

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
 Data File : BF103938.D
 Acq On : 26 Mar 2018 21:14
 Operator : JU/SJ
 Sample : J1758-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

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

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	35	40	48	rVB	108876	155824	3.68%	0.368%
2	5.222	526	533	540	rBV	114160	152992	3.61%	0.362%
3	5.604	592	598	601	rBV	2715000	2966102	69.96%	7.010%
4	6.598	761	767	770	rBV	2858719	3003869	70.85%	7.100%
5	6.745	787	792	795	rBV	3260376	3062067	72.23%	7.237%
6	6.975	827	831	834	rBV	1061150	880128	20.76%	2.080%
7	7.128	853	857	861	rBV	2516059	2375498	56.03%	5.614%
8	7.540	922	927	934	rBV	1815228	1948398	45.96%	4.605%
9	8.222	1040	1043	1045	rBV	66469	59197	1.40%	0.140%
10	8.257	1045	1049	1051	rVV	1380600	1171534	27.63%	2.769%
11	8.275	1051	1052	1054	rVB	107319	61143	1.44%	0.145%
12	8.828	1143	1146	1151	rVB	92134	80455	1.90%	0.190%
13	8.969	1167	1170	1173	rBV	64696	54547	1.29%	0.129%
14	9.069	1184	1187	1192	rBV2	74224	67009	1.58%	0.158%
15	9.328	1226	1231	1235	rBV	3547440	3341054	78.81%	7.897%
16	9.398	1240	1243	1246	rVB	115369	103918	2.45%	0.246%
17	9.598	1272	1277	1281	rBV5	30324	50092	1.18%	0.118%
18	9.669	1281	1289	1292	rVV4	58844	97258	2.29%	0.230%
19	9.716	1294	1297	1305	rVB	250917	261715	6.17%	0.619%
20	9.928	1328	1333	1337	rBV	222475	179672	4.24%	0.425%
21	10.016	1344	1348	1351	rBV	1500171	1277418	30.13%	3.019%
22	10.045	1351	1353	1356	rVB	95338	73466	1.73%	0.174%
23	10.216	1379	1382	1385	rVB2	82786	75559	1.78%	0.179%
24	10.251	1385	1388	1392	rVB2	52486	55578	1.31%	0.131%
25	10.428	1415	1418	1421	rVB	266036	205240	4.84%	0.485%
26	10.563	1438	1441	1447	rVB	142141	126620	2.99%	0.299%
27	10.645	1450	1455	1458	rBV2	88199	109047	2.57%	0.258%
28	10.810	1478	1483	1486	rBV	2095276	2043401	48.20%	4.830%
29	10.898	1495	1498	1508	rVB	213180	276863	6.53%	0.654%
30	11.145	1538	1540	1545	rVB4	36774	44787	1.06%	0.106%
31	11.351	1572	1575	1577	rBV	173300	133321	3.14%	0.315%
32	11.380	1577	1580	1582	rVB2	49441	43442	1.02%	0.103%
33	11.510	1598	1602	1604	rBV	1427492	1259338	29.70%	2.976%
34	11.533	1604	1606	1610	rVV	578281	463016	10.92%	1.094%

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Integration Parameters: rteint.p
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35	11.580	1610	1614	1618	rVB	122167	117056	2.76%	0.277%
36	11.733	1638	1640	1645	rVB	46536	44288	1.04%	0.105%
37	11.780	1645	1648	1650	rBV	95988	77899	1.84%	0.184%
38	11.880	1661	1665	1669	rVB	1479959	1185907	27.97%	2.803%
39	12.016	1684	1688	1690	rBV2	116868	137345	3.24%	0.325%
40	12.039	1690	1692	1696	rVB2	89380	82300	1.94%	0.195%
41	12.122	1703	1706	1709	rBV2	89566	106114	2.50%	0.251%
42	12.186	1713	1717	1720	rBV2	80329	78483	1.85%	0.185%
43	12.574	1778	1783	1785	rBV	4029333	4239579	100.00%	10.020%
44	12.592	1785	1786	1793	rVV2	2861243	2566540	60.54%	6.066%
45	12.674	1797	1800	1804	rVB	644618	498077	11.75%	1.177%
46	12.727	1805	1809	1812	rBV	421694	363796	8.58%	0.860%
47	12.910	1837	1840	1843	rBV2	74382	77613	1.83%	0.183%
48	12.957	1843	1848	1854	rVB	349094	405858	9.57%	0.959%
49	13.098	1867	1872	1875	rBV	2806010	2654311	62.61%	6.273%
50	13.169	1881	1884	1887	rBV3	30105	42916	1.01%	0.101%
51	13.239	1893	1896	1898	rBV	123957	111242	2.62%	0.263%
52	13.286	1901	1904	1906	rVB	69412	69355	1.64%	0.164%
53	13.351	1913	1915	1918	rVB	62121	47994	1.13%	0.113%
54	13.398	1920	1923	1926	rVB2	65070	76448	1.80%	0.181%
55	13.974	2017	2021	2025	rBV2	432671	420776	9.92%	0.994%
56	14.074	2035	2038	2042	rBV2	56919	64266	1.52%	0.152%
57	14.116	2042	2045	2048	rVV	220600	193011	4.55%	0.456%
58	14.163	2048	2053	2056	rVV	1045687	1113475	26.26%	2.632%
59	14.186	2056	2057	2065	rVB	136129	101796	2.40%	0.241%
60	14.298	2074	2076	2080	rBV2	37757	46202	1.09%	0.109%
61	15.015	2196	2198	2202	rVB	103192	104777	2.47%	0.248%
62	15.245	2234	2237	2241	rBV	77442	132565	3.13%	0.313%
63	15.639	2301	2304	2310	rVB	68535	86892	2.05%	0.205%
64	15.710	2311	2316	2321	rBV	596988	803891	18.96%	1.900%

