

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
 Data File : BF103767.D
 Acq On : 19 Mar 2018 20:30
 Operator : JU/SJ
 Sample : J1971-05
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ES-6-SA-3-WW@4.5

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:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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	2.334	35	42	51	rVB	169584	227518	5.22%	0.457%
2	5.228	530	534	547	rVB	167743	231451	5.31%	0.465%
3	5.604	592	598	601	rBV	3819388	4245742	97.39%	8.532%
4	5.922	648	652	660	rVB	64437	60703	1.39%	0.122%
5	6.598	760	767	770	rBV	3492212	4359389	100.00%	8.760%
6	6.745	787	792	795	rBV	4012940	4271138	97.98%	8.583%
7	6.975	827	831	834	rVB	1091119	908167	20.83%	1.825%
8	7.133	853	858	861	rBV	3364781	3224307	73.96%	6.479%
9	7.539	921	927	930	rBV	2555976	2729084	62.60%	5.484%
10	7.598	932	937	945	rVB	38546	43890	1.01%	0.088%
11	8.257	1045	1049	1051	rBV	1297234	1245242	28.56%	2.502%
12	8.275	1051	1052	1055	rVB	334290	185535	4.26%	0.373%
13	8.963	1166	1169	1173	rBV	216028	185470	4.25%	0.373%
14	9.063	1182	1186	1190	rBV	194728	174101	3.99%	0.350%
15	9.327	1225	1231	1234	rBV	4008929	3940598	90.39%	7.919%
16	9.427	1245	1248	1251	rBV	58078	45467	1.04%	0.091%
17	9.598	1272	1277	1280	rBV2	51197	68727	1.58%	0.138%
18	9.669	1284	1289	1291	rBV	108829	109033	2.50%	0.219%
19	9.692	1291	1293	1295	rVV	55661	51270	1.18%	0.103%
20	9.716	1295	1297	1300	rVB	413766	293491	6.73%	0.590%
21	9.786	1305	1309	1310	rBV2	58064	52180	1.20%	0.105%
22	9.927	1330	1333	1336	rVB	134959	97188	2.23%	0.195%
23	10.016	1344	1348	1351	rVV	1263953	1161032	26.63%	2.333%
24	10.045	1351	1353	1356	rVB	185993	129476	2.97%	0.260%
25	10.216	1379	1382	1385	rVB2	88801	80180	1.84%	0.161%
26	10.251	1385	1388	1392	rBV2	55267	46330	1.06%	0.093%
27	10.404	1411	1414	1416	rBV	48987	52584	1.21%	0.106%
28	10.427	1416	1418	1421	rVB	167531	136288	3.13%	0.274%
29	10.563	1437	1441	1444	rVB	168602	149981	3.44%	0.301%
30	10.645	1453	1455	1458	rVB2	68150	65742	1.51%	0.132%
31	10.804	1478	1482	1486	rBV	2231704	2327518	53.39%	4.677%
32	10.904	1496	1499	1508	rVB	140794	161912	3.71%	0.325%
33	11.110	1528	1534	1536	rBV4	53347	81583	1.87%	0.164%
34	11.139	1536	1539	1542	rVV3	40102	48823	1.12%	0.098%

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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-BF031518.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.310	1564	1568	1571	rVB	57194	52530	1.20%	0.106%
36	11.351	1571	1575	1578	rBV2	104102	101140	2.32%	0.203%
37	11.404	1582	1584	1587	rVB	102963	83034	1.90%	0.167%
38	11.504	1597	1601	1603	rBV	1079761	1023445	23.48%	2.057%
39	11.533	1603	1606	1609	rVV	1477160	1264036	29.00%	2.540%
40	11.580	1609	1614	1617	rVV	331598	294193	6.75%	0.591%
41	11.610	1617	1619	1625	rVB	70274	73765	1.69%	0.148%
42	11.733	1633	1640	1643	rBV2	206036	213249	4.89%	0.429%
43	12.010	1681	1687	1689	rBV2	277568	330670	7.59%	0.664%
44	12.033	1689	1691	1696	rVV2	279387	277546	6.37%	0.558%
45	12.080	1696	1699	1702	rVV	87128	79530	1.82%	0.160%
46	12.121	1702	1706	1708	rVV2	282098	305453	7.01%	0.614%
47	12.145	1708	1710	1713	rVB	135151	107380	2.46%	0.216%
48	12.304	1734	1737	1739	rBV	121584	111679	2.56%	0.224%
49	12.327	1739	1741	1747	rVB	167902	151200	3.47%	0.304%
50	12.451	1757	1762	1764	rBV2	53889	55596	1.28%	0.112%
51	12.557	1777	1780	1783	rBV	108460	111095	2.55%	0.223%
52	12.645	1792	1795	1800	rBV2	79591	88090	2.02%	0.177%
53	12.727	1805	1809	1812	rVV	1631955	1402785	32.18%	2.819%
54	12.786	1817	1819	1826	rVB5	46511	69095	1.58%	0.139%
55	12.851	1826	1830	1832	rBV3	35982	49246	1.13%	0.099%
56	12.957	1841	1848	1853	rBV	1361205	1466347	33.64%	2.947%
57	12.998	1853	1855	1860	rVB2	68333	75653	1.74%	0.152%
58	13.098	1864	1872	1875	rBV	3043113	3202508	73.46%	6.435%
59	13.168	1879	1884	1888	rBV2	89167	158097	3.63%	0.318%
60	13.257	1895	1899	1900	rBV3	66977	76656	1.76%	0.154%
61	13.280	1900	1903	1906	rVB	225281	223139	5.12%	0.448%
62	13.351	1911	1915	1917	rBV	137725	129453	2.97%	0.260%
63	13.386	1917	1921	1924	rBV2	119890	135458	3.11%	0.272%
64	13.480	1934	1937	1940	rBV2	71870	61852	1.42%	0.124%
65	13.509	1940	1942	1945	rVV	77325	71889	1.65%	0.144%
66	13.792	1987	1990	1994	rVB	94125	102919	2.36%	0.207%
67	13.904	2005	2009	2012	rBV3	127292	151969	3.49%	0.305%
68	13.933	2012	2014	2016	rVB	108017	78701	1.81%	0.158%
69	13.974	2016	2021	2024	rBV4	553493	600353	13.77%	1.206%
70	14.157	2047	2052	2055	rBV2	1157909	1618756	37.13%	3.253%
71	14.180	2055	2056	2062	rVB	643175	519911	11.93%	1.045%

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72	14.262	2068	2070	2073	rVB	60636	51873	1.19%	0.104%
73	14.380	2087	2090	2093	rBV2	75117	84286	1.93%	0.169%
74	14.545	2116	2118	2121	rVB2	101362	79321	1.82%	0.159%
75	14.580	2122	2124	2127	rBV2	67257	59337	1.36%	0.119%
76	14.615	2128	2130	2134	rBV5	72092	86355	1.98%	0.174%
77	14.915	2178	2181	2191	rVB	567064	681426	15.63%	1.369%
78	15.239	2231	2236	2239	rBV	399887	663568	15.22%	1.333%
79	15.351	2253	2255	2258	rVB	60877	50656	1.16%	0.102%
80	15.562	2287	2291	2297	rVB	228062	317514	7.28%	0.638%
81	15.627	2298	2302	2308	rBV	307115	373309	8.56%	0.750%
82	15.692	2309	2313	2317	rBV	514147	678430	15.56%	1.363%
83	17.262	2576	2580	2588	rVB2	144333	302068	6.93%	0.607%
84	17.739	2657	2661	2673	rVB2	99697	225831	5.18%	0.454%

