

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
 Data File : BF103935.D
 Acq On : 26 Mar 2018 19:55
 Operator : JU/SJ
 Sample : J1758-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-02-032318-A

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

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.322	33	40	46	rBV	106509	143916	4.58%	0.344%
2	5.216	528	532	542	rVB	101933	137024	4.36%	0.328%
3	5.598	592	597	600	rBV	2793838	2812940	89.49%	6.730%
4	6.598	761	767	782	rVB	2637543	2866155	91.18%	6.857%
5	6.745	787	792	795	rBV	2924633	2914426	92.72%	6.972%
6	6.969	827	830	834	rBV	903727	822477	26.17%	1.968%
7	7.128	853	857	861	rBV	2576889	2252827	71.67%	5.390%
8	7.534	921	926	929	rBV	1878019	1811074	57.62%	4.333%
9	8.222	1040	1043	1045	rBV	55919	53324	1.70%	0.128%
10	8.257	1045	1049	1054	rVV	1146006	1163043	37.00%	2.782%
11	8.298	1054	1056	1060	rVB2	41506	35387	1.13%	0.085%
12	8.534	1092	1096	1100	rVB4	30136	35091	1.12%	0.084%
13	8.751	1129	1133	1136	rBV	40078	37332	1.19%	0.089%
14	8.828	1141	1146	1151	rVV	87211	79179	2.52%	0.189%
15	8.963	1166	1169	1172	rVB	51190	46067	1.47%	0.110%
16	9.069	1181	1187	1190	rBV2	79833	86132	2.74%	0.206%
17	9.328	1226	1231	1234	rBV	3574588	3143300	100.00%	7.520%
18	9.392	1239	1242	1246	rVB	107866	91572	2.91%	0.219%
19	9.598	1272	1277	1280	rBV3	40315	61214	1.95%	0.146%
20	9.669	1280	1289	1291	rBV4	69428	100603	3.20%	0.241%
21	9.716	1294	1297	1305	rVB	228936	238050	7.57%	0.570%
22	9.786	1305	1309	1310	rBV2	38895	45755	1.46%	0.109%
23	9.869	1320	1323	1327	rBV	41784	43441	1.38%	0.104%
24	9.928	1329	1333	1336	rVB	182423	156673	4.98%	0.375%
25	10.016	1344	1348	1351	rBV	1238331	1188535	37.81%	2.843%
26	10.045	1351	1353	1357	rVB	124117	93259	2.97%	0.223%
27	10.163	1369	1373	1375	rBV4	28043	43284	1.38%	0.104%
28	10.216	1379	1382	1385	rVV2	95563	90989	2.89%	0.218%
29	10.245	1385	1387	1392	rVV3	57742	61051	1.94%	0.146%
30	10.328	1397	1401	1403	rBV2	33992	40560	1.29%	0.097%
31	10.428	1411	1418	1421	rBV	250885	239695	7.63%	0.573%
32	10.563	1437	1441	1448	rVB	180156	168259	5.35%	0.403%
33	10.645	1450	1455	1458	rBV3	89728	112123	3.57%	0.268%
34	10.804	1478	1482	1486	rBV	1937042	1917204	60.99%	4.587%

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Filtering: 5
 Min Area: 3 % of largest Peak
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35	10.898	1495	1498	1508	rVB	202103	264392	8.41%	0.633%
36	11.098	1528	1532	1533	rBV2	36003	44408	1.41%	0.106%
37	11.186	1545	1547	1551	rVB2	37211	40801	1.30%	0.098%
38	11.222	1551	1553	1559	rBV7	26567	54173	1.72%	0.130%
39	11.351	1572	1575	1577	rBV	152965	123769	3.94%	0.296%
40	11.380	1577	1580	1582	rVV2	57850	60870	1.94%	0.146%
41	11.404	1582	1584	1586	rVB	66005	52288	1.66%	0.125%
42	11.510	1597	1602	1604	rBV	1188297	1248803	39.73%	2.988%
43	11.528	1604	1605	1609	rVV	843104	676055	21.51%	1.617%
44	11.580	1609	1614	1618	rVV	240440	211928	6.74%	0.507%
45	11.733	1637	1640	1644	rVB	47090	39539	1.26%	0.095%
46	11.775	1644	1647	1650	rBV	80731	65372	2.08%	0.156%
47	11.880	1661	1665	1668	rVB	1068133	842618	26.81%	2.016%
48	12.010	1684	1687	1689	rBV2	155768	166682	5.30%	0.399%
49	12.033	1689	1691	1696	rVB	150053	142772	4.54%	0.342%
50	12.086	1696	1700	1703	rBV	53368	55030	1.75%	0.132%
51	12.122	1703	1706	1708	rVV	188216	188012	5.98%	0.450%
52	12.139	1708	1709	1713	rVB	65606	53947	1.72%	0.129%
53	12.186	1714	1717	1720	rBV2	82326	68724	2.19%	0.164%
54	12.304	1734	1737	1739	rVV	71662	53165	1.69%	0.127%
55	12.333	1739	1742	1747	rVB2	40039	42793	1.36%	0.102%
56	12.574	1777	1783	1784	rBV	2849737	2835388	90.20%	6.783%
57	12.592	1784	1786	1792	rVB2	2374024	2181302	69.40%	5.219%
58	12.674	1797	1800	1805	rVB	439625	369499	11.76%	0.884%
59	12.722	1805	1808	1812	rBV	820555	786031	25.01%	1.880%
60	12.910	1837	1840	1842	rBV3	57307	52618	1.67%	0.126%
61	12.957	1842	1848	1853	rVB	767294	832305	26.48%	1.991%
62	13.004	1853	1856	1860	rVB2	48598	51950	1.65%	0.124%
63	13.092	1864	1871	1874	rBV	2682710	2558246	81.39%	6.120%
64	13.169	1880	1884	1885	rBV2	45930	55622	1.77%	0.133%
65	13.239	1893	1896	1898	rBV2	82675	85401	2.72%	0.204%
66	13.280	1900	1903	1906	rVB	128717	124578	3.96%	0.298%
67	13.351	1912	1915	1917	rVB	119264	100626	3.20%	0.241%
68	13.386	1917	1921	1929	rBV4	74136	150838	4.80%	0.361%
69	13.510	1940	1942	1945	rBV	41209	41804	1.33%	0.100%
70	13.645	1962	1965	1968	rBV2	45171	52690	1.68%	0.126%
71	13.974	2016	2021	2026	rBV4	331894	399562	12.71%	0.956%

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72	14.074	2035	2038	2041	rBV3	49125	60377	1.92%	0.144%
73	14.157	2047	2052	2055	rBV2	1029534	1217512	38.73%	2.913%
74	14.186	2055	2057	2064	rVB	271041	251498	8.00%	0.602%
75	14.269	2068	2071	2073	rBV	48368	44999	1.43%	0.108%
76	14.298	2073	2076	2082	rBV4	60444	83745	2.66%	0.200%
77	14.610	2127	2129	2134	rVB4	43328	48995	1.56%	0.117%
78	14.704	2143	2145	2150	rVB3	45798	44180	1.41%	0.106%
79	15.016	2195	2198	2202	rVB	102621	106811	3.40%	0.256%
80	15.245	2233	2237	2244	rBV	183844	381602	12.14%	0.913%
81	15.568	2289	2292	2299	rVB	110146	153594	4.89%	0.367%
82	15.633	2299	2303	2309	rBV	186246	267134	8.50%	0.639%
83	15.704	2310	2315	2319	rBV	630347	832100	26.47%	1.991%

