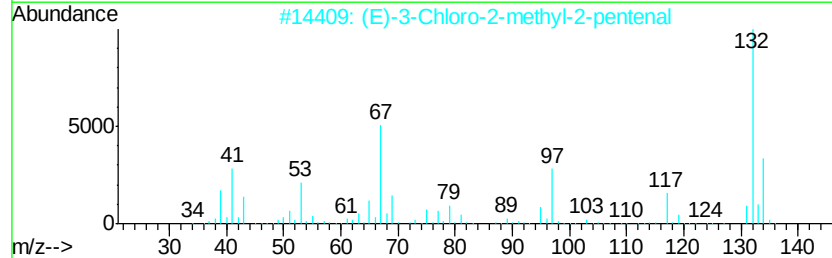
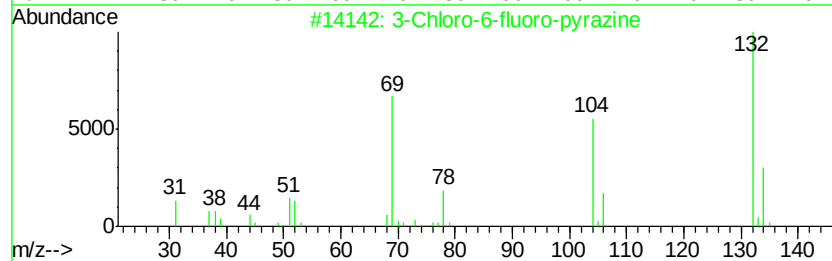
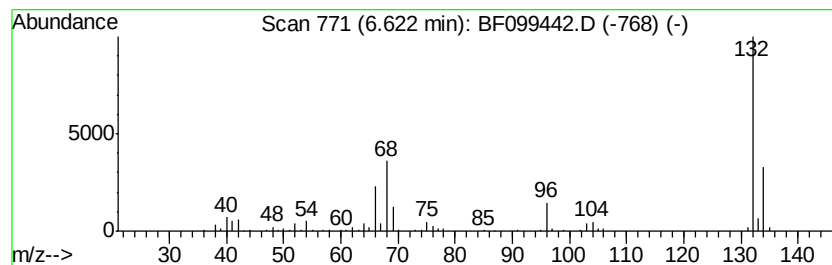
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.62 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	8.20 ng	227494	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	10
4		5-Aminoindole	132	C8H8N2	005192-03-0	10
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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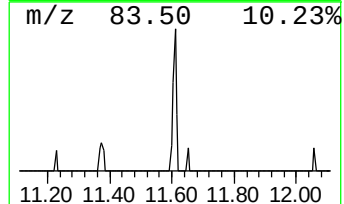
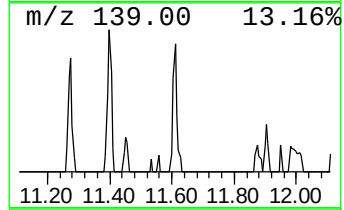
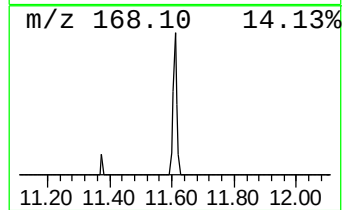
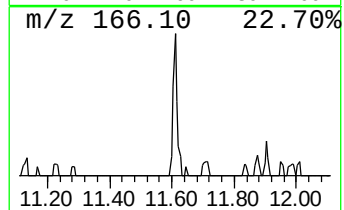
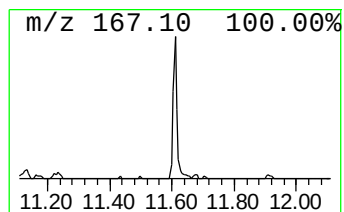
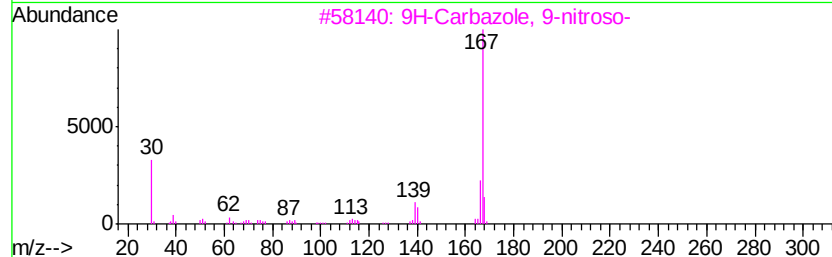
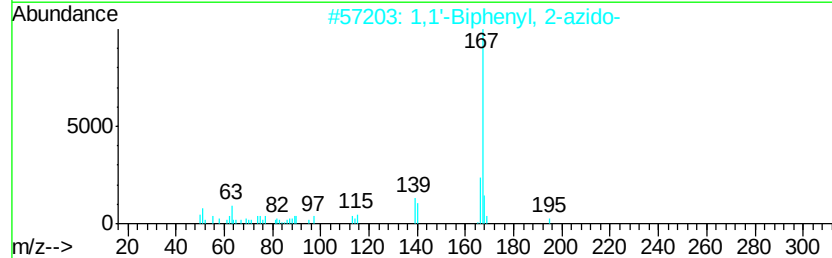
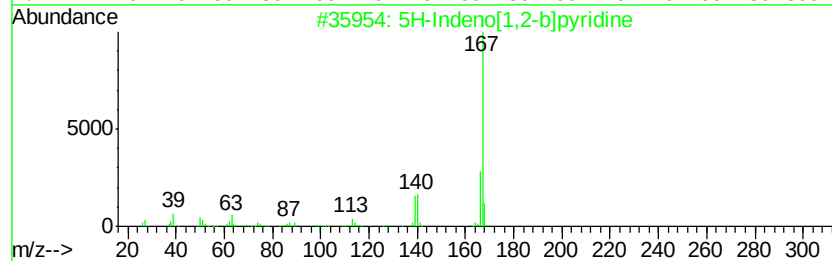
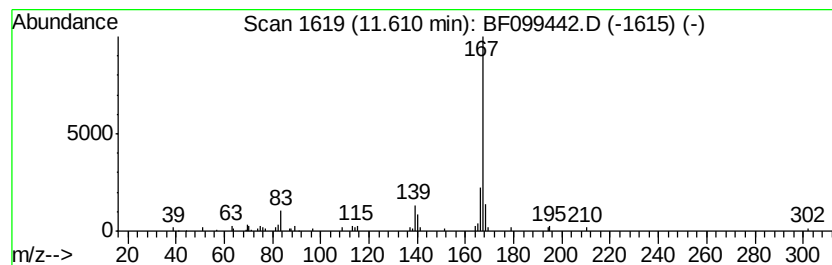