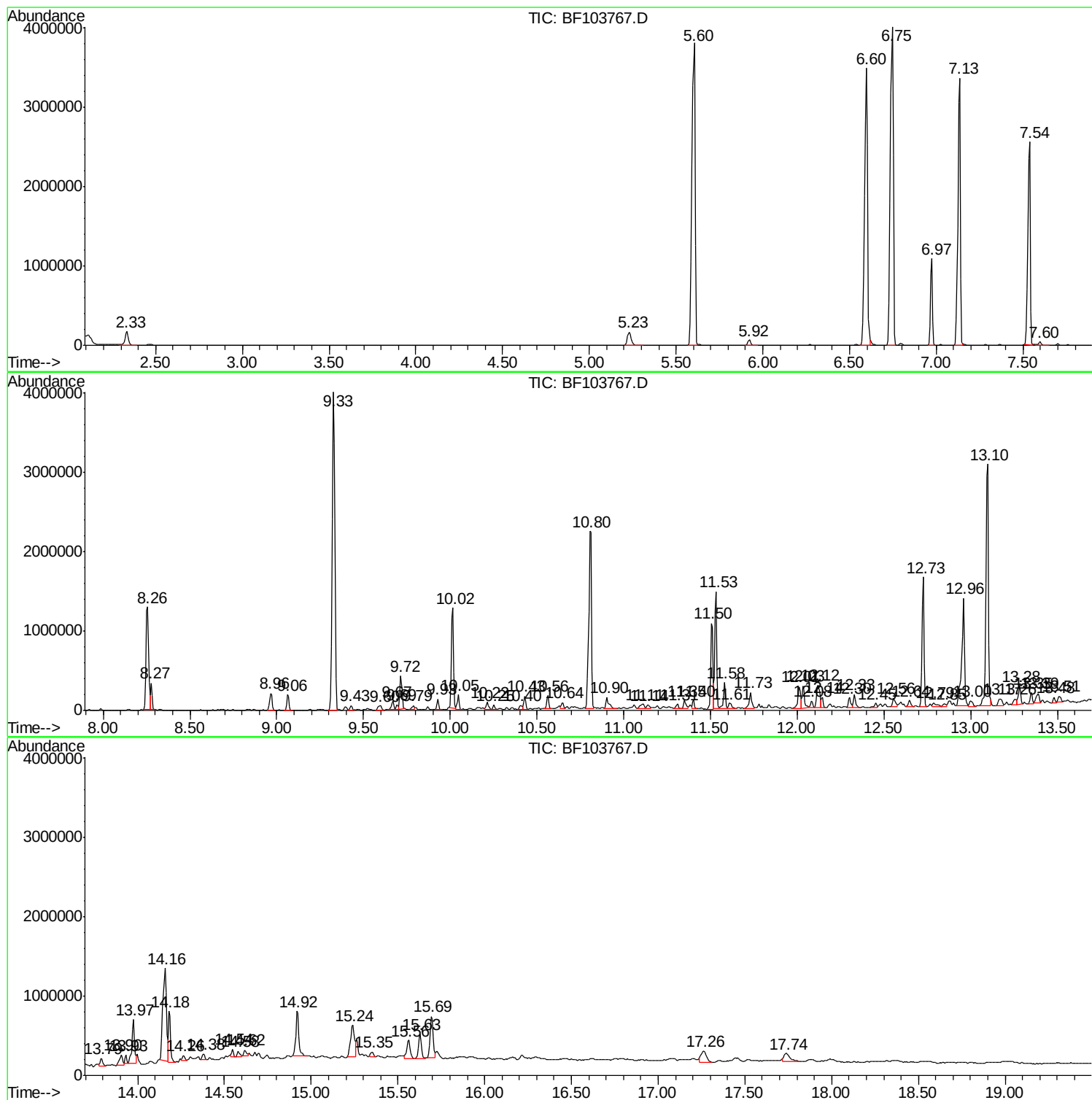
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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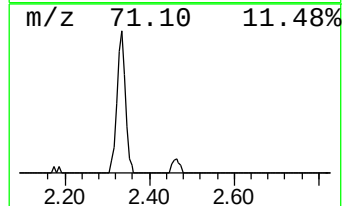
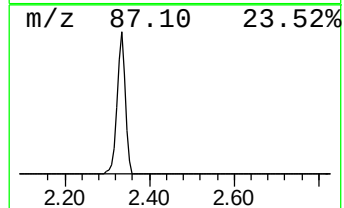
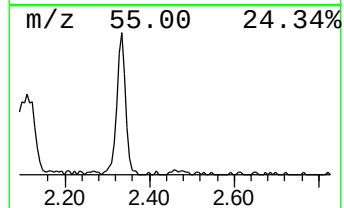
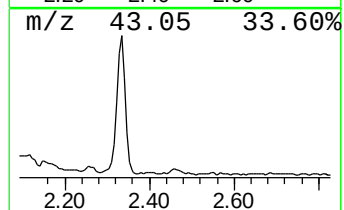
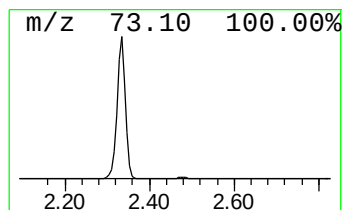
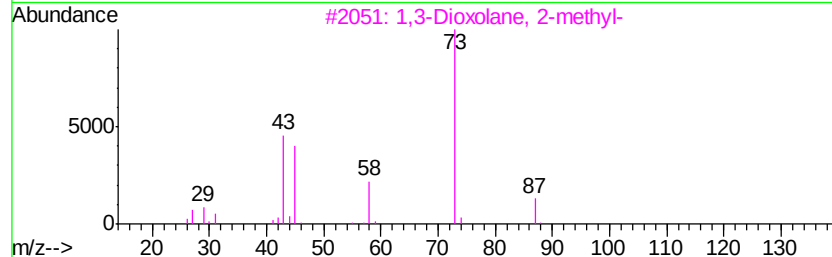
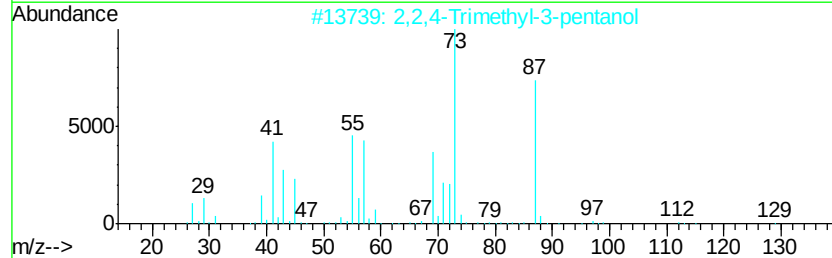
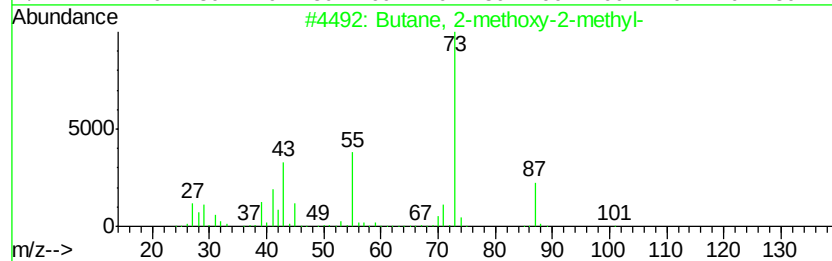
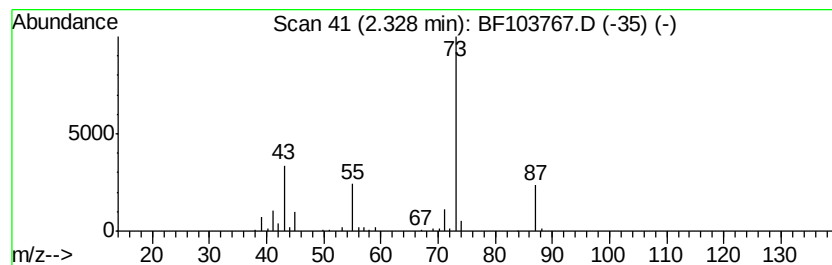
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.33	5.01 ng	227518	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
4		3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	17
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17



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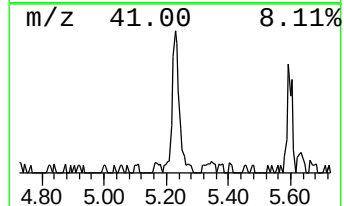
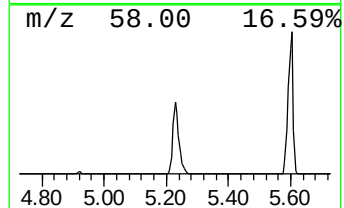
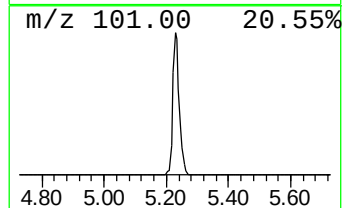
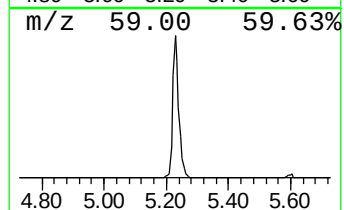
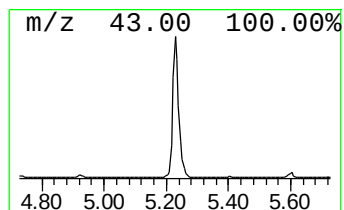
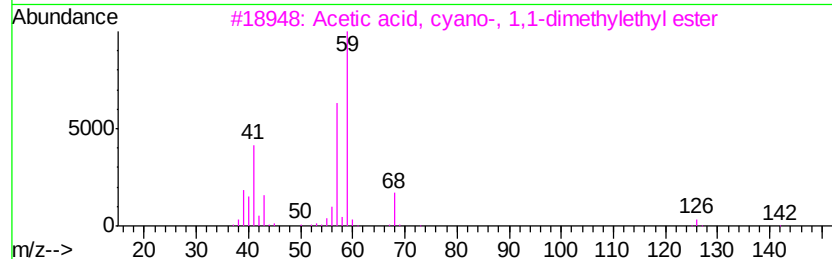
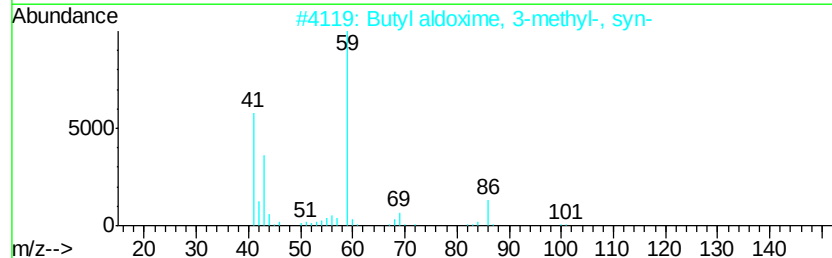
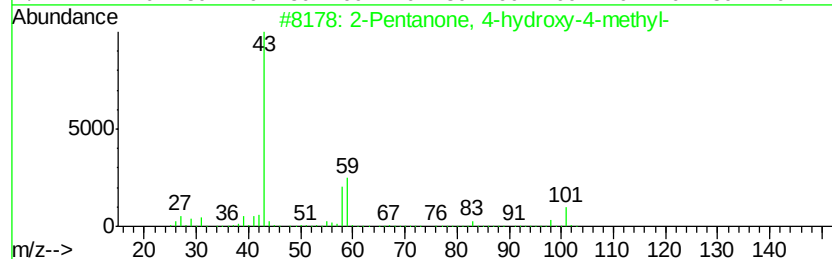
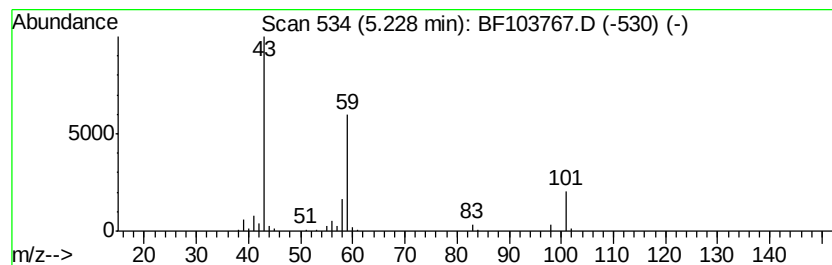
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.23	5.10 ng	231451	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4			Guanidine	59	CH5N3	000113-00-8	9
5			Acetone	58	C3H6O	000067-64-1	9