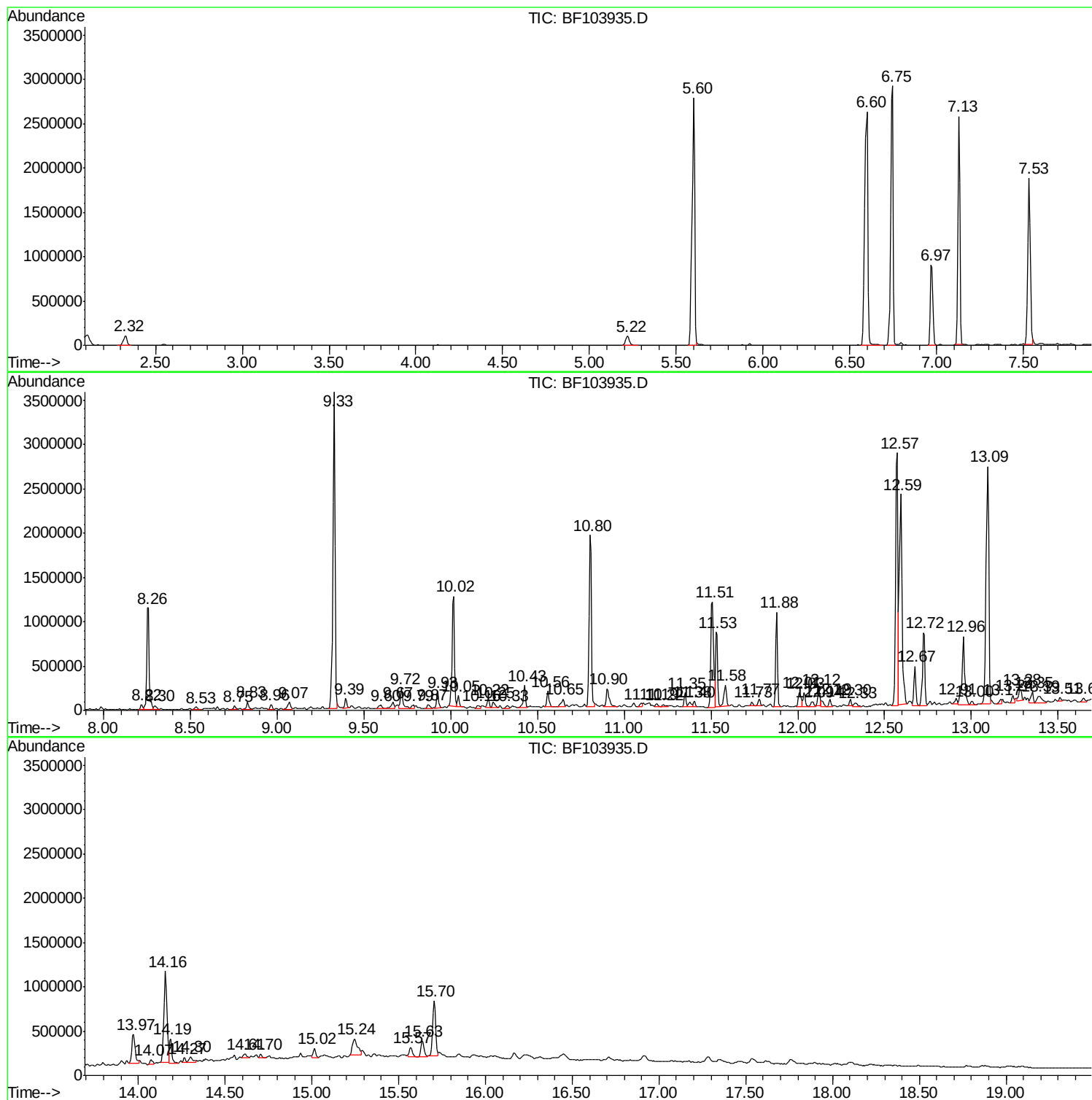
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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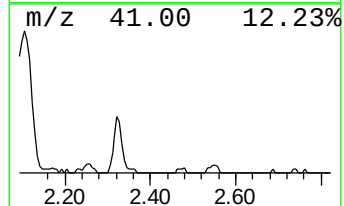
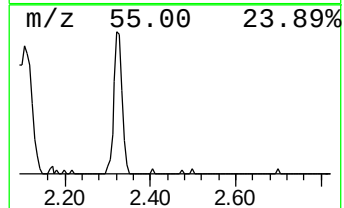
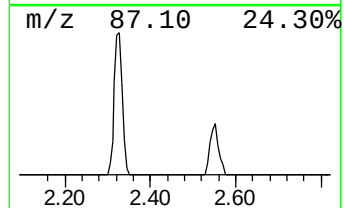
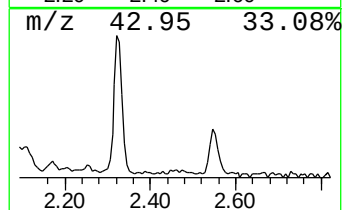
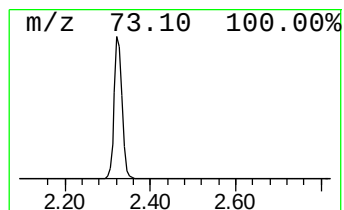
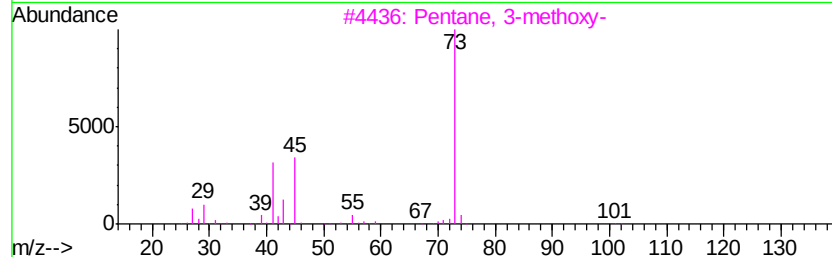
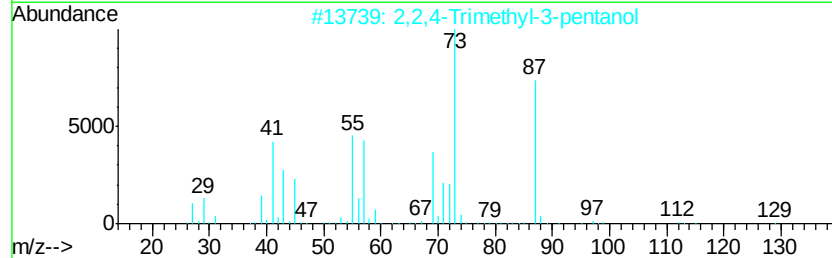
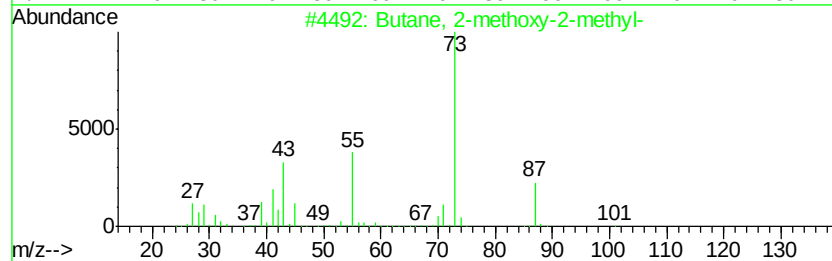
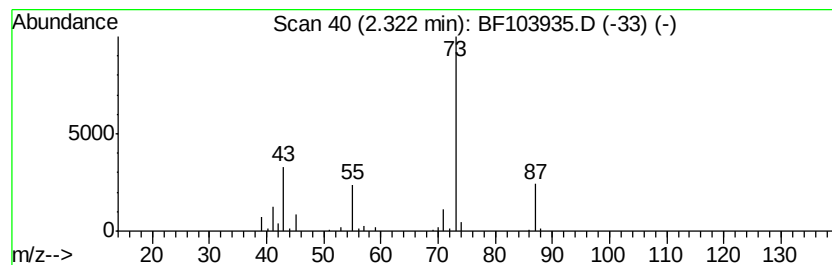
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.32	3.50 ng	143916	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
5			Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9



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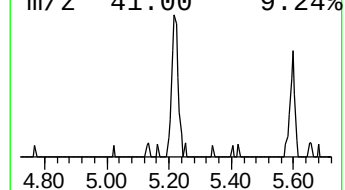
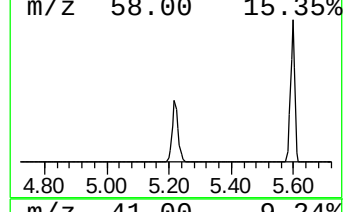
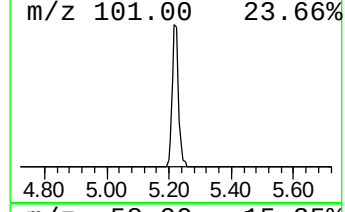
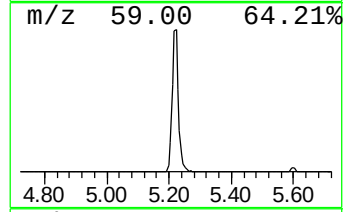
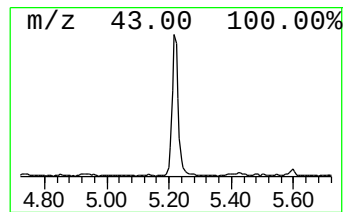
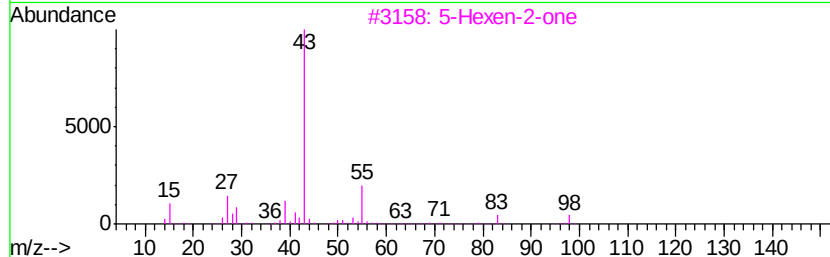
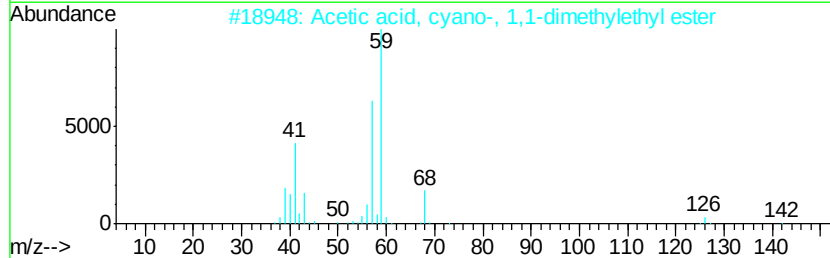
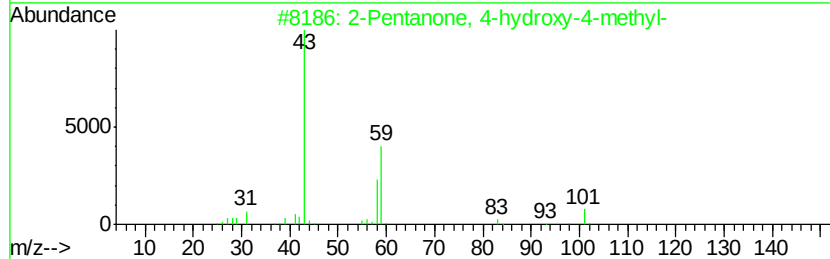
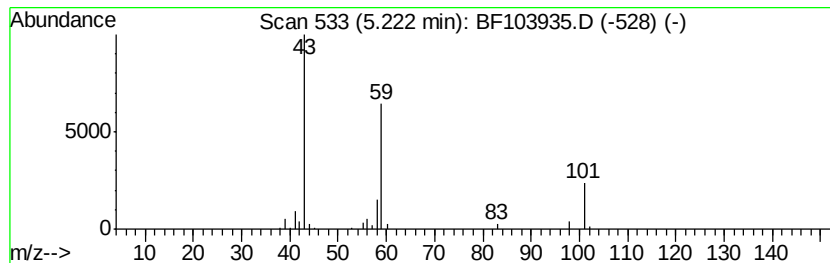
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown5.22 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	3.33 ng	137024	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	40
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3			5-Hexen-2-one	98	C6H10O	000109-49-9	9
4			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
5			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9