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 5H-Indeno[1,2-b]pyridine Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.61	2.10 ng	57924	Phenanthrene-d10	11.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	90
2		1,1'-Biphenyl, 2-azido-	195	C12H9N3	007599-23-7	90
3		9H-Carbazole, 9-nitroso-	196	C12H8N2O	002788-23-0	90
4		Carbazole	167	C12H9N	000086-74-8	87
5		D-Galactitol, 2-(acetylmethylami...	363	C16H29NO8	074410-42-7	80



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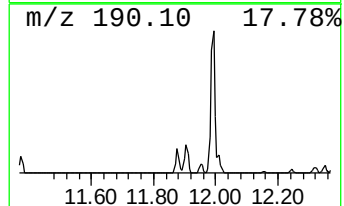
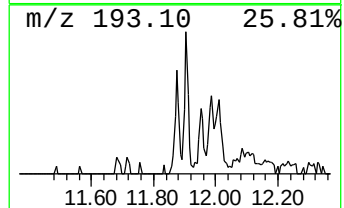
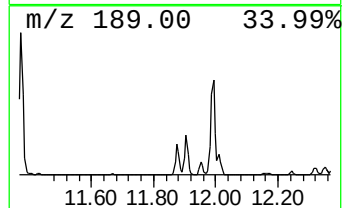
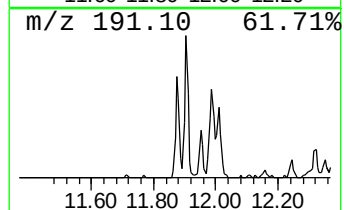
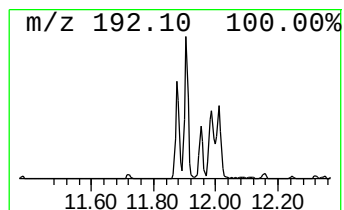
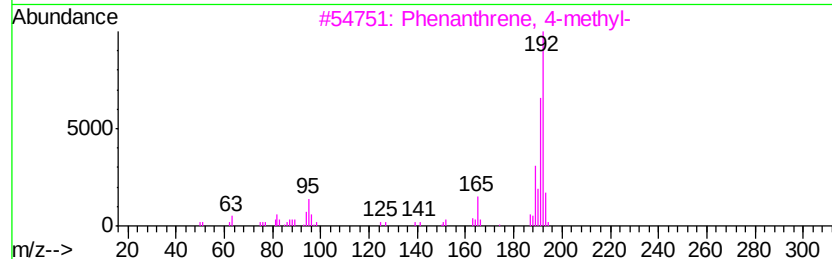
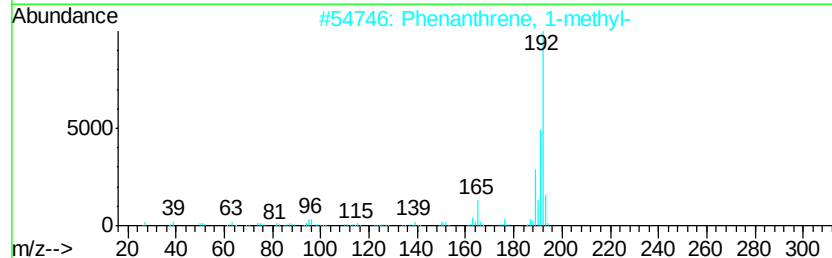
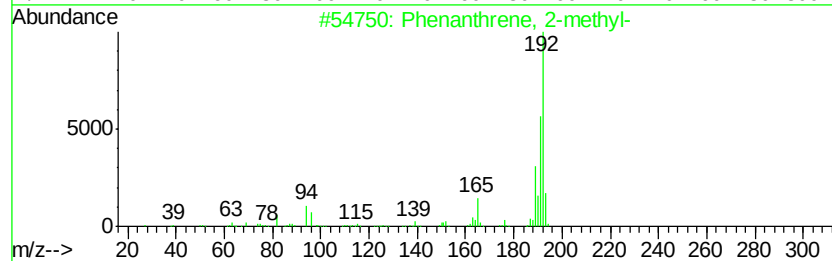
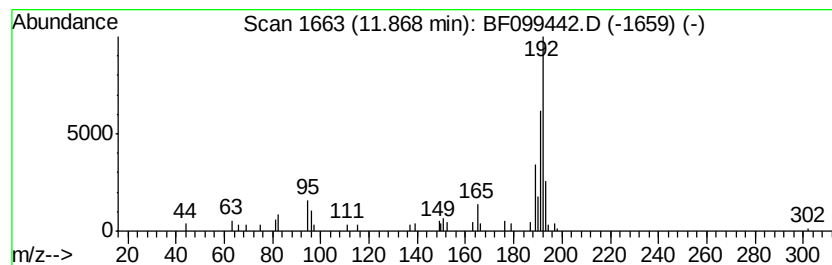