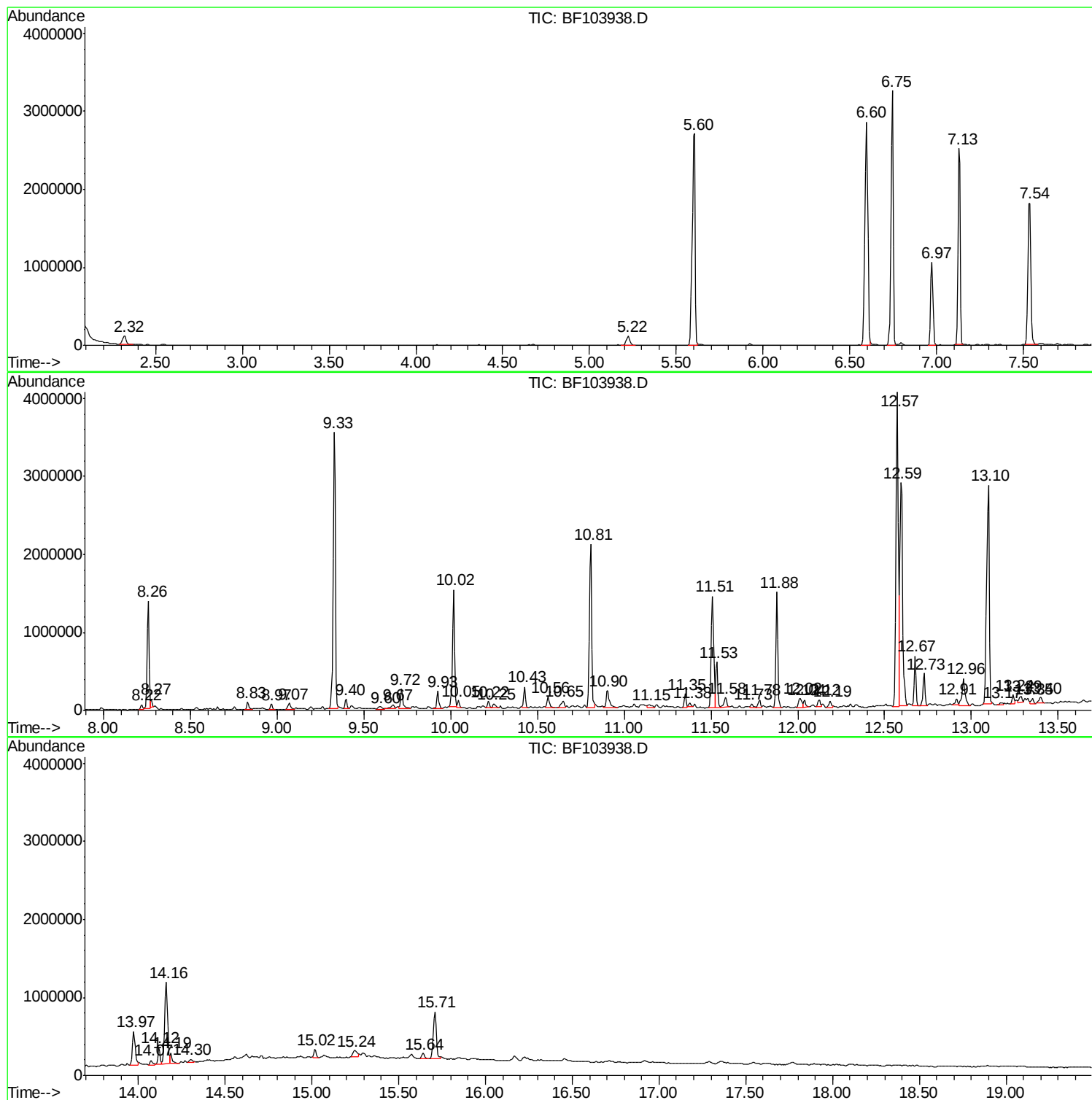
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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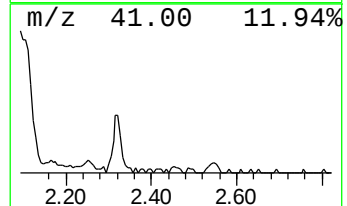
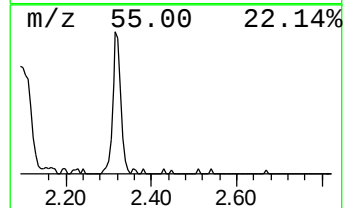
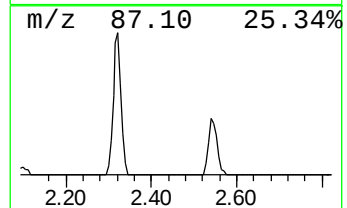
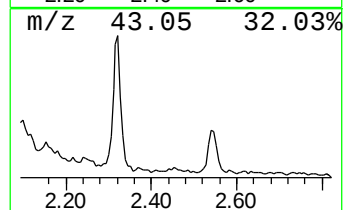
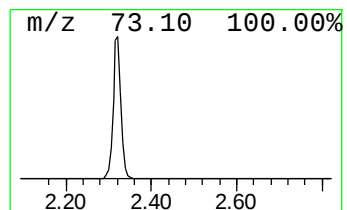
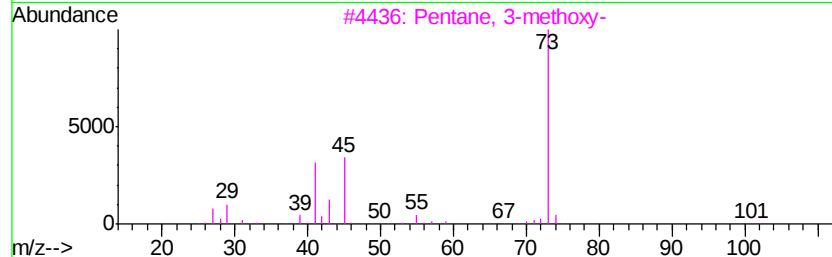
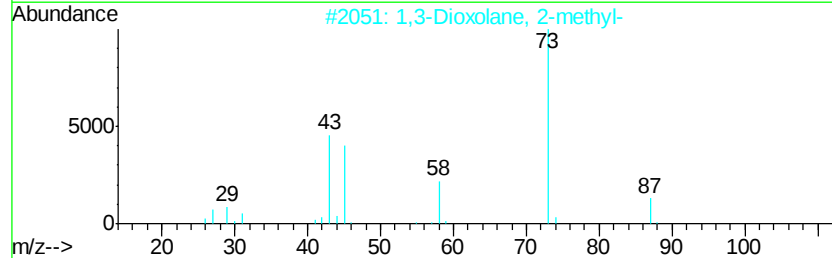
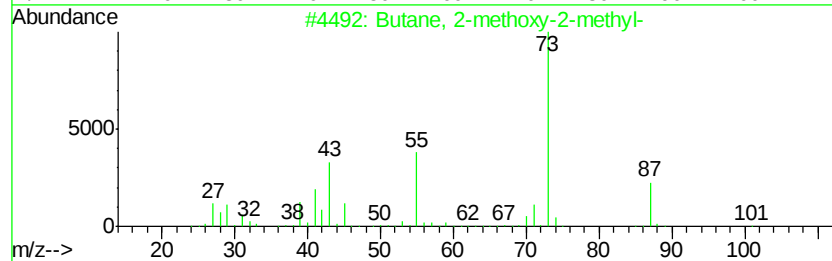
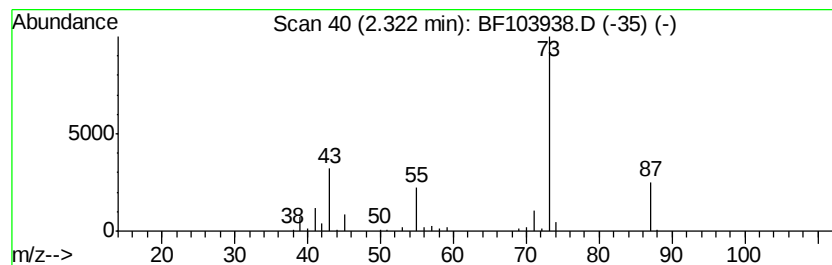
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.32	3.54 ng	155824	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	64
2			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	17
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	17



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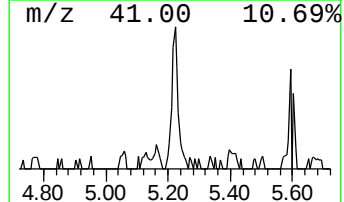
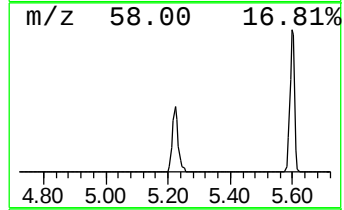
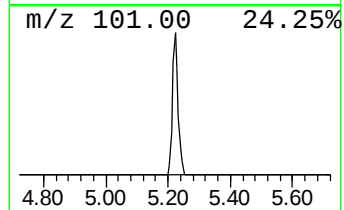
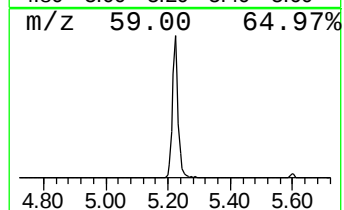
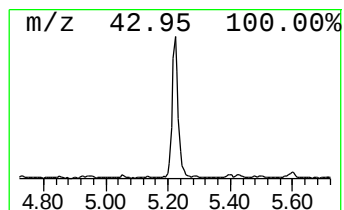
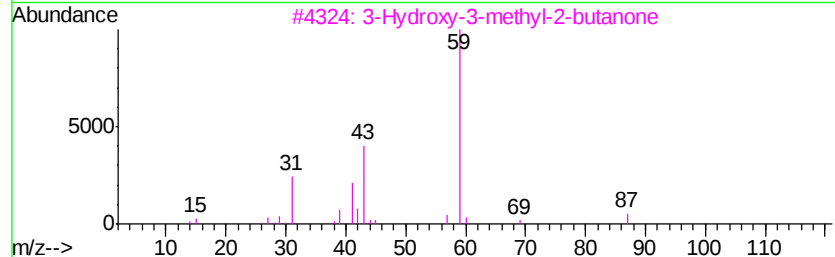
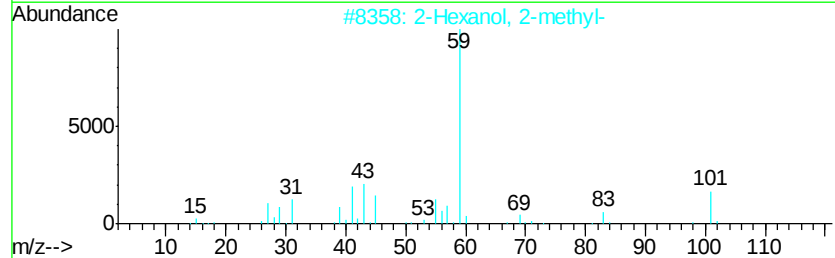
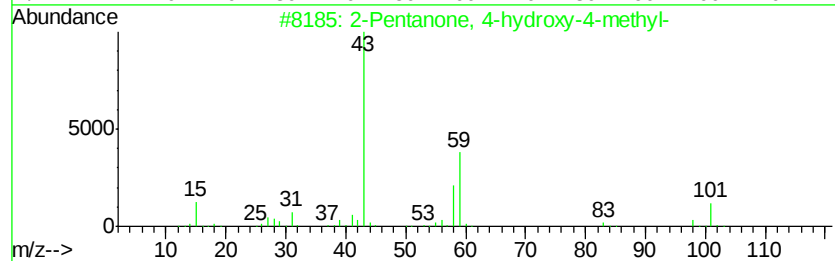
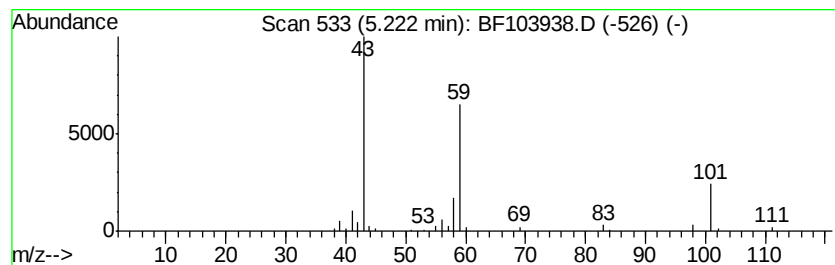
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	3.48 ng	152992	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	38
3		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	35
4		1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	32
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25