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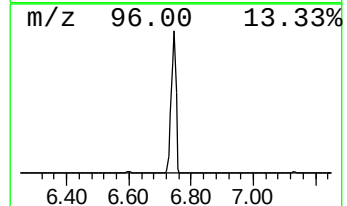
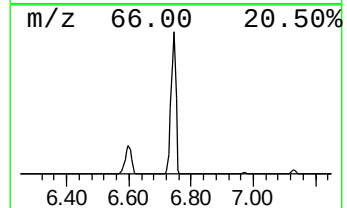
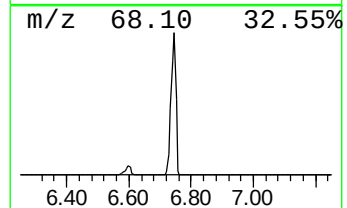
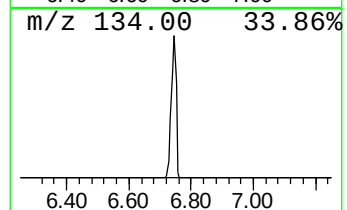
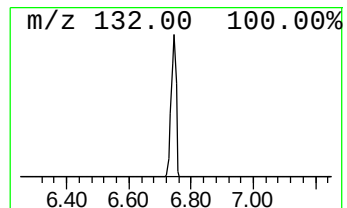
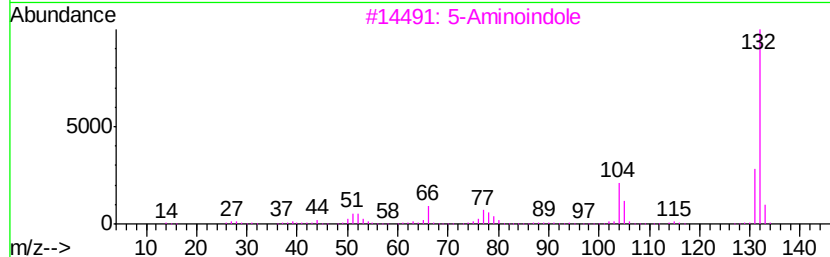
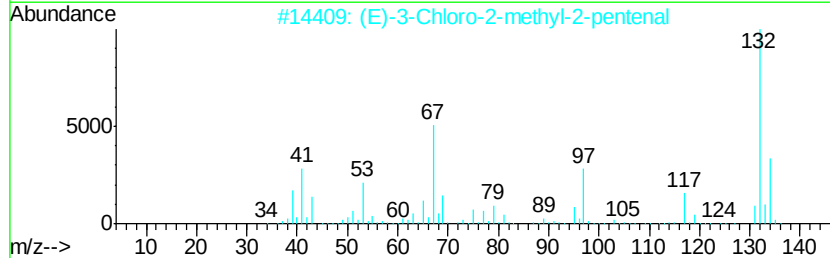
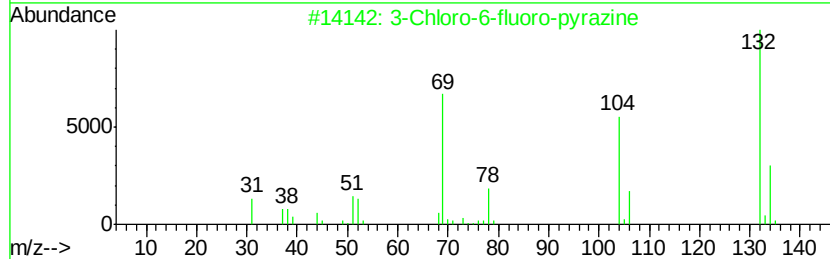
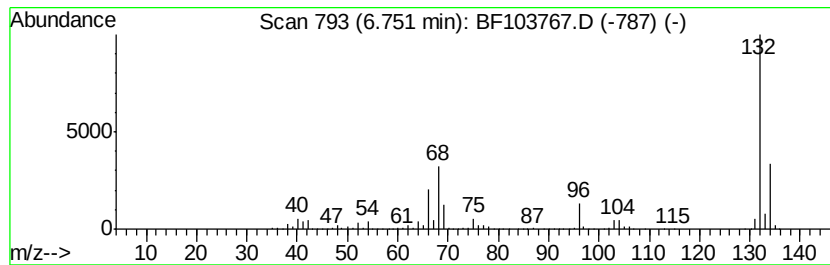
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.75 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	94.06 ng	4271140	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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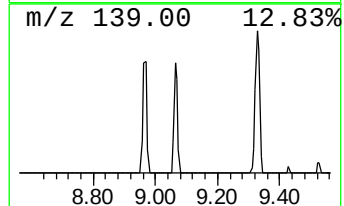
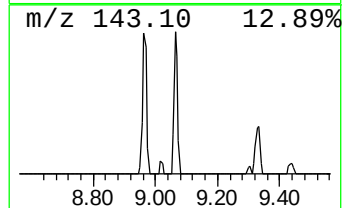
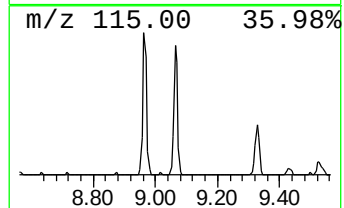
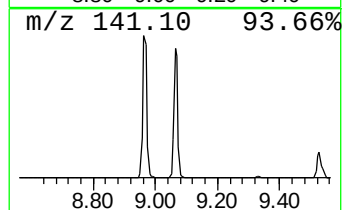
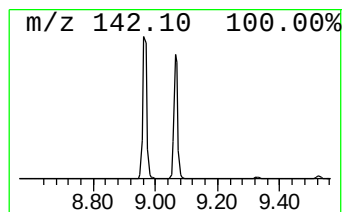
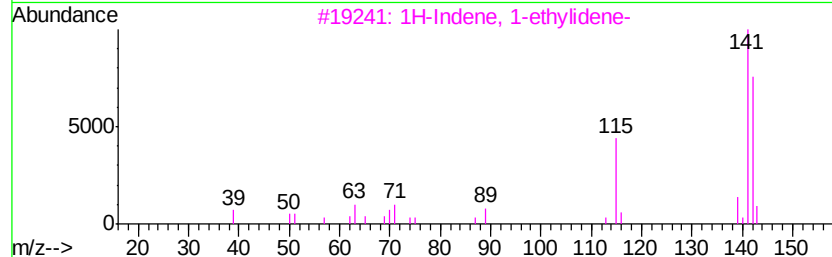
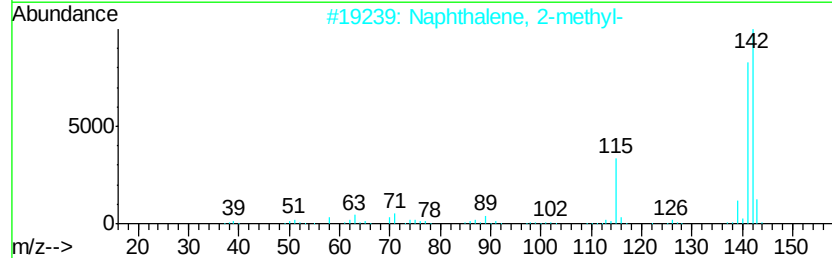
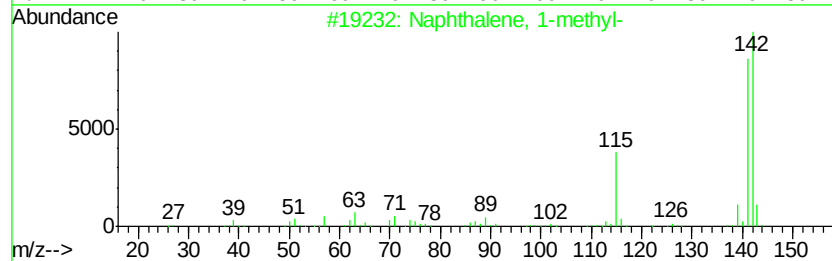
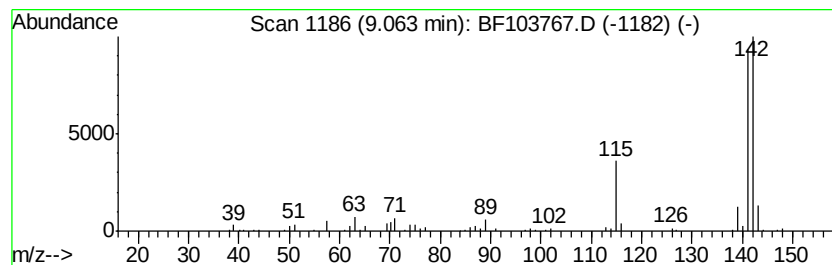
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Naphthalene, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.06	2.80 ng	174101	Naphthalene-d8	8.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	95
4		Benzocycloheptatriene	142	C11H10	000264-09-5	94
5		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
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Acq On : 19 Mar 2018 20:30
Operator : JU/SJ
Sample : J1971-05
Misc :
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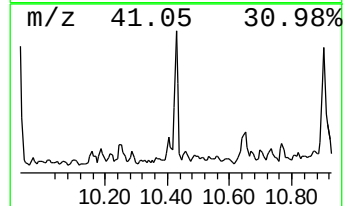
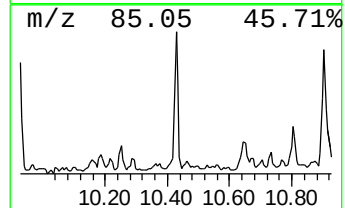
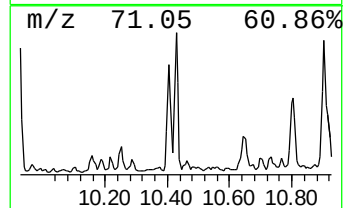
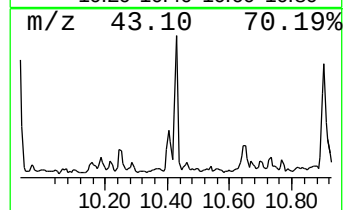
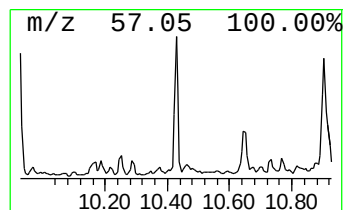
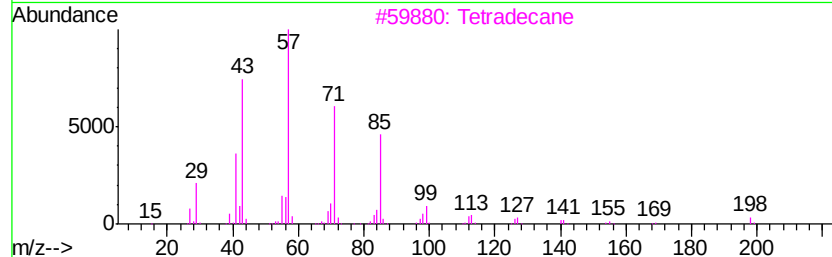
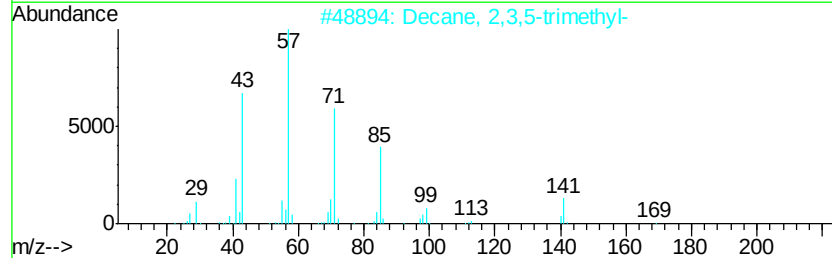
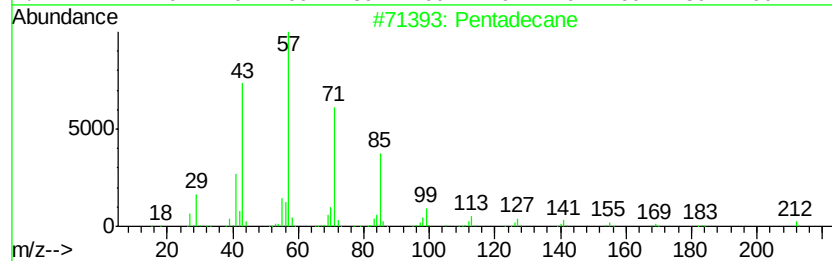
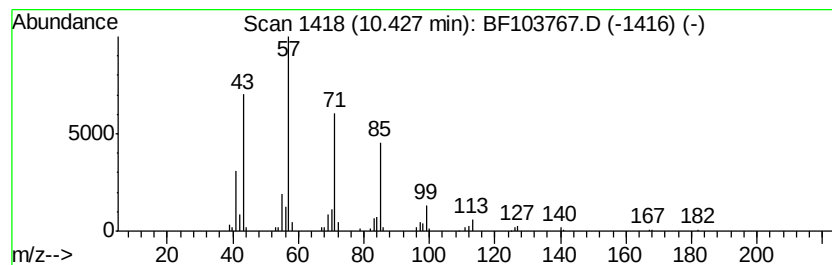
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Pentadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.43	2.35 ng	136288	Acenaphthene-d10	10.02
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Pentadecane	212 C15H32	000629-62-9	90
2	Decane, 2,3,5-trimethyl-	184 C13H28	062238-11-3	90
3	Tetradecane	198 C14H30	000629-59-4	86
4	Hexadecane	226 C16H34	000544-76-3	83
5	Heptadecane	240 C17H36	000629-78-7	83