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Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.87	3.00 ng	82960	Phenanthrene-d10		11.37
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
3	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	94
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101317\
Data File : BF099442.D
Acq On : 13 Oct 2017 16:30
Operator : SJ/JU
Sample : I5780-05 5X
Misc :
ALS Vial : 5 Sample Multiplier: 1

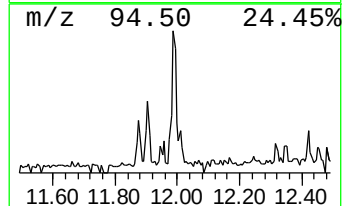
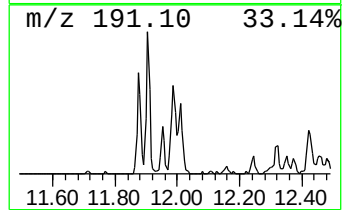
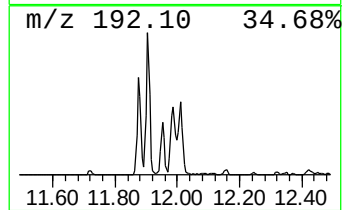
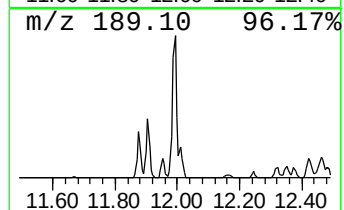
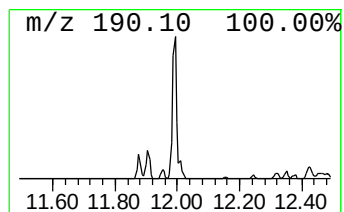
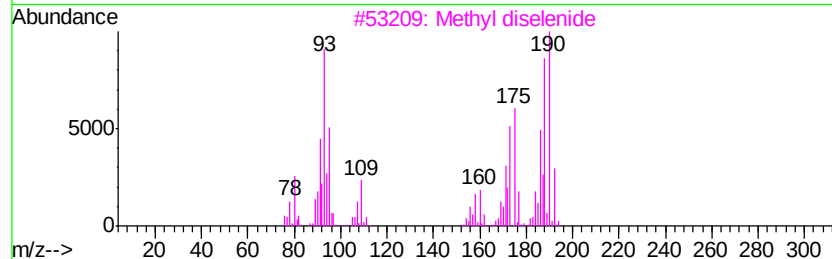
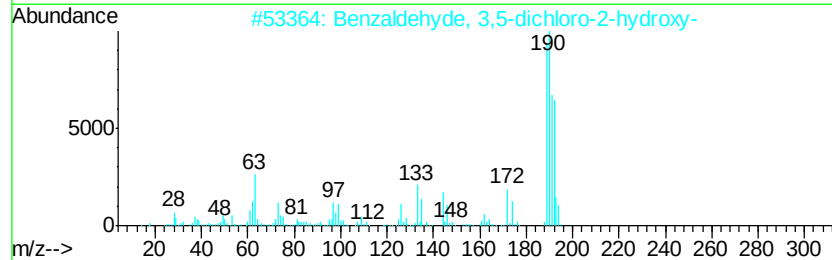
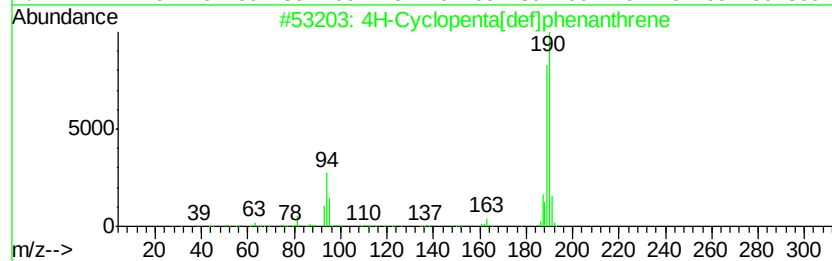
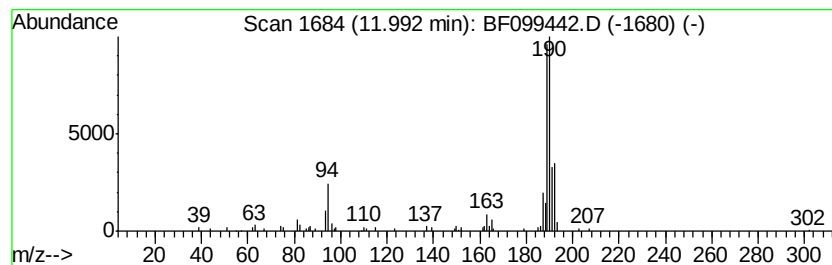
Instrument :
BNA_F
ClientSampleId :