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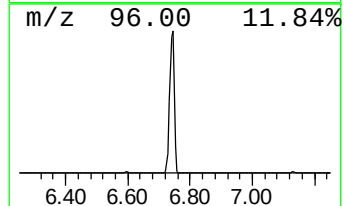
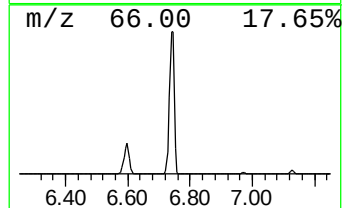
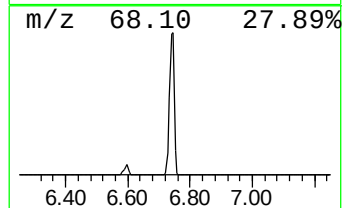
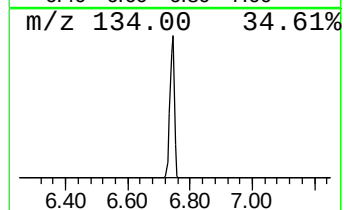
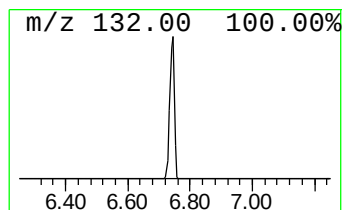
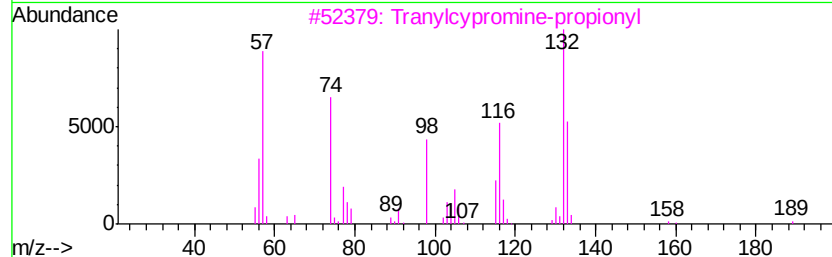
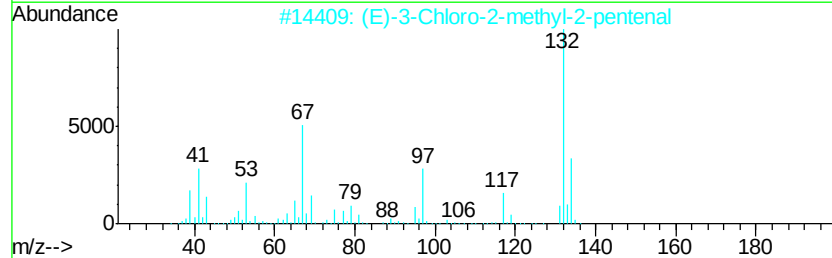
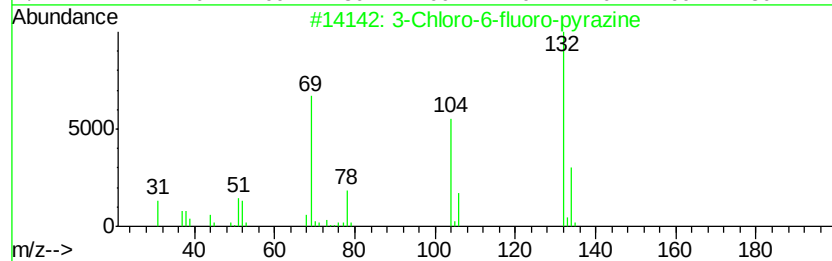
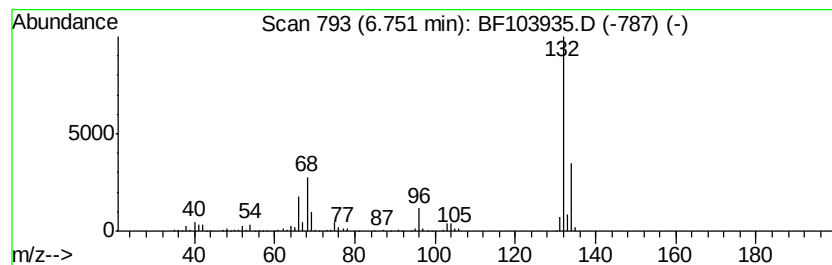
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 unknown6.75 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	70.87 ng	2914430	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	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10



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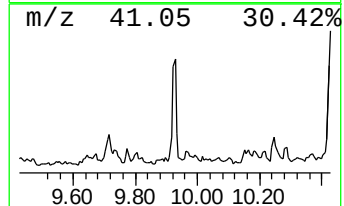
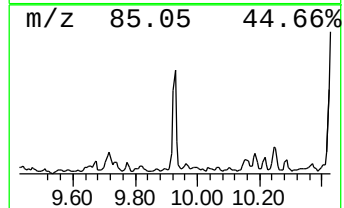
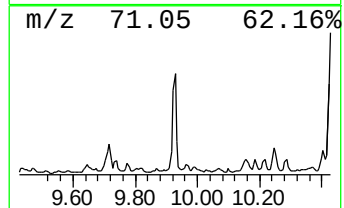
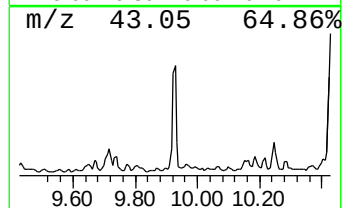
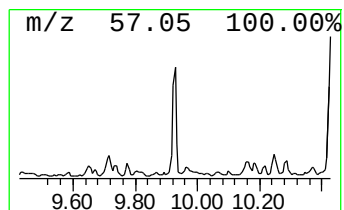
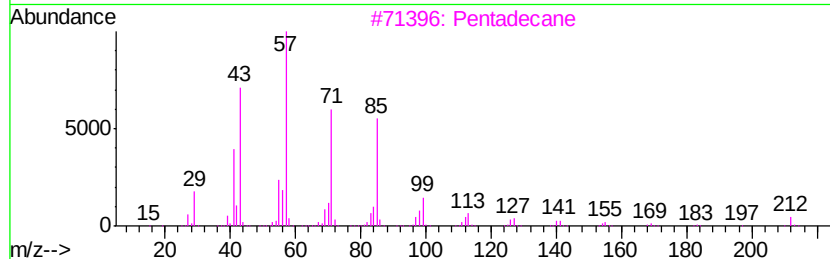
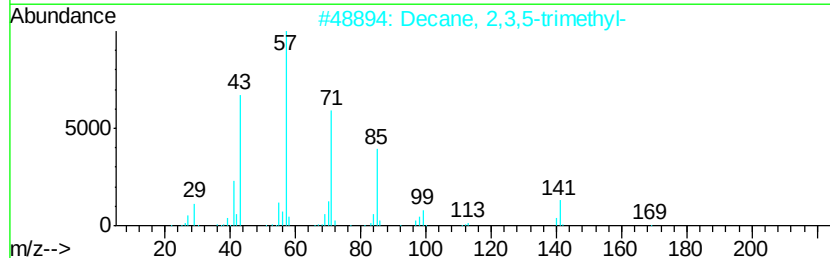
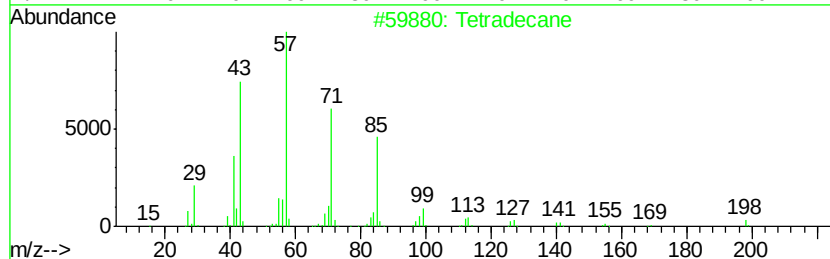
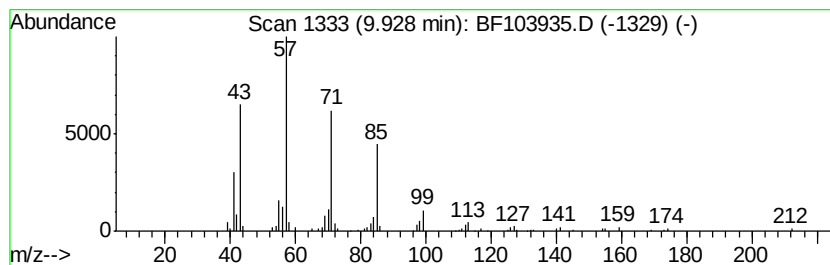
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Tetradecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.93	2.64 ng	156673	Acenaphthene-d10	10.02

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	91
2		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	90
3		Pentadecane	212	C15H32	000629-62-9	80
4		Hexadecane	226	C16H34	000544-76-3	72
5		Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
Data File : BF103935.D
Acq On : 26 Mar 2018 19:55
Operator : JU/SJ
Sample : J1758-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

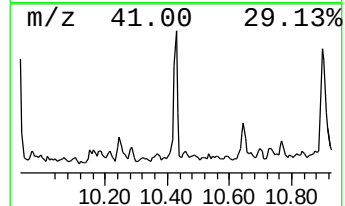
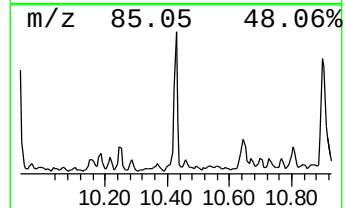
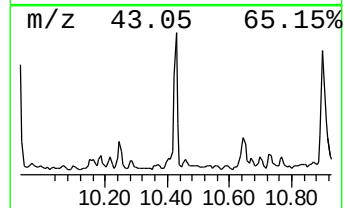
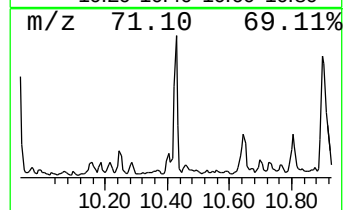
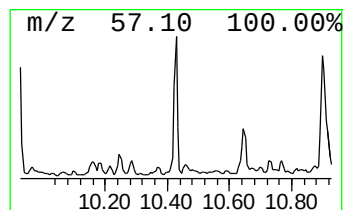
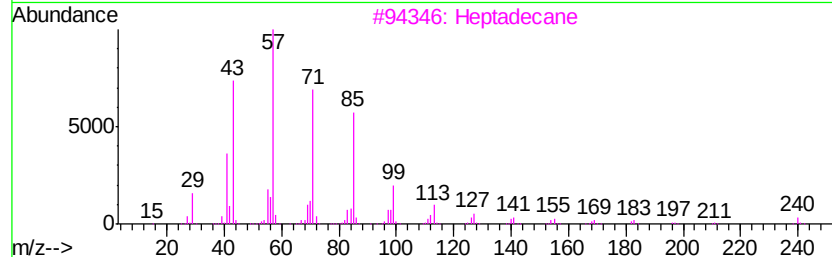
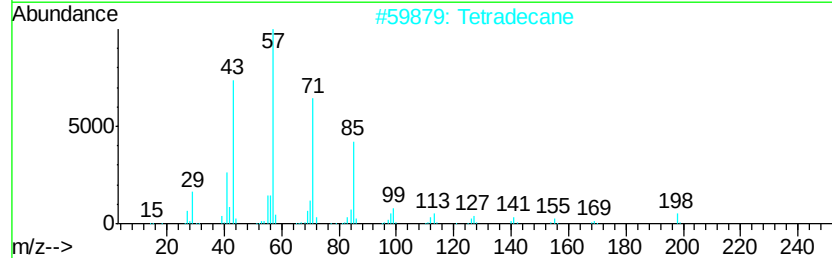
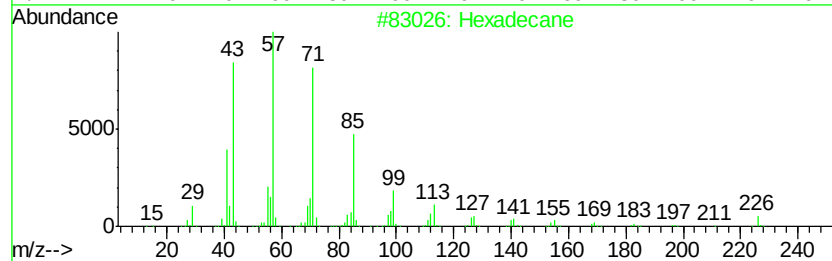
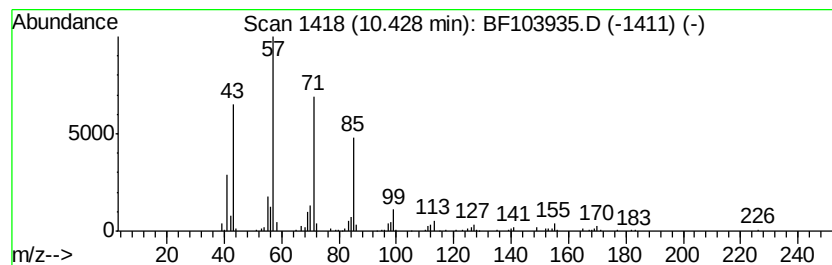
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Hexadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.43	4.03 ng	239695	Acenaphthene-d10		10.02
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	95
2	Tetradecane	198	C14H30	000629-59-4	91
3	Heptadecane	240	C17H36	000629-78-7	91
4	Undecane, 2-methyl-	170	C12H26	007045-71-8	86
5	Dodecane	170	C12H26	000112-40-3	81



