

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092418\
 Data File : BF109276.D
 Acq On : 24 Sep 2018 19:03
 Operator : JU/SJ
 Sample : J4770-15 2X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-092118-H

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	5.310	545	548	561	rVB	739885	662839	16.51%	3.005%
2	6.339	720	723	734	rVB	842024	659090	16.41%	2.988%
3	6.475	742	746	750	rBV	870547	684767	17.05%	3.105%
4	6.710	782	786	789	rBV	731910	564553	14.06%	2.560%
5	6.863	809	812	816	rBV	618408	519991	12.95%	2.358%
6	7.263	877	880	884	rBV	464323	376634	9.38%	1.708%
7	7.992	1000	1004	1007	rBV	818717	671530	16.72%	3.045%
8	9.063	1183	1186	1190	rBV	896854	660109	16.44%	2.993%
9	9.457	1250	1253	1259	rBV	200773	157441	3.92%	0.714%
10	9.739	1297	1301	1305	rBV2	883157	726979	18.10%	3.296%
11	10.186	1373	1377	1379	rBV	67276	57622	1.43%	0.261%
12	10.404	1410	1414	1416	rBV	41332	43028	1.07%	0.195%
13	10.522	1430	1434	1439	rBV	544110	436885	10.88%	1.981%
14	10.657	1453	1457	1464	rBV	119567	159988	3.98%	0.725%
15	11.104	1530	1533	1536	rBV2	119141	118235	2.94%	0.536%
16	11.133	1536	1538	1543	rVB	78808	71245	1.77%	0.323%
17	11.216	1549	1552	1555	rBV	923256	820114	20.42%	3.719%
18	11.239	1555	1556	1560	rVV	242431	136187	3.39%	0.618%
19	11.292	1562	1565	1568	rVB	79294	72242	1.80%	0.328%
20	11.527	1602	1605	1607	rBV	135877	100728	2.51%	0.457%
21	11.621	1618	1621	1625	rVB	1082463	906085	22.56%	4.108%
22	11.757	1640	1644	1647	rBV2	149446	163517	4.07%	0.741%
23	11.827	1653	1656	1659	rBV2	88359	98268	2.45%	0.446%
24	11.933	1671	1674	1680	rVB	140540	138155	3.44%	0.626%
25	12.268	1728	1731	1734	rBV2	49526	47983	1.19%	0.218%
26	12.310	1734	1738	1740	rBV	3928435	4015722	100.00%	18.208%
27	12.333	1740	1742	1749	rVB	3452189	2885594	71.86%	13.084%
28	12.416	1752	1756	1760	rBV2	858135	1133889	28.24%	5.141%
29	12.468	1762	1765	1772	rVB3	158045	230583	5.74%	1.046%
30	12.651	1791	1796	1799	rBV	743726	724114	18.03%	3.283%
31	12.692	1800	1803	1807	rVV2	88318	87660	2.18%	0.397%
32	12.798	1818	1821	1826	rVB	902938	738452	18.39%	3.348%
33	12.968	1847	1850	1852	rBV3	139681	155703	3.88%	0.706%
34	13.045	1860	1863	1867	rBV3	149913	181174	4.51%	0.821%

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 Sampling : 1 Min Area: 3 % of largest Peak
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 Stop Thrs : 0 Peak Location: TOP

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35	13.080	1867	1869	1872	rVB	60246	54348	1.35%	0.246%
36	13.139	1876	1879	1881	rBV2	103038	92446	2.30%	0.419%
37	13.174	1883	1885	1888	rVB	54340	41596	1.04%	0.189%
38	13.204	1888	1890	1894	rBV3	54021	68191	1.70%	0.309%
39	13.386	1918	1921	1924	rBV	81839	82074	2.04%	0.372%
40	13.551	1947	1949	1952	rVB3	51398	58029	1.45%	0.263%
41	13.710	1974	1976	1981	rVB	78083	78348	1.95%	0.355%
42	13.804	1989	1992	1994	rBV	77527	70279	1.75%	0.319%
43	13.845	1994	1999	2002	rVV2	860791	1005851	25.05%	4.561%
44	13.868	2002	2003	2006	rVB	239269	141262	3.52%	0.641%
45	14.027	2028	2030	2033	rVB2	58991	45947	1.14%	0.208%
46	14.727	2146	2149	2154	rBV	146369	178982	4.46%	0.812%
47	14.851	2167	2170	2179	rVB	188747	351019	8.74%	1.592%
48	15.192	2225	2228	2234	rVB	136423	152121	3.79%	0.690%
49	15.251	2234	2238	2241	rBV	384915	426689	10.63%	1.935%

