

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
 Data File : BF109347.D
 Acq On : 26 Sep 2018 19:55
 Operator : JU/SJ
 Sample : J4773-13
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-092518-B

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.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	4.887	472	476	485	rBV	31043	39259	3.08%	0.218%
2	5.310	544	548	558	rBV	1176530	1027405	80.58%	5.707%
3	6.339	719	723	726	rBV	1362256	1095866	85.95%	6.087%
4	6.475	742	746	749	rBV	1302055	1115706	87.50%	6.198%
5	6.704	782	785	789	rBV	756911	596881	46.81%	3.316%
6	6.863	808	812	815	rBV	1076317	849263	66.61%	4.718%
7	7.263	876	880	883	rBV	861927	682313	53.51%	3.790%
8	7.986	999	1003	1009	rBV	997476	807370	63.32%	4.485%
9	8.492	1086	1089	1093	rVB2	19126	16524	1.30%	0.092%
10	8.592	1102	1106	1109	rBV	50090	42720	3.35%	0.237%
11	8.810	1137	1143	1145	rBV3	21637	25607	2.01%	0.142%
12	9.022	1177	1179	1182	rVB3	23626	21867	1.72%	0.121%
13	9.063	1182	1186	1189	rBV	1673775	1275030	100.00%	7.083%
14	9.157	1196	1202	1205	rBV	90492	78910	6.19%	0.438%
15	9.322	1226	1230	1237	rBV10	15009	27709	2.17%	0.154%
16	9.457	1250	1253	1255	rBV	391998	323118	25.34%	1.795%
17	9.527	1263	1265	1268	rVB3	15206	16326	1.28%	0.091%
18	9.686	1289	1292	1296	rVB	107444	88254	6.92%	0.490%
19	9.739	1296	1301	1304	rBV	950355	872634	68.44%	4.847%
20	9.922	1327	1332	1334	rBV5	17798	28681	2.25%	0.159%
21	9.945	1334	1336	1339	rVB4	28058	24464	1.92%	0.136%
22	10.004	1344	1346	1350	rVB3	28641	31203	2.45%	0.173%
23	10.180	1373	1376	1379	rVB	141543	105119	8.24%	0.584%
24	10.233	1383	1385	1391	rVB5	12757	20398	1.60%	0.113%
25	10.280	1391	1393	1395	rBV	26921	19930	1.56%	0.111%
26	10.404	1409	1414	1416	rBV	49229	56779	4.45%	0.315%
27	10.451	1420	1422	1425	rVB	22123	18654	1.46%	0.104%
28	10.480	1425	1427	1430	rBV2	18142	18592	1.46%	0.103%
29	10.522	1430	1434	1438	rBV	726805	714187	56.01%	3.967%
30	10.651	1453	1456	1463	rBV	170696	179649	14.09%	0.998%
31	10.857	1487	1491	1493	rBV4	18665	22402	1.76%	0.124%
32	10.974	1509	1511	1515	rVB3	21050	22390	1.76%	0.124%
33	11.022	1515	1519	1524	rVB6	17361	24726	1.94%	0.137%
34	11.098	1529	1532	1535	rVV	153554	130926	10.27%	0.727%

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35	11.127	1535	1537	1542	rVB	60380	57865	4.54%	0.321%
36	11.216	1548	1552	1554	rBV	844059	770462	60.43%	4.280%
37	11.233	1554	1555	1559	rVB	86922	62557	4.91%	0.347%
38	11.286	1559	1564	1567	rBV2	34149	40184	3.15%	0.223%
39	11.474	1594	1596	1600	rVB4	15418	16889	1.32%	0.094%
40	11.521	1601	1604	1607	rBV	115420	102557	8.04%	0.570%
41	11.621	1617	1621	1624	rBV	518206	414641	32.52%	2.303%
42	11.751	1639	1643	1647	rBV2	83656	85526	6.71%	0.475%
43	11.821	1653	1655	1658	rBV2	34259	35634	2.79%	0.198%
44	11.927	1670	1673	1679	rBV	104442	87625	6.87%	0.487%
45	12.080	1697	1699	1705	rBV6	16709	22405	1.76%	0.124%
46	12.263	1727	1730	1733	rBV2	18576	19354	1.52%	0.108%
47	12.304	1733	1737	1739	rBV	1262634	1250935	98.11%	6.949%
48	12.321	1739	1740	1747	rVB	1288219	1065917	83.60%	5.921%
49	12.410	1752	1755	1760	rVV2	259671	277067	21.73%	1.539%
50	12.457	1760	1763	1771	rVB9	18845	35934	2.82%	0.200%
51	12.533	1773	1776	1781	rBV5	16317	24306	1.91%	0.135%
52	12.645	1790	1795	1799	rBV	114647	120577	9.46%	0.670%
53	12.686	1799	1802	1805	rVB	61033	47247	3.71%	0.262%
54	12.792	1817	1820	1824	rBV	1183190	1023059	80.24%	5.683%
55	12.974	1849	1851	1855	rVB2	28567	23492	1.84%	0.130%
56	13.039	1859	1862	1866	rBV2	70290	75112	5.89%	0.417%
57	13.080	1866	1869	1872	rVV2	18723	17788	1.40%	0.099%
58	13.168	1882	1884	1887	rVB2	21126	16918	1.33%	0.094%
59	13.198	1887	1889	1893	rBV2	17438	19928	1.56%	0.111%
60	13.245	1894	1897	1900	rBV4	20247	27722	2.17%	0.154%
61	13.380	1917	1920	1925	rBV	52573	50350	3.95%	0.280%
62	13.551	1946	1949	1952	rVB4	23800	28151	2.21%	0.156%
63	13.710	1972	1976	1984	rVB	122441	169058	13.26%	0.939%
64	13.839	1994	1998	2001	rBV	866325	803957	63.05%	4.466%
65	14.021	2027	2029	2032	rVB2	46811	40710	3.19%	0.226%
66	14.327	2078	2081	2083	rBV	57691	56658	4.44%	0.315%
67	14.621	2129	2131	2134	rBV	61400	62090	4.87%	0.345%
68	14.692	2140	2143	2146	rVB	75878	74430	5.84%	0.413%
69	15.245	2232	2237	2241	rBV	489238	554368	43.48%	3.079%