DUCT-BANK-1-8-COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.99	5.13 ng	141803	Phenanthrene-d10	11.37		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
2		Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	59
3		Methyl diselenide	190	C2H6Se2	007101-31-7	38
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	30
5		8,9-Dihydrocyclopenta[def]phenan...	192	C15H12	027410-55-5	20



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101317\
Data File : BF099442.D
Acq On : 13 Oct 2017 16:30
Operator : SJ/JU
Sample : I5780-05 5X
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DUCT-BANK-1-8-COMP

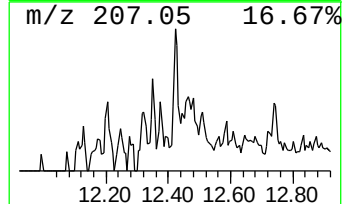
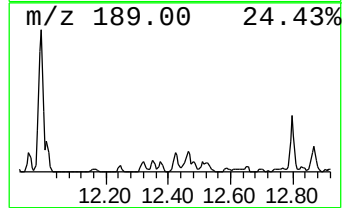
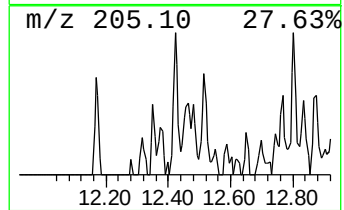
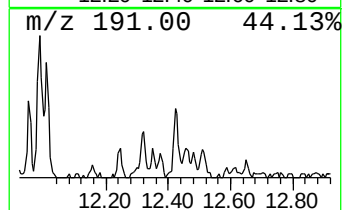
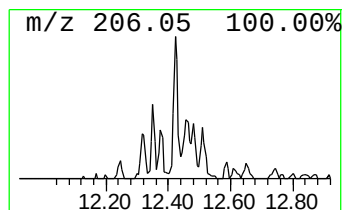