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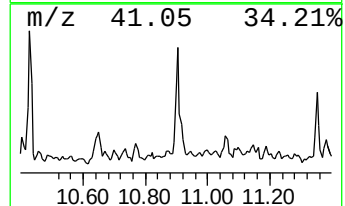
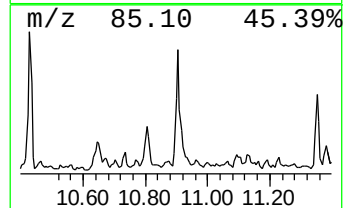
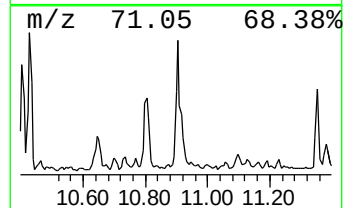
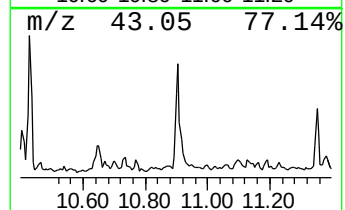
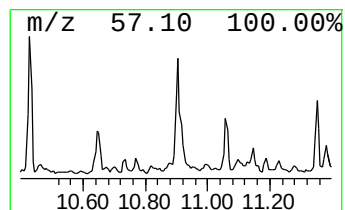
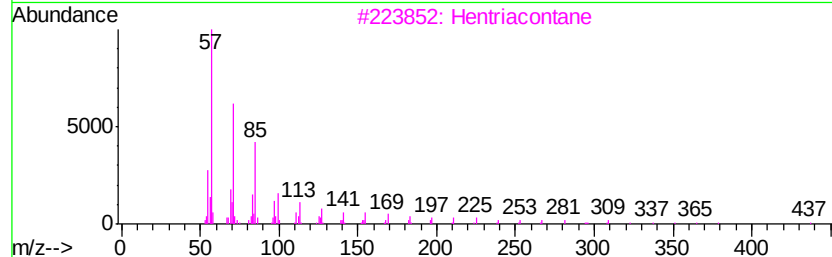
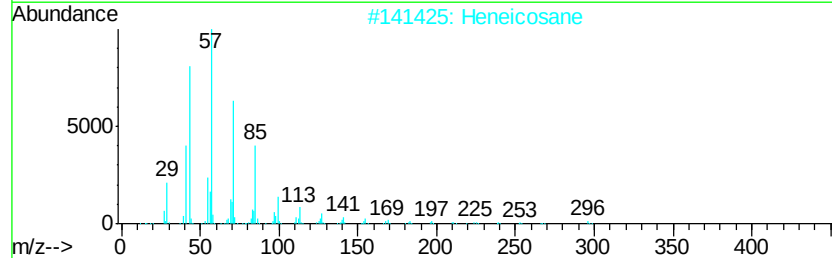
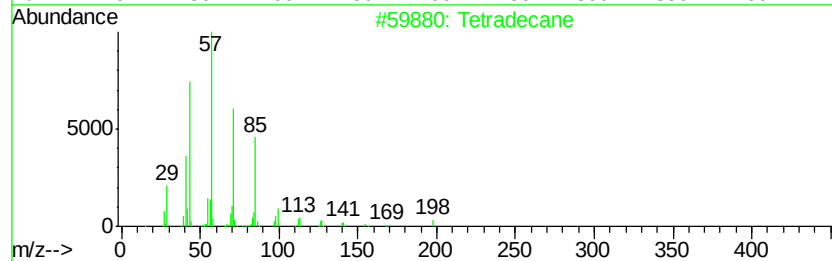
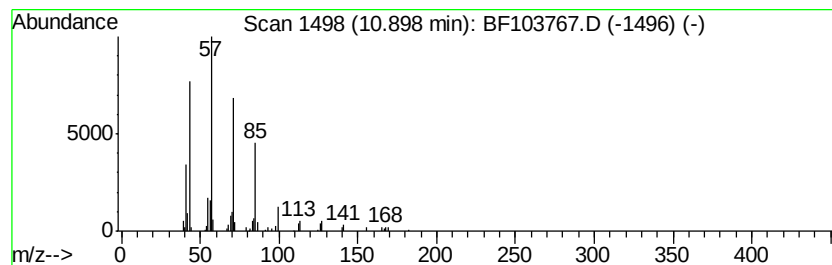
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tetradecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.90	3.16 ng	161912	Phenanthrene-d10	11.50

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	90
2		Heneicosane	296	C21H44	000629-94-7	80
3		Hentriacontane	437	C31H64	000630-04-6	78
4		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	78
5		Pentadecane	212	C15H32	000629-62-9	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
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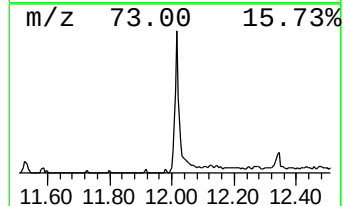
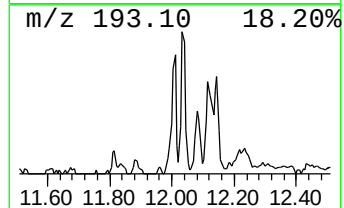
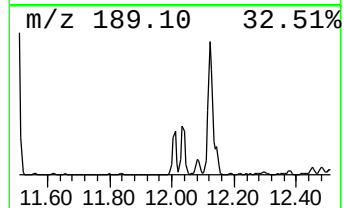
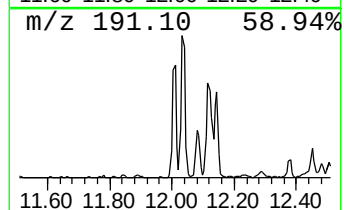
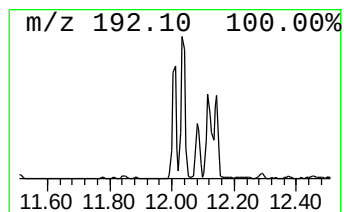
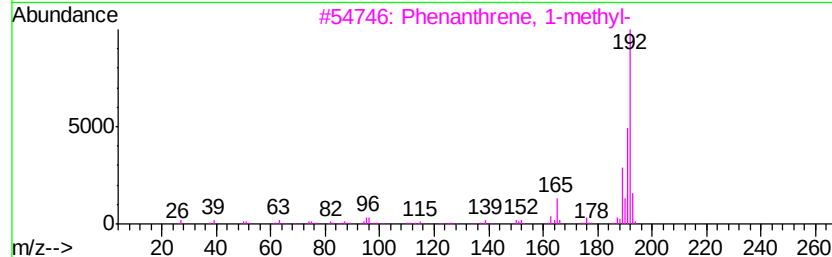
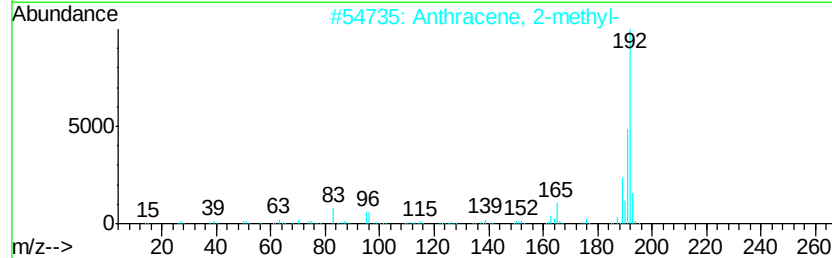
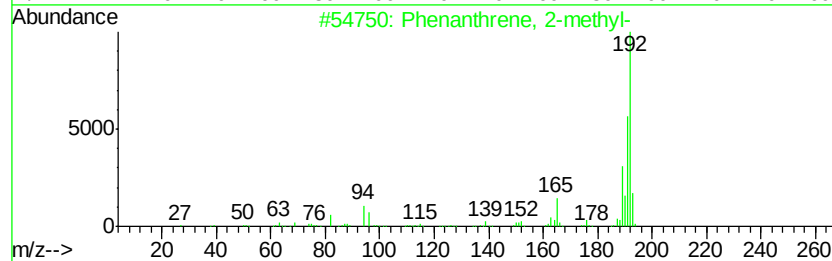
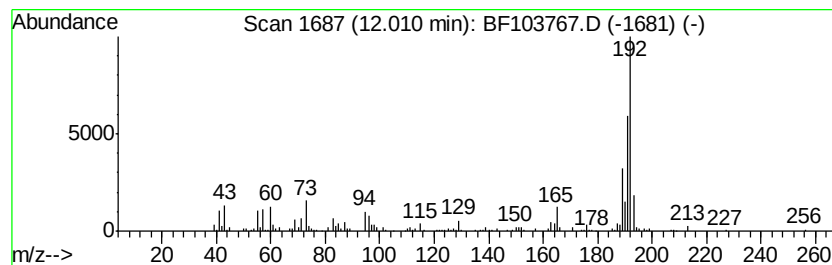
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phenanthrene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	6.46 ng	330670	Phenanthrene-d10	11.50

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		5H-Dibenzofa,dlcycloheptene	192	C15H12	000256-81-5	94
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	94