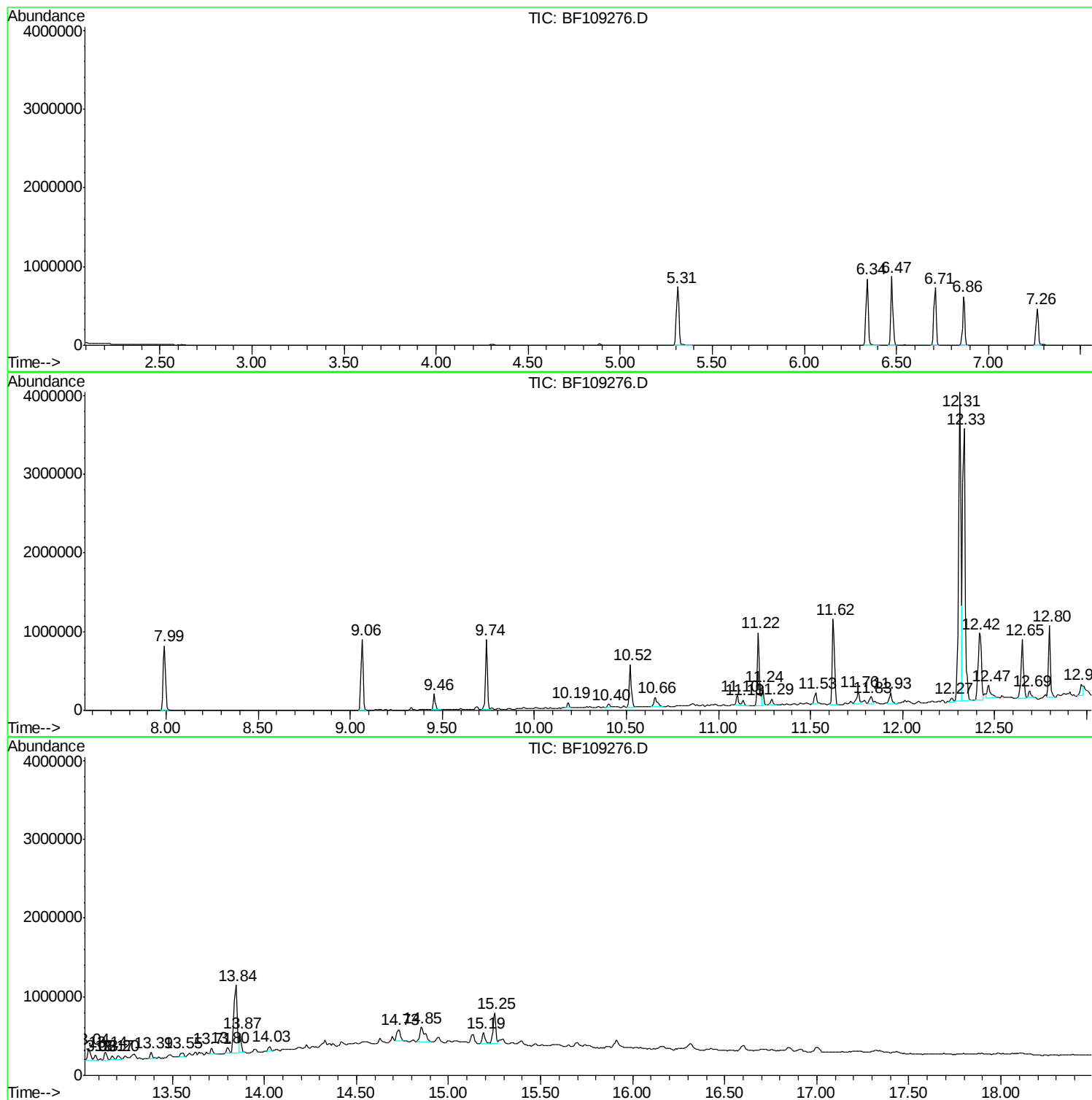
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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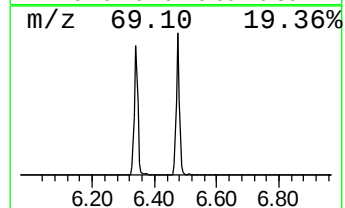
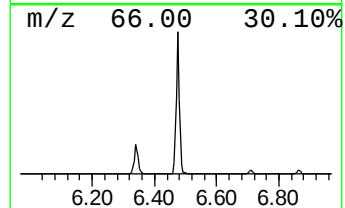
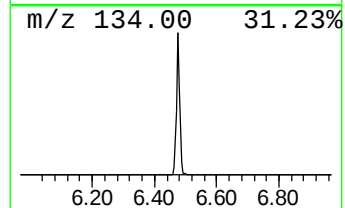
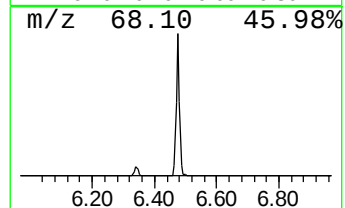
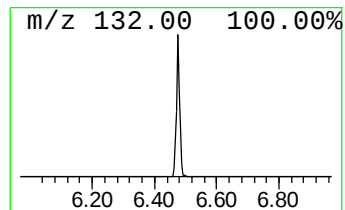
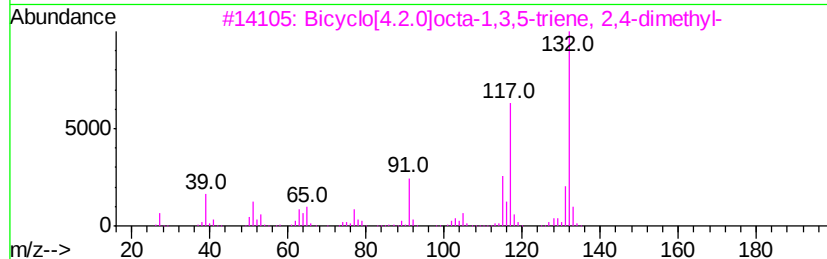
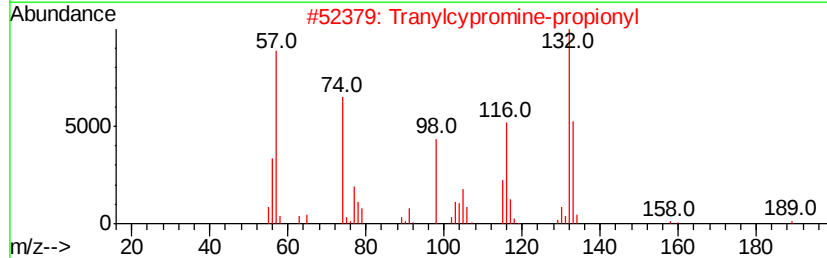
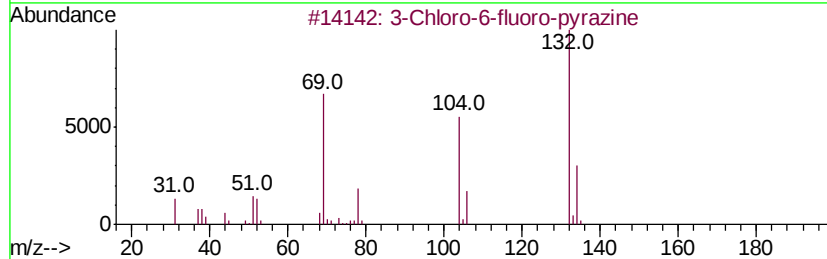
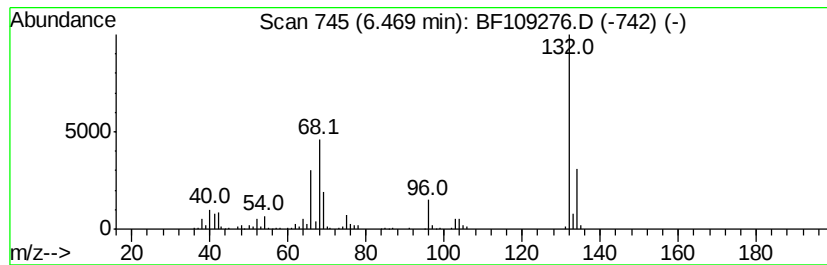
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.47 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	24.26 ng	684767	1,4-Dichlorobenzene-d4	6.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	16
2	Tranylcypromine-propionyl			189	C12H15NO	1000123-86-3	10
3	Bicvclo[4.2.0]octa-1,3,5-triene, ...			132	C10H12	028749-81-7	9
4	(E)-3-Chloro-2-methyl-2-pentenal			132	C6H9ClO	031357-76-3	9
5	4-Chloro-2-fluoro-pyrimidine			132	C4H2ClFN2	051422-00-5	9



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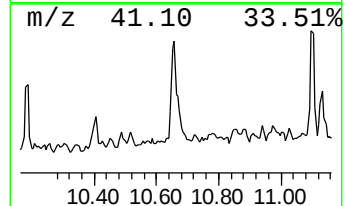
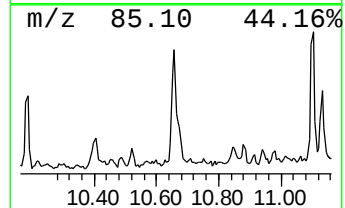
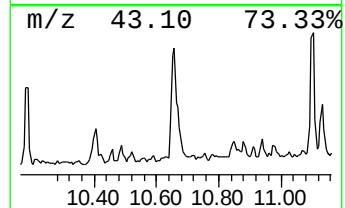
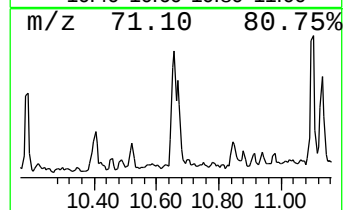
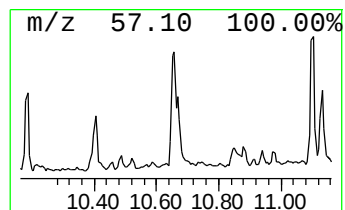
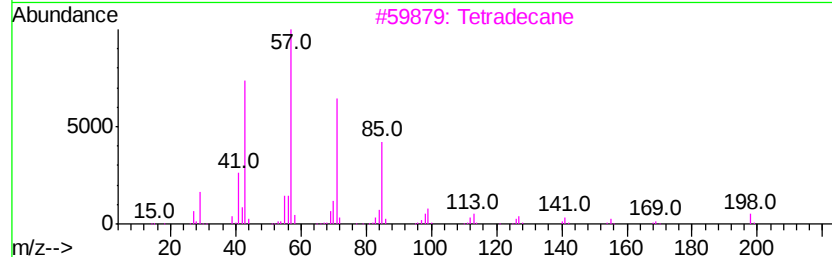
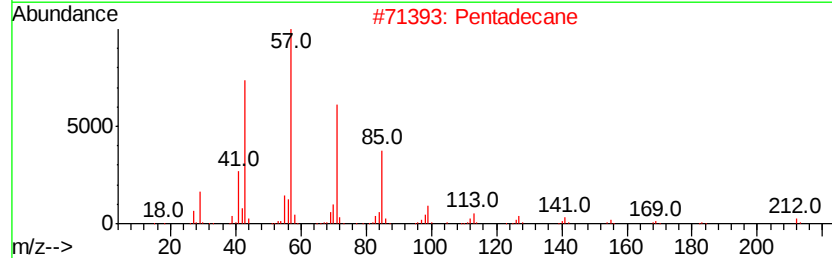
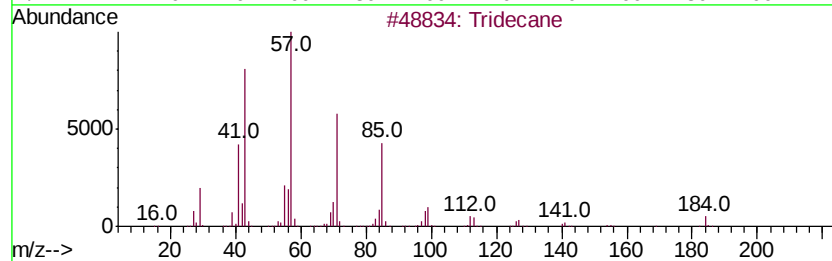
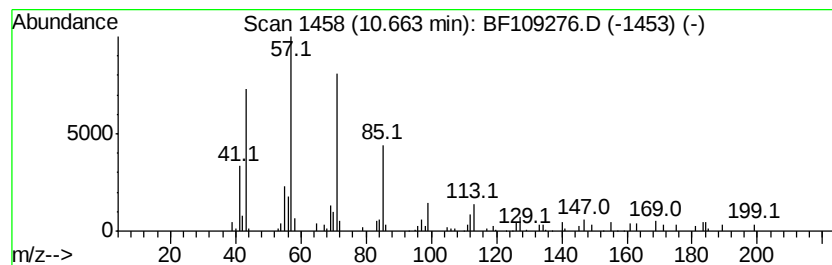
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Tridecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.66	3.90 ng	159988	Phenanthrene-d10	11.22