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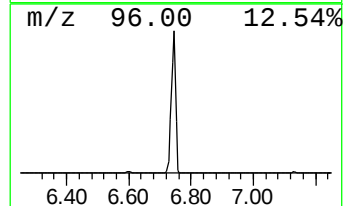
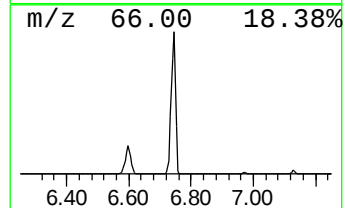
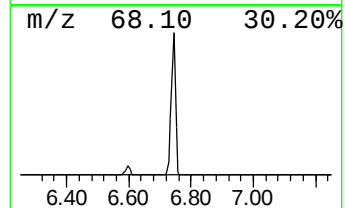
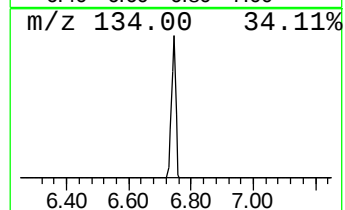
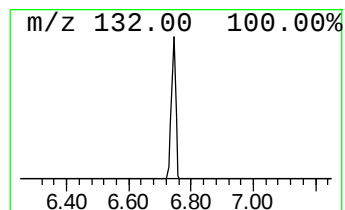
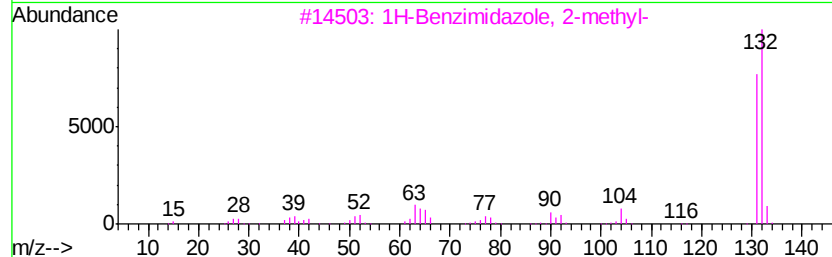
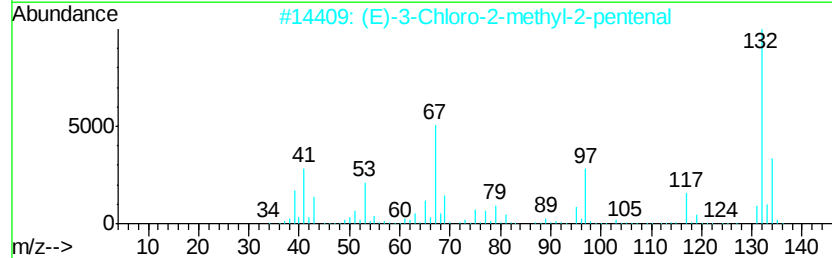
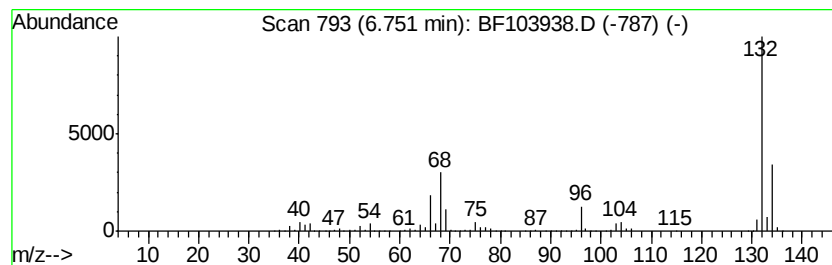
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.75 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	69.58 ng	3062070	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3			1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
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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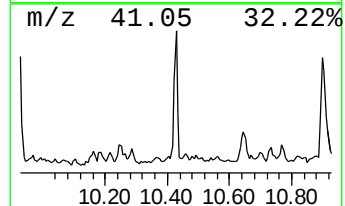
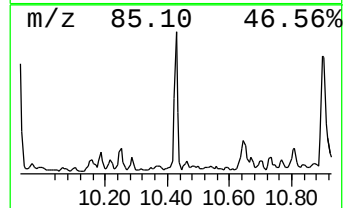
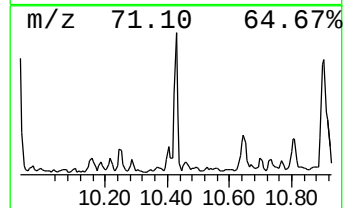
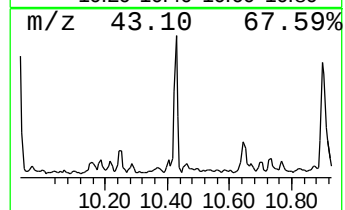
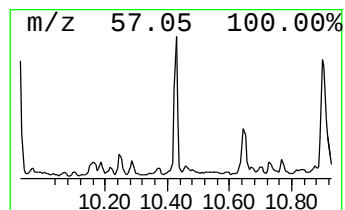
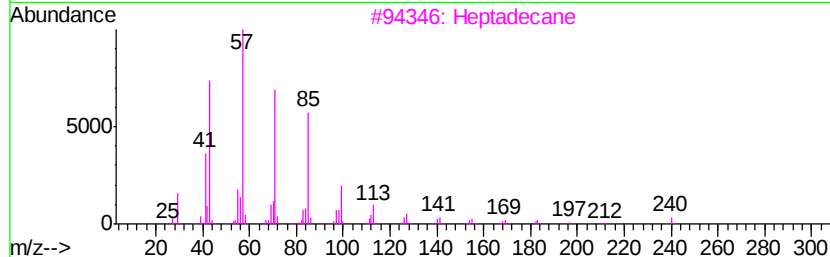
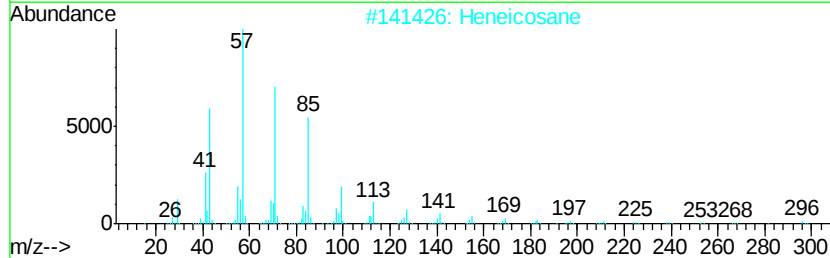
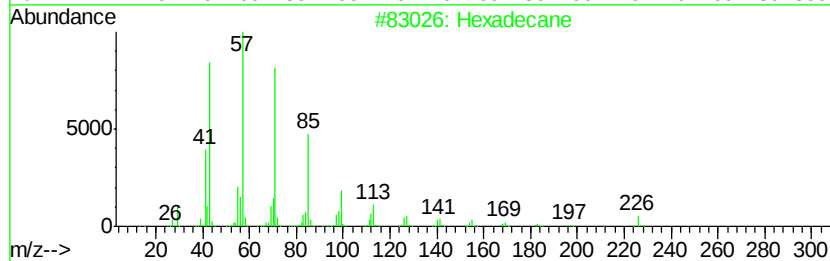
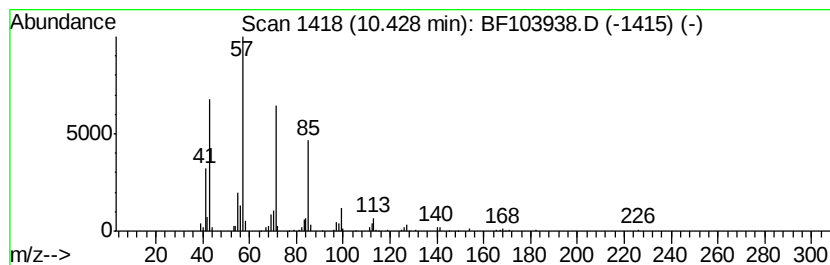
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.43	3.21 ng	205240	Acenaphthene-d10	10.02
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane	226 C16H34	000544-76-3	95
2	Heneicosane	296 C21H44	000629-94-7	90
3	Heptadecane	240 C17H36	000629-78-7	90
4	Tetradecane	198 C14H30	000629-59-4	87
5	Pentadecane	212 C15H32	000629-62-9	86