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
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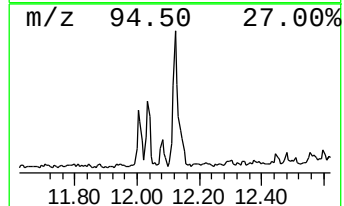
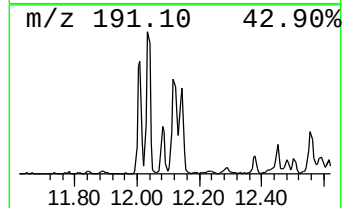
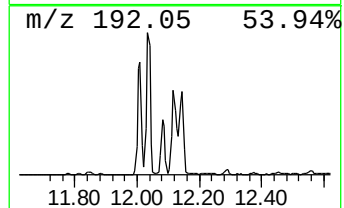
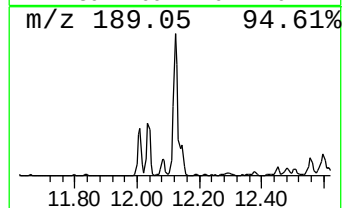
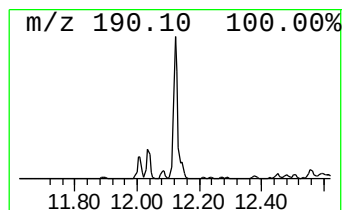
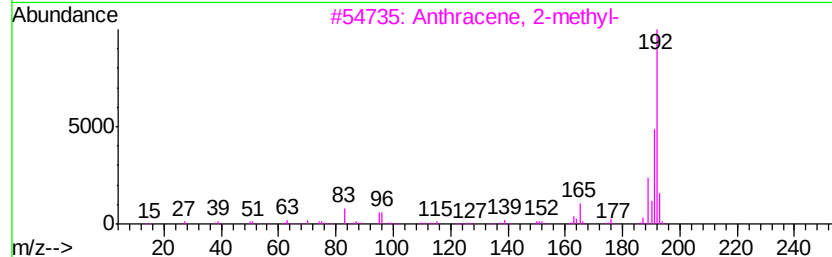
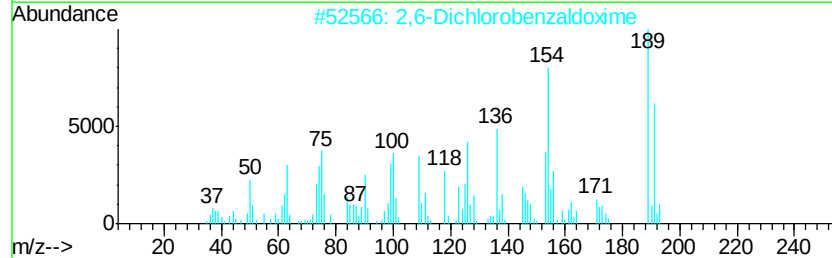
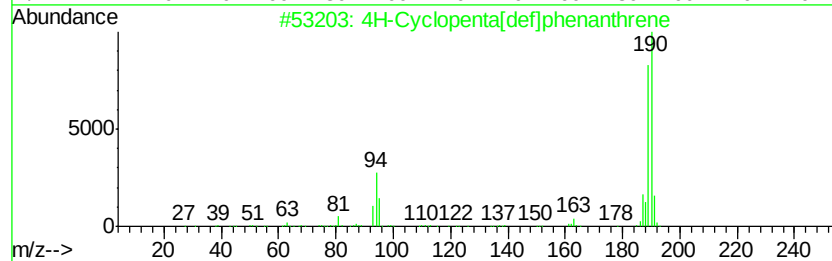
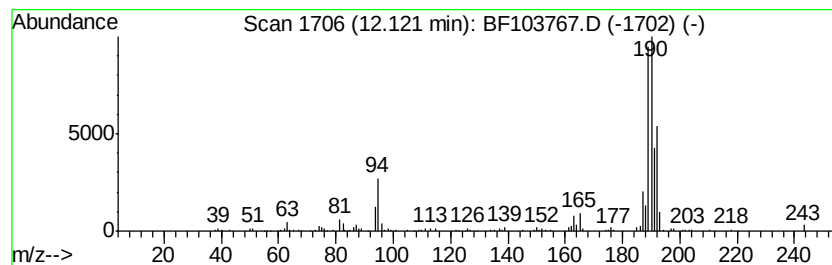
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 4H-Cyclopenta[def]phenanthrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.12	5.97 ng	305453	Phenanthrene-d10	11.50	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2	2,6-Dichlorobenzaloxime	189	C7H5Cl2NO	025185-95-9	42
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	25
4	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	25
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031918\
Data File : BF103767.D
Acq On : 19 Mar 2018 20:30
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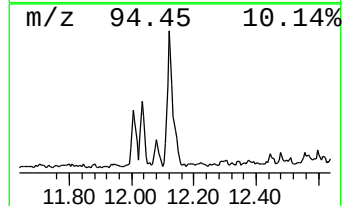
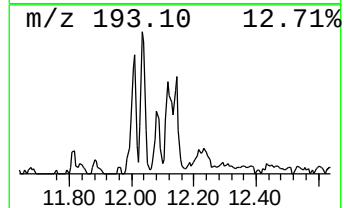
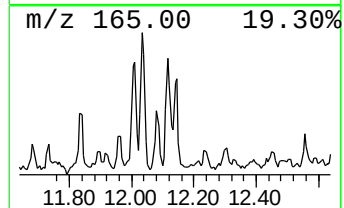
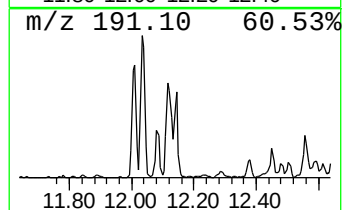
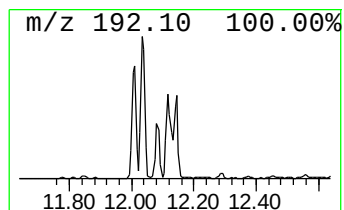
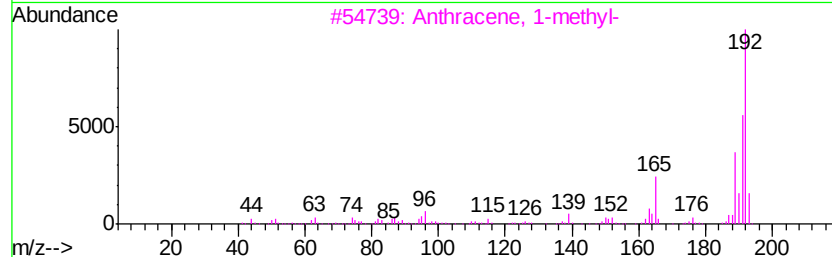
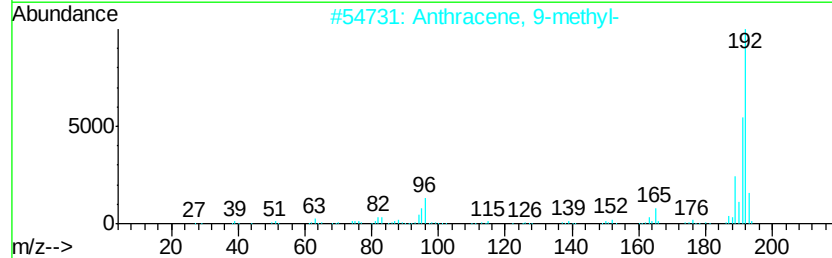
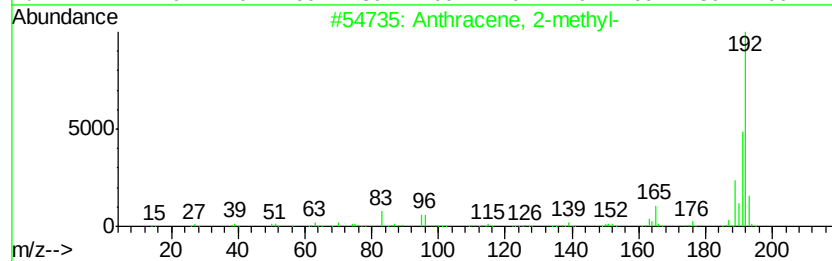
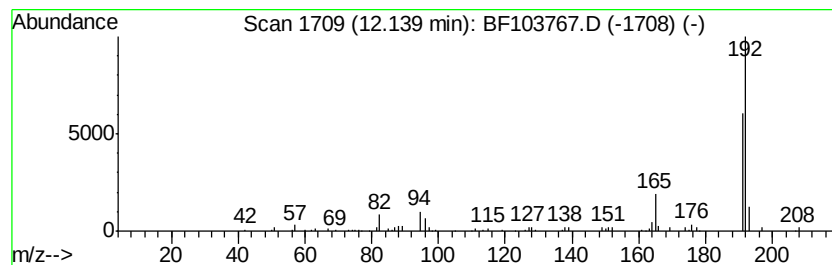
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Anthracene, 2-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	2.10 ng	107380	Phenanthrene-d10	11.50