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane		184	C13H28	000629-50-5	93
2	Pentadecane		212	C15H32	000629-62-9	86
3	Tetradecane		198	C14H30	000629-59-4	86
4	Heptadecane		240	C17H36	000629-78-7	86
5	Heneicosane		296	C21H44	000629-94-7	86



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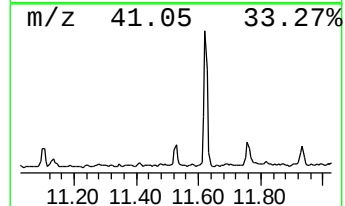
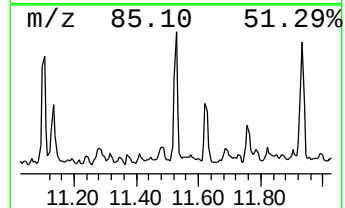
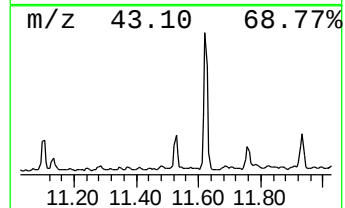
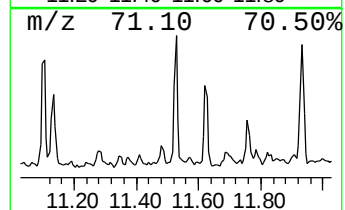
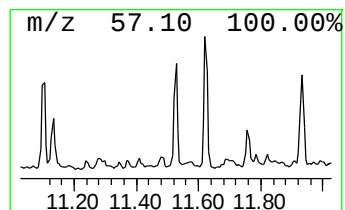
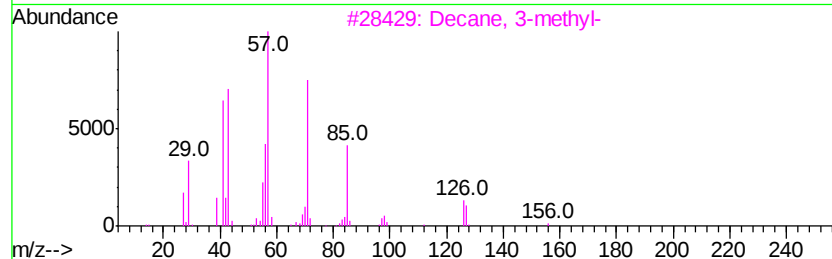
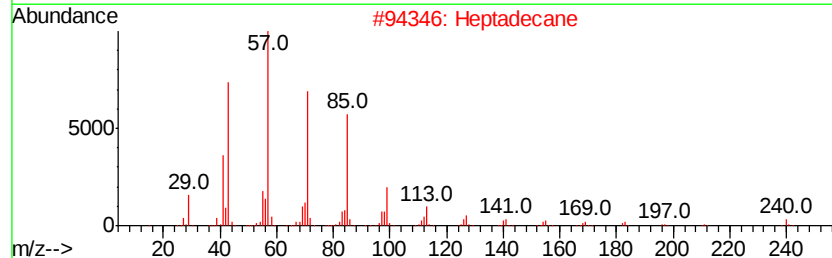
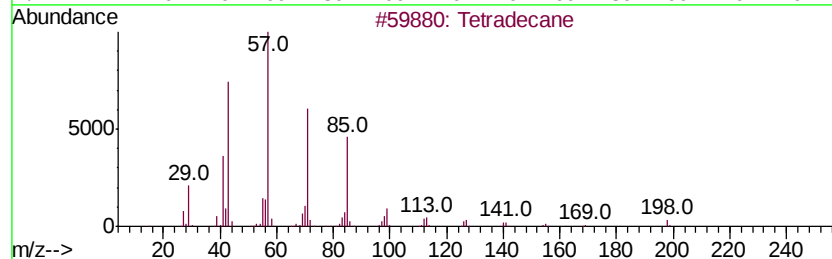
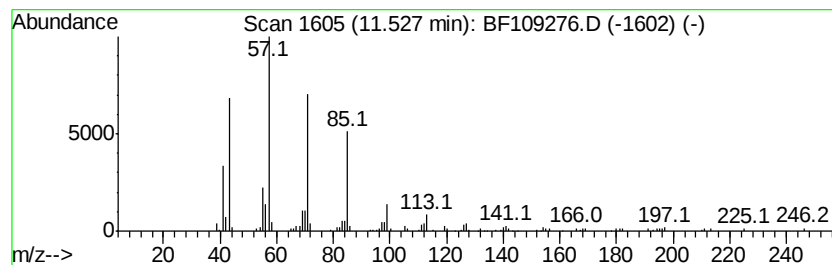
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Tetradecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.53	2.46 ng	100728	Phenanthrene-d10	11.22

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	86
2		Heptadecane	240	C17H36	000629-78-7	83
3		Decane, 3-methyl-	156	C11H24	013151-34-3	74
4		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	68
5		Tetracosane	338	C24H50	000646-31-1	68