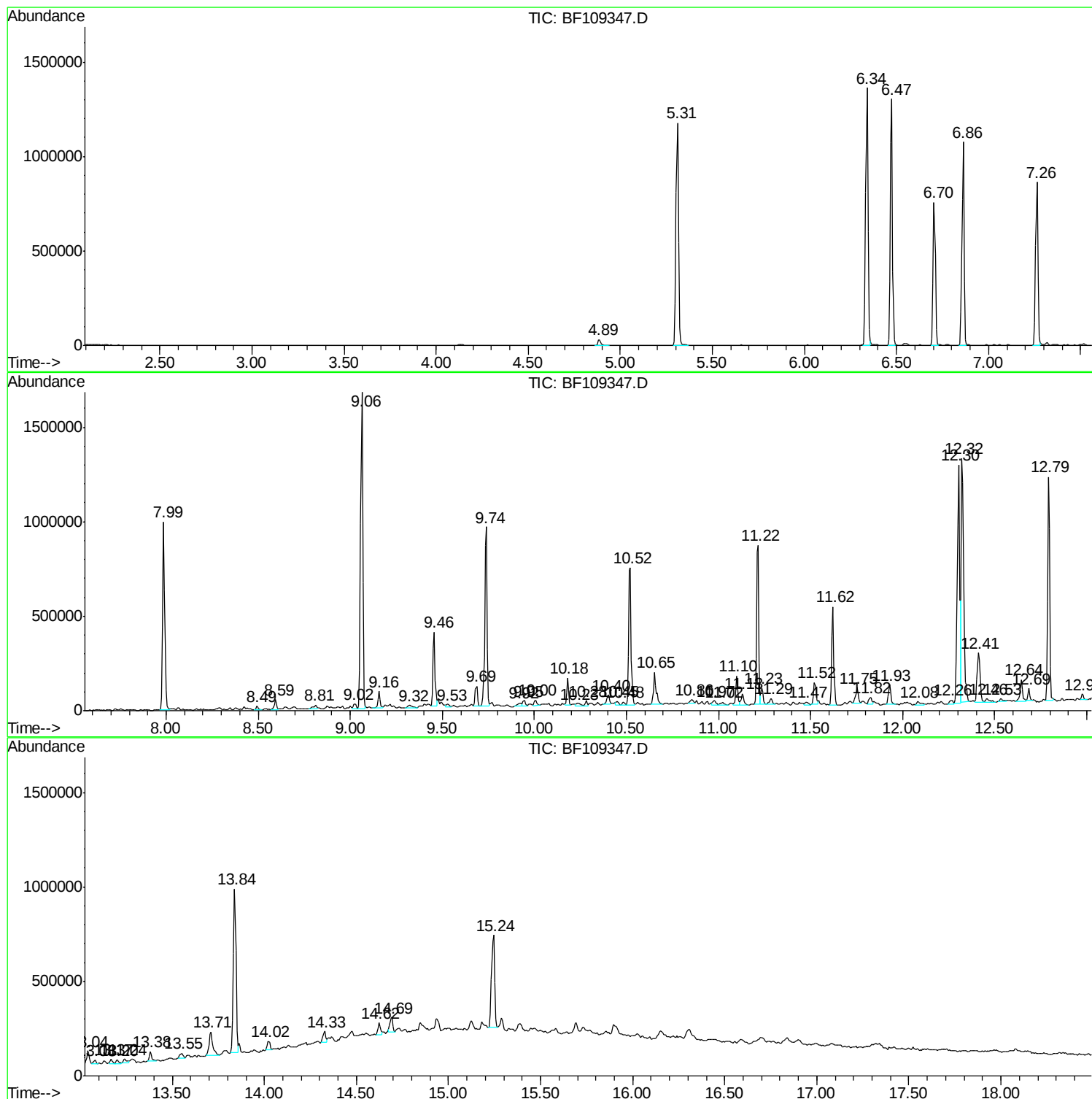
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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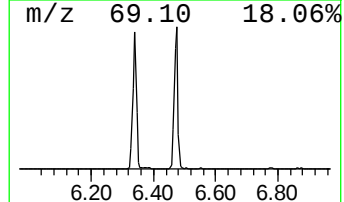
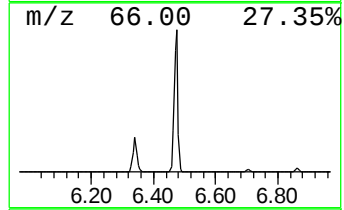
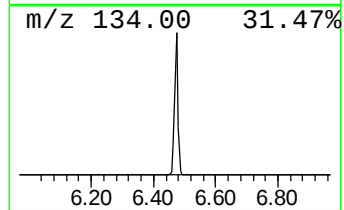
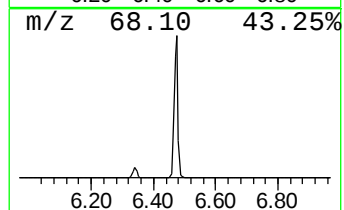
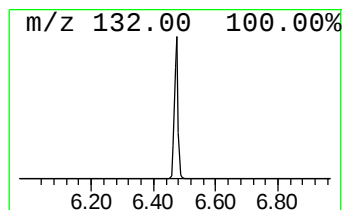
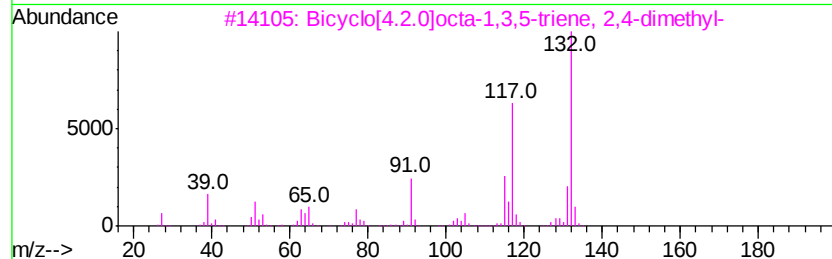
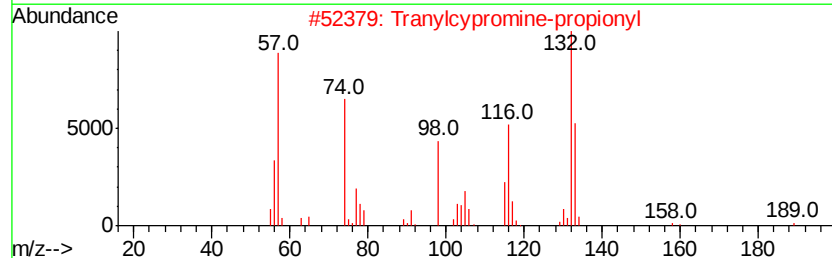
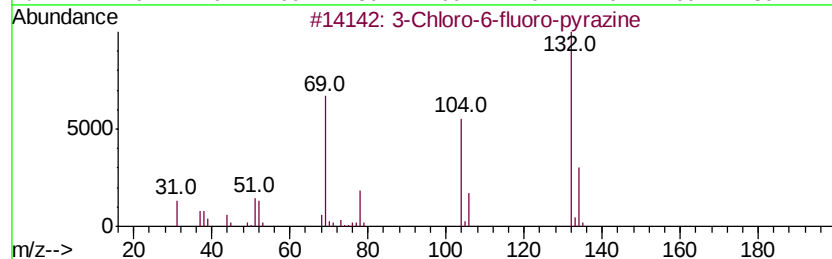
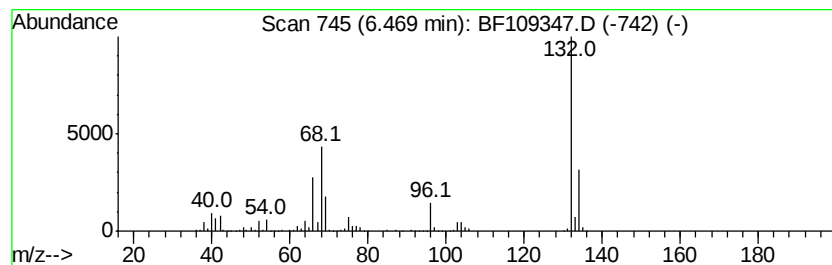
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.47 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	37.38 ng	1115710	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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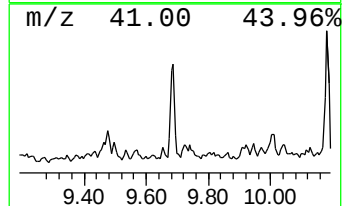
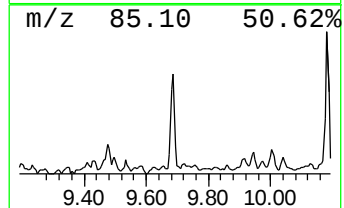
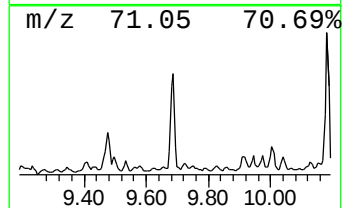
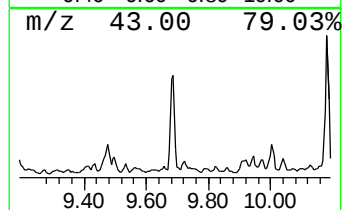
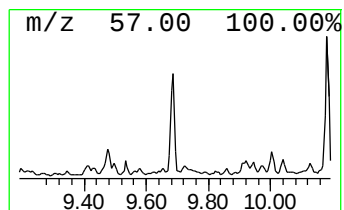
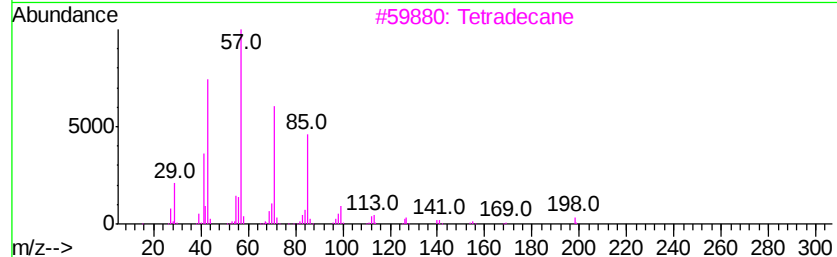
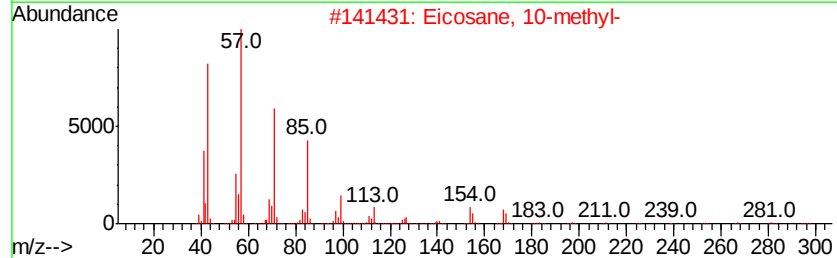
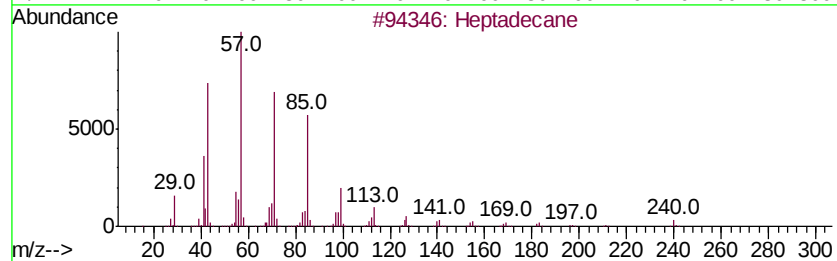
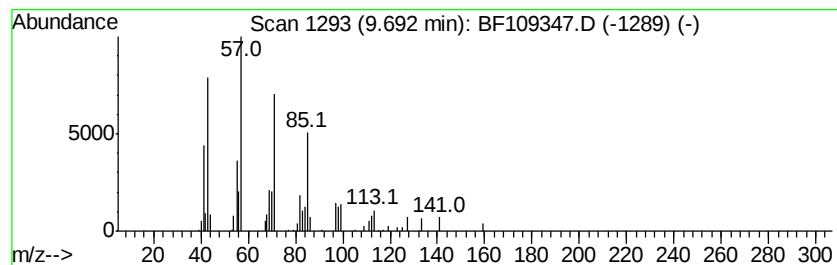
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Heptadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.69	2.02 ng	88254	Acenaphthene-d10	9.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	90
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	86
3		Tetradecane	198	C14H30	000629-59-4	86
4		Tridecane, 7-propyl-	226	C16H34	055045-09-5	80
5		9-methylheptadecane	254	C18H38	026741-18-4	78



