

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092418\
 Data File : BF109275.D
 Acq On : 24 Sep 2018 18:37
 Operator : JU/SJ
 Sample : J4770-13 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

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

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	544	548	558	rBV	982996	841795	11.92%	2.791%
2	6.339	719	723	726	rBV	1118585	887513	12.57%	2.943%
3	6.475	742	746	749	rBV	1143841	906515	12.84%	3.006%
4	6.710	782	786	789	rBV	868316	696936	9.87%	2.311%
5	6.863	809	812	816	rVB	879369	693532	9.82%	2.300%
6	7.263	876	880	888	rBV	629244	543365	7.70%	1.802%
7	7.992	1000	1004	1007	rBV	1111157	925793	13.11%	3.070%
8	9.063	1183	1186	1190	rBV	1181208	945100	13.39%	3.134%
9	9.457	1250	1253	1256	rBV	289810	223114	3.16%	0.740%
10	9.739	1297	1301	1305	rBV2	1156350	987377	13.99%	3.274%
11	10.404	1410	1414	1417	rBV	66110	70716	1.00%	0.234%
12	10.522	1431	1434	1439	rBV	677012	544959	7.72%	1.807%
13	10.657	1454	1457	1458	rBV	153512	132829	1.88%	0.440%
14	11.104	1530	1533	1536	rBV2	169744	155295	2.20%	0.515%
15	11.133	1536	1538	1544	rVB	94424	87238	1.24%	0.289%
16	11.216	1549	1552	1555	rBV	977159	902421	12.78%	2.992%
17	11.239	1555	1556	1560	rVV	287783	180062	2.55%	0.597%
18	11.292	1560	1565	1568	rVB	86398	90602	1.28%	0.300%
19	11.527	1602	1605	1607	rBV	183575	135045	1.91%	0.448%
20	11.627	1618	1622	1625	rVB	1417614	1209889	17.14%	4.012%
21	11.757	1640	1644	1647	rBV2	137976	149771	2.12%	0.497%
22	11.827	1653	1656	1659	rBV2	78267	86629	1.23%	0.287%
23	11.933	1671	1674	1679	rVB	180991	163443	2.32%	0.542%
24	12.316	1733	1739	1741	rBV	5877511	7059325	100.00%	23.409%
25	12.339	1741	1743	1748	rVB	5356581	4422582	62.65%	14.665%
26	12.416	1752	1756	1760	rBV2	1161026	1201907	17.03%	3.986%
27	12.468	1762	1765	1767	rVB	124361	110943	1.57%	0.368%
28	12.651	1791	1796	1799	rBV	552977	551037	7.81%	1.827%
29	12.692	1800	1803	1810	rVB2	161244	199141	2.82%	0.660%
30	12.798	1818	1821	1827	rVB	823969	693652	9.83%	2.300%
31	12.968	1847	1850	1860	rBV2	228189	379849	5.38%	1.260%
32	13.045	1860	1863	1867	rBV2	241091	265102	3.76%	0.879%
33	13.139	1876	1879	1882	rBV	131616	123103	1.74%	0.408%
34	13.386	1918	1921	1924	rBV	173123	138630	1.96%	0.460%

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JC-02-092118-G

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.715	1974	1977	1982	rVB	160066	142079	2.01%	0.471%
36	13.804	1989	1992	1994	rBV	96536	96444	1.37%	0.320%
37	13.845	1994	1999	2002	rVV2	899337	996639	14.12%	3.305%
38	13.868	2002	2003	2006	rVB	181131	107534	1.52%	0.357%
39	14.027	2027	2030	2033	rVB2	188301	147206	2.09%	0.488%
40	14.333	2079	2082	2084	rVB	172192	160240	2.27%	0.531%
41	14.627	2129	2132	2136	rBV	175802	163656	2.32%	0.543%
42	14.727	2146	2149	2158	rVB	180794	278614	3.95%	0.924%
43	14.851	2167	2170	2180	rVB	158156	284968	4.04%	0.945%
44	14.939	2183	2185	2190	rVB2	171297	193777	2.74%	0.643%
45	15.127	2215	2217	2223	rVB	97427	110368	1.56%	0.366%
46	15.186	2225	2227	2233	rVB	121103	126466	1.79%	0.419%
47	15.251	2234	2238	2241	rBV	426946	495343	7.02%	1.643%
48	15.698	2311	2314	2318	rVB	115839	148185	2.10%	0.491%