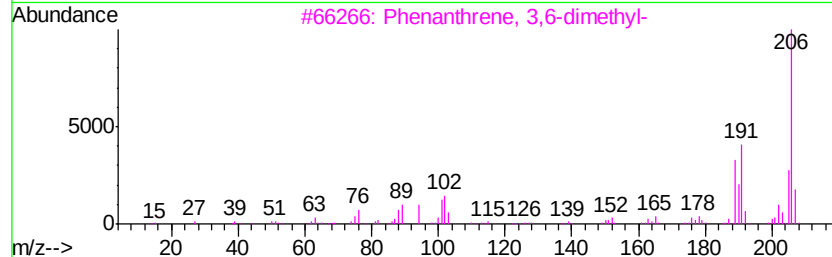
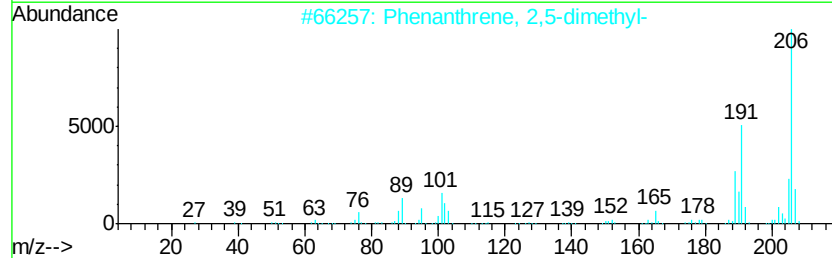
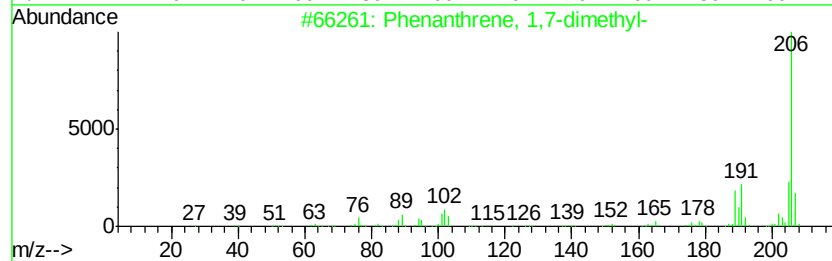
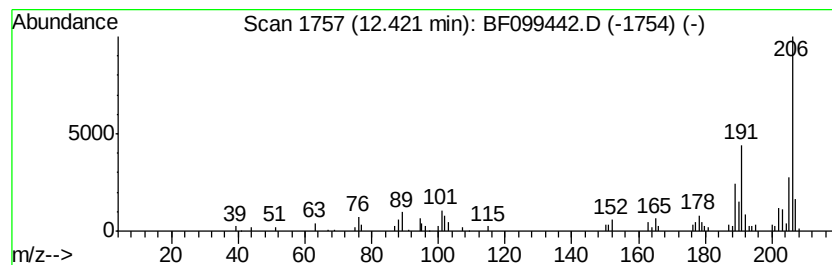
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenanthrene, 1,7-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.42	2.07 ng	57102	Phenanthrene-d10	11.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	96
2		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	95
3		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	94
4		9,10-Dimethylantracene	206	C16H14	000781-43-1	93
5		Naphthalene, 1,2-dihydro-1-phenyl-	206	C16H14	016606-46-5	92



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101317\
Data File : BF099442.D
Acq On : 13 Oct 2017 16:30
Operator : SJ/JU
Sample : I5780-05 5X
Misc :
ALS Vial : 5 Sample Multiplier: 1

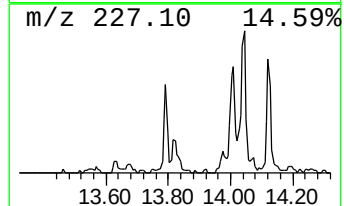