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
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Acq On : 26 Mar 2018 19:55
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ClientSampleId :
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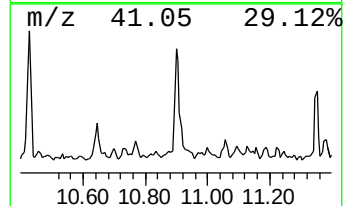
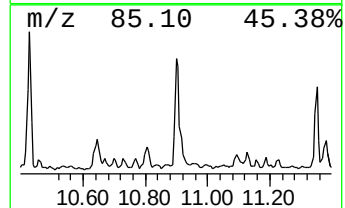
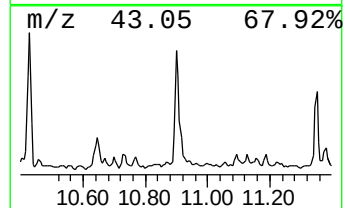
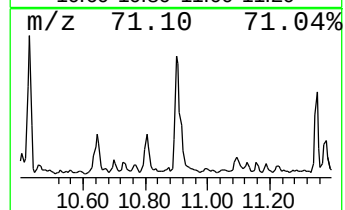
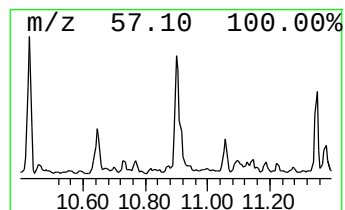
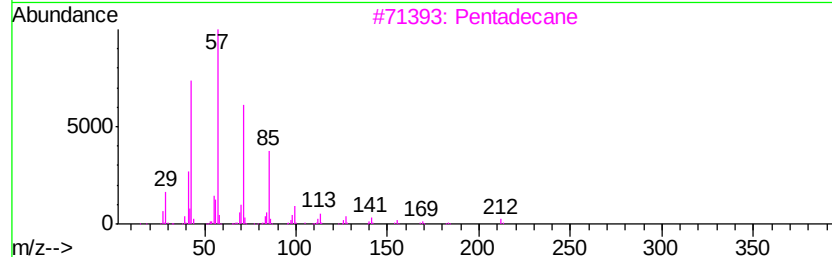
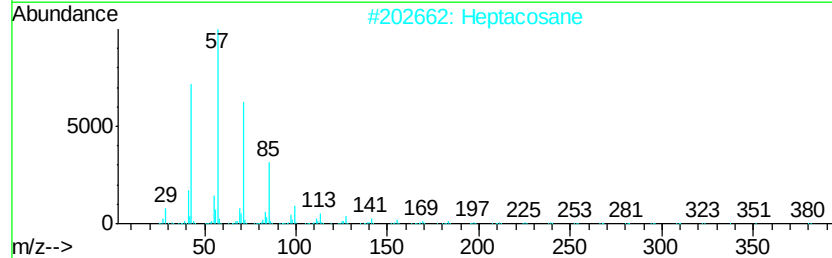
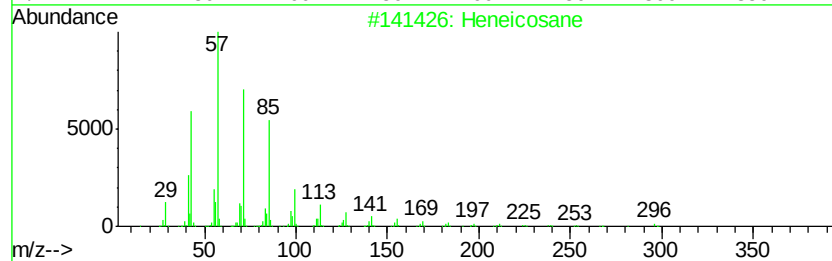
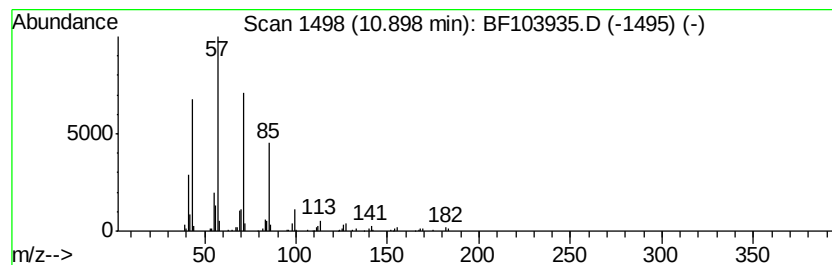
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heneicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.90	4.23 ng	264392	Phenanthrene-d10	11.51

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Heptacosane		380	C27H56	000593-49-7	91
3	Pentadecane		212	C15H32	000629-62-9	90
4	Tridecane, 1-iodo-		310	C13H27I	035599-77-0	86
5	Nonadecane		268	C19H40	000629-92-5	86



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
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Misc :
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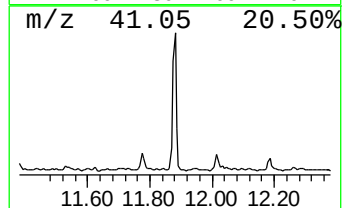
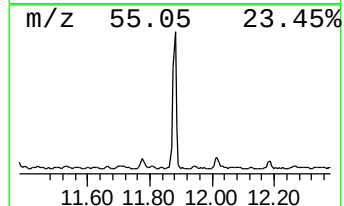
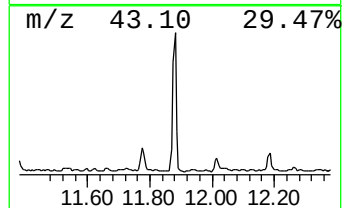
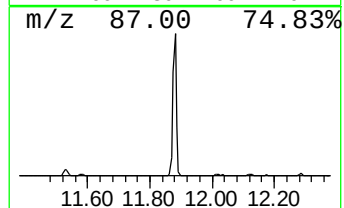
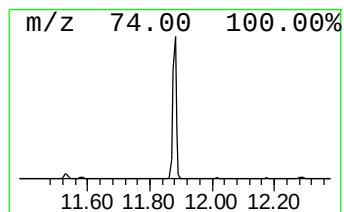
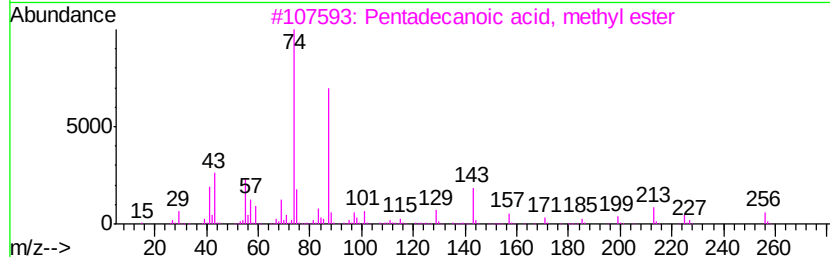
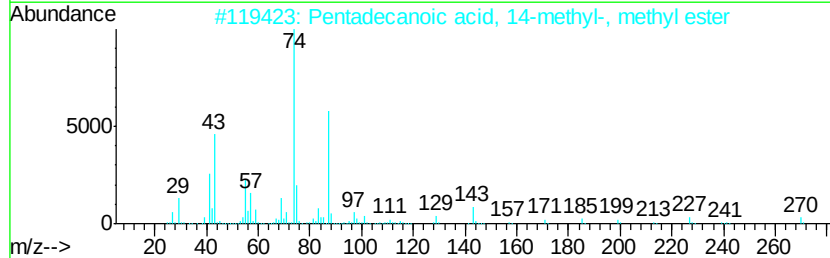
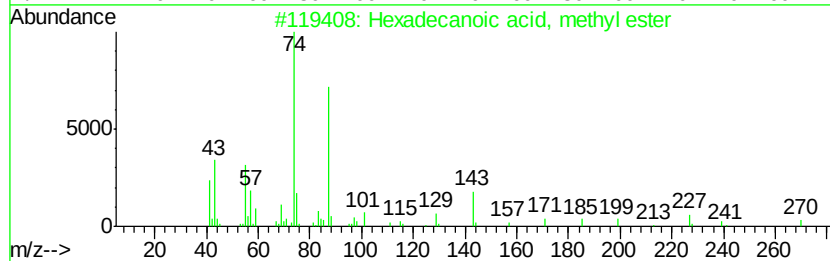
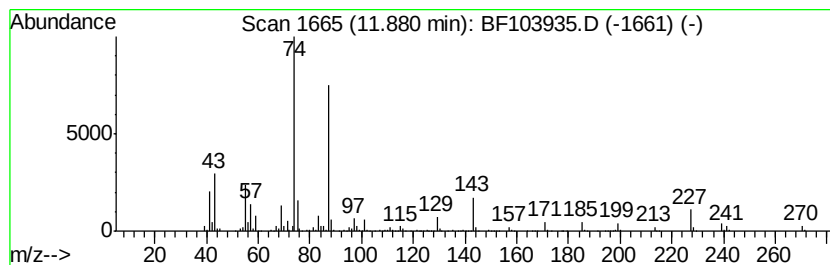
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexadecanoic acid, methyl e... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.88	13.49 ng	842618	Phenanthrene-d10	11.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	92
3		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86
4		Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	86
5		Tridecanoic acid, 12-methyl-, me...	242	C15H30O2	005129-58-8	80