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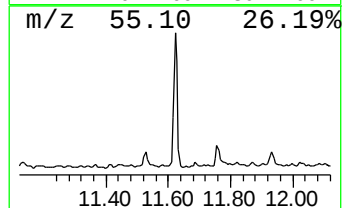
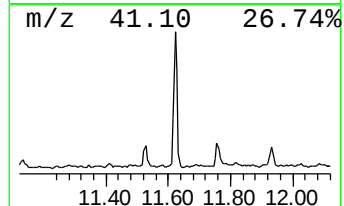
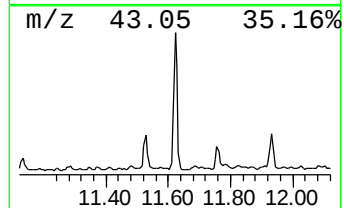
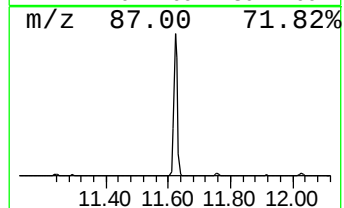
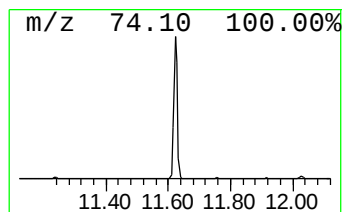
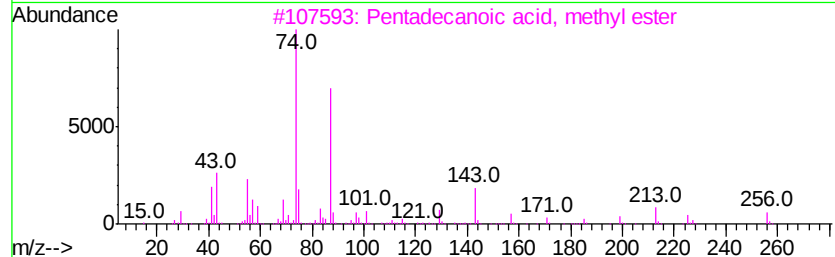
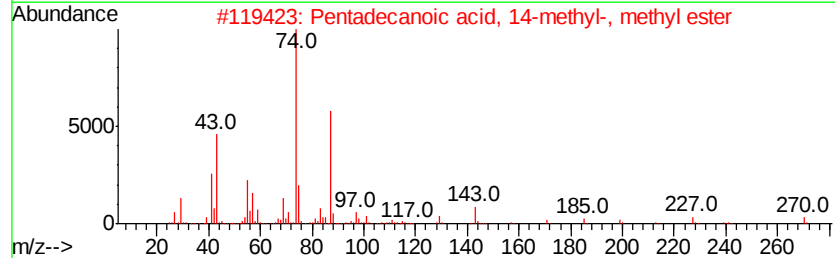
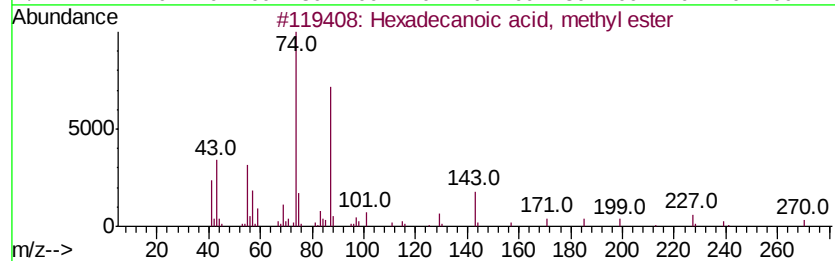
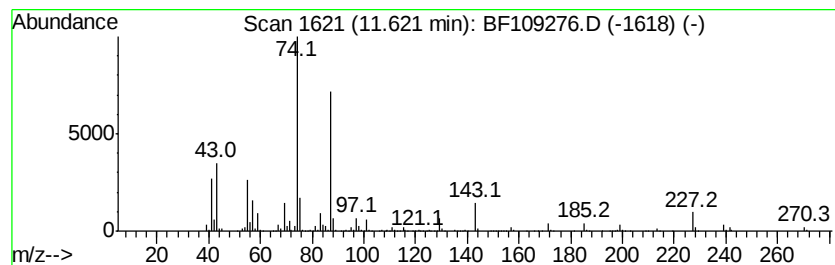
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Hexadecanoic acid, methyl e... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.62	22.10 ng	906085	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	87
3	Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86
4	Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	80
5	Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	80



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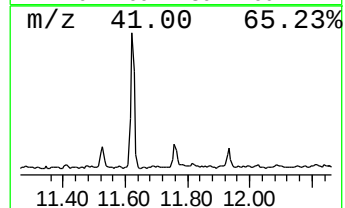
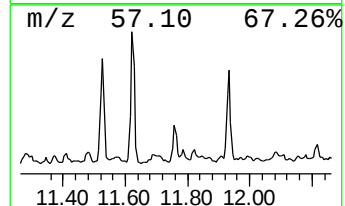
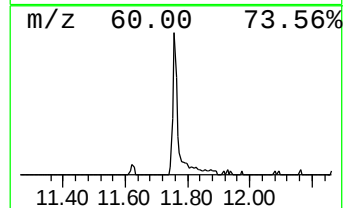
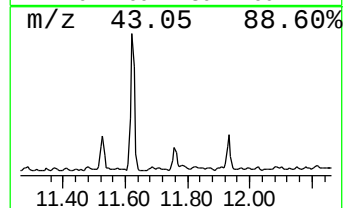
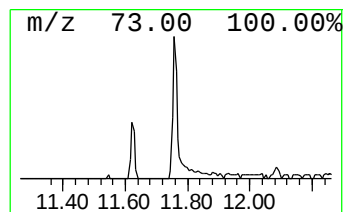
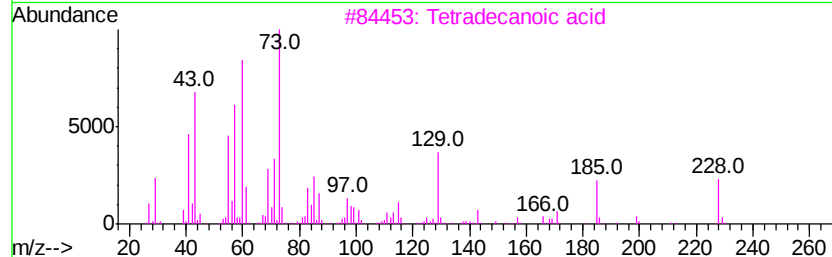
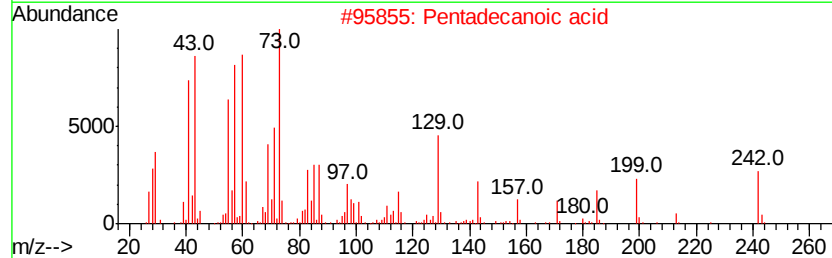
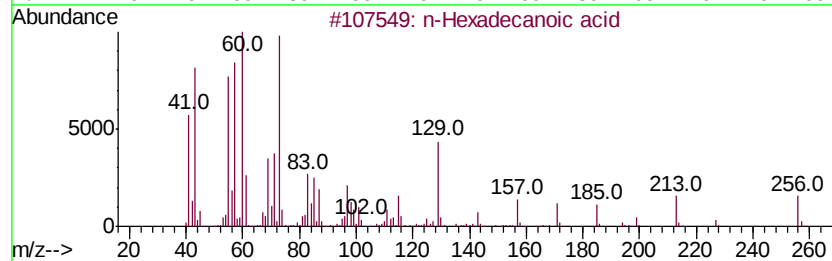
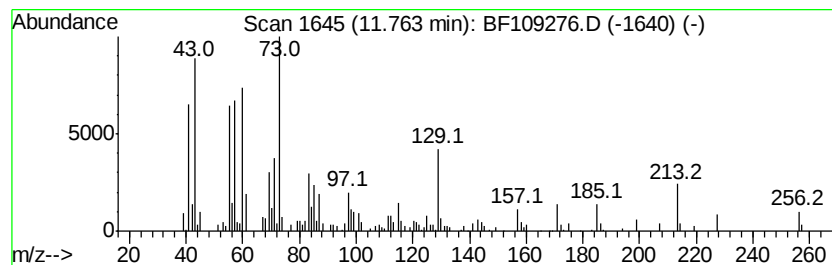
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.76	3.99 ng	163517	Phenanthrene-d10	11.22

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4		Tridecanoic acid	214	C13H26O2	000638-53-9	72
5		n-Decanoic acid	172	C10H20O2	000334-48-5	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092418\
Data File : BF109276.D
Acq On : 24 Sep 2018 19:03
Operator : JU/SJ
Sample : J4770-15 2X
Misc :
ALS Vial : 23 Sample Multiplier: 1