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
2			Anthracene, 9-methyl-	192	C15H12	000779-02-2	90
3			Anthracene, 1-methyl-	192	C15H12	000610-48-0	87
4			Phenanthrene, 9-methyl-	192	C15H12	000883-20-5	86
5			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031918\
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Acq On : 19 Mar 2018 20:30
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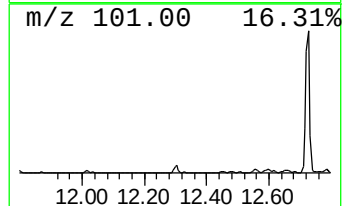
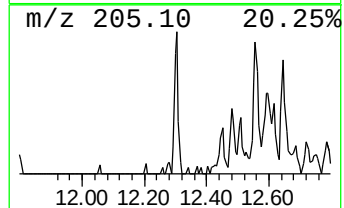
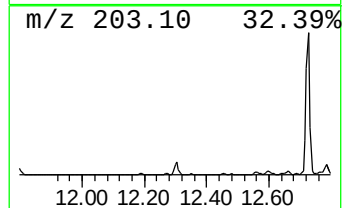
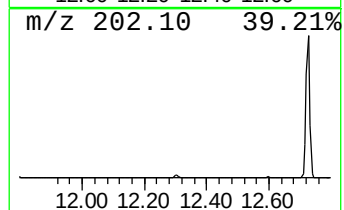
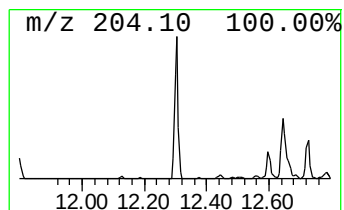
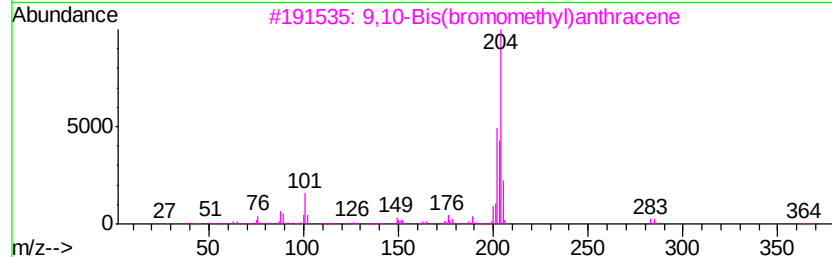
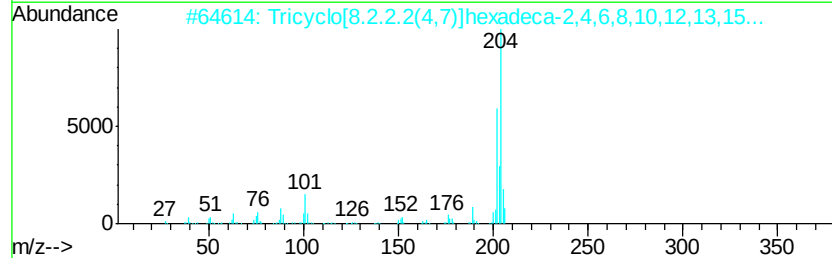
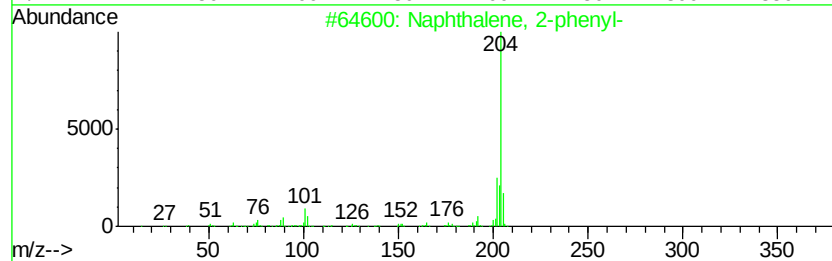
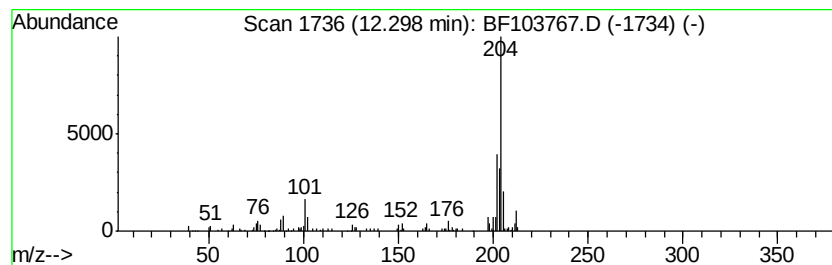
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Naphthalene, 2-phenyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	2.18 ng	111679	Phenanthrene-d10	11.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	95
2		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	70
3		9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	62
4		Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	55
5		3,4-Biphenyldicarbonitrile	204	C14H8N2	004128-63-6	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031918\
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Acq On : 19 Mar 2018 20:30
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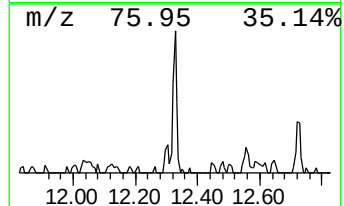
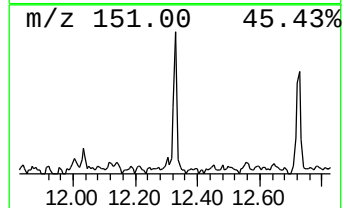
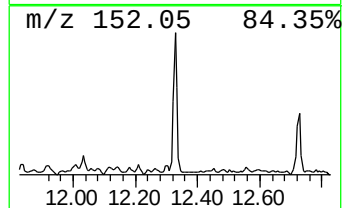
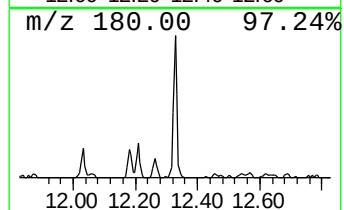
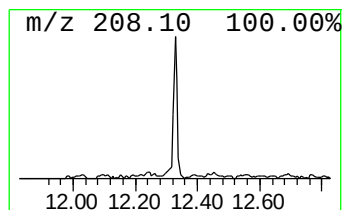
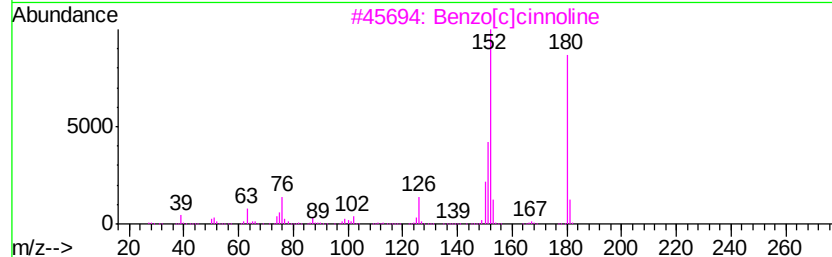
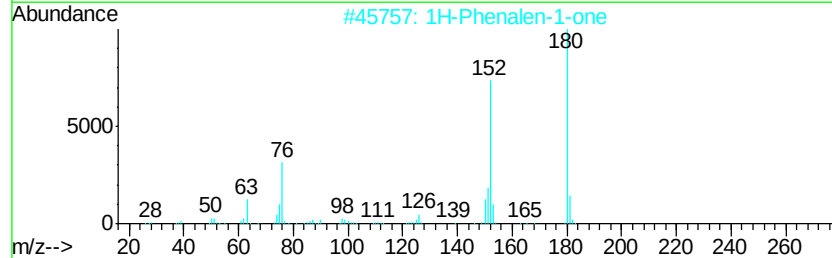
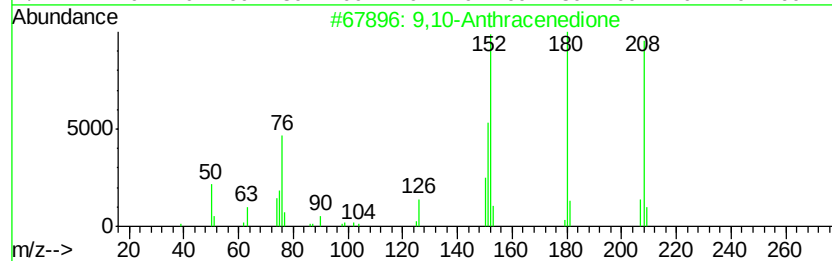
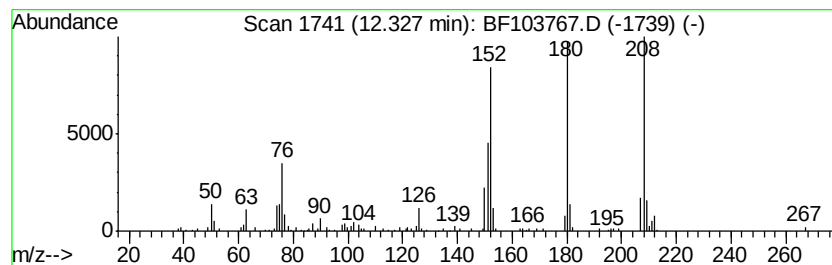
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 9,10-Anthracenedione Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.33	2.95 ng	151200	Phenanthrene-d10	11.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Anthracenedione	208	C14H8O2	000084-65-1	97
2		1H-Phenalen-1-one	180	C13H8O	000548-39-0	90
3		Benzo[c]cinnoline	180	C12H8N2	000230-17-1	86
4		9H-Fluoren-9-one	180	C13H8O	000486-25-9	86
5		1,4-Anthracenedione	208	C14H8O2	000635-12-1	49



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
Data File : BF103767.D
Acq On : 19 Mar 2018 20:30
Operator : JU/SJ
Sample : J1971-05
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ES-6-SA-3-WW@4.5

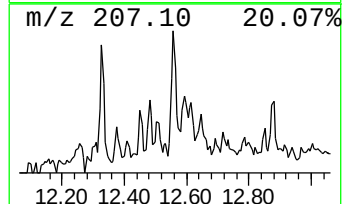
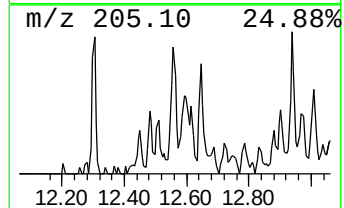
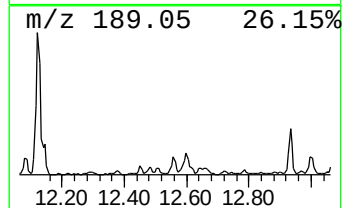
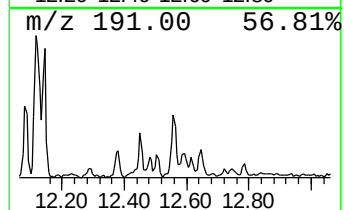
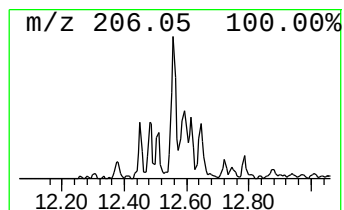
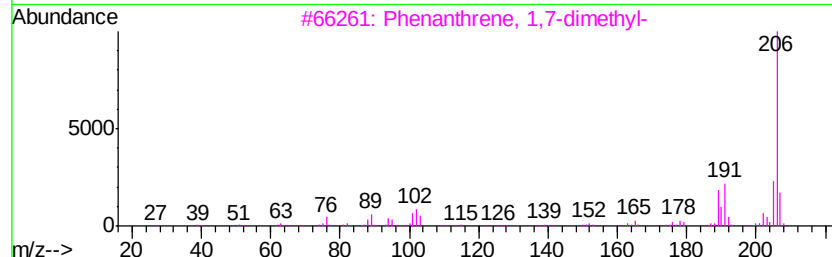
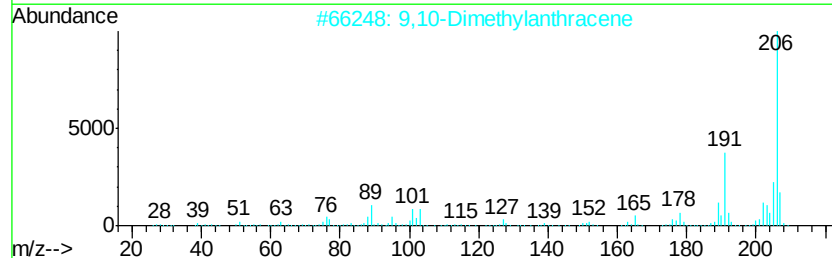
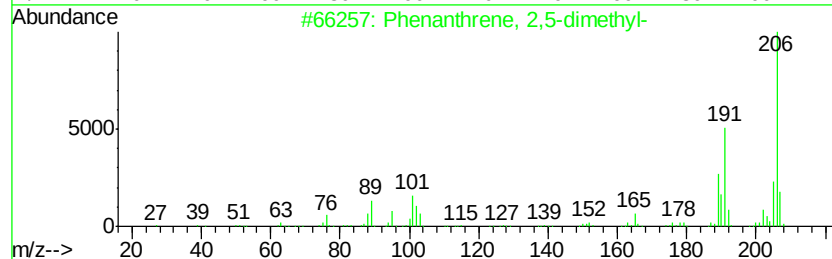
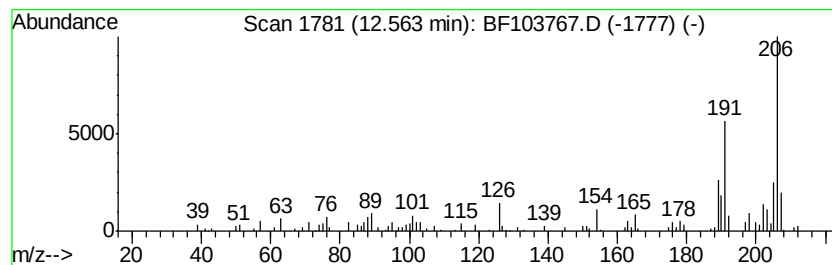
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Phenanthrene, 2,5-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.56	2.17 ng	111095	Phenanthrene-d10	11.50