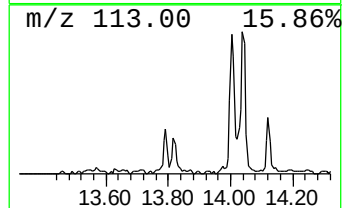
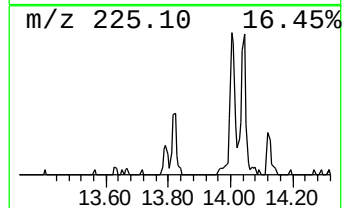
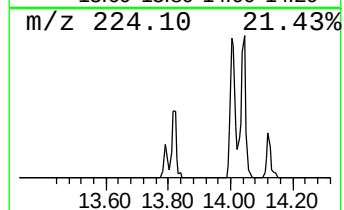
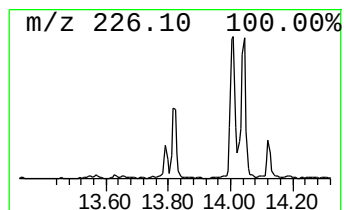
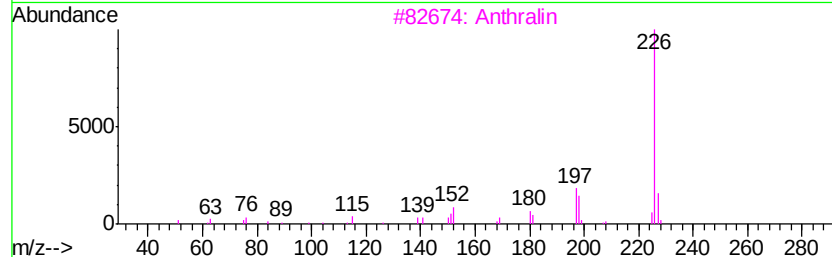
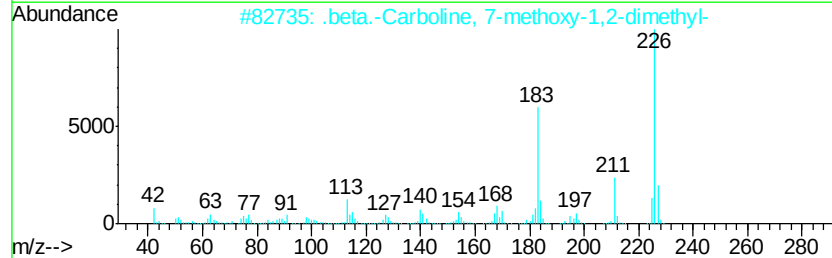
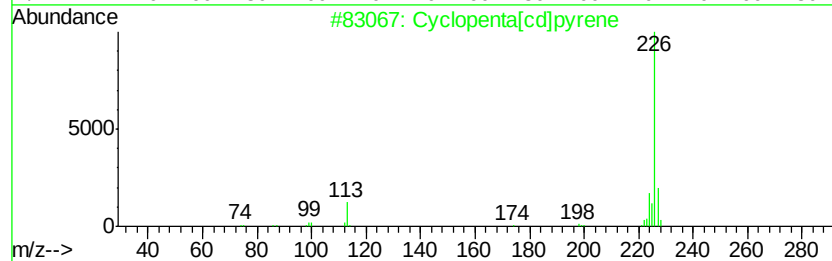
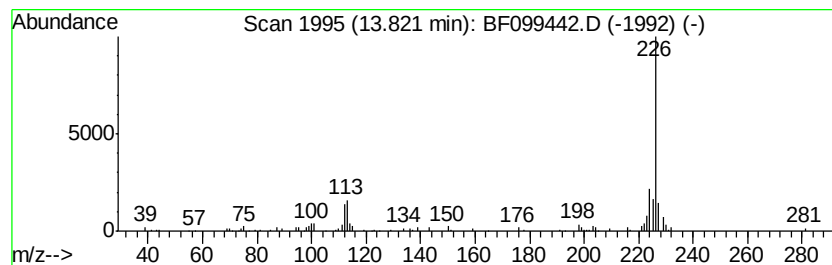
Instrument :
BNA_F
ClientSampleId :
DUCT-BANK-1-8-COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Cyclopenta[cd]pyrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.82	2.10 ng	113490	Chrysene-d12	14.02		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	58
2		.beta.-Carboline, 7-methoxy-1,2-...	226	C14H14N2O	006519-18-2	53
3		Anthralin	226	C14H10O3	001143-38-0	47
4		4-Naphthalen-2-yl-thiazol-2-ylamine	226	C13H10N2S	1000300-53-4	42
5		1,3-Diamino-5,6-dihydro-9-methyl...	226	C13H14N4	037436-39-8	42