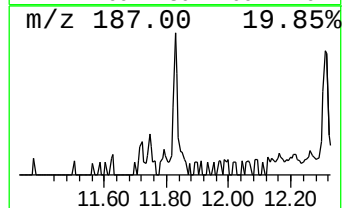
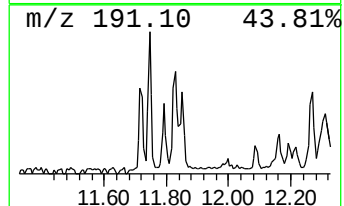
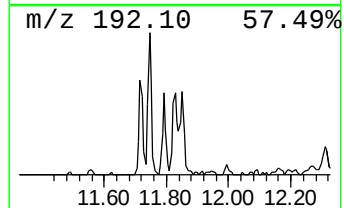
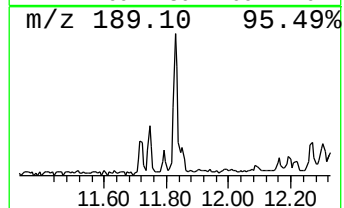
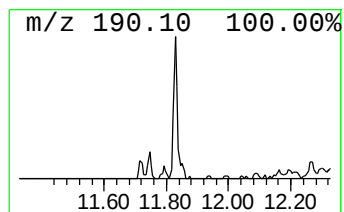
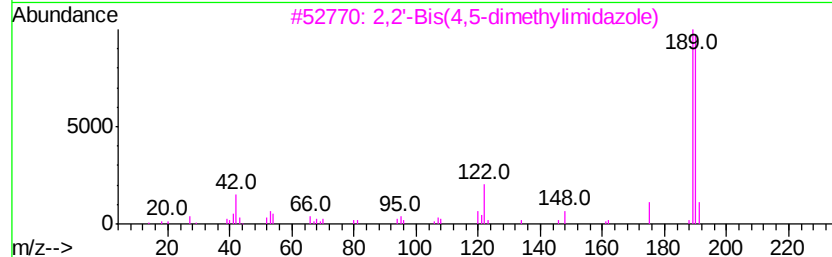
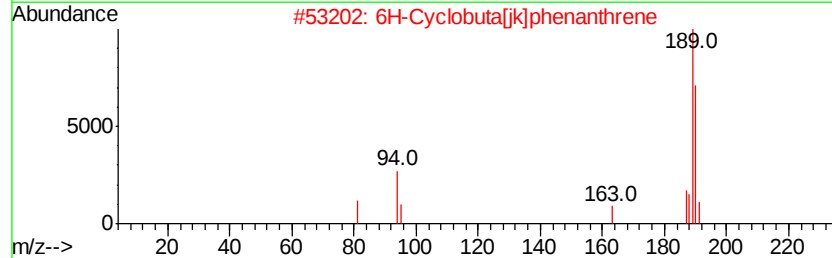
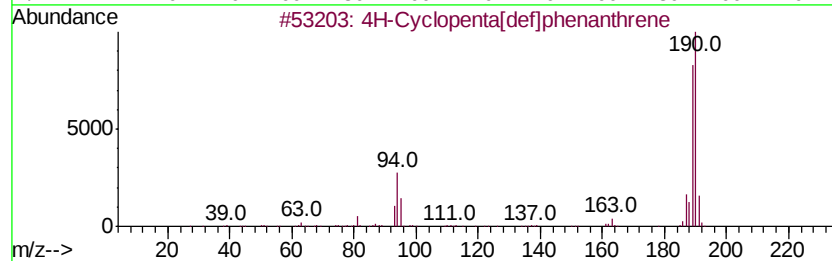
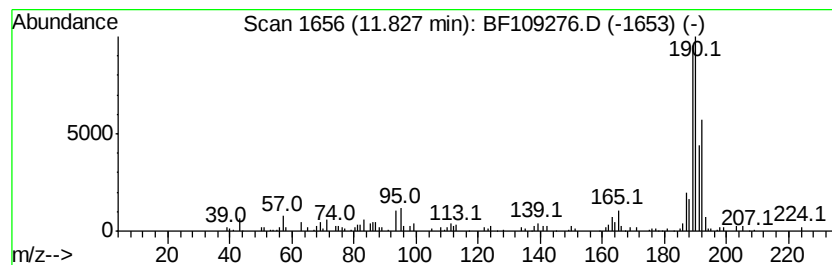
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 4H-Cyclopenta[def]phenanthrene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.83	2.40 ng	98268	Phenanthrene-d10	11.22		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2		6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	53
3		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	47
4		Benzene, 1,1'-(1,2-propadienylid...	192	C15H12	014251-57-1	25
5		Propyne, 1,3-diphenyl-	192	C15H12	004980-70-5	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092418\
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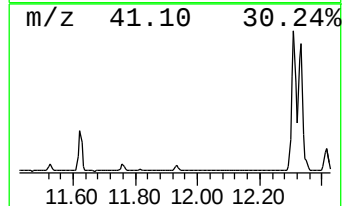
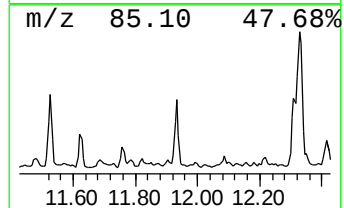
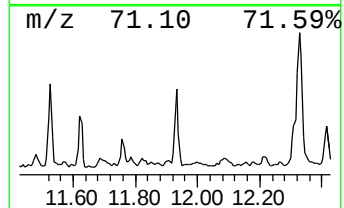
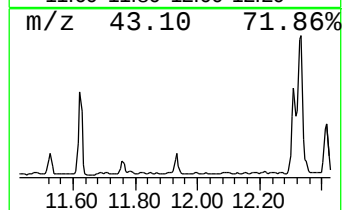
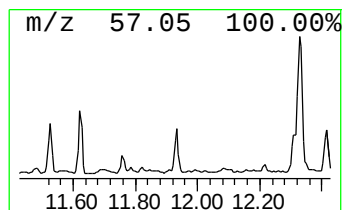
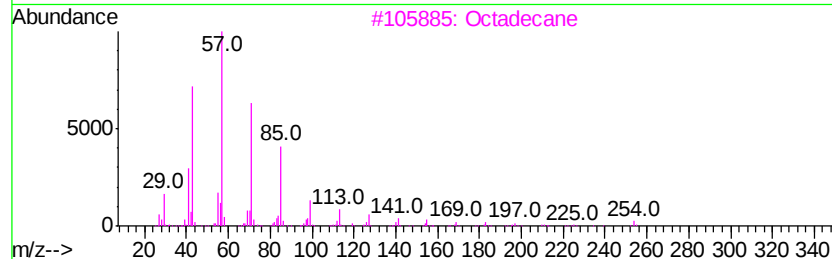
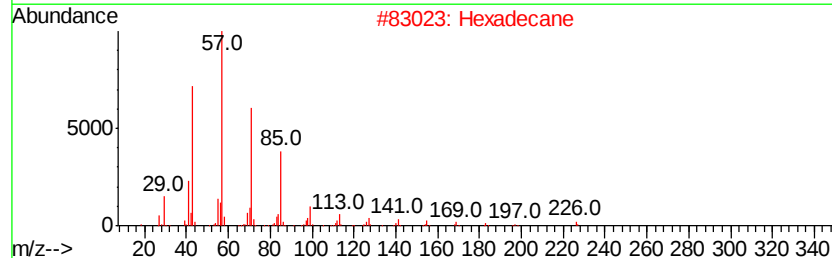
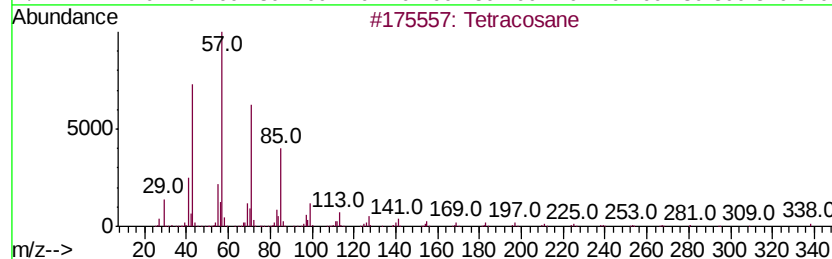
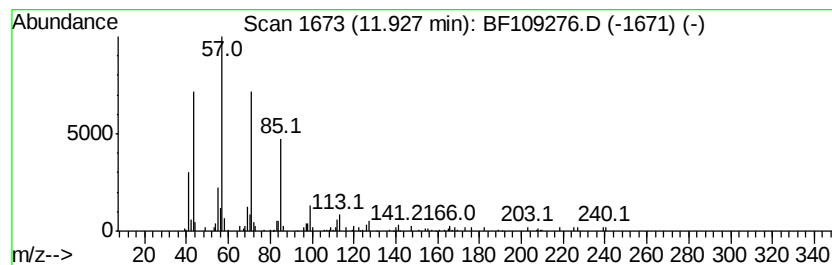
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetracosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	3.37 ng	138155	Phenanthrene-d10	11.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	91
2		Hexadecane	226	C16H34	000544-76-3	90
3		Octadecane	254	C18H38	000593-45-3	90
4		Tetradecane	198	C14H30	000629-59-4	90
5		Pentadecane	212	C15H32	000629-62-9	90



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Misc :
ALS Vial : 23 Sample Multiplier: 1