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2			9,10-Dimethylantracene	206	C16H14	000781-43-1	93
3			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	90
4			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	89
5			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
Data File : BF103767.D
Acq On : 19 Mar 2018 20:30
Operator : JU/SJ
Sample : J1971-05
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
ES-6-SA-3-WW@4.5

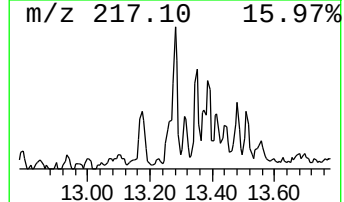
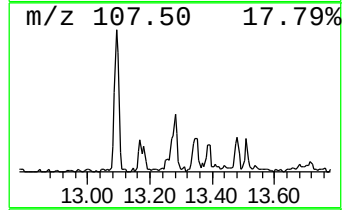
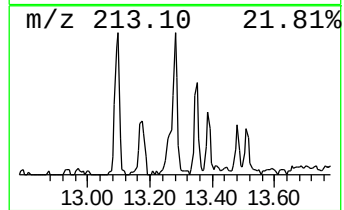
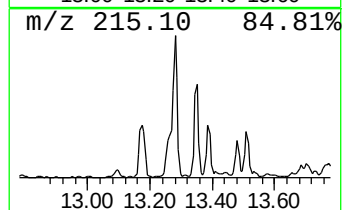
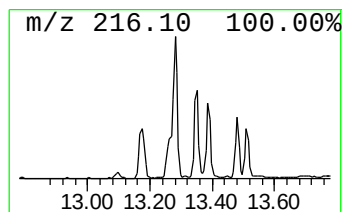
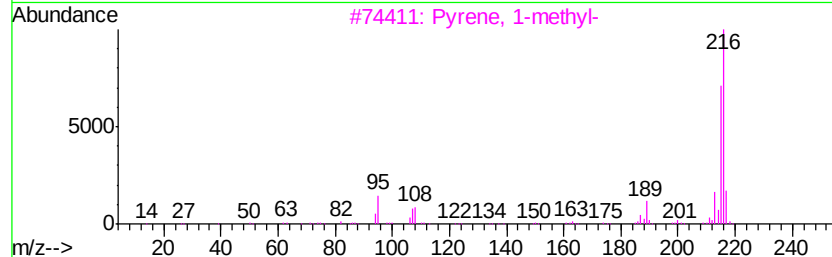
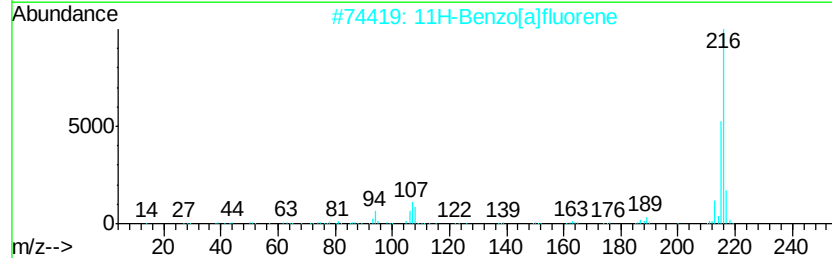
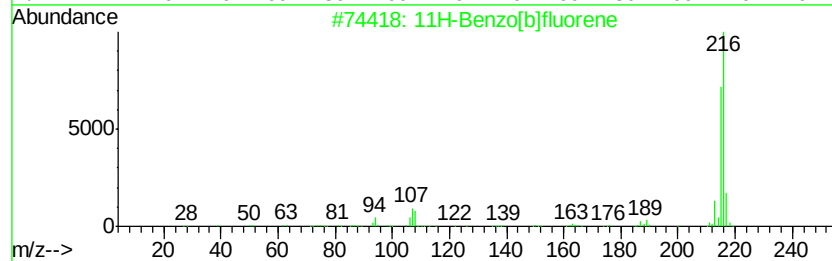
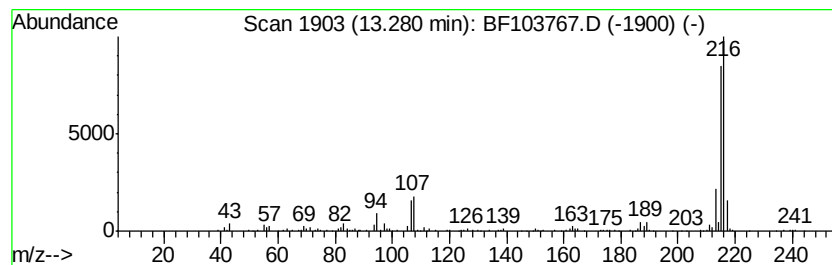
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 11H-Benzo[b]fluorene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.28	2.76 ng	223139	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
2		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
3		Pvrene, 1-methyl-	216	C17H12	002381-21-7	87
4		Pvrene, 2-methyl-	216	C17H12	003442-78-2	76
5		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	76



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
Data File : BF103767.D
Acq On : 19 Mar 2018 20:30
Operator : JU/SJ
Sample : J1971-05
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
ES-6-SA-3-WW@4.5

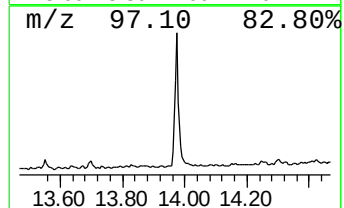
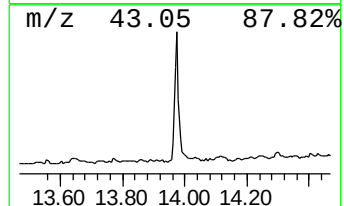
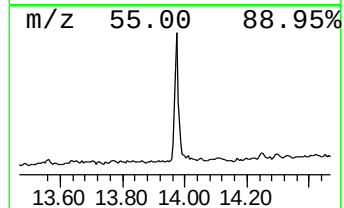
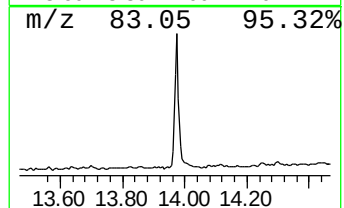
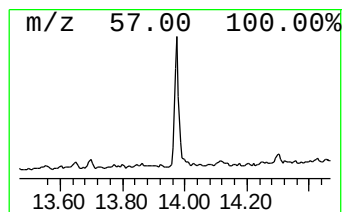
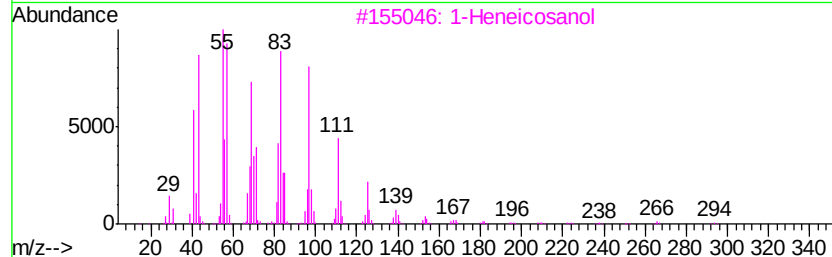
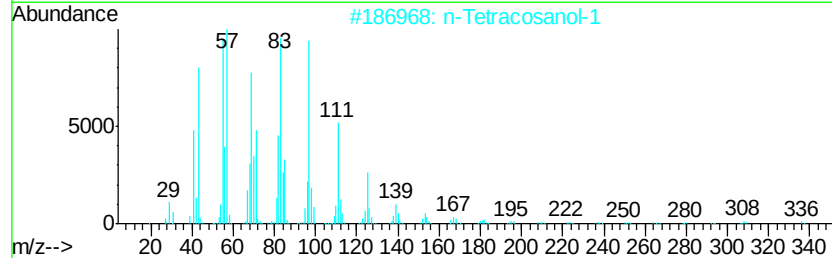
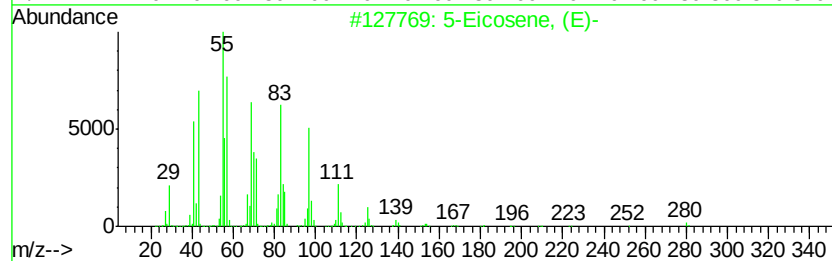
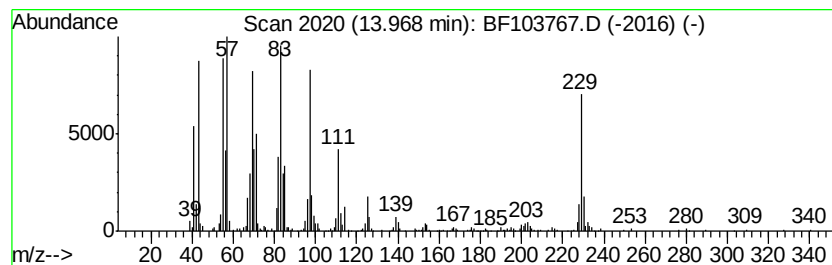
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 5-Eicosene, (E)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	7.42 ng	600353	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Eicosene, (E)-	280	C20H40	074685-30-6	95
2		n-Tetracosanol-1	354	C24H50O	000506-51-4	95
3		1-Heneicosanol	312	C21H44O	015594-90-8	95
4		n-Nonadecanol-1	284	C19H40O	001454-84-8	95
5		1-Heneicosyl formate	340	C22H44O2	077899-03-7	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
Data File : BF103767.D
Acq On : 19 Mar 2018 20:30
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Sample : J1971-05
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
ES-6-SA-3-WW@4.5