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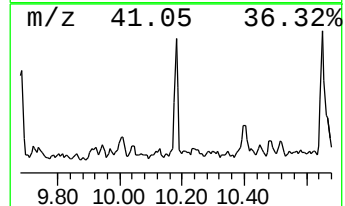
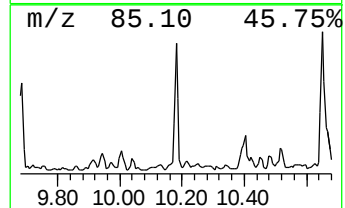
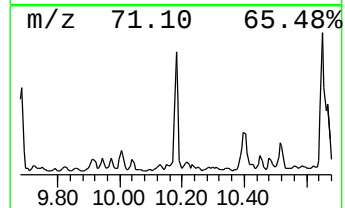
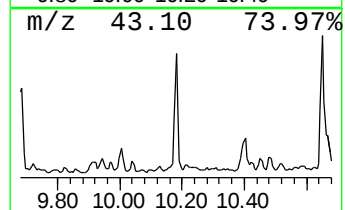
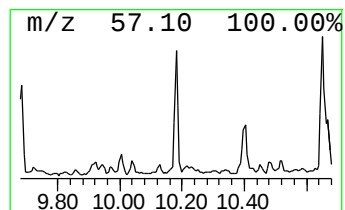
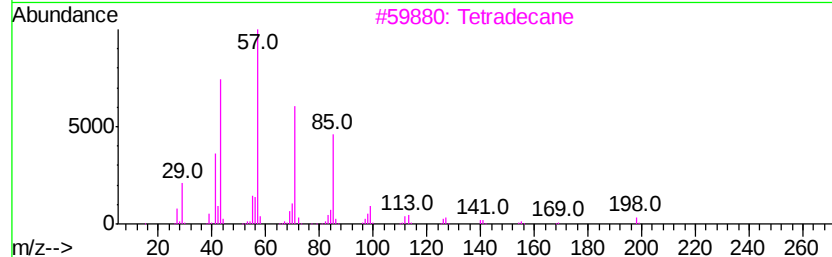
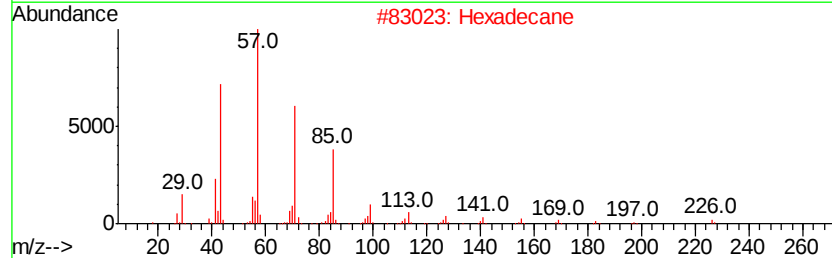
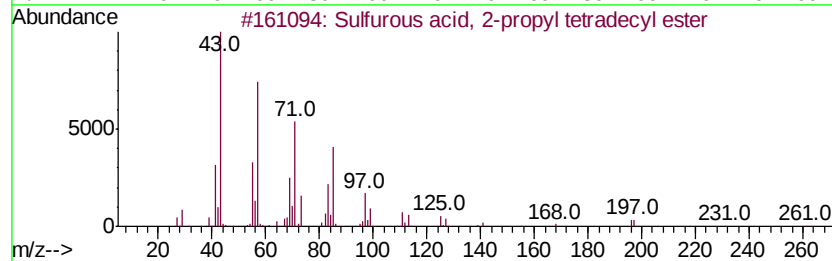
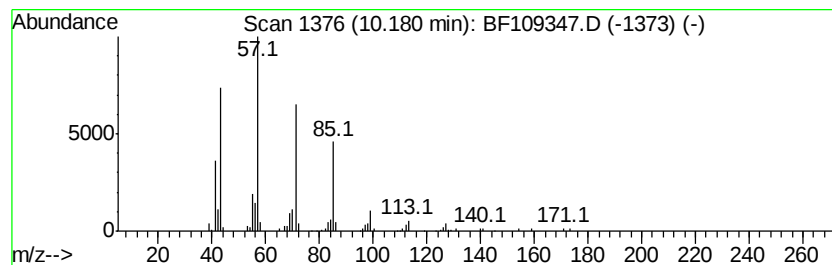
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Peak Number 4 Sulfurous acid, 2-propyl te... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.18	2.41 ng	105119	Acenaphthene-d10	9.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, 2-propyl tetradec...	320	C17H36O3S	1000309-12-5	90
2		Hexadecane	226	C16H34	000544-76-3	90
3		Tetradecane	198	C14H30	000629-59-4	90
4		Pentadecane	212	C15H32	000629-62-9	86
5		Eicosane	282	C20H42	000112-95-8	86