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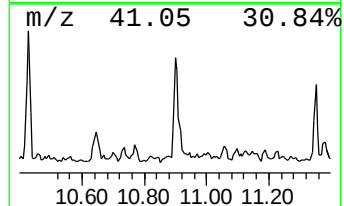
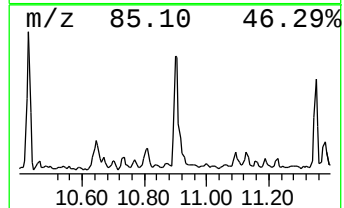
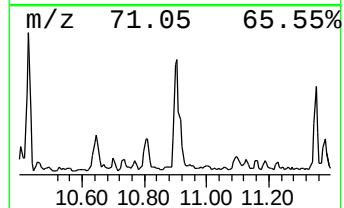
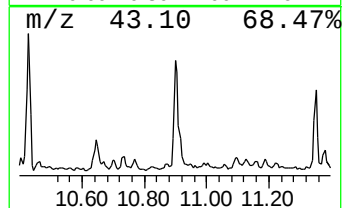
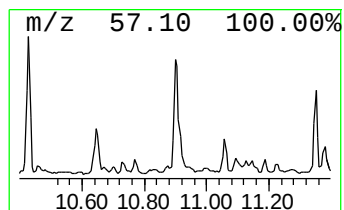
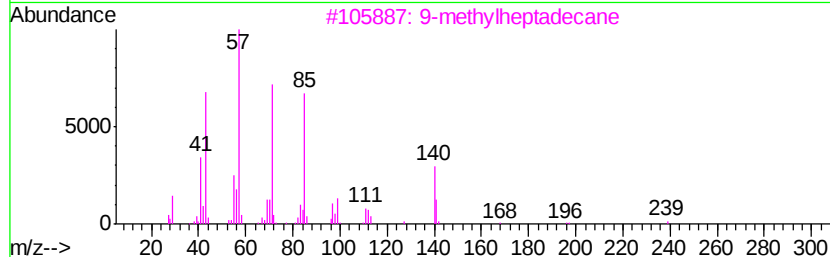
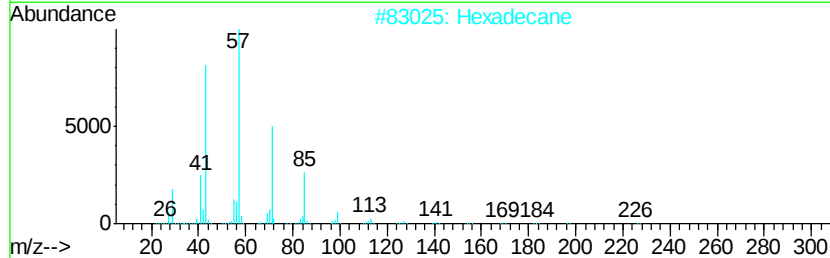
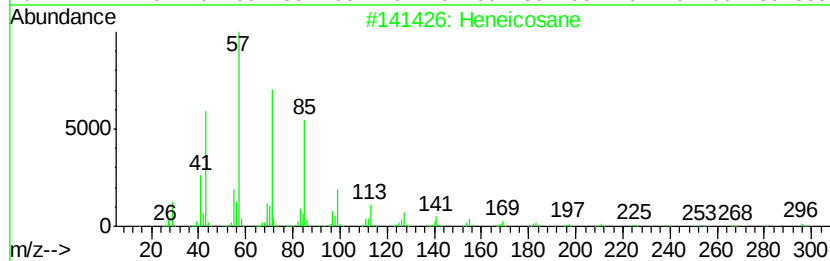
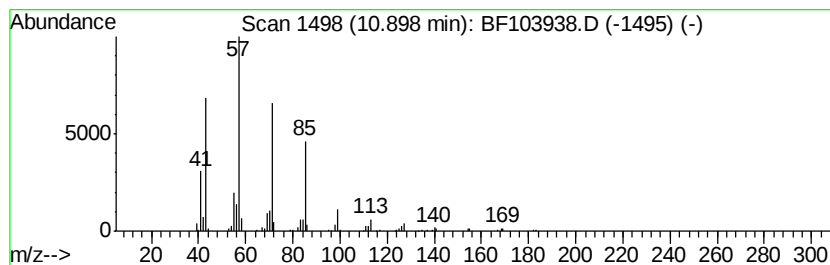
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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	4.40 ng	276863	Phenanthrene-d10	11.51

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	90
2	Hexadecane		226	C16H34	000544-76-3	90
3	9-methylheptadecane		254	C18H38	026741-18-4	90
4	Pentadecane		212	C15H32	000629-62-9	90
5	Dodecane, 1-iodo-		296	C12H25I	004292-19-7	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
Data File : BF103938.D
Acq On : 26 Mar 2018 21:14
Operator : JU/SJ
Sample : J1758-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

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

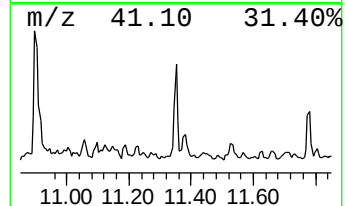
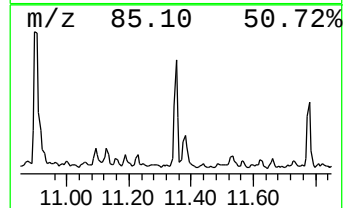
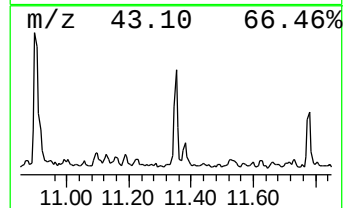
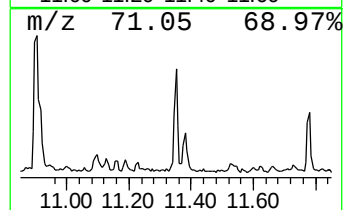
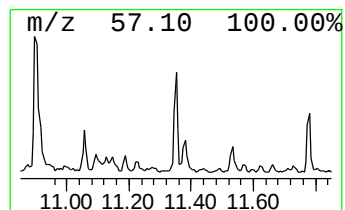
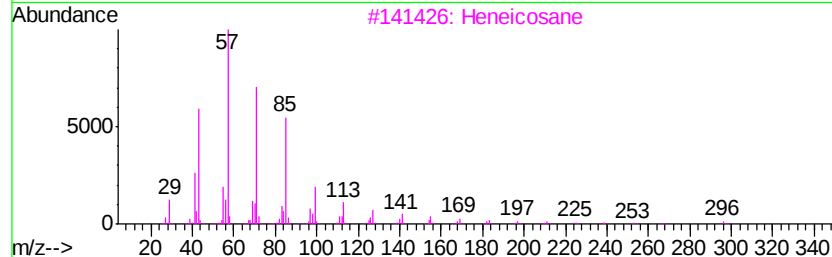
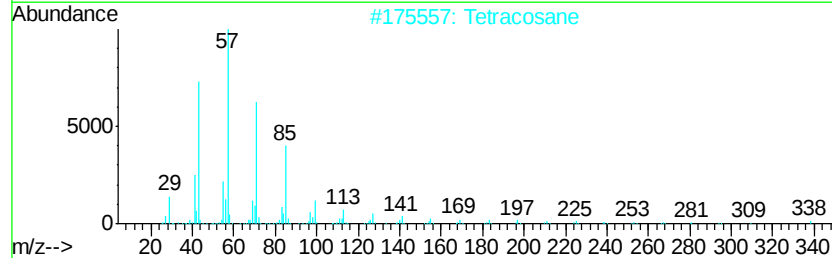
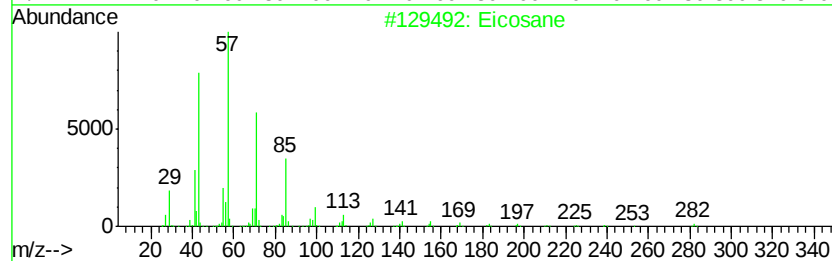
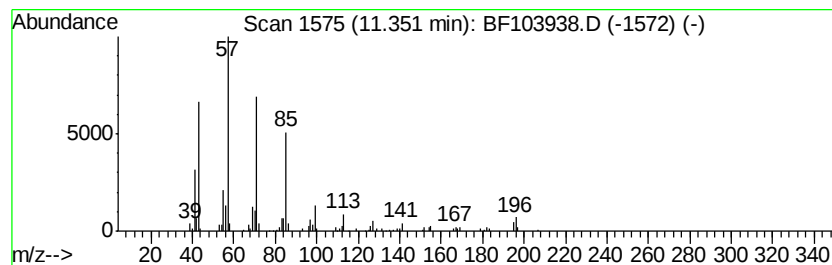
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Eicosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.35	2.12 ng	133321	Phenanthrene-d10	11.51