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101317\
Data File : BF099442.D
Acq On : 13 Oct 2017 16:30
Operator : SJ/JU
Sample : I5780-05 5X
Misc :
ALS Vial : 5 Sample Multiplier: 1

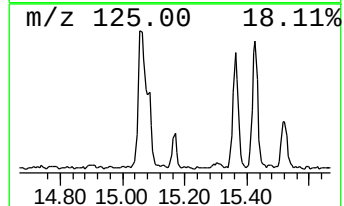
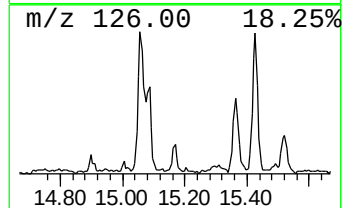
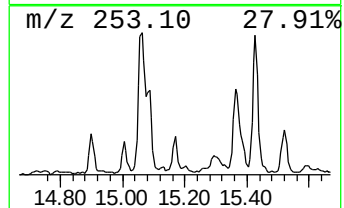
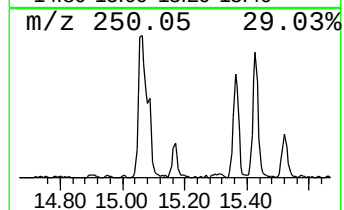
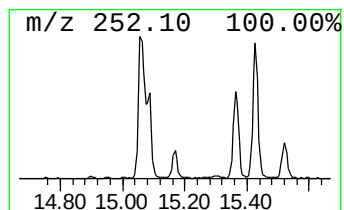
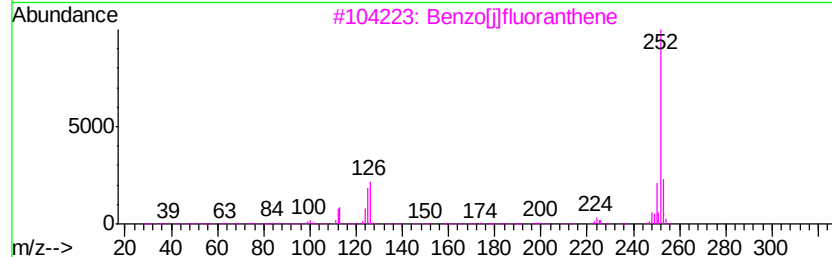
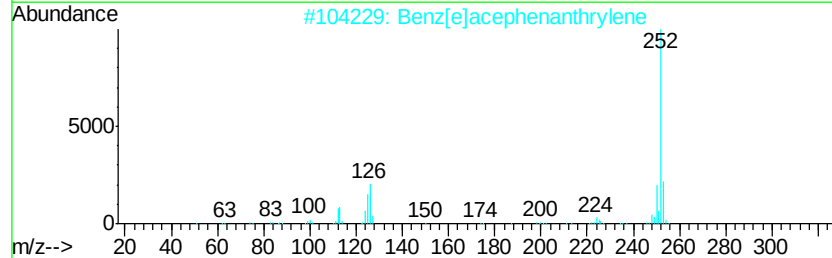
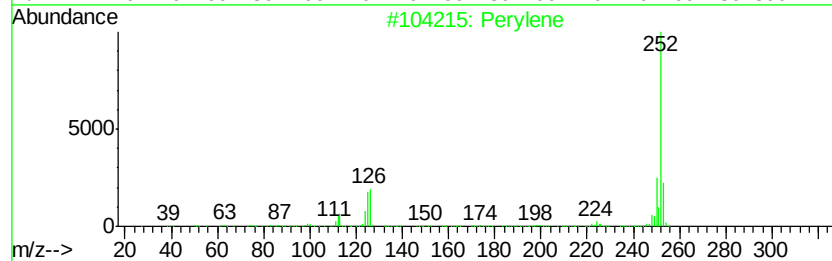
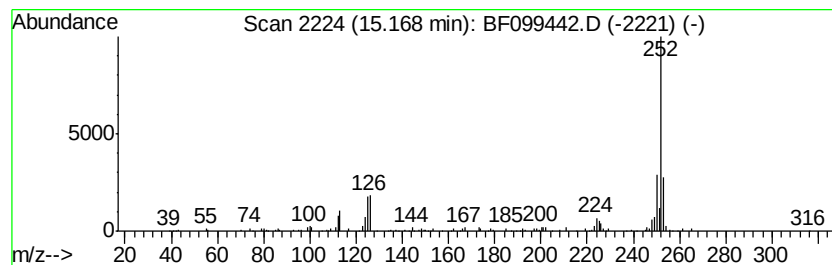
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Perylene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.17	2.50 ng	79870	Perylene-d12	15.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Perylene	252 C20H12	000198-55-0	96
2	Benz[e]acephenanthrylene	252 C20H12	000205-99-2	95
3	Benzo[j]fluoranthene	252 C20H12	000205-82-3	94
4	Benzo[e]pyrene	252 C20H12	000192-97-2	94
5	Benzo[k]fluoranthene	252 C20H12	000207-08-9	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101317\
Data File : BF099442.D
Acq On : 13 Oct 2017 16:30
Operator : SJ/JU
Sample : I5780-05 5X
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DUCT-BANK-1-8-COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.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
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5H-Indeno[1,2-b]p...	11.61	2.1	ng	57924	4	11.37	552895	20.0
Phenanthrene, 2-m...	11.87	3.0	ng	82960	4	11.37	552895	20.0
4H-Cyclopenta[def...	11.99	5.1	ng	141803	4	11.37	552895	20.0
Phenanthrene, 1,7...	12.42	2.1	ng	57102	4	11.37	552895	20.0
Cyclopenta[cd]pyrene	13.82	2.1	ng	113490	5	14.02	1082380	20.0
Perylene	15.17	2.5	ng	79870	6	15.49	639661	20.0