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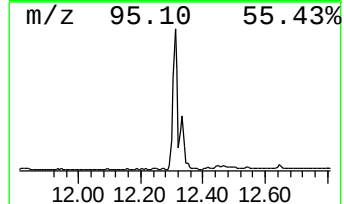
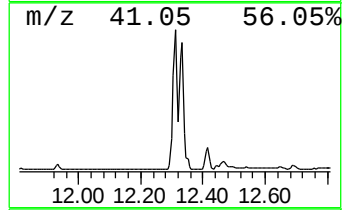
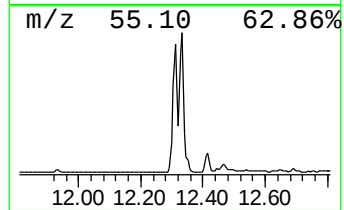
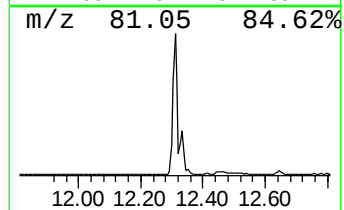
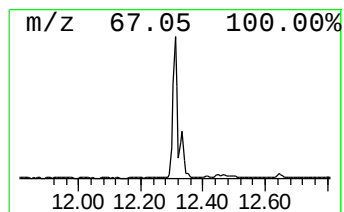
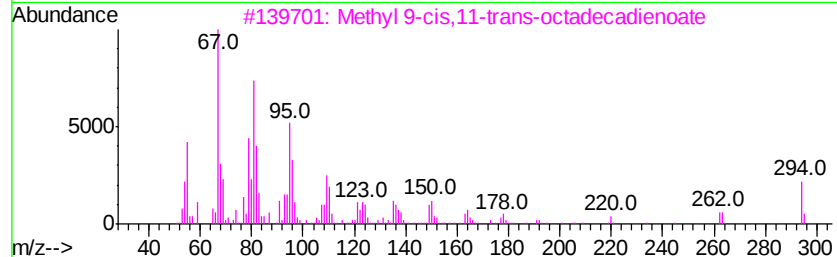
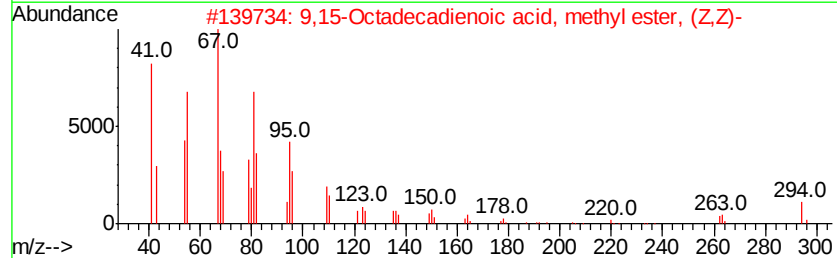
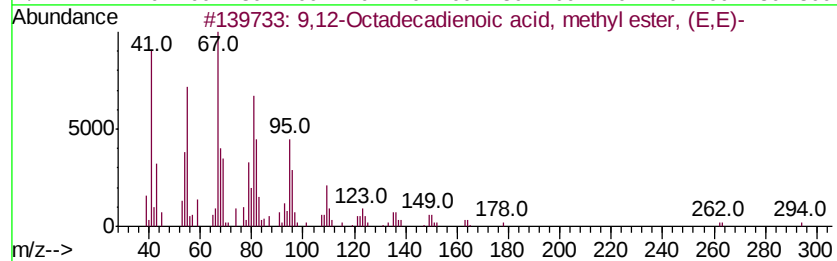
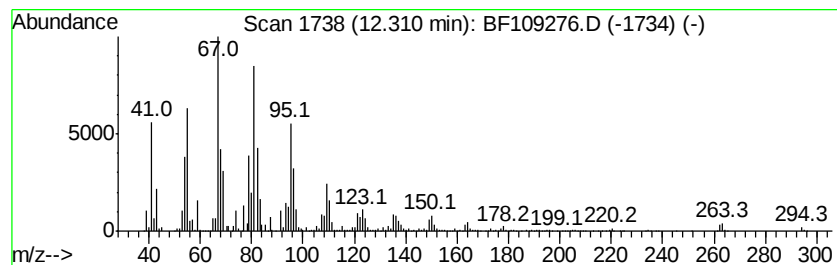
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.31	97.93 ng	4015720	Phenanthrene-d10	11.22

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		Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	1000336-44-0	98
4		10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	97
5		11,14-Octadecadienoic acid, meth...	294	C19H34O2	056554-61-1	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092418\
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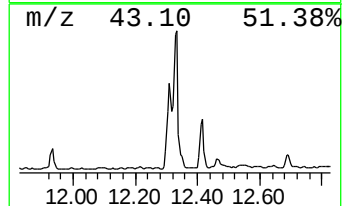
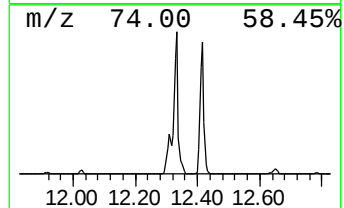
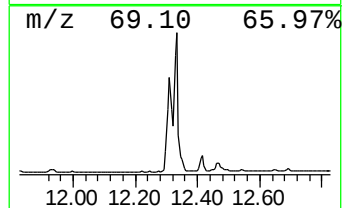
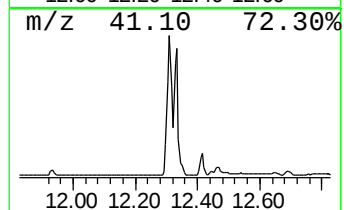
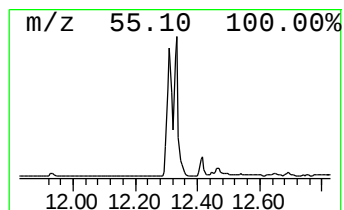
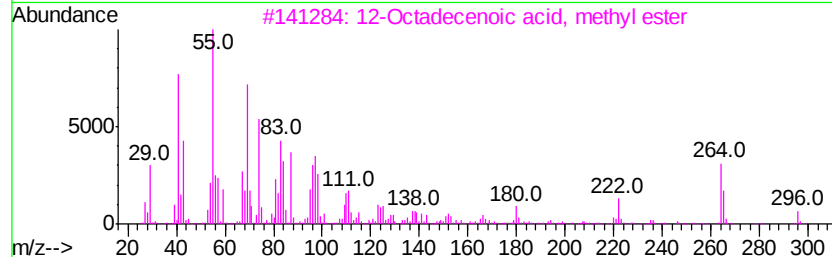
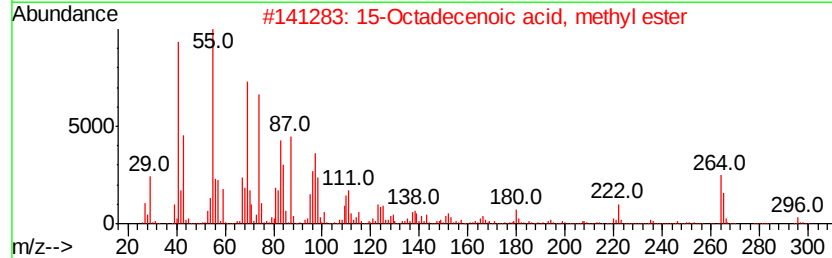
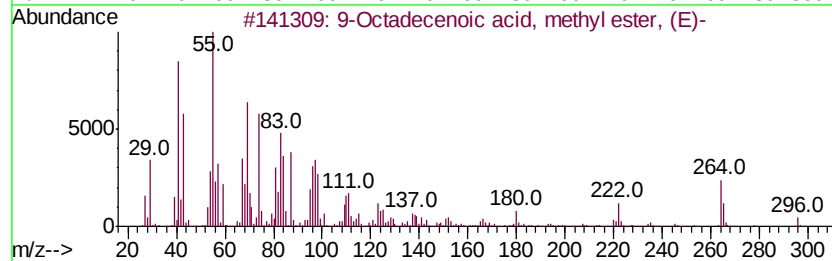
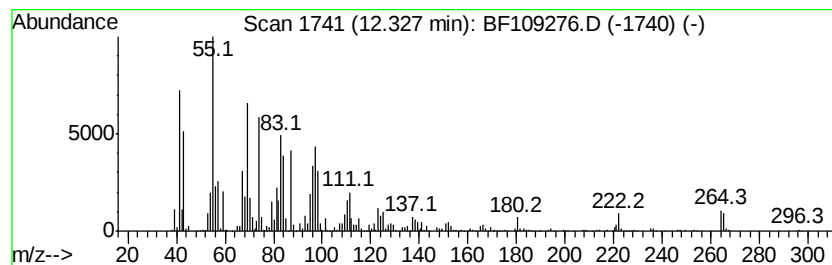
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 9-Octadecenoic acid, methyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.33	70.37 ng	2885590	Phenanthrene-d10	11.22	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	99
2	15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	99
3	12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	99
4	16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	99
5	6-Octadecenoic acid, methyl este...	296	C19H36O2	002777-58-4	99