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	90
2	Tetracosane		338	C24H50	000646-31-1	90
3	Heneicosane		296	C21H44	000629-94-7	87
4	Pentadecane		212	C15H32	000629-62-9	87
5	Hexadecane		226	C16H34	000544-76-3	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
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Acq On : 26 Mar 2018 21:14
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ClientSampleId :
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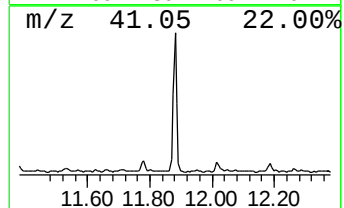
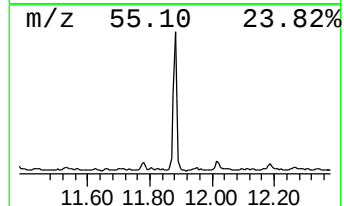
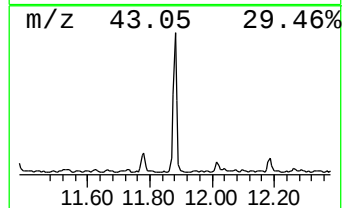
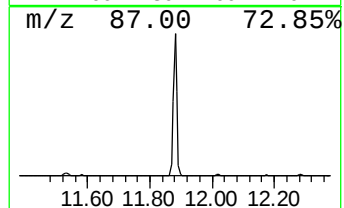
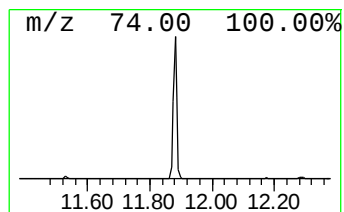
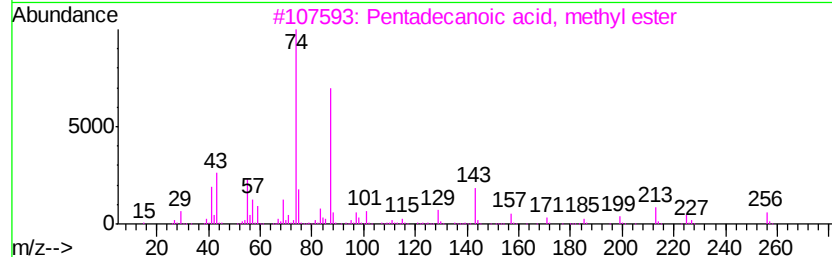
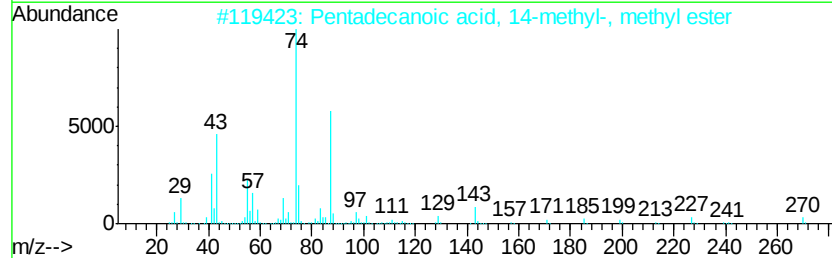
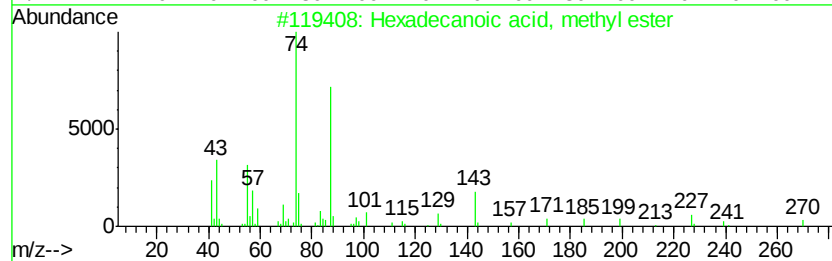
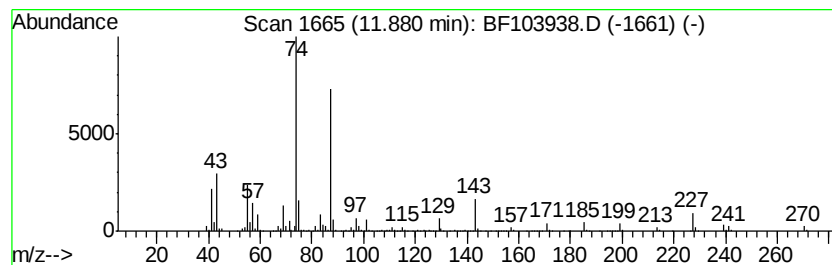
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexadecanoic acid, methyl e... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.88	18.83 ng	1185910	Phenanthrene-d10	11.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	99
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	94
3		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86
4		Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	86
5		Methyl tetradecanoate	242	C15H30O2	000124-10-7	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
Data File : BF103938.D
Acq On : 26 Mar 2018 21:14
Operator : JU/SJ
Sample : J1758-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

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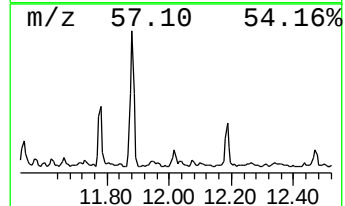
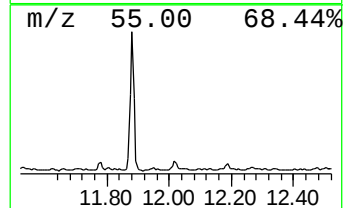
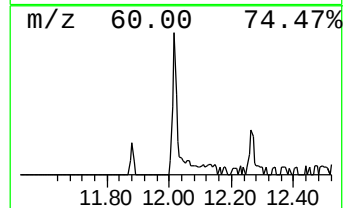
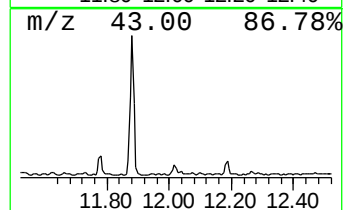
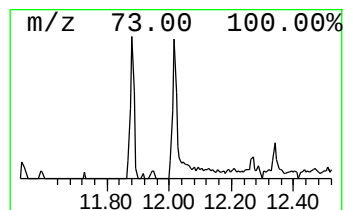
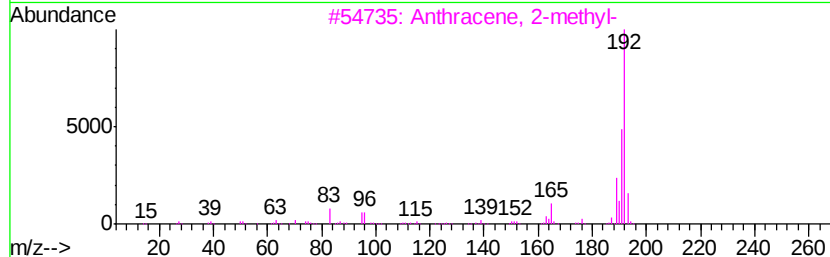
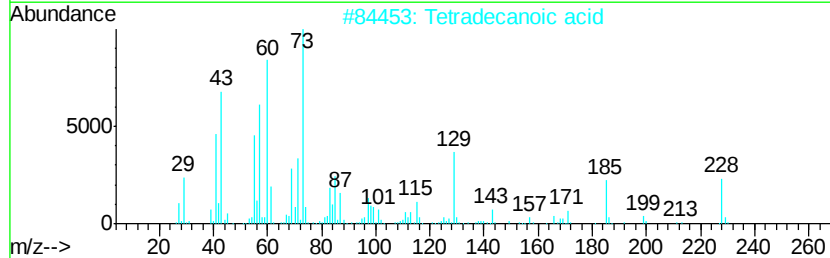
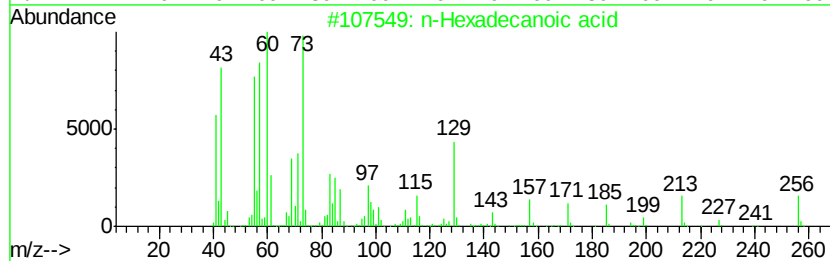
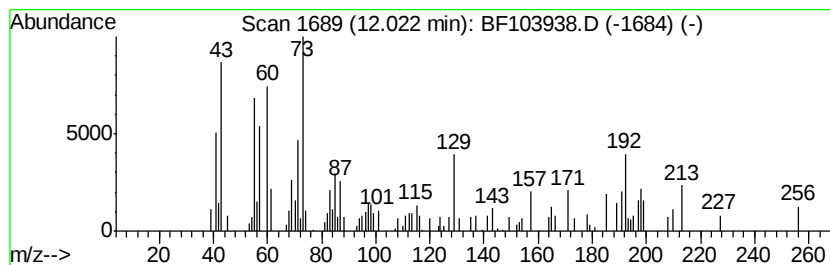
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 n-Hexadecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.02	2.18 ng	137345	Phenanthrene-d10	11.51