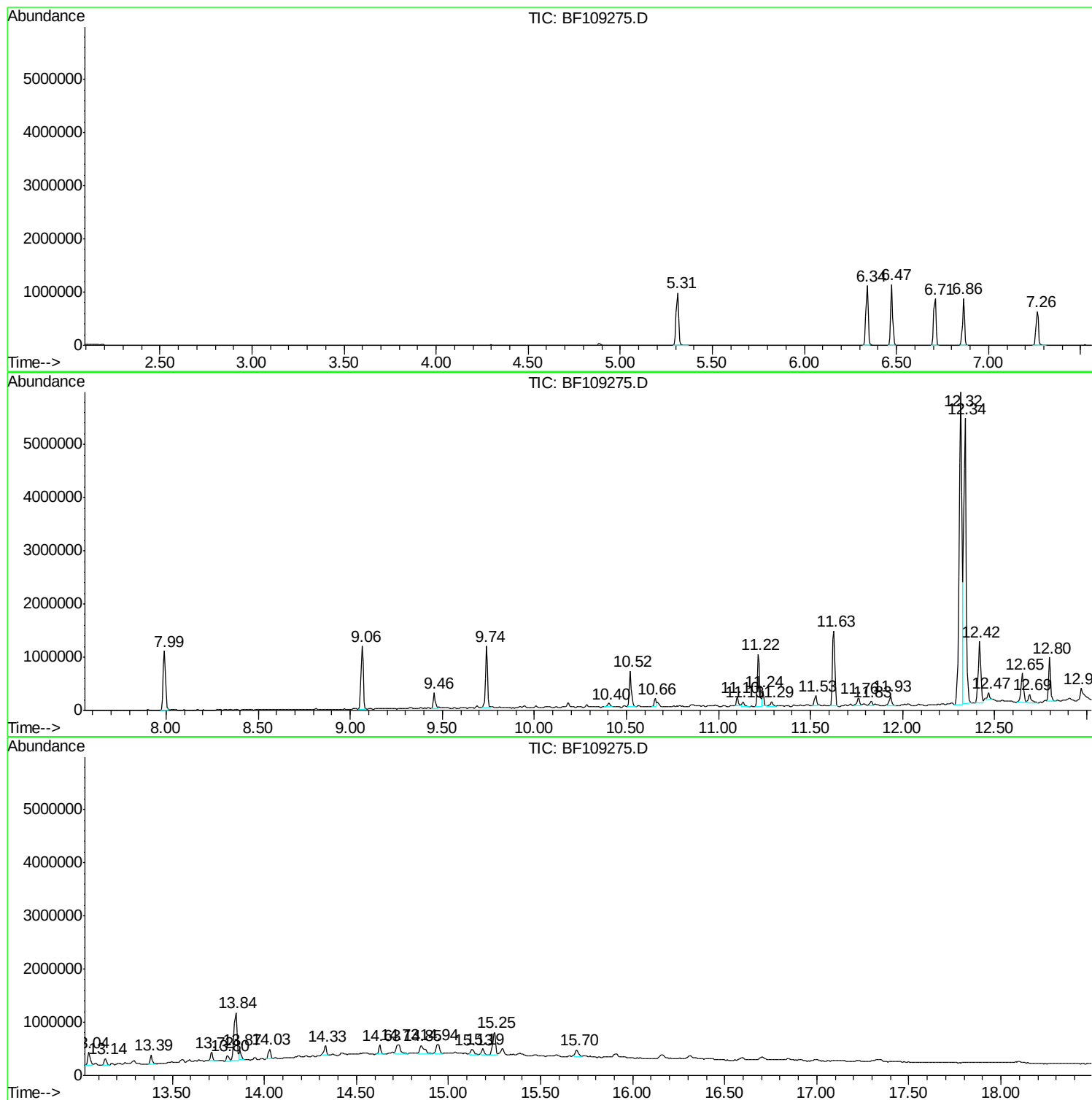
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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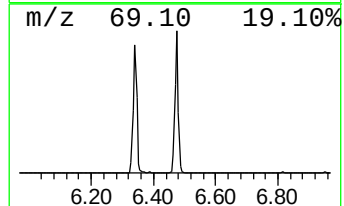
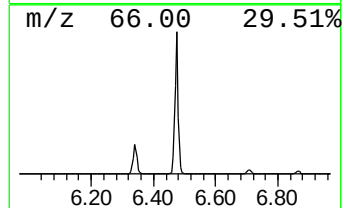
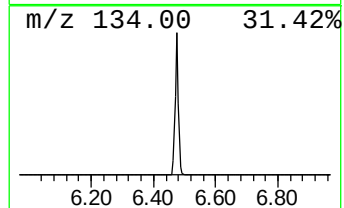
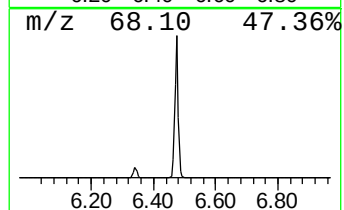
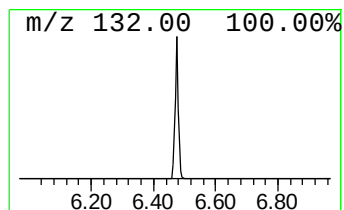
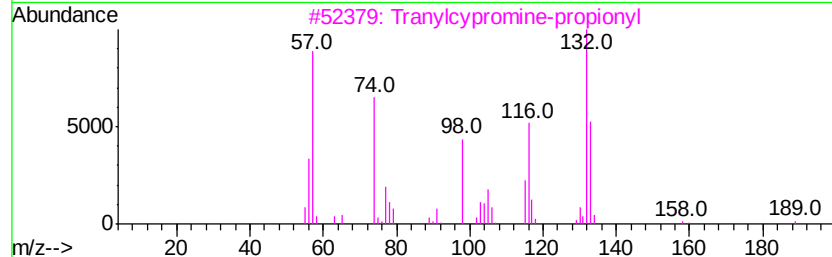
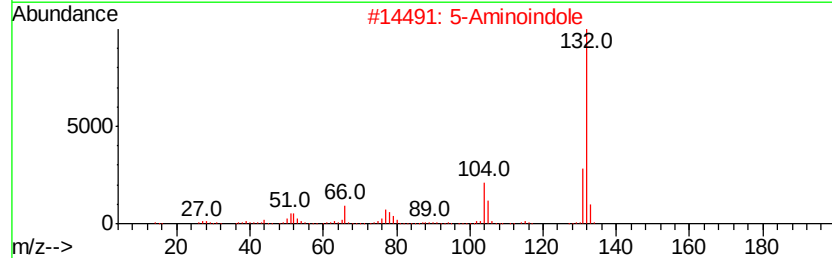
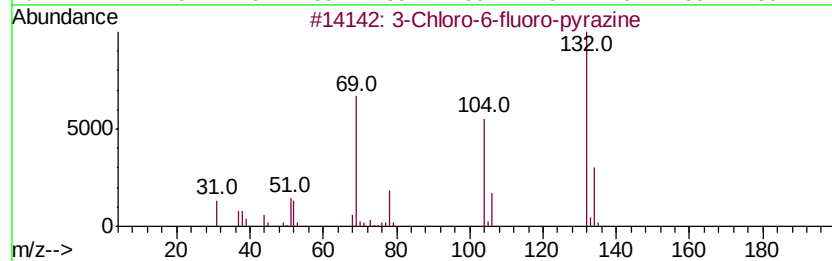
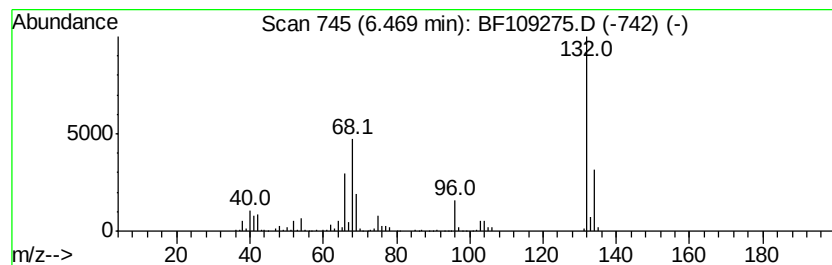
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.47 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	26.01 ng	906515	