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
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Acq On : 26 Mar 2018 19:55
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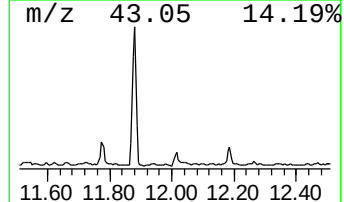
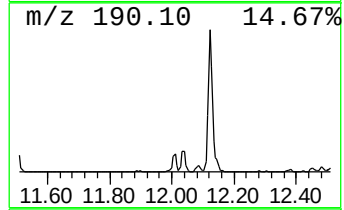
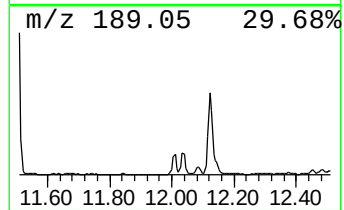
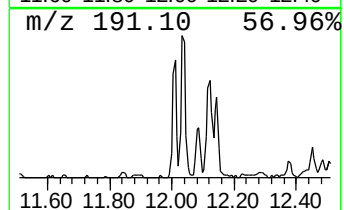
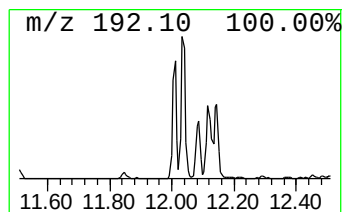
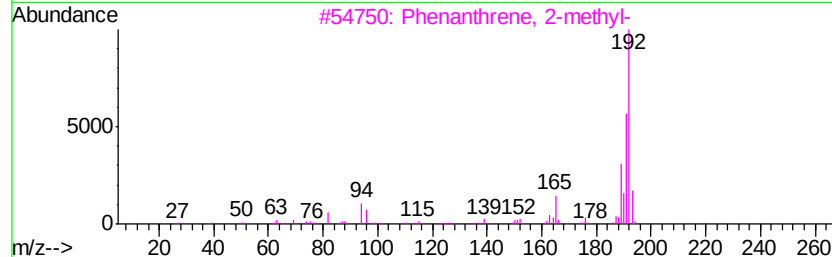
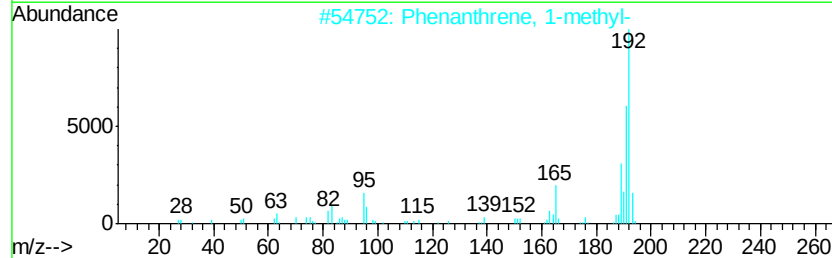
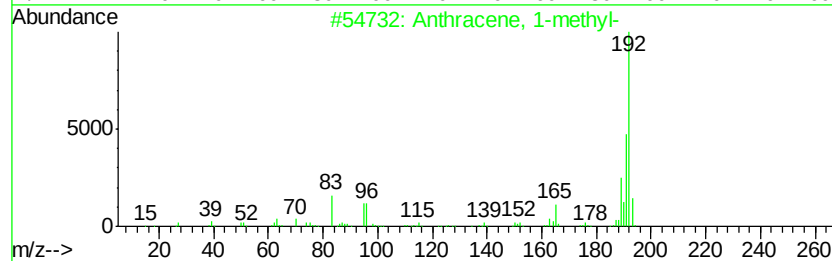
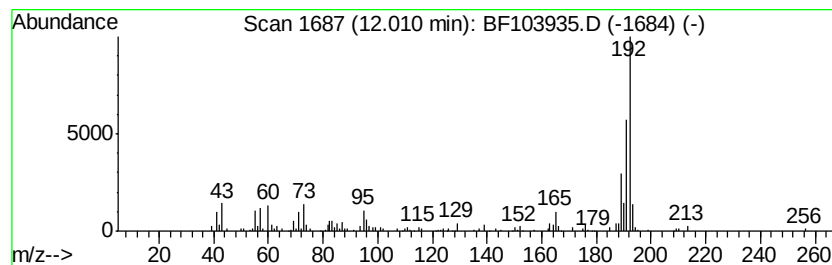
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Anthracene, 1-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	2.67 ng	166682	Phenanthrene-d10	11.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	92
4		Anthracene, 9-methyl-	192	C15H12	000779-02-2	90
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	89



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
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Acq On : 26 Mar 2018 19:55
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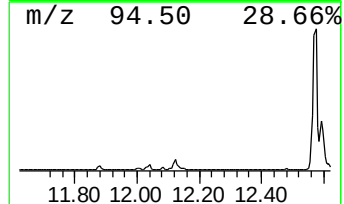
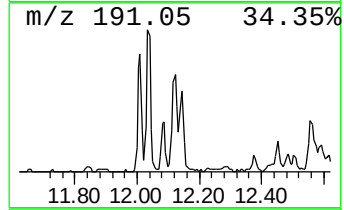
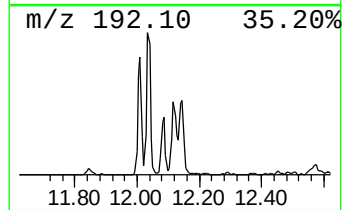
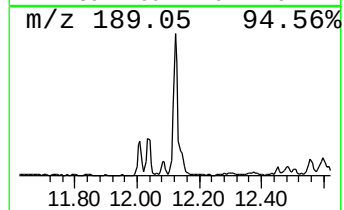
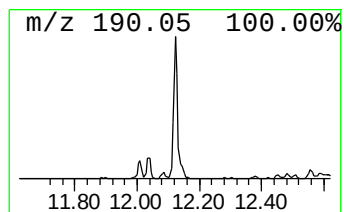
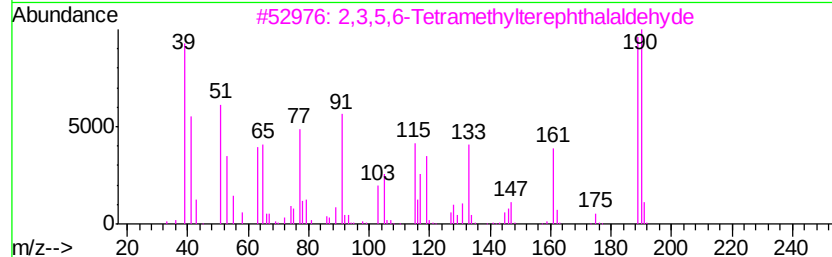
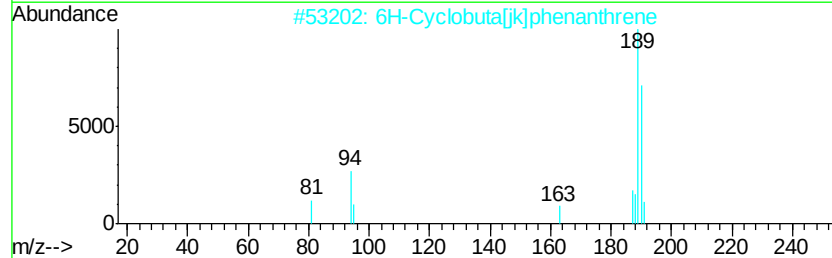
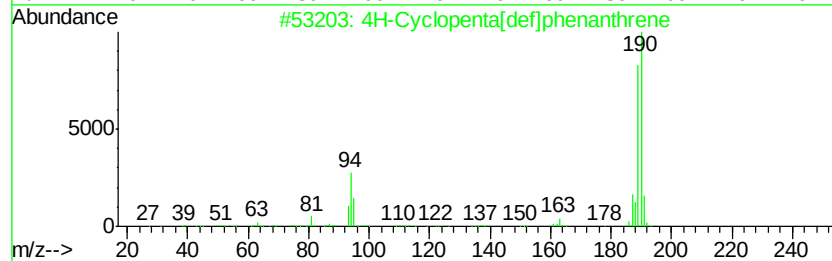
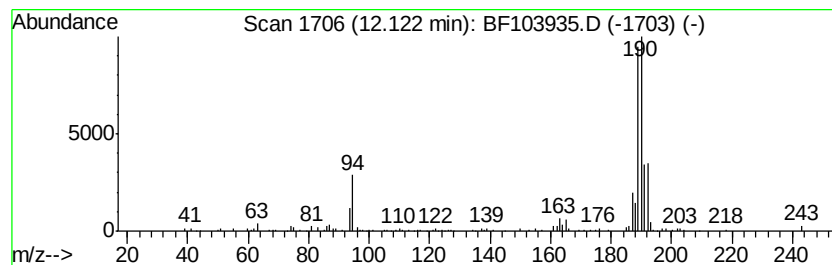
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 4H-Cyclopenta[def]phenanthrene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.12	3.01 ng	188012	Phenanthrene-d10	11.51

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2			6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	59
3			2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	47
4			2,6-Dichlorobenzaloxime	189	C7H5Cl2NO	025185-95-9	38
5			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	38