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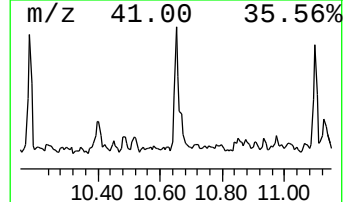
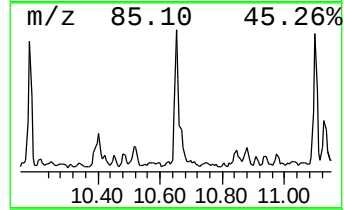
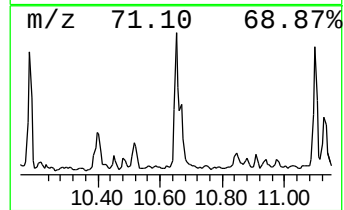
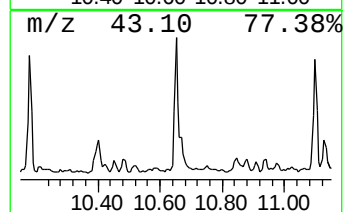
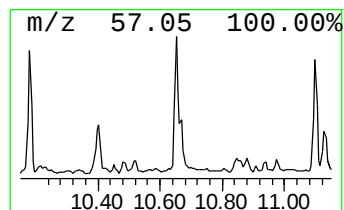
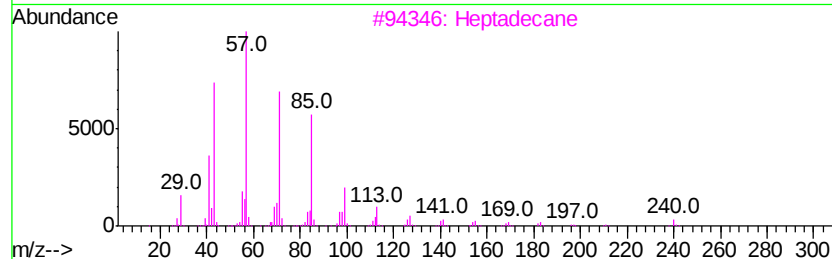
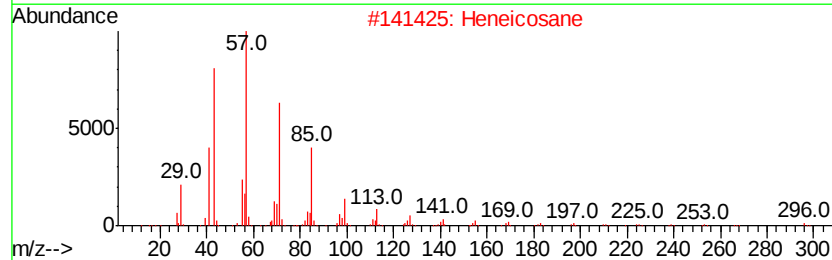
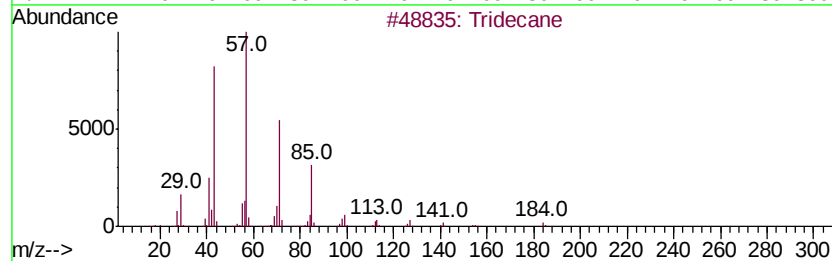
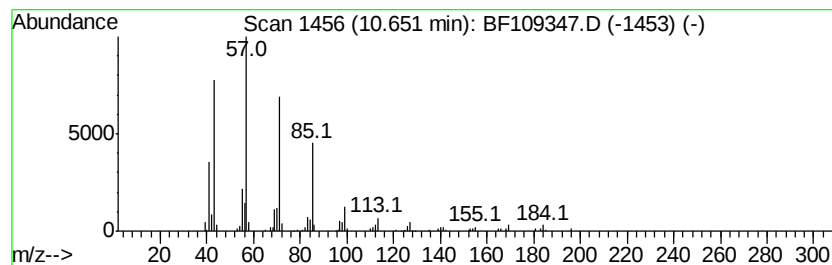
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Peak Number 5 Tridecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.65	4.66 ng	179649	Phenanthrene-d10	11.22	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	93
2	Heneicosane	296	C21H44	000629-94-7	91
3	Heptadecane	240	C17H36	000629-78-7	91
4	Tetracosane	338	C24H50	000646-31-1	91
5	Heptacosane	380	C27H56	000593-49-7	91



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Sample : J4773-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CL-02-092518-B

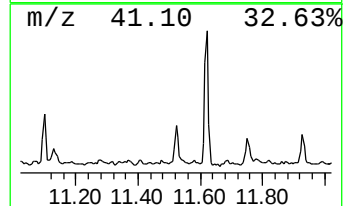
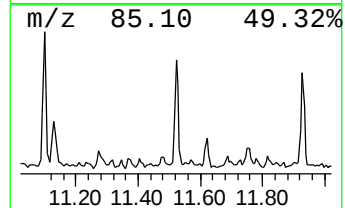
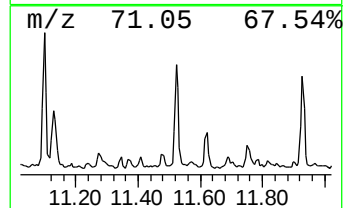
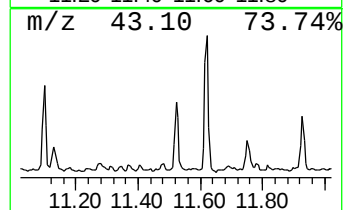
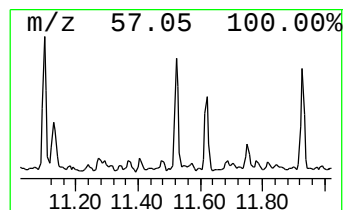
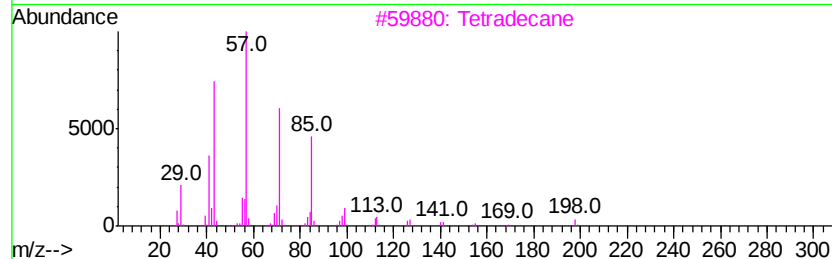
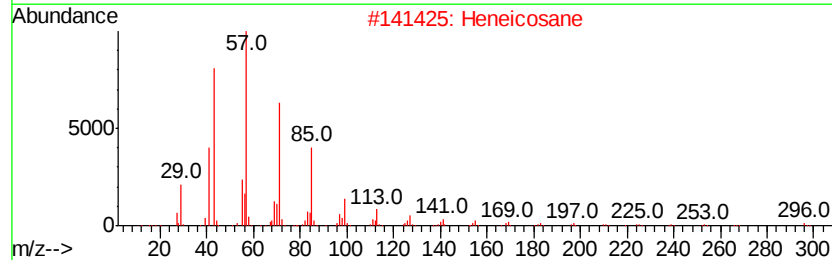
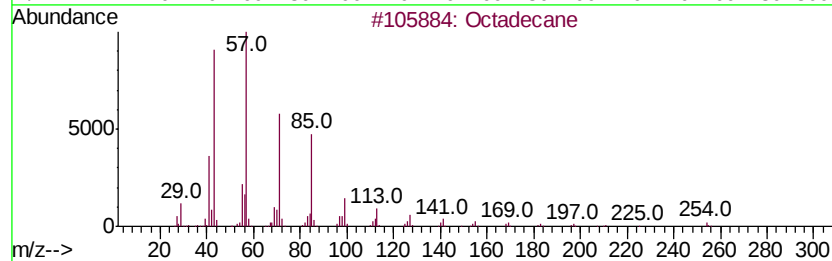
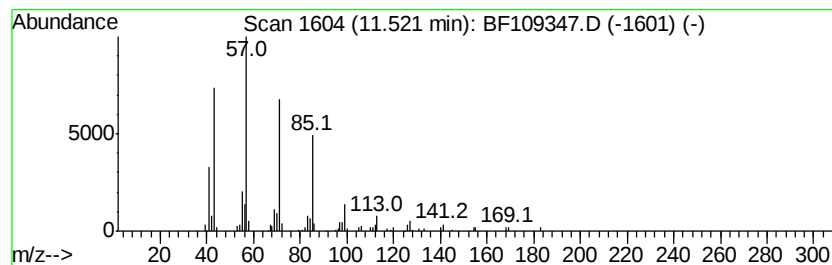
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.52	2.66 ng	102557	Phenanthrene-d10	11.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Tetradecane	198	C14H30	000629-59-4	91
4		Hexadecane	226	C16H34	000544-76-3	91
5		Heptadecane	240	C17H36	000629-78-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
Data File : BF109347.D
Acq On : 26 Sep 2018 19:55
Operator : JU/SJ
Sample : J4773-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