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	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	10
3			Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4			4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5			2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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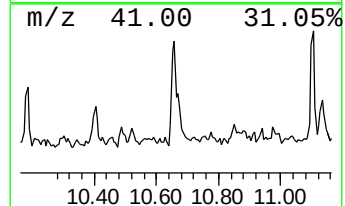
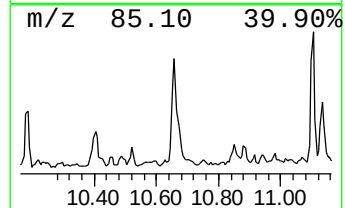
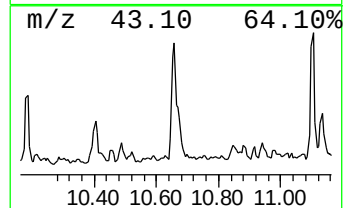
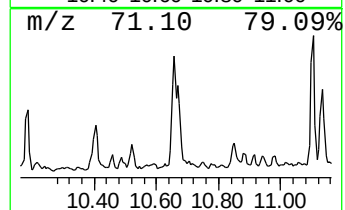
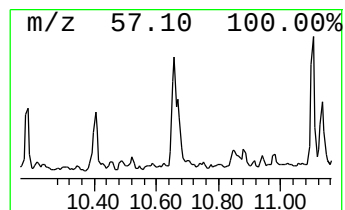
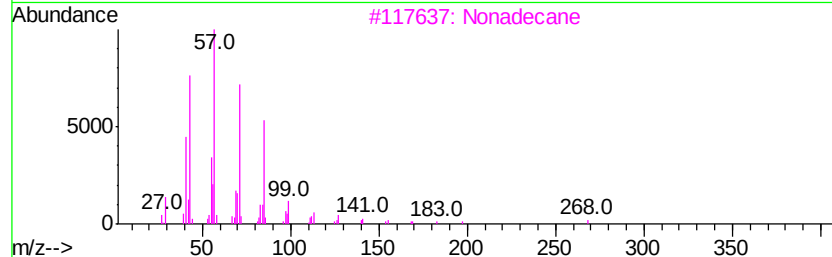
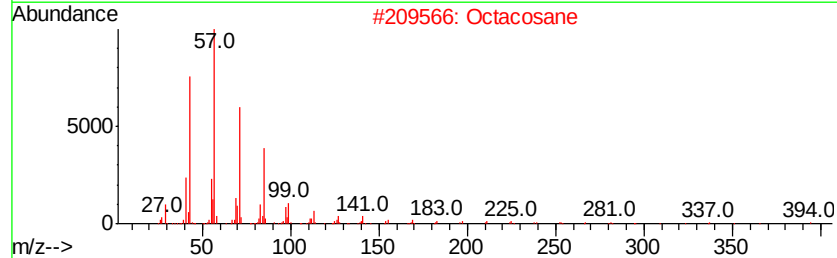
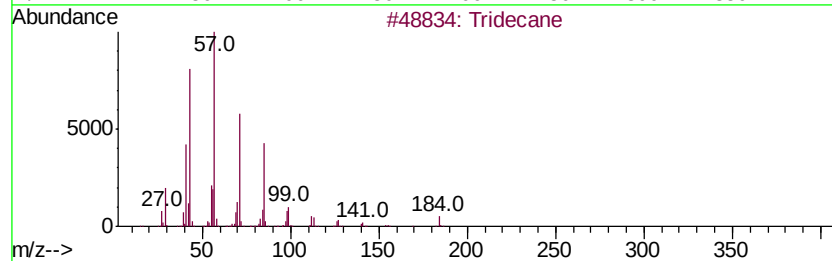