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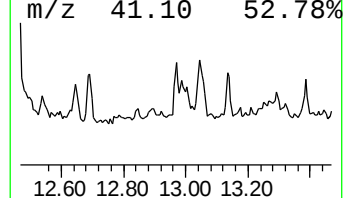
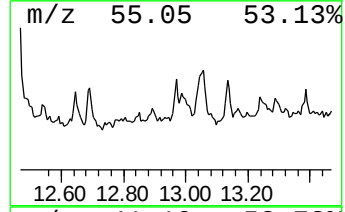
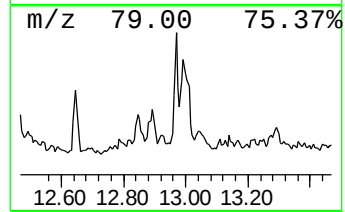
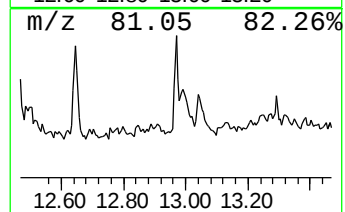
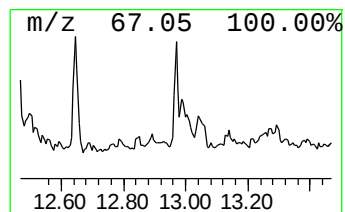
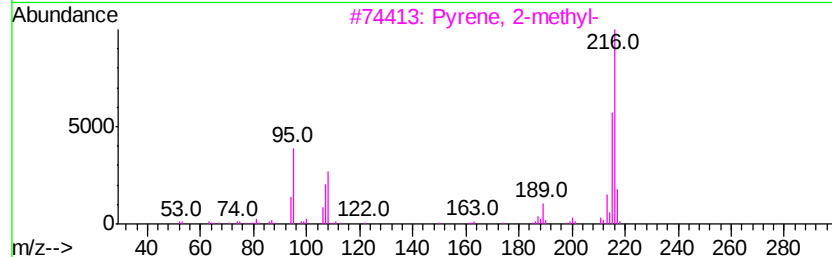
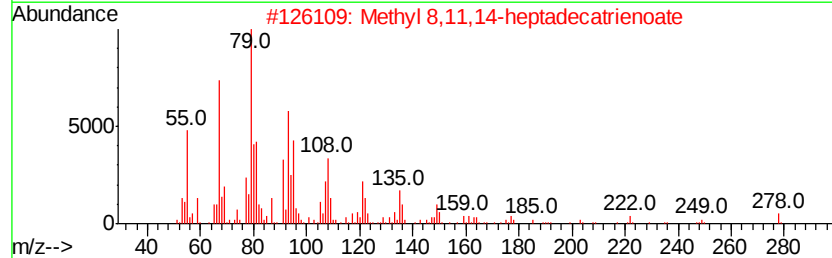
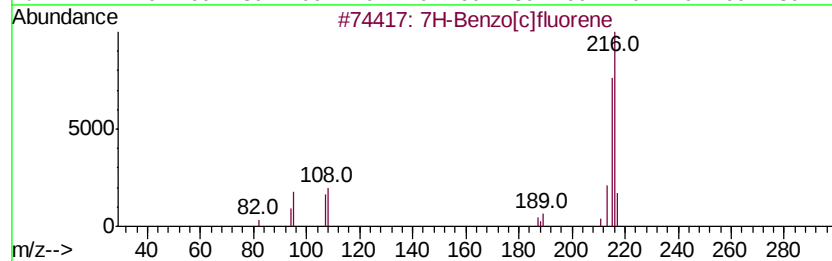
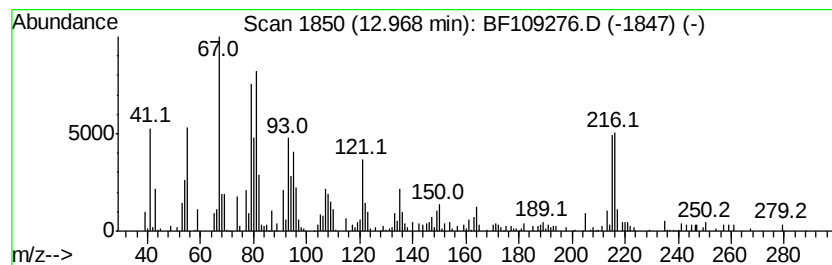
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown12.97 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.97	3.10 ng	155703	Chrysene-d12	13.84	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	46
2	Methyl 8,11,14-heptadecatrienoate	278	C18H30O2	1000336-35-1	38
3	Pvrene, 2-methyl-	216	C17H12	003442-78-2	35
4	4-Hexadecen-6-yne, (E)-	220	C16H28	074744-51-7	30
5	Spiro[2.5]octane	110	C8H14	000185-65-9	30



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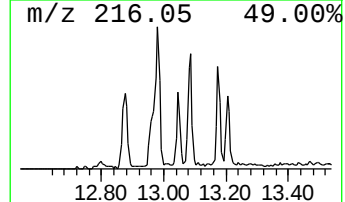
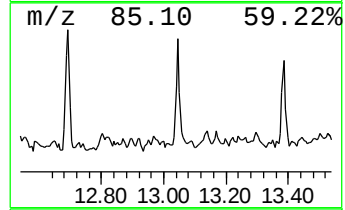
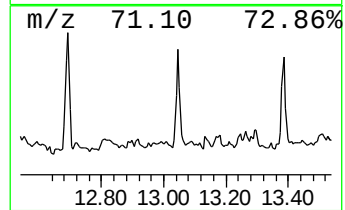
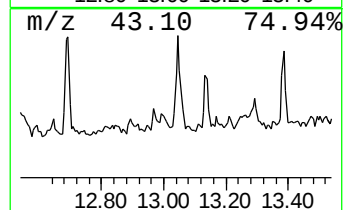
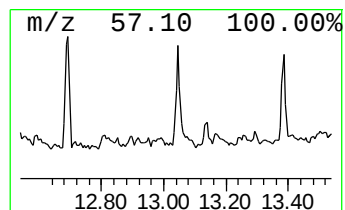
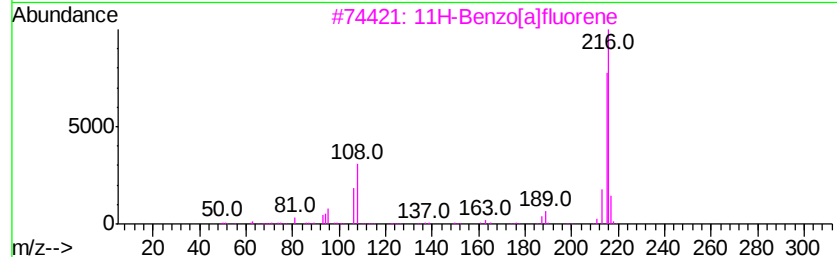
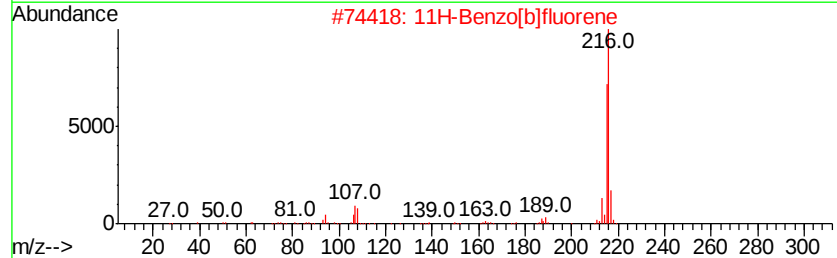
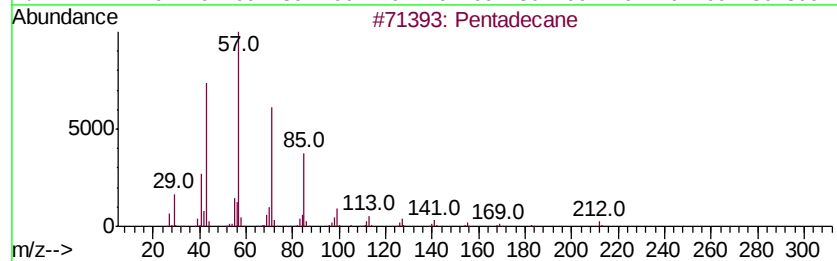
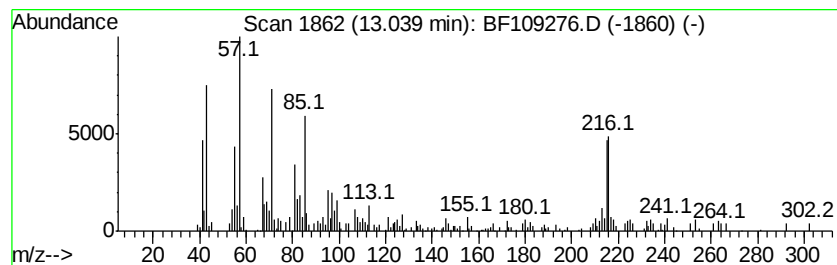
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Pentadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.04	3.60 ng	181174	Chrysene-d12		13.84
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane	212	C15H32	000629-62-9	83
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	50
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	50
4	Hexadecane	226	C16H34	000544-76-3	46
5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	45