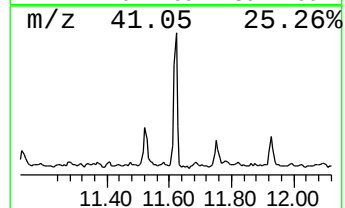
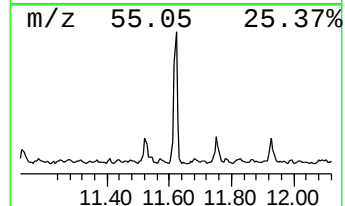
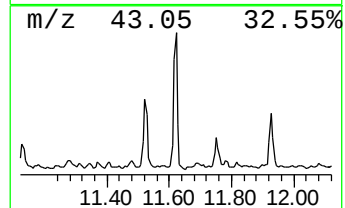
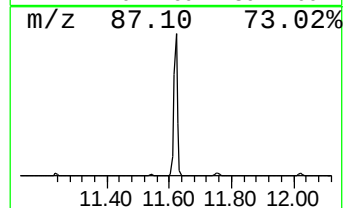
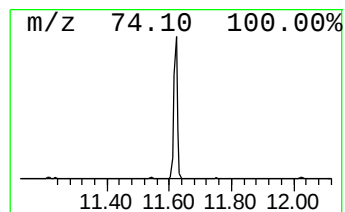
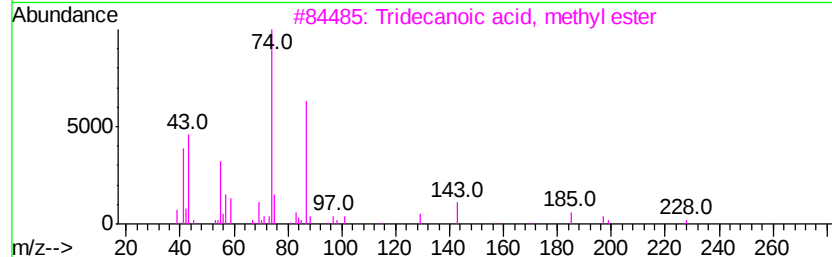
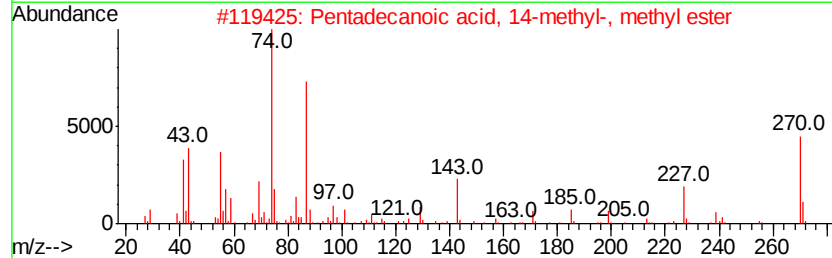
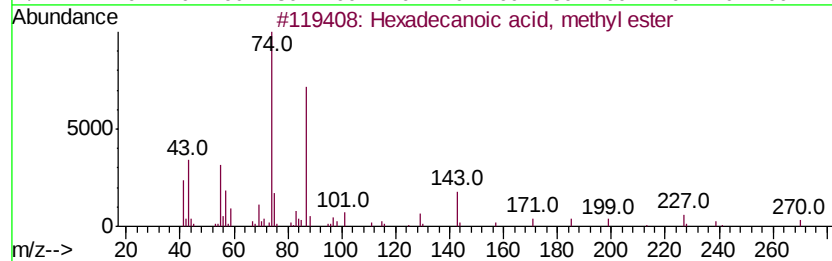
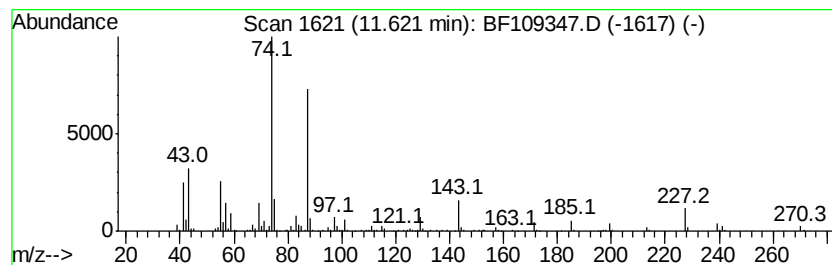
Instrument :
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ClientSampleId :
CL-02-092518-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
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.	
11.62	10.76 ng	414641	Phenanthrene-d10		11.22	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	94
3		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	90
4		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86
5		Tridecanoic acid, 12-methyl-, me...	242	C15H30O2	005129-58-8	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
Data File : BF109347.D
Acq On : 26 Sep 2018 19:55
Operator : JU/SJ
Sample : J4773-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

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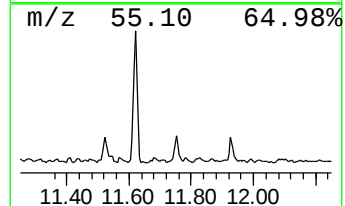
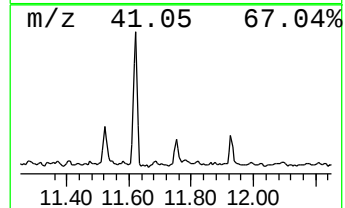
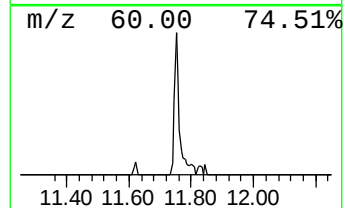
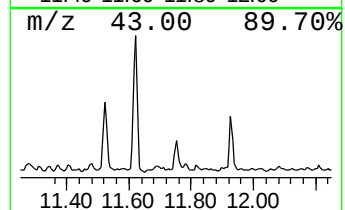
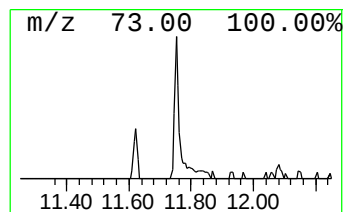
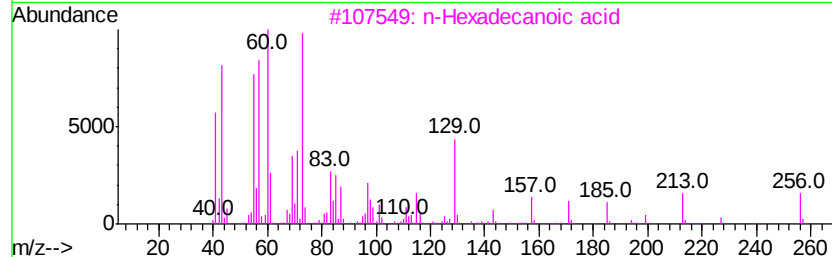
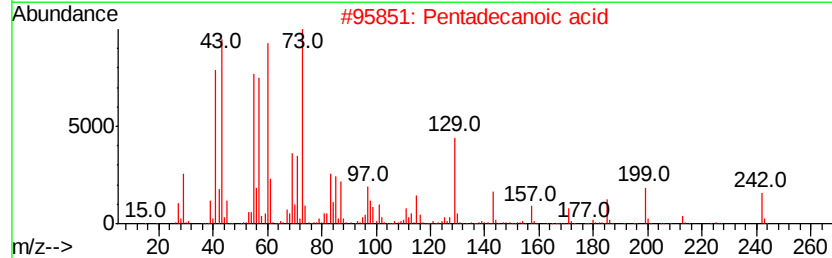
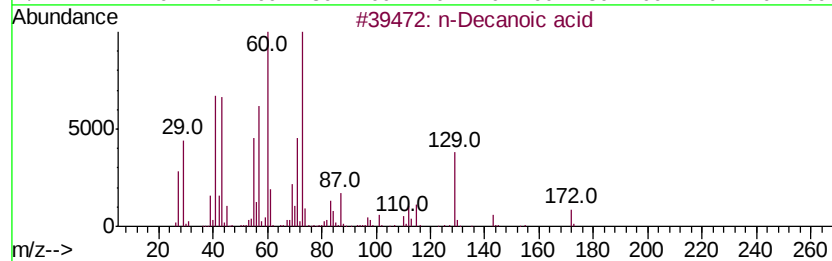
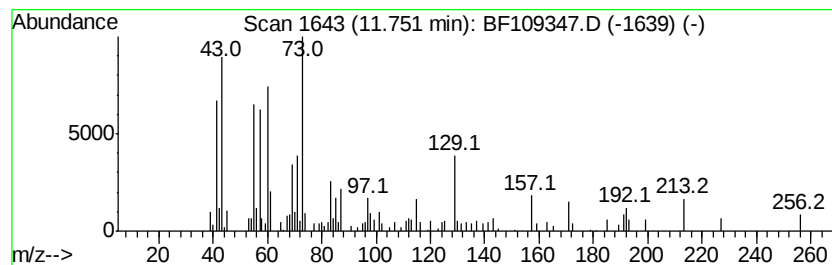
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 n-Decanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.75	2.22 ng	85526	Phenanthrene-d10	11.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Decanoic acid	172	C10H20O2	000334-48-5	87
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	64
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	64
4		Tridecanoic acid	214	C13H26O2	000638-53-9	53
5		Dodecanoic acid	200	C12H24O2	000143-07-7	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
Data File : BF109347.D
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Operator : JU/SJ
Sample : J4773-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