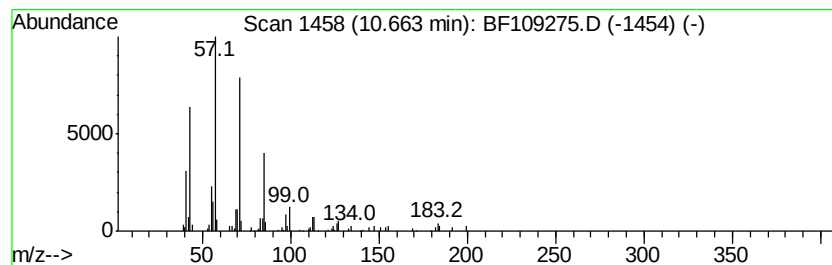
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Tridecane Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.66	2.94 ng	132829	Phenanthrene-d10		11.22
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	94
2	Octacosane	394	C28H58	000630-02-4	90
3	Nonadecane	268	C19H40	000629-92-5	90
4	Hexadecane	226	C16H34	000544-76-3	90
5	Heptacosane	380	C27H56	000593-49-7	86



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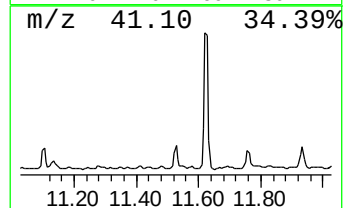
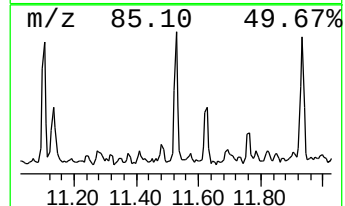
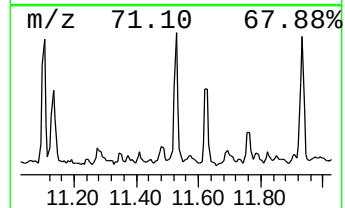
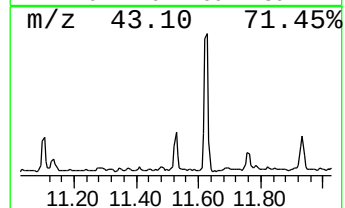