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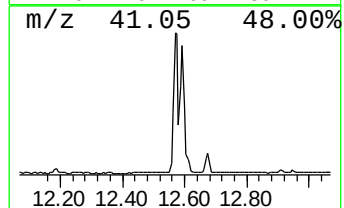
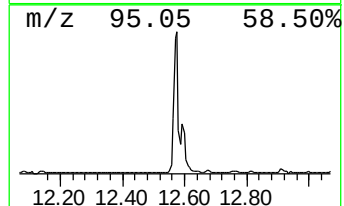
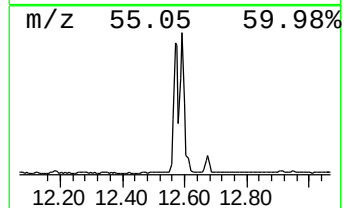
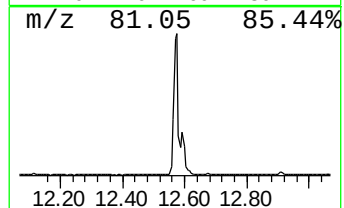
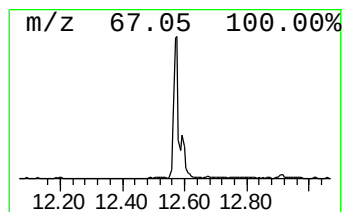
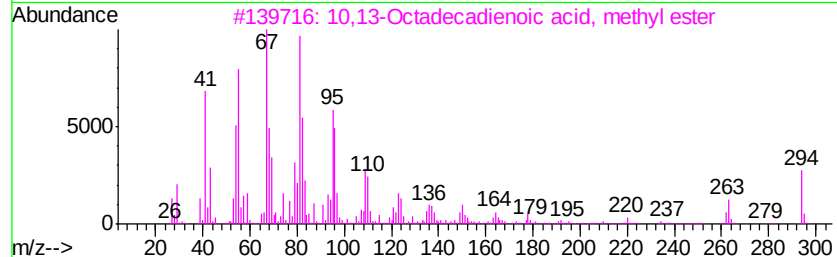
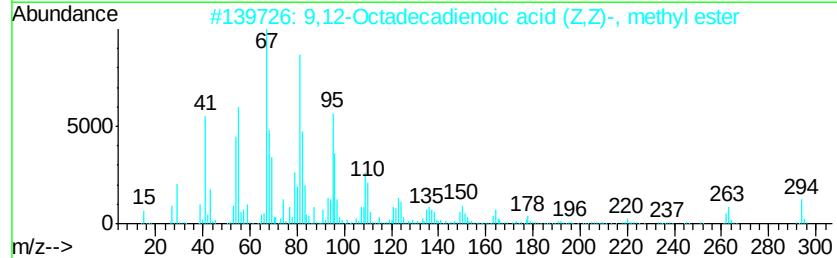
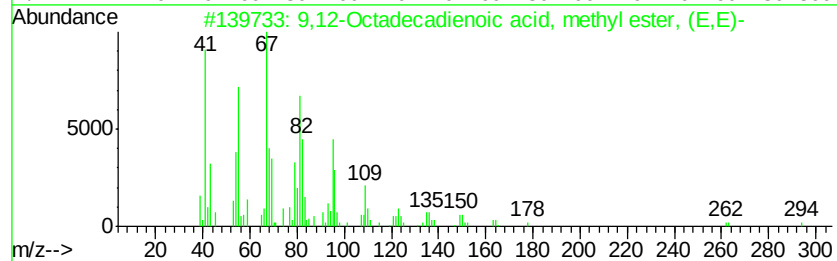
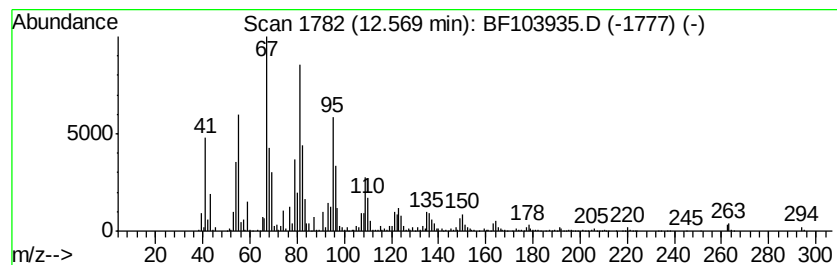
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 9,12-Octadecadienoic acid, ... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.57	45.41 ng	2835390	Phenanthrene-d10	11.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
3		10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99
4		9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99
5		Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	1000336-44-0	99



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
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ClientSampleId :
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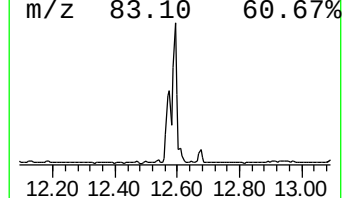
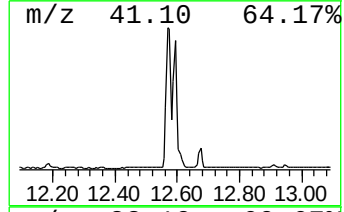
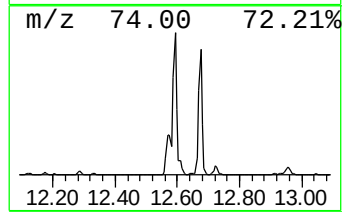
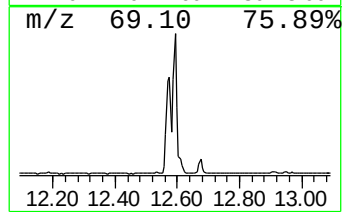
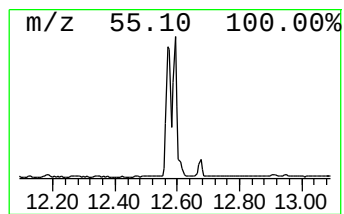
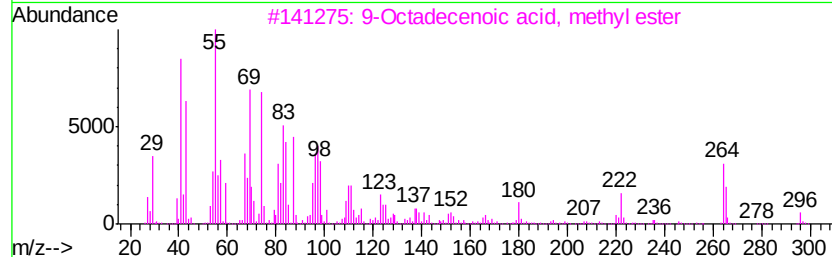
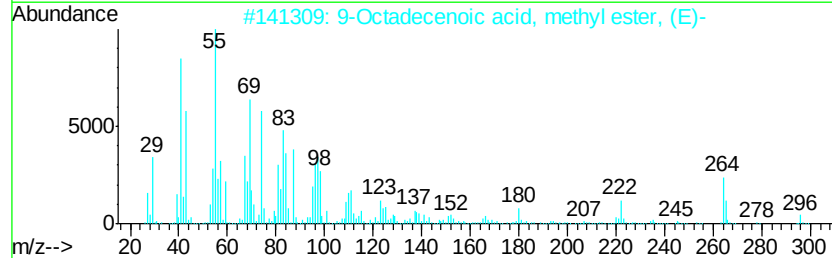
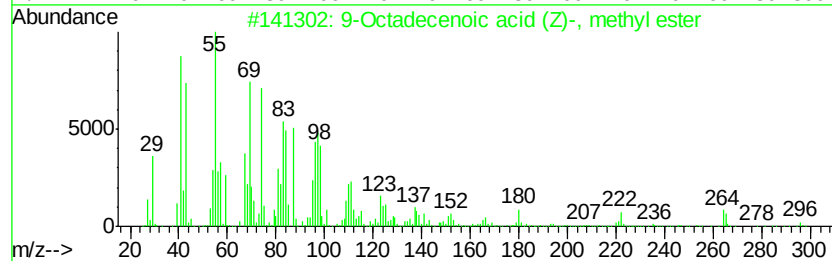
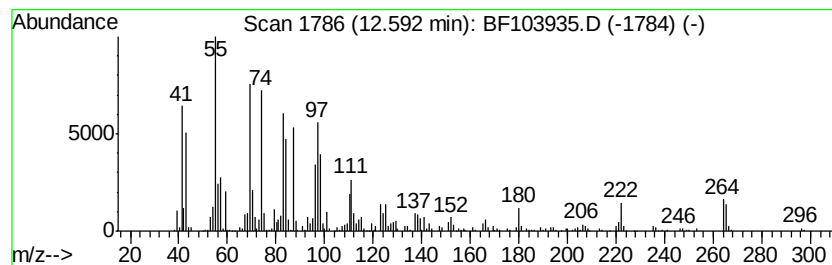
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 9-Octadecenoic acid (Z)-, m... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.59	34.93 ng	2181300	Phenanthrene-d10	11.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	95
2		9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	92
3		9-Octadecenoic acid, methyl ester	296	C19H36O2	002462-84-2	92
4		8-Octadecenoic acid, methyl este...	296	C19H36O2	026528-50-7	89
5		11-Octadecenoic acid, methyl est...	296	C19H36O2	001937-63-9	89



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103935.D
Acq On : 26 Mar 2018 19:55
Operator : JU/SJ
Sample : J1758-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PL-02-032318-A

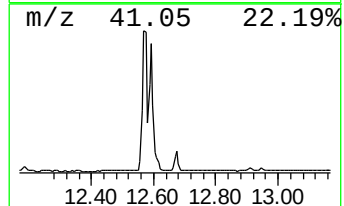
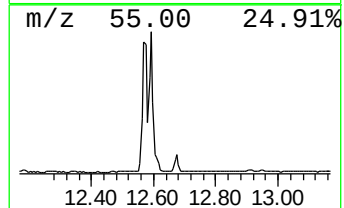
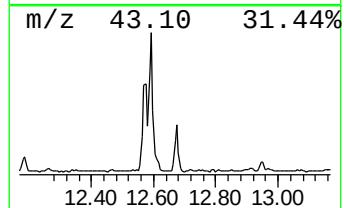
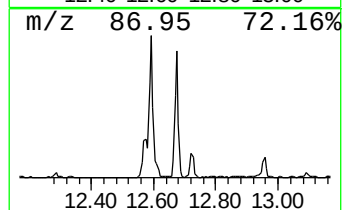
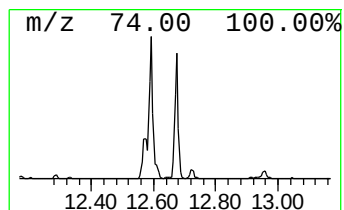
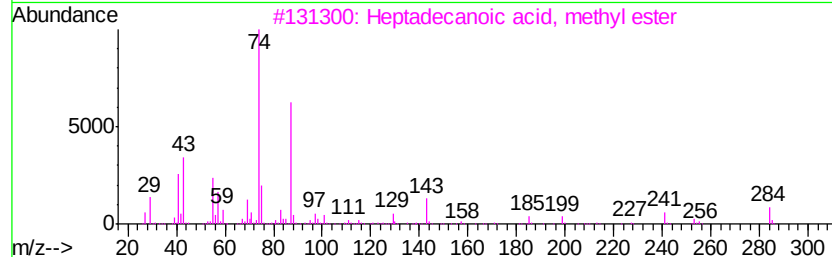
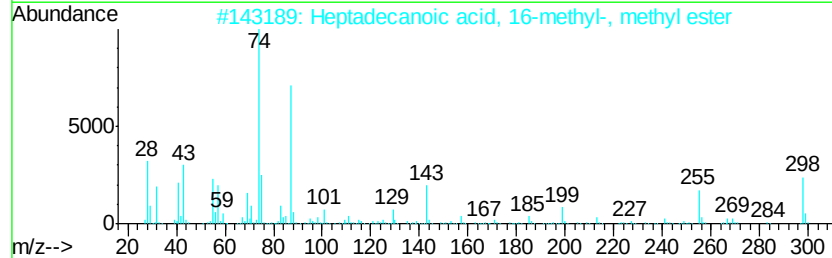
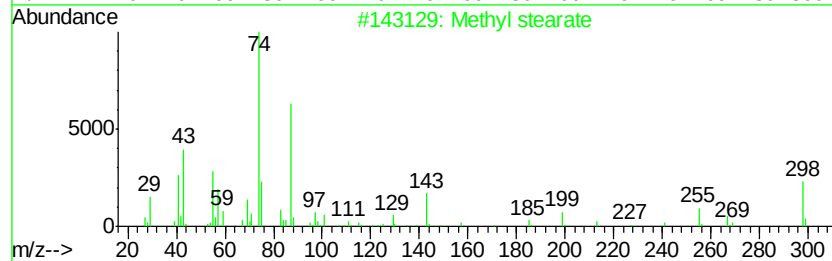
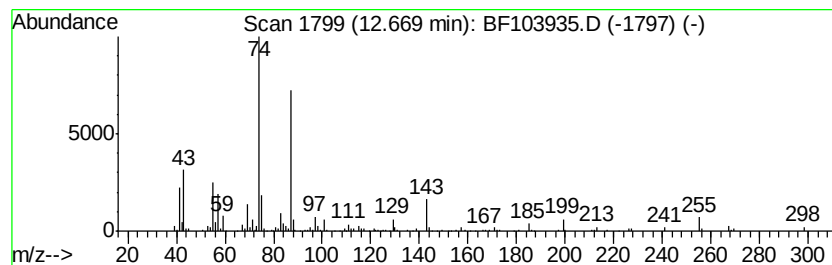
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Methyl stearate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	5.92 ng	369499	Phenanthrene-d10	11.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyl stearate	298	C19H38O2	000112-61-8	97
2		Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	96
3		Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	91
4		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	90
5		Methyl tetradecanoate	242	C15H30O2	000124-10-7	86