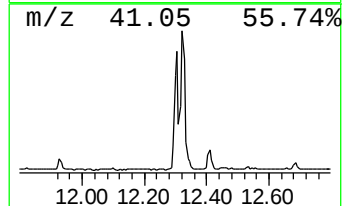
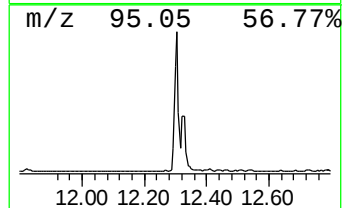
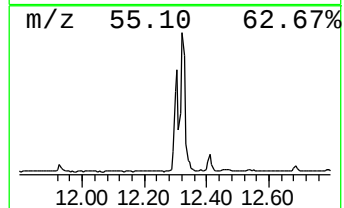
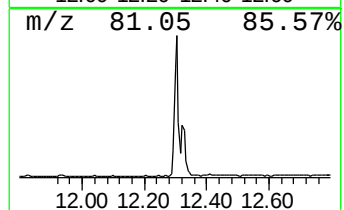
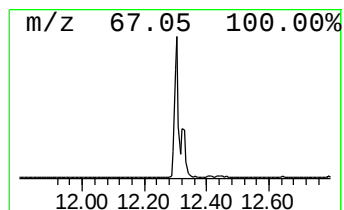
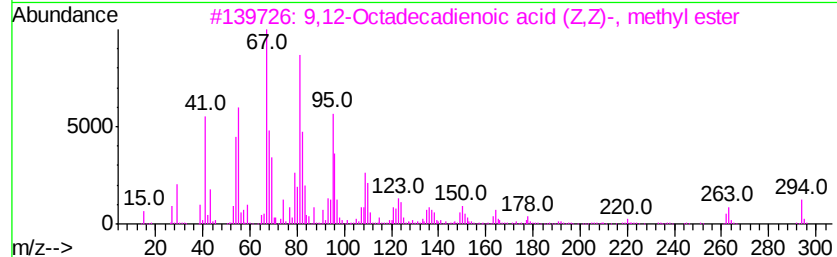
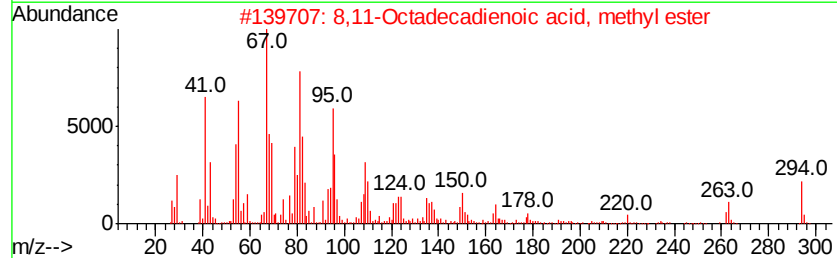
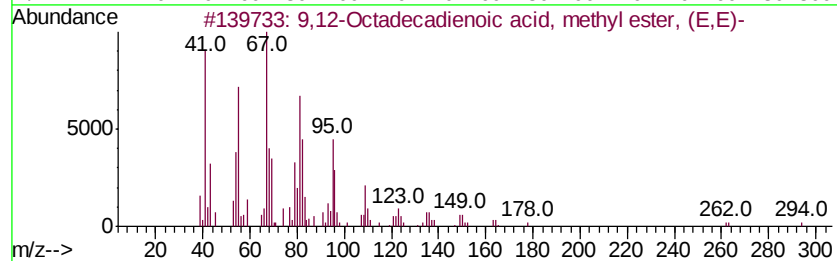
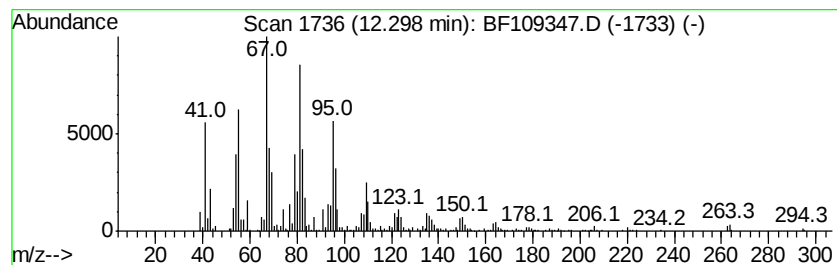
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.30	32.47 ng	1250940	Phenanthrene-d10	11.22		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99	
2	8,11-Octadecadienoic acid, methy...	294	C19H34O2	056599-58-7	99	
3	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99	
4	10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	98	
5	9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	95	



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
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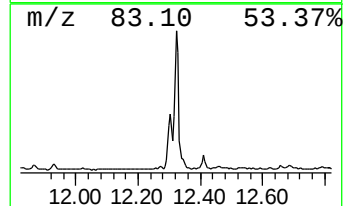
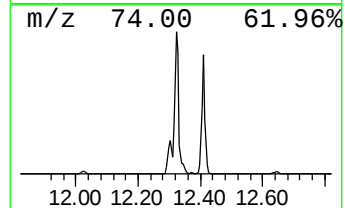
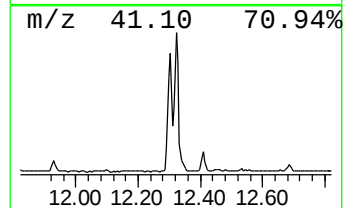
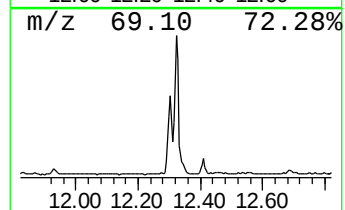
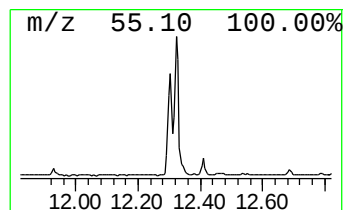
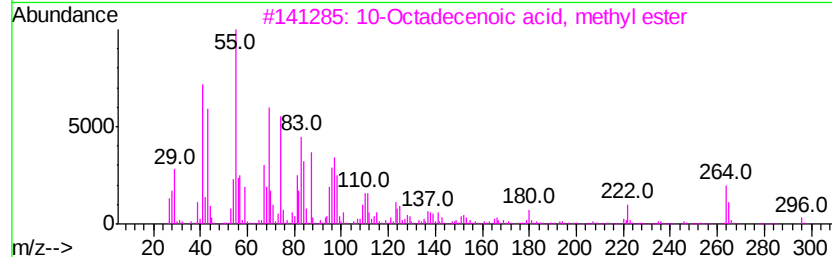
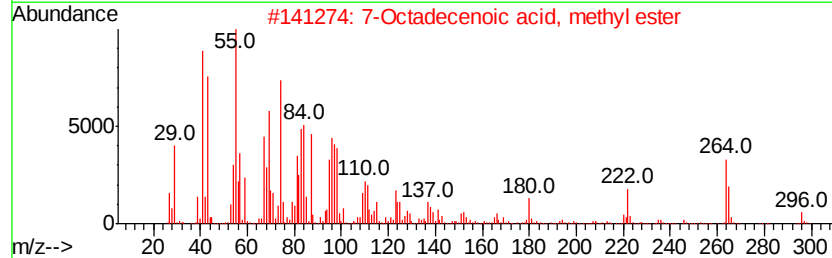
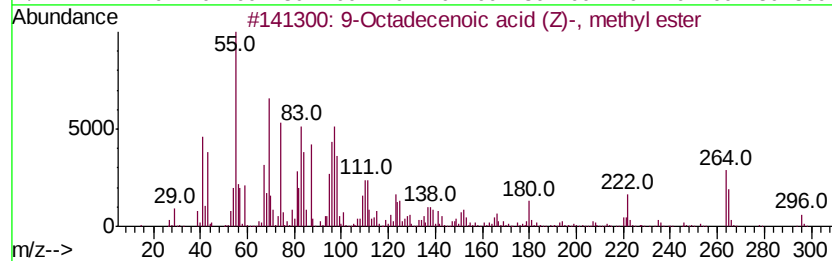
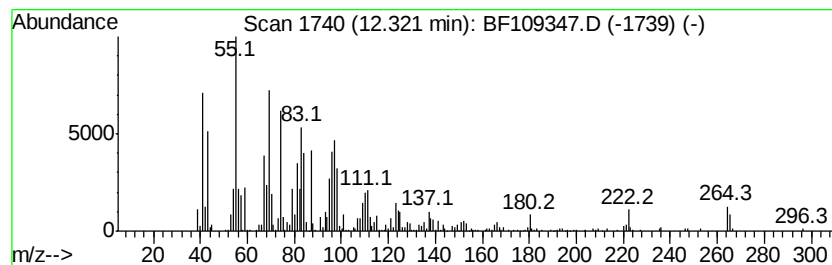
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 9-Octadecenoic acid (Z)-, m... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.32	27.67 ng	1065920	Phenanthrene-d10	11.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2		7-Octadecenoic acid, methyl ester	296	C19H36O2	057396-98-2	99
3		10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	99
4		9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	99
5		8-Octadecenoic acid, methyl ester	296	C19H36O2	002345-29-1	99