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	41
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	38
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	38
5			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103938.D
Acq On : 26 Mar 2018 21:14
Operator : JU/SJ
Sample : J1758-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
PL-02-032318-D

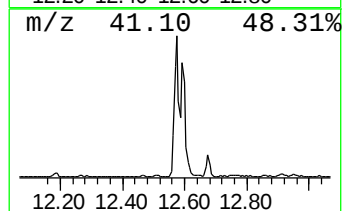
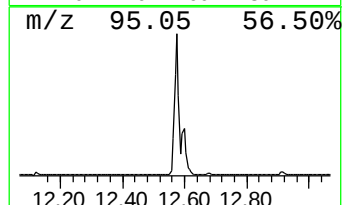
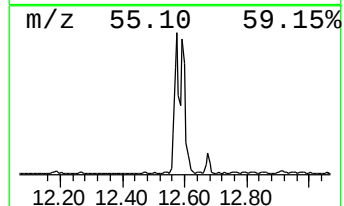
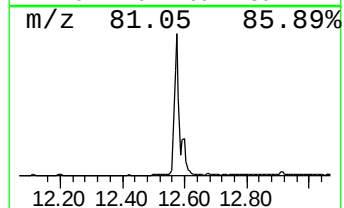
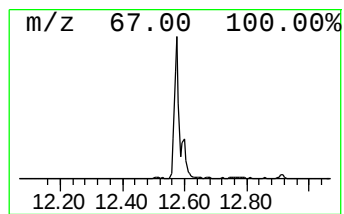
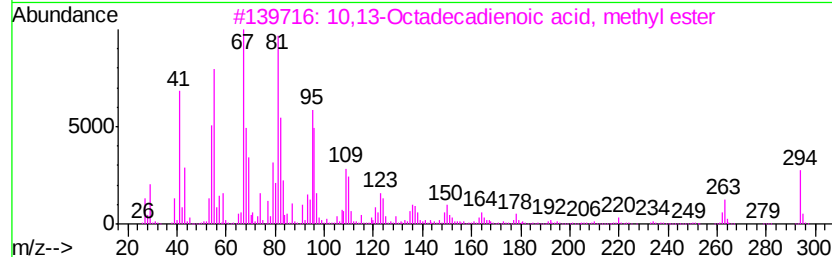
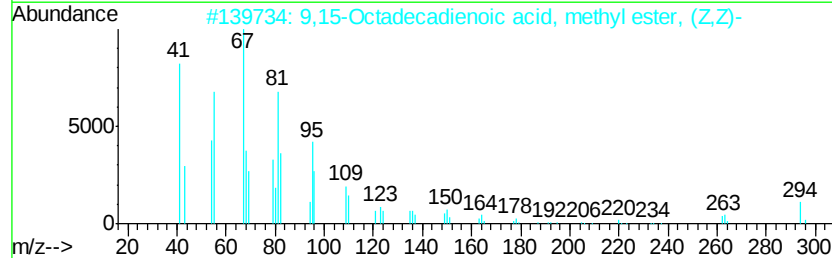
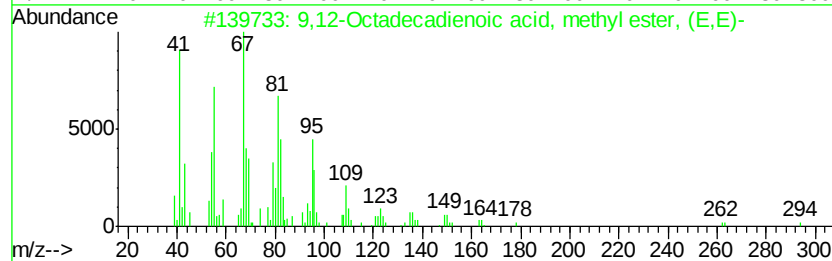
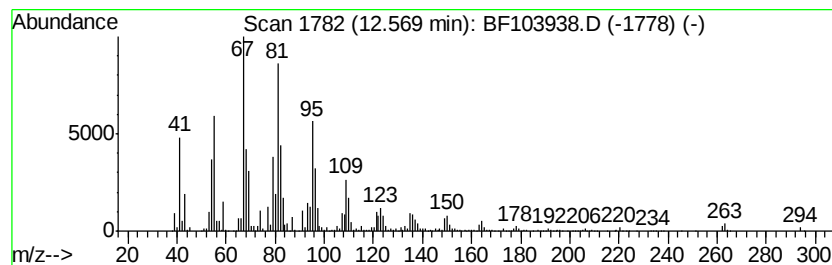
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 9,12-Octadecadienoic acid, ... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.57	67.33 ng	4239580	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		9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
3		10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99
4		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
5		11,14-Octadecadienoic acid, meth...	294	C19H34O2	056554-61-1	98



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103938.D
Acq On : 26 Mar 2018 21:14
Operator : JU/SJ
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Misc :
ALS Vial : 11 Sample Multiplier: 1

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PL-02-032318-D

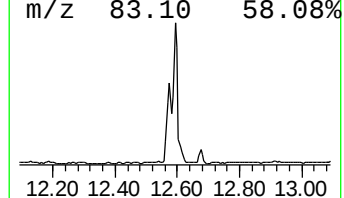
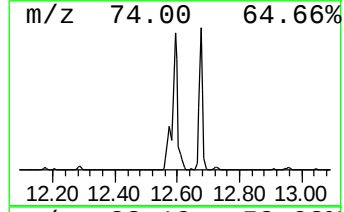
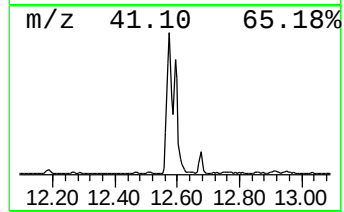
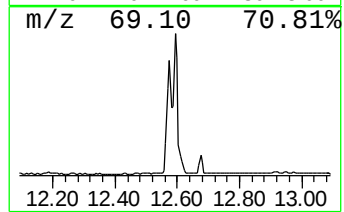
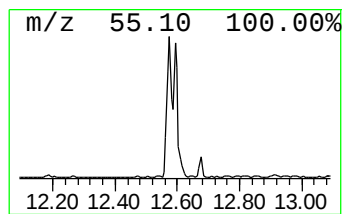
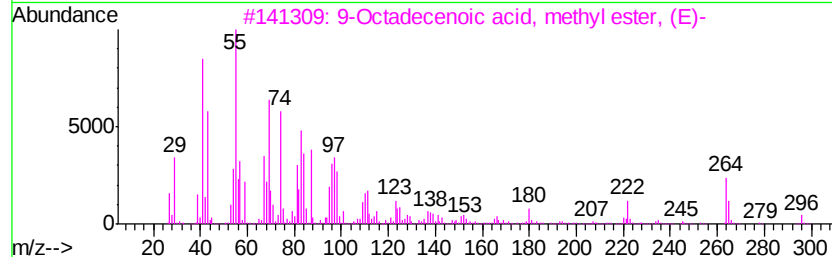
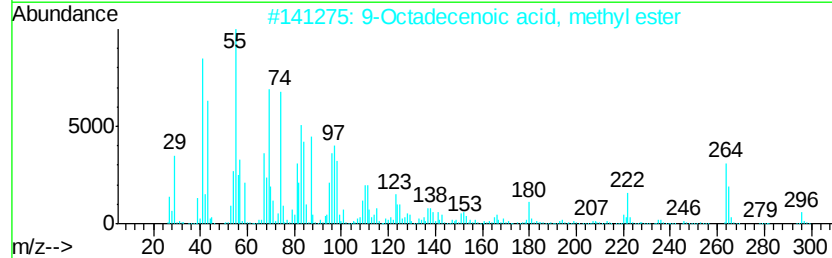
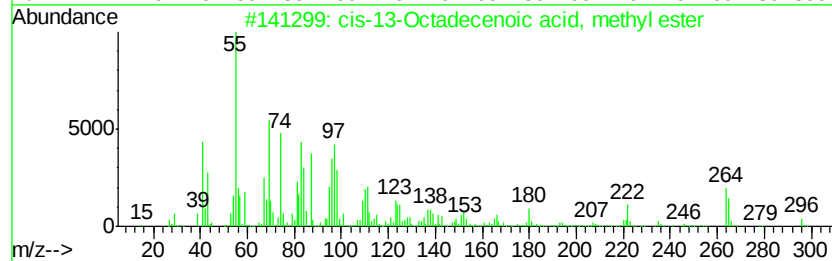
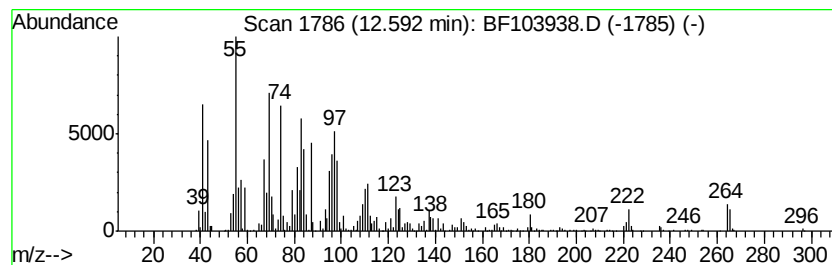
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 cis-13-Octadecenoic acid, m... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.59	40.76 ng	2566540	Phenanthrene-d10	11.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	cis-13-Octadecenoic acid, methyl...	296	C19H36O2	1000333-58-3	99
2		9-Octadecenoic acid, methyl ester	296	C19H36O2	002462-84-2	99
3		9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	99
4		10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	99
5		trans-13-Octadecenoic acid, meth...	296	C19H36O2	1000333-61-3	99