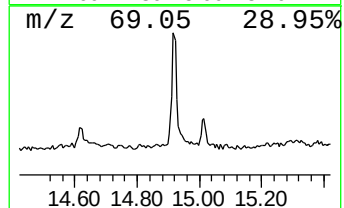
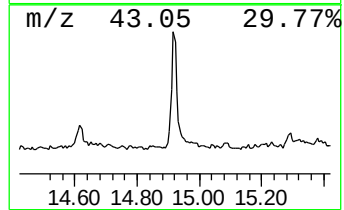
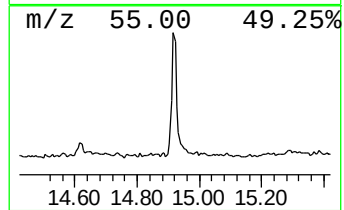
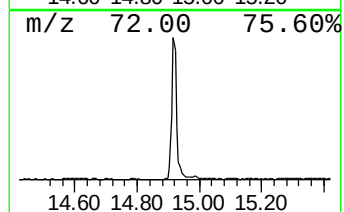
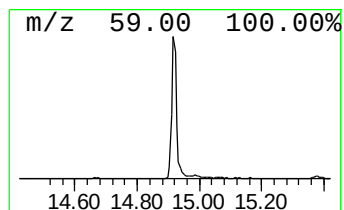
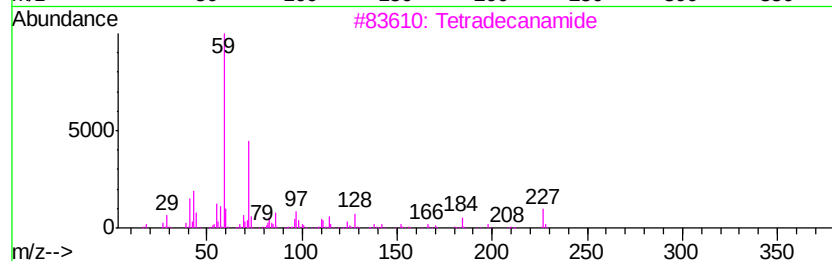
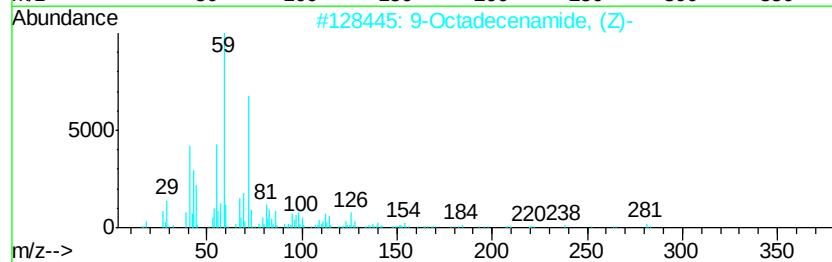
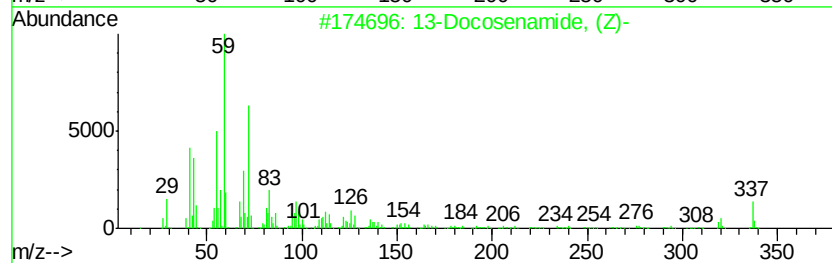
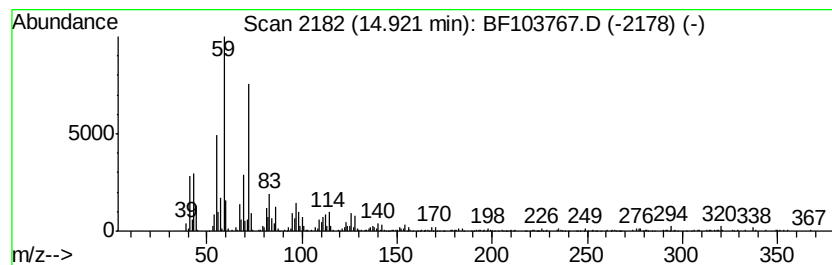
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 13-Docosenamide, (Z)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.92	8.42 ng	681426	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	96
2		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	91
3		Tetradecanamide	227	C14H29NO	000638-58-4	50
4		Decanamide-	171	C10H21NO	002319-29-1	47
5		Dodecanamide	199	C12H25NO	001120-16-7	42



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
Data File : BF103767.D
Acq On : 19 Mar 2018 20:30
Operator : JU/SJ
Sample : J1971-05
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ES-6-SA-3-WW@4.5

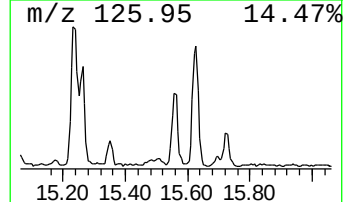
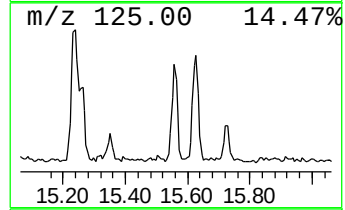
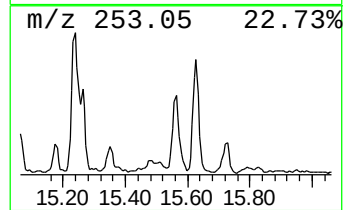
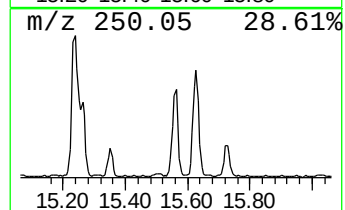
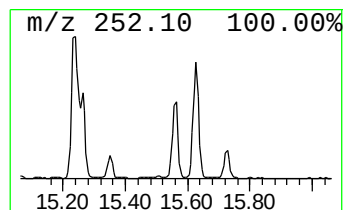
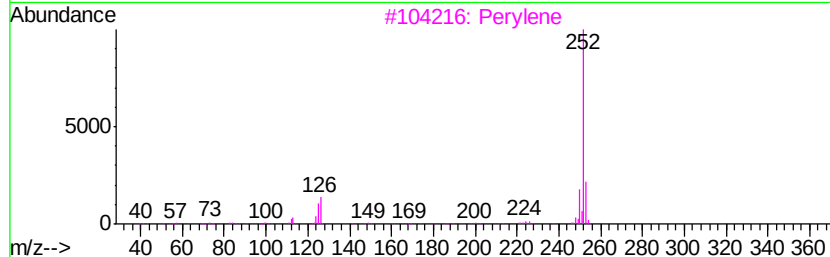
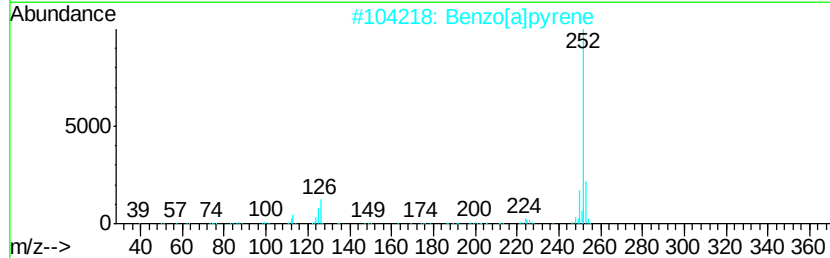
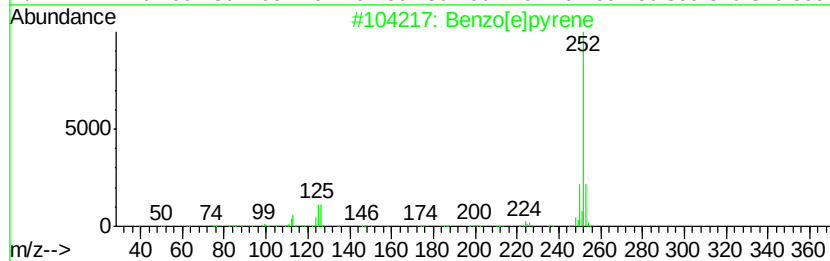
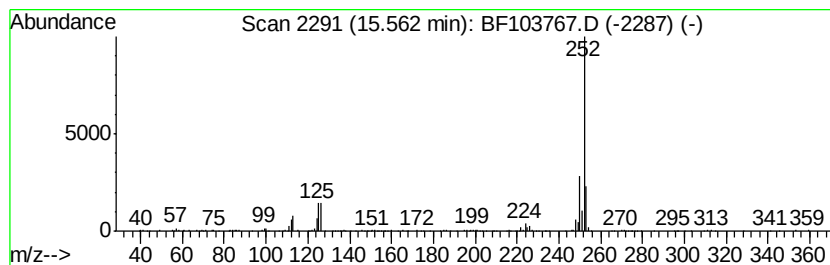
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.56	9.36 ng	317514	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2		Benzo[a]pyrene	252	C20H12	000050-32-8	96
3		Perylene	252	C20H12	000198-55-0	95
4		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	93
5		Benzo[k]fluoranthene	252	C20H12	000207-08-9	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031918\
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 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
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 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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2-Pentanone, 4-hy...	5.23	5.1	ng	231451	1	6.97	908167	20.0
unknown6.75	6.75	94.1	ng	4271140	1	6.97	908167	20.0
Naphthalene, 1-me...	9.06	2.8	ng	174101	2	8.26	1245240	20.0
Pentadecane	10.43	2.4	ng	136288	3	10.02	1161030	20.0
Tetradecane	10.90	3.2	ng	161912	4	11.50	1023450	20.0
Phenanthrene, 2-m...	12.01	6.5	ng	330670	4	11.50	1023450	20.0
4H-Cyclopenta[def...	12.12	6.0	ng	305453	4	11.50	1023450	20.0
Anthracene, 2-met...	12.14	2.1	ng	107380	4	11.50	1023450	20.0
Naphthalene, 2-ph...	12.30	2.2	ng	111679	4	11.50	1023450	20.0
9,10-Anthracenedione	12.33	3.0	ng	151200	4	11.50	1023450	20.0
Phenanthrene, 2,5...	12.56	2.2	ng	111095	4	11.50	1023450	20.0
11H-Benzo[b]fluorene	13.28	2.8	ng	223139	5	14.16	1618760	20.0
5-Eicosene, (E)-	13.97	7.4	ng	600353	5	14.16	1618760	20.0
13-Docosenamide, ...	14.92	8.4	ng	681426	5	14.16	1618760	20.0
Benzo[e]pyrene	15.56	9.4	ng	317514	6	15.69	678430	20.0