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103935.D
Acq On : 26 Mar 2018 19:55
Operator : JU/SJ
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Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
PL-02-032318-A

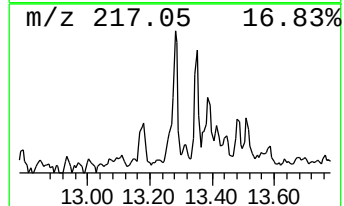
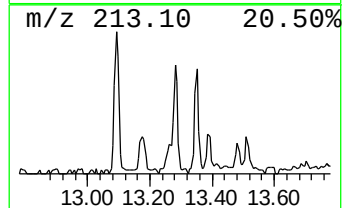
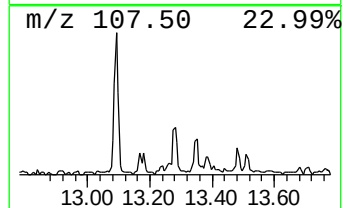
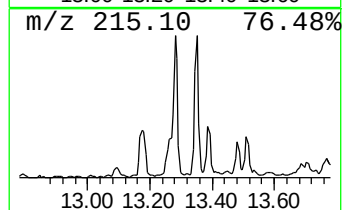
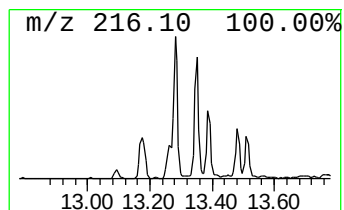
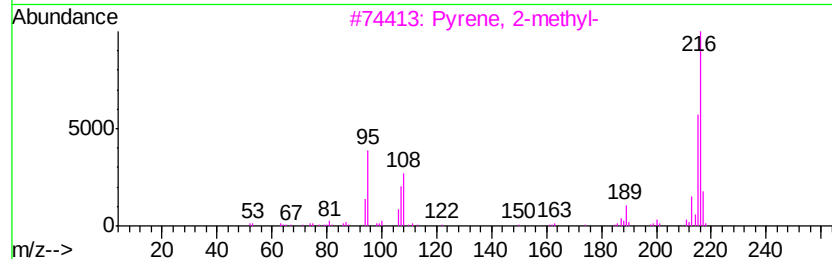
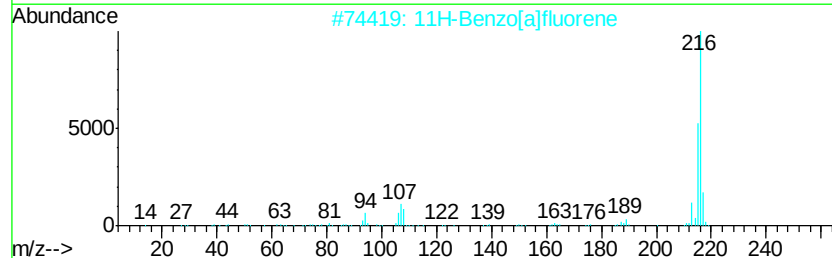
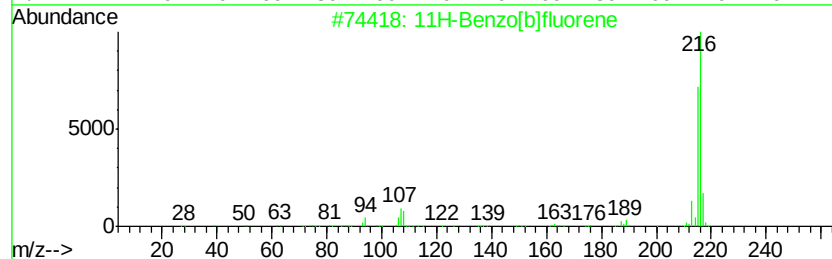
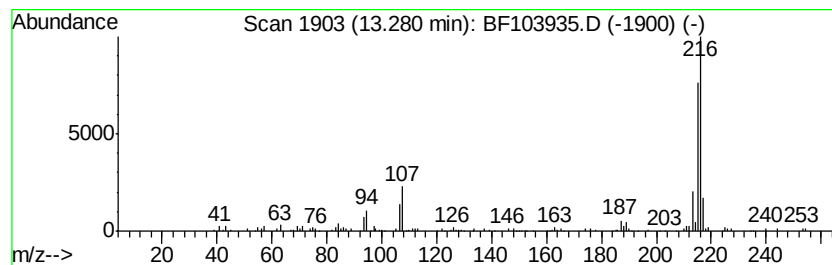
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 11H-Benzo[b]fluorene Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.28	2.05 ng	124578	Chrysene-d12	14.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
3			Pvrene, 2-methyl-	216	C17H12	003442-78-2	87
4			Pvrene, 1-methyl-	216	C17H12	002381-21-7	80
5			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
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Acq On : 26 Mar 2018 19:55
Operator : JU/SJ
Sample : J1758-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

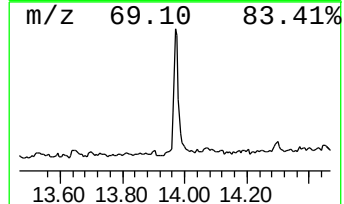
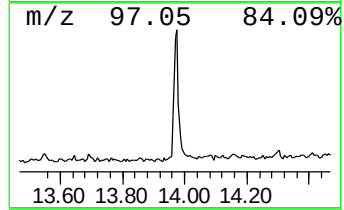
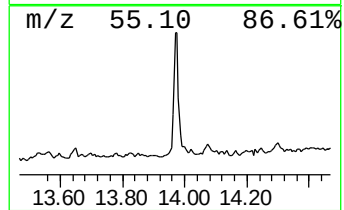
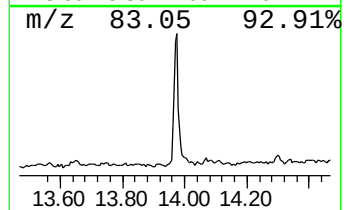
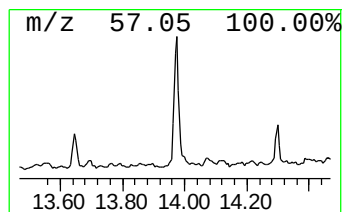
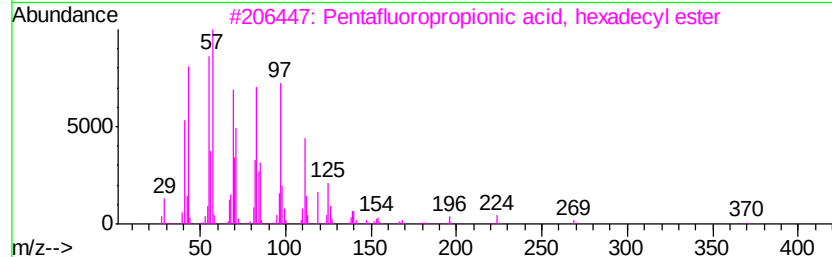
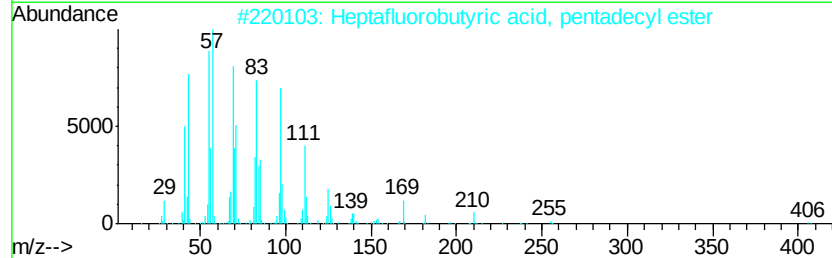
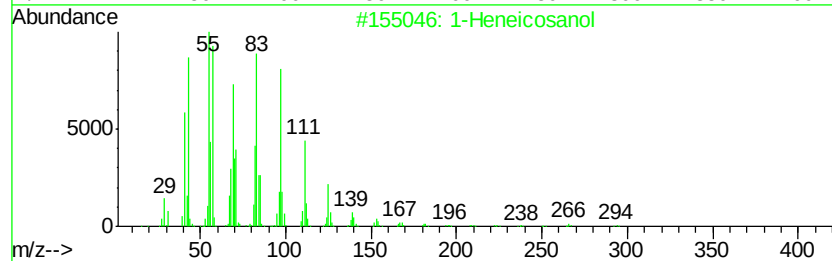
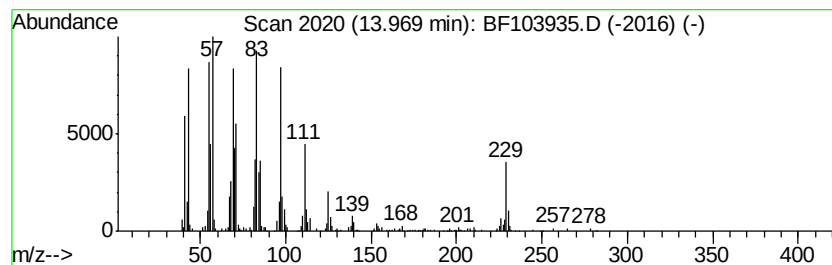
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 1-Heneicosanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.97	6.56 ng	399562	Chrysene-d12		14.16
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	93
2	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	91
3	Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	91
4	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	91
5	Cyclopentadecane	210	C15H30	000295-48-7	90



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103935.D
Acq On : 26 Mar 2018 19:55
Operator : JU/SJ
Sample : J1758-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
PL-02-032318-A