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
Data File : BF103938.D
Acq On : 26 Mar 2018 21:14
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Misc :
ALS Vial : 11 Sample Multiplier: 1

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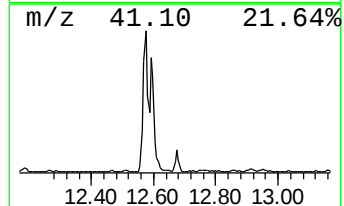
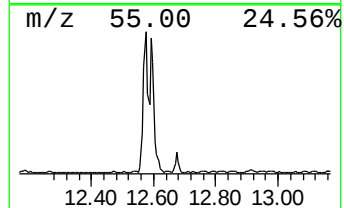
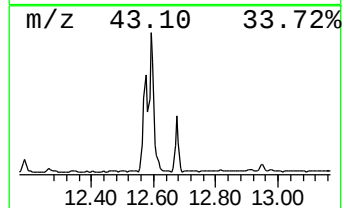
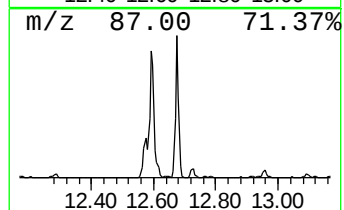
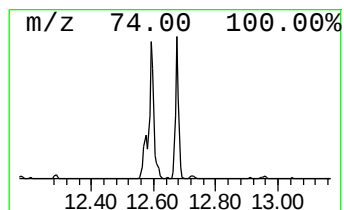
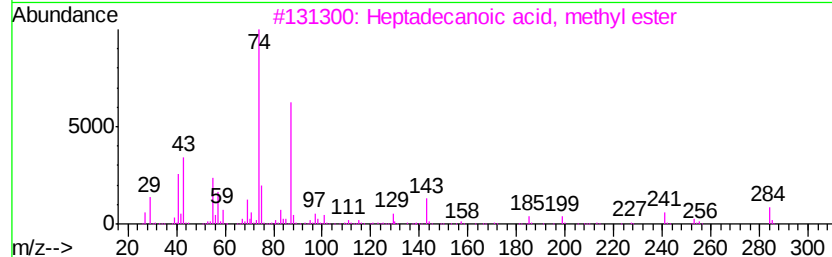
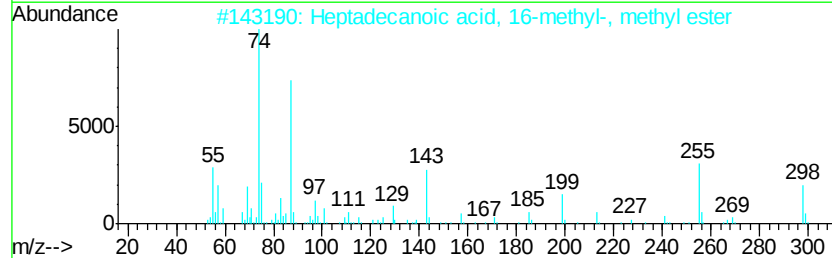
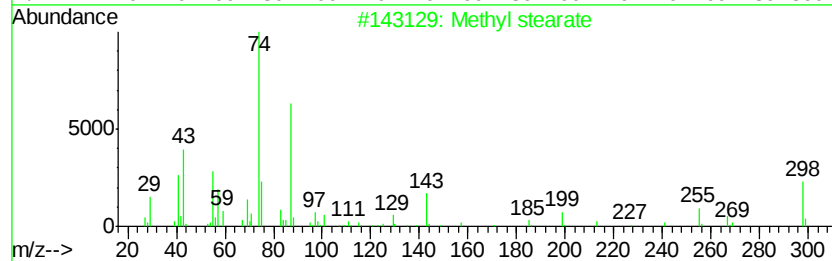
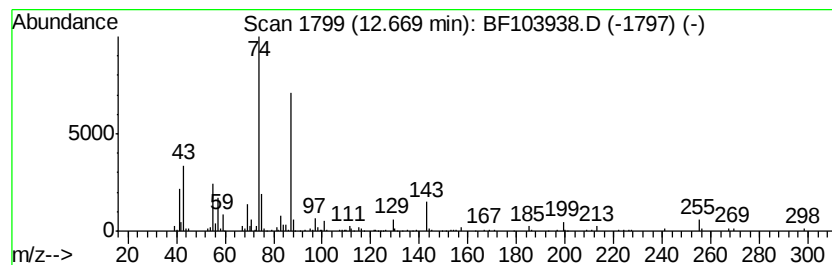
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Methyl stearate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	7.91 ng	498077	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	94
3		Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	91
4		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	90
5		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
Data File : BF103938.D
Acq On : 26 Mar 2018 21:14
Operator : JU/SJ
Sample : J1758-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