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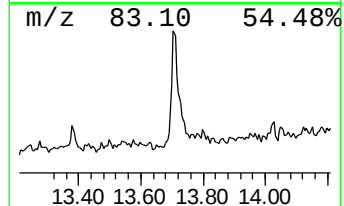
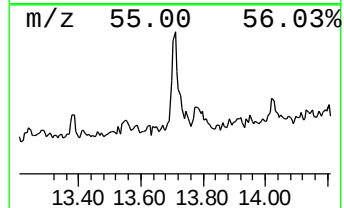
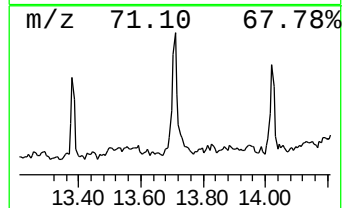
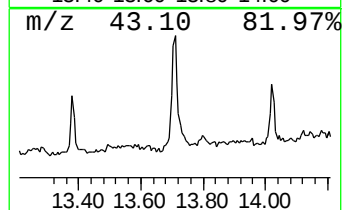
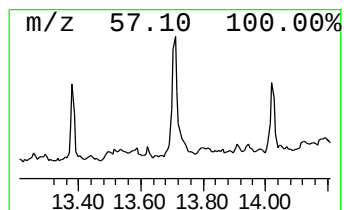
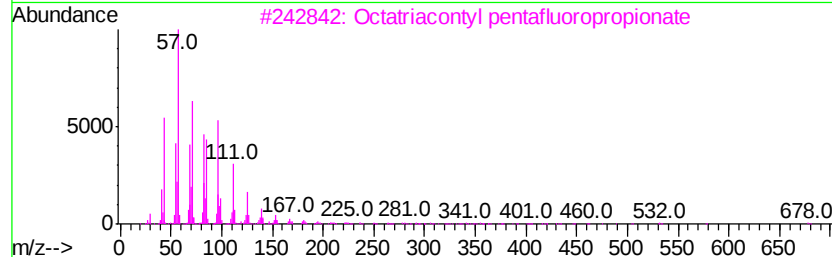
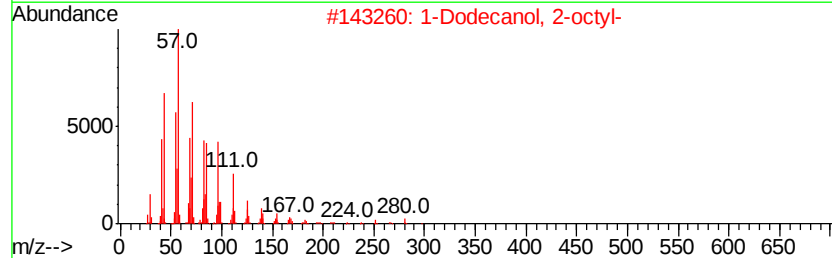
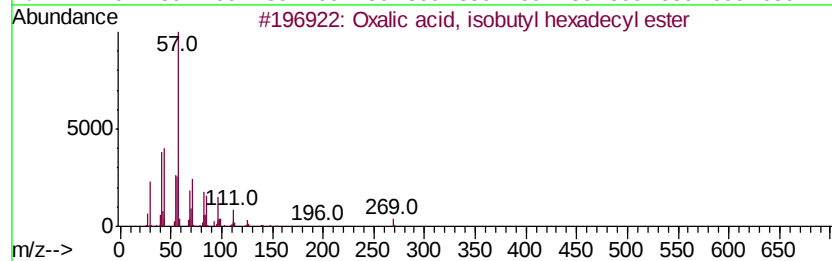
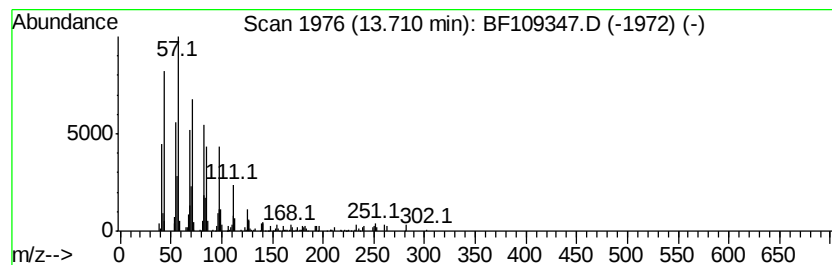
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Oxalic acid, isobutyl hexad... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.71	4.21 ng	169058	Chrysene-d12	13.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	90
2		1-Dodecanol, 2-octyl-	298	C20H42O	005333-42-6	90
3		Octatriacontyl pentafluoropropio...	697	C41H77F5O2	1000351-89-1	90
4		Cyclotetradecane	196	C14H28	000295-17-0	86
5		1-Dodecanol, 2-hexyl-	270	C18H38O	110225-00-8	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
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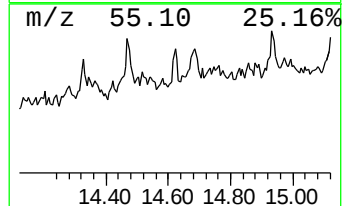
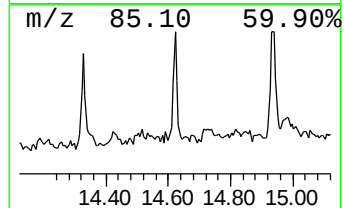
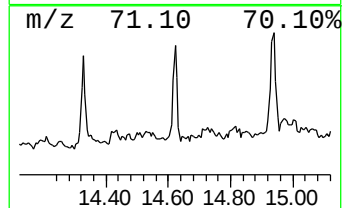
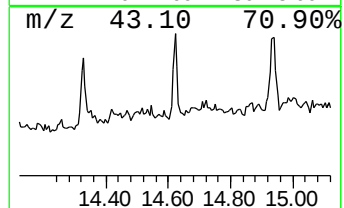
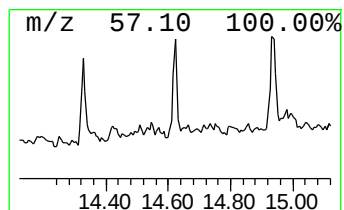
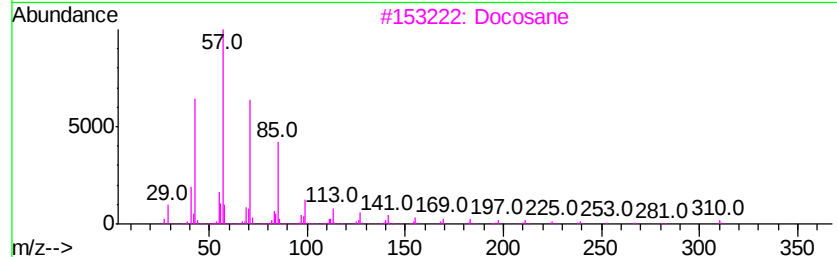
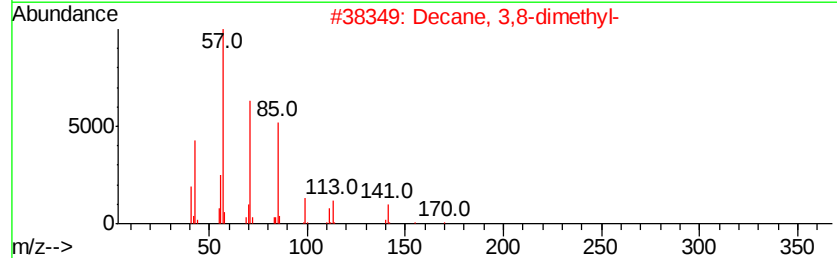
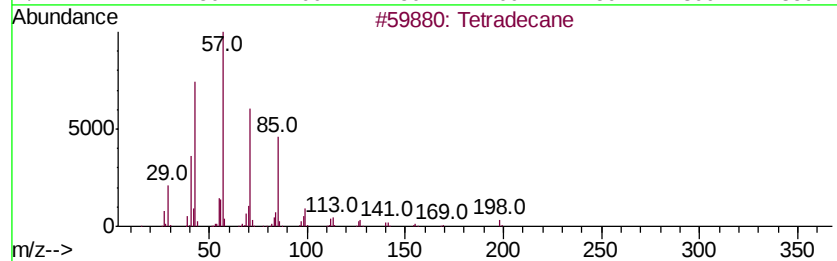
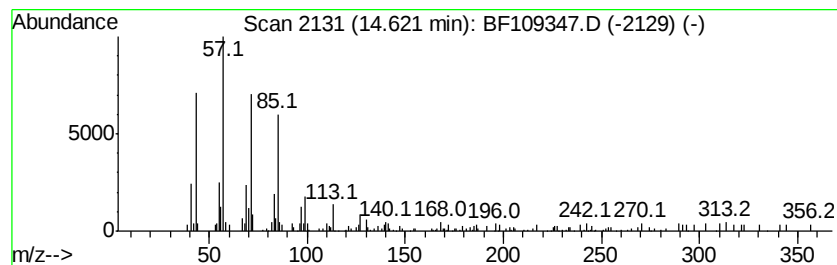
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Tetradecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.
14.62	2.24 ng	62090	Perylene-d12		15.24
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	70
2	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	64
3	Docosane	310	C22H46	000629-97-0	58
4	Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	58
5	Hexacosane	366	C26H54	000630-01-3	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092618\
Data File : BF109347.D
Acq On : 26 Sep 2018 19:55
Operator : JU/SJ
Sample : J4773-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CL-02-092518-B