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ALS Vial : 23 Sample Multiplier: 1

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ClientSampleId :
JC-02-092118-H

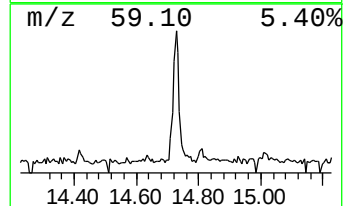
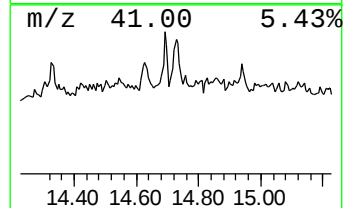
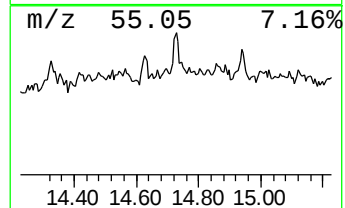
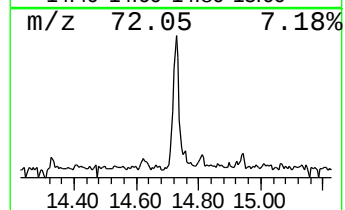
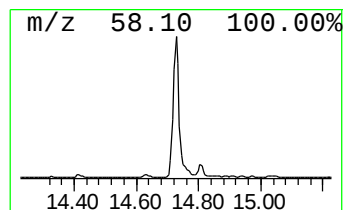
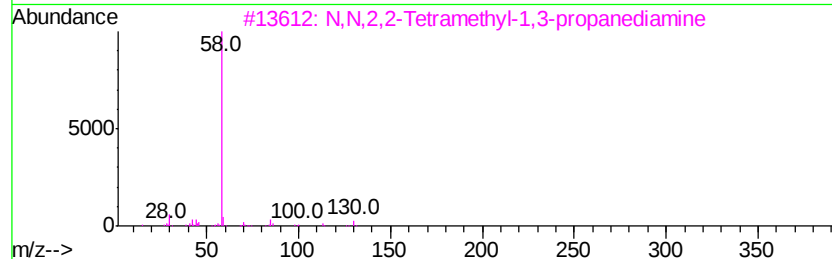
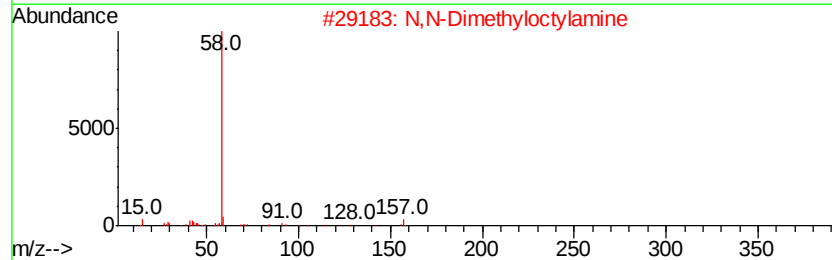
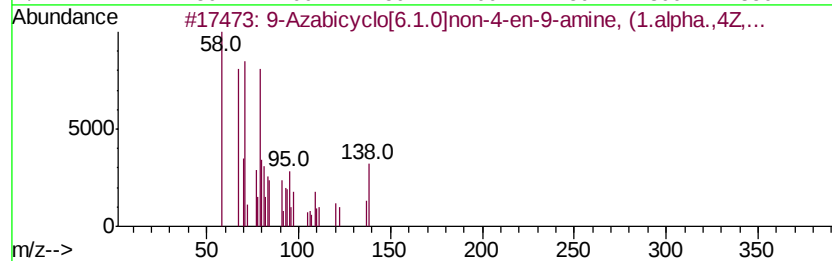
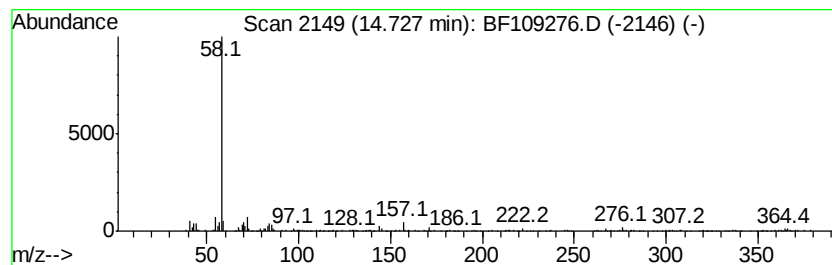
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 9-Azabicyclo[6.1.0]non-4-en... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.73	8.39 ng	178982	Perylene-d12	15.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Azabicyclo[6.1.0]non-4-en-9-am...	138	C8H14N2	066387-78-8	76
2		N,N-Dimethyloctylamine	157	C10H23N	007378-99-6	64
3		N,N,2,2-Tetramethyl-1,3-propaned...	130	C7H18N2	053369-71-4	59
4		5H-Dibenzo[a,d]cyclohepten-5-ol,...	295	C20H25NO	001159-03-1	59
5		1-Dodecanamine, N,N-dimethyl-	213	C14H31N	000112-18-5	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092418\
 Data File : BF109276.D
 Acq On : 24 Sep 2018 19:03
 Operator : JU/SJ
 Sample : J4770-15 2X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-092118-H

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	24.3	ng	684767	1	6.71	564553	20.0
Tridecane	10.66	3.9	ng	159988	4	11.22	820114	20.0
Tetradecane	11.53	2.5	ng	100728	4	11.22	820114	20.0
Hexadecanoic acid...	11.62	22.1	ng	906085	4	11.22	820114	20.0
n-Hexadecanoic acid	11.76	4.0	ng	163517	4	11.22	820114	20.0
4H-Cyclopenta[def...	11.83	2.4	ng	98268	4	11.22	820114	20.0
Tetracosane	11.93	3.4	ng	138155	4	11.22	820114	20.0
9,12-Octadecadien...	12.31	97.9	ng	4015720	4	11.22	820114	20.0
9-Octadecenoic ac...	12.33	70.4	ng	2885590	4	11.22	820114	20.0
unknown12.97	12.97	3.1	ng	155703	5	13.84	1005850	20.0
Pentadecane	13.04	3.6	ng	181174	5	13.84	1005850	20.0
9-Azabicyclo[6.1....	14.73	8.4	ng	178982	6	15.25	426689	20.0