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

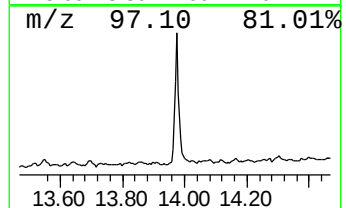
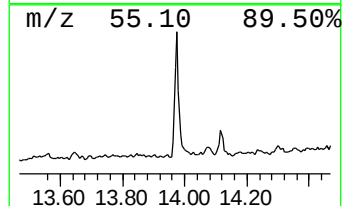
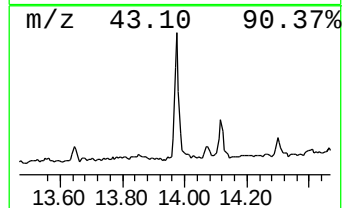
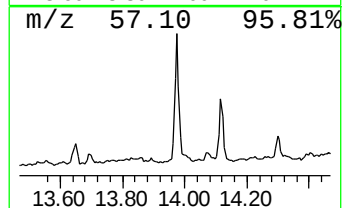
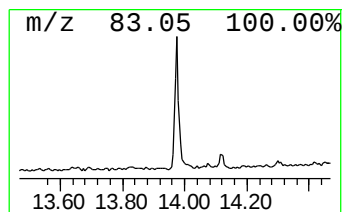
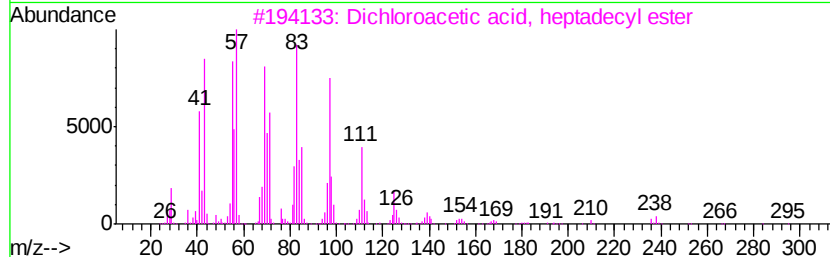
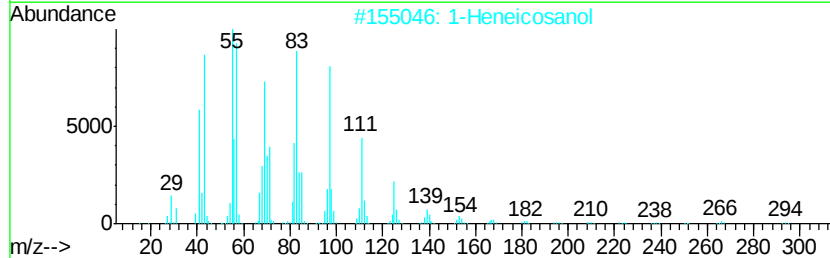
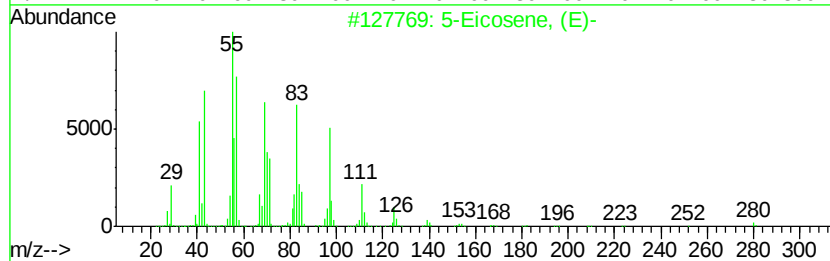
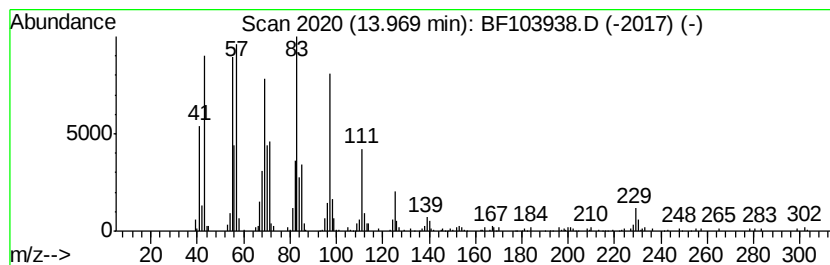
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 5-Eicosene, (E)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	7.56 ng	420776	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Eicosene, (E)-	280	C20H40	074685-30-6	98
2		1-Heneicosanol	312	C21H44O	015594-90-8	95
3		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	94
4		n-Tetracosanol-1	354	C24H50O	000506-51-4	93
5		Cyclopentadecane	210	C15H30	000295-48-7	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
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Acq On : 26 Mar 2018 21:14
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Sample : J1758-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

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

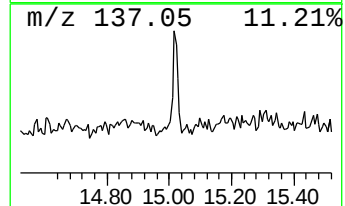
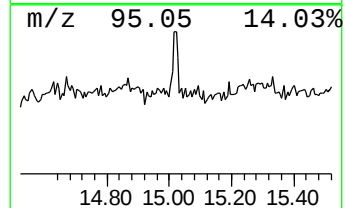
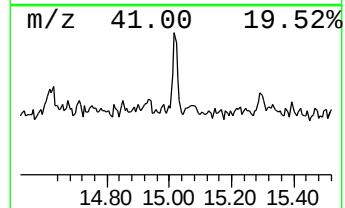
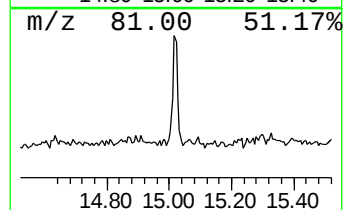
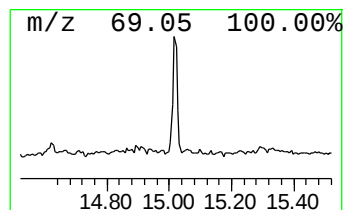
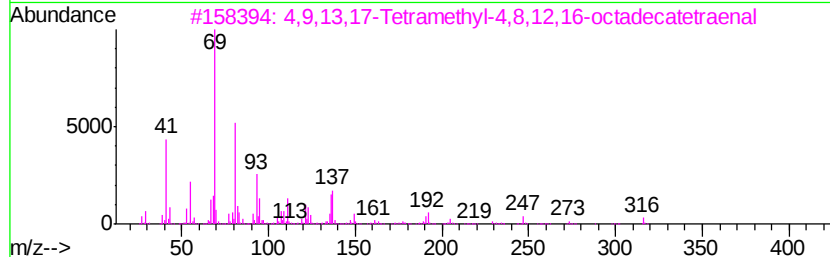
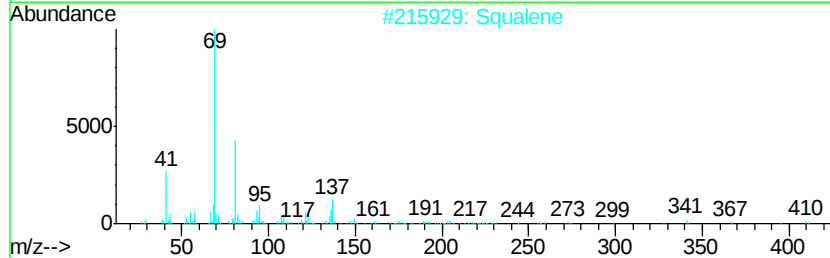
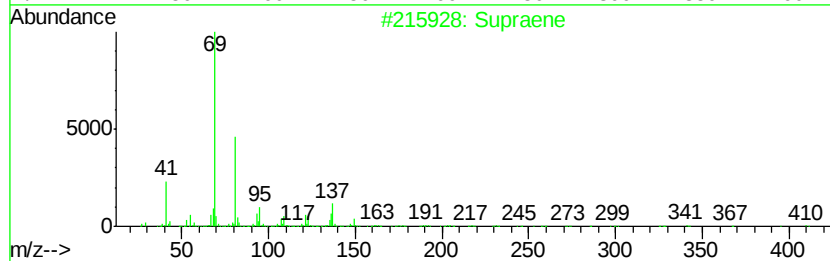
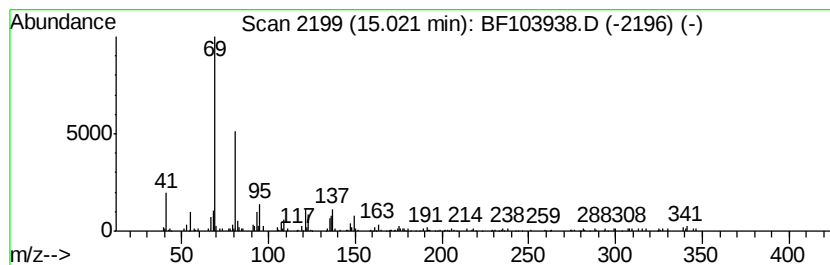
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Supraene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.02	2.61 ng	104777	Perylene-d12	15.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Supraene	410	C30H50	007683-64-9	87
2		Squalene	410	C30H50	000111-02-4	80
3		4,9,13,17-Tetramethyl-4,8,12,16-...	316	C22H36O	056882-09-8	72
4		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	72
5		2,6,10,14-Hexadecatetraenoic aci...	332	C22H36O2	024035-35-6	68



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 Misc :
 ALS Vial : 11 Sample Multiplier: 1

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.32	3.5	ng	155824	1	6.97	880128	20.0
2-Pentanone, 4-hy...	5.22	3.5	ng	152992	1	6.97	880128	20.0
unknown6.75	6.75	69.6	ng	3062070	1	6.97	880128	20.0
Hexadecane	10.43	3.2	ng	205240	3	10.02	1277420	20.0
Heneicosane	10.90	4.4	ng	276863	4	11.51	1259340	20.0
Eicosane	11.35	2.1	ng	133321	4	11.51	1259340	20.0
Hexadecanoic acid...	11.88	18.8	ng	1185910	4	11.51	1259340	20.0
n-Hexadecanoic acid	12.02	2.2	ng	137345	4	11.51	1259340	20.0
9,12-Octadecadien...	12.57	67.3	ng	4239580	4	11.51	1259340	20.0
cis-13-Octadeceno...	12.59	40.8	ng	2566540	4	11.51	1259340	20.0
Methyl stearate	12.67	7.9	ng	498077	4	11.51	1259340	20.0
5-Eicosene, (E)-	13.97	7.6	ng	420776	5	14.16	1113480	20.0
Supraene	15.02	2.6	ng	104777	6	15.71	803891	20.0