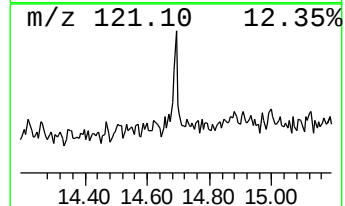
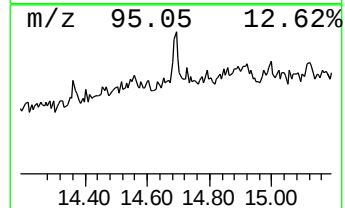
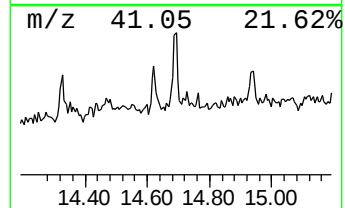
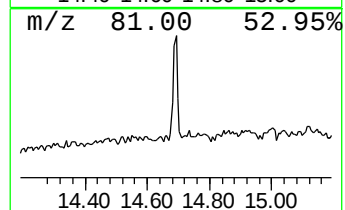
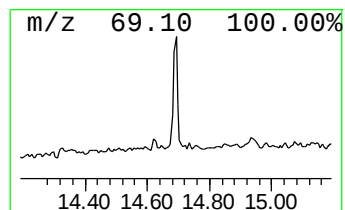
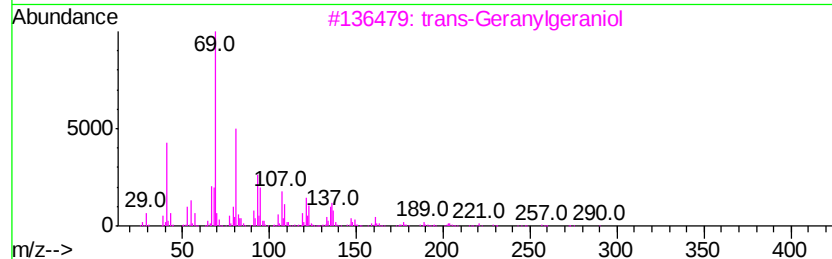
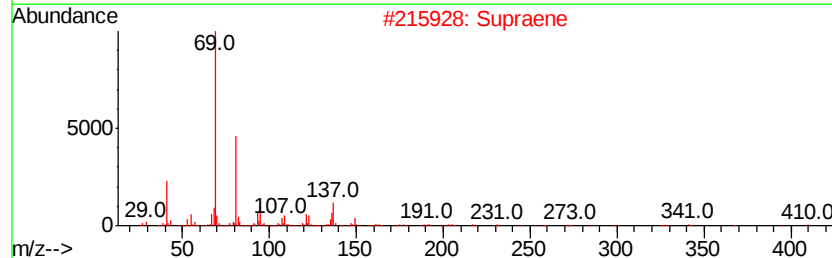
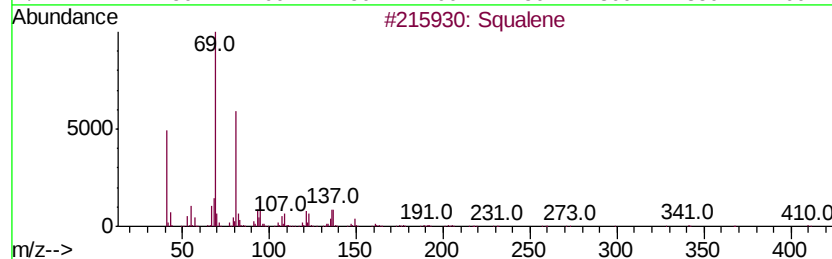
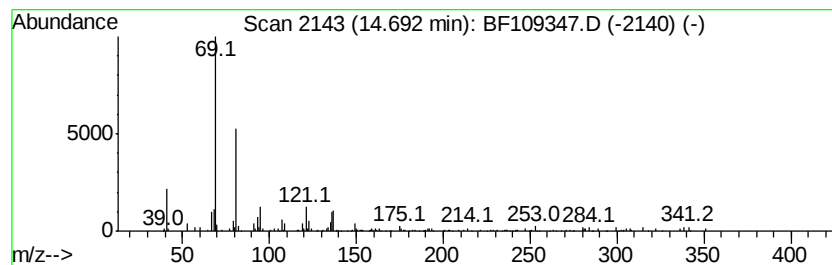
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Squalene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	2.69 ng	74430	Perylene-d12	15.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	86
2		Supraene	410	C30H50	007683-64-9	80
3		trans-Geranylgeraniol	290	C20H34O	024034-73-9	64
4		Cyclopropanecarboxylic acid, 4-i...	204	C13H16O2	1000331-01-9	58
5		4,9,13,17-Tetramethyl-4,8,12,16-...	316	C22H36O	056882-09-8	56



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092618\
 Data File : BF109347.D
 Acq On : 26 Sep 2018 19:55
 Operator : JU/SJ
 Sample : J4773-13
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-092518-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.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
unknown6.47	6.47	37.4	ng	1115710	1	6.70	596881	20.0
Heptadecane	9.69	2.0	ng	88254	3	9.74	872634	20.0
Sulfurous acid, 2...	10.18	2.4	ng	105119	3	9.74	872634	20.0
Tridecane	10.65	4.7	ng	179649	4	11.22	770462	20.0
Octadecane	11.52	2.7	ng	102557	4	11.22	770462	20.0
Hexadecanoic acid...	11.62	10.8	ng	414641	4	11.22	770462	20.0
n-Decanoic acid	11.75	2.2	ng	85526	4	11.22	770462	20.0
9,12-Octadecadien...	12.30	32.5	ng	1250940	4	11.22	770462	20.0
9-Octadecenoic ac...	12.32	27.7	ng	1065920	4	11.22	770462	20.0
Oxalic acid, isob...	13.71	4.2	ng	169058	5	13.84	803957	20.0
Tetradecane	14.62	2.2	ng	62090	6	15.24	554368	20.0
Squalene	14.69	2.7	ng	74430	6	15.24	554368	20.0