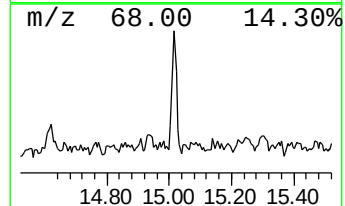
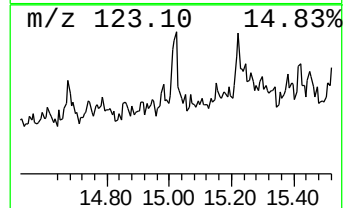
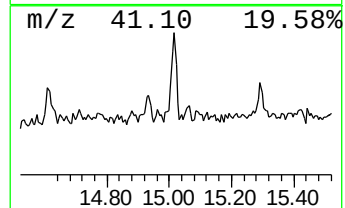
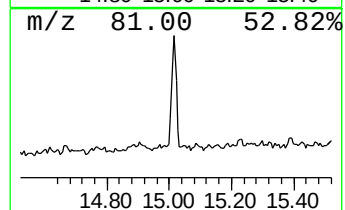
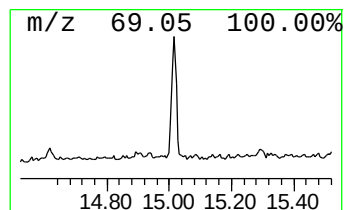
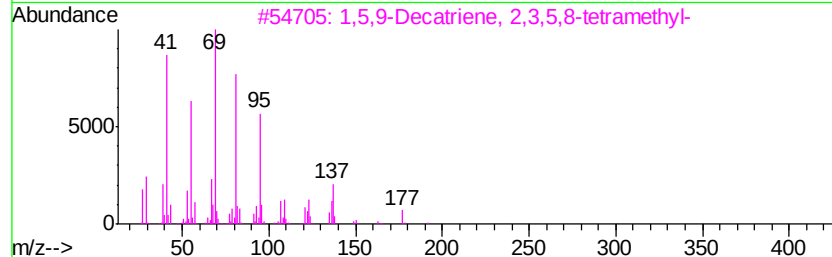
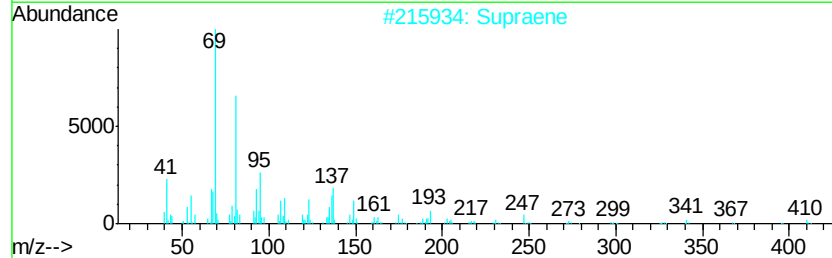
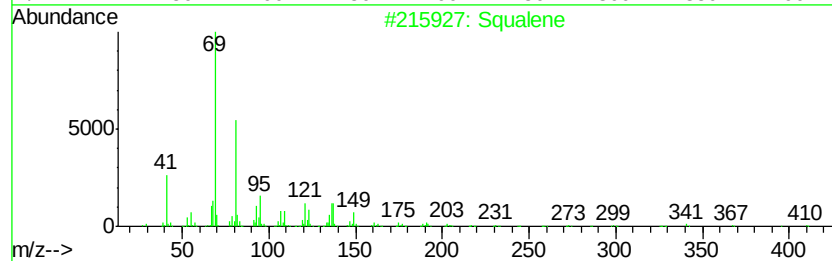
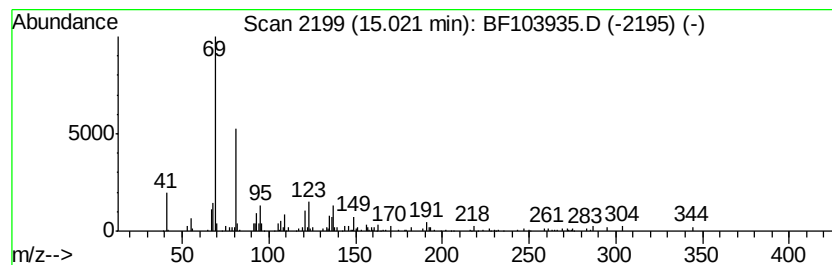
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Squalene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.02	2.57 ng	106811	Perylene-d12	15.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	90
2		Supraene	410	C30H50	007683-64-9	72
3		1,5,9-Decatriene, 2,3,5,8-tetram...	192	C14H24	230646-72-7	72
4		6,10,14-Hexadecatrien-1-ol, 3,7,...	292	C20H36O	036237-66-8	72
5		2,6,10-Dodecatrien-1-ol, 3,7,11-...	264	C17H28O2	004128-17-0	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103935.D
Acq On : 26 Mar 2018 19:55
Operator : JU/SJ
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Misc :
ALS Vial : 8 Sample Multiplier: 1

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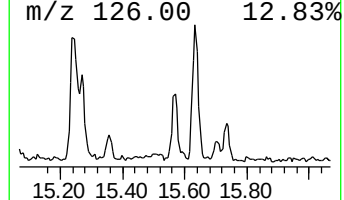
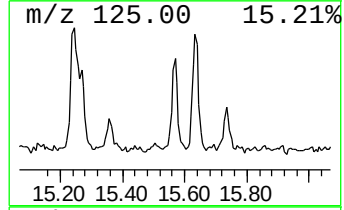
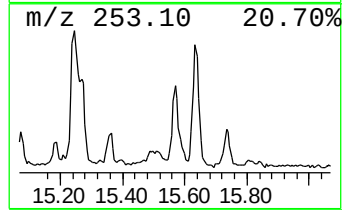
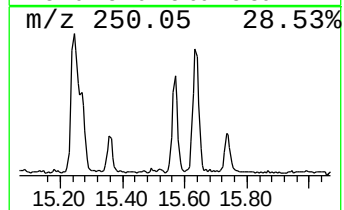
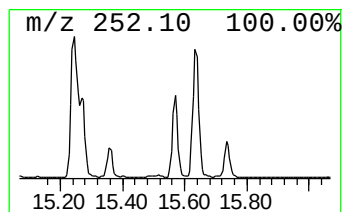
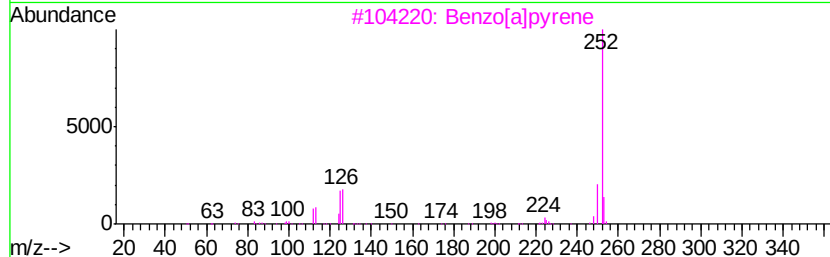
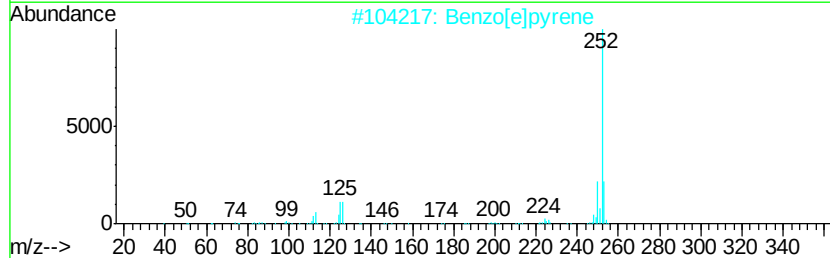
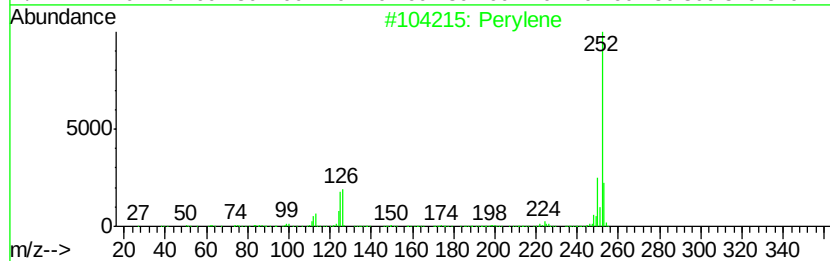
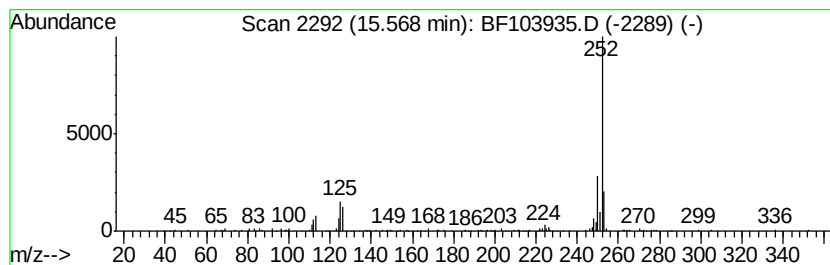
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Perylene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.57	3.69 ng	153594	Perylene-d12	15.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Perylene	252	C20H12	000198-55-0	99
2		Benzo[e]pyrene	252	C20H12	000192-97-2	96
3		Benzo[a]pyrene	252	C20H12	000050-32-8	95
4		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	93
5		Benzo[k]fluoranthene	252	C20H12	000207-08-9	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
 Data File : BF103935.D
 Acq On : 26 Mar 2018 19:55
 Operator : JU/SJ
 Sample : J1758-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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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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unknown5.22	5.22	3.3	ng	137024	1	6.97	822477	20.0
unknown6.75	6.75	70.9	ng	2914430	1	6.97	822477	20.0
Tetradecane	9.93	2.6	ng	156673	3	10.02	1188540	20.0
Hexadecane	10.43	4.0	ng	239695	3	10.02	1188540	20.0
Heneicosane	10.90	4.2	ng	264392	4	11.51	1248800	20.0
Hexadecanoic acid...	11.88	13.5	ng	842618	4	11.51	1248800	20.0
Anthracene, 1-met...	12.01	2.7	ng	166682	4	11.51	1248800	20.0
4H-Cyclopenta[def...	12.12	3.0	ng	188012	4	11.51	1248800	20.0
9,12-Octadecadien...	12.57	45.4	ng	2835390	4	11.51	1248800	20.0
9-Octadecenoic ac...	12.59	34.9	ng	2181300	4	11.51	1248800	20.0
Methyl stearate	12.67	5.9	ng	369499	4	11.51	1248800	20.0
11H-Benzo[b]fluorene	13.28	2.0	ng	124578	5	14.16	1217510	20.0
1-Heneicosanol	13.97	6.6	ng	399562	5	14.16	1217510	20.0
Squalene	15.02	2.6	ng	106811	6	15.70	832100	20.0
Perylene	15.57	3.7	ng	153594	6	15.70	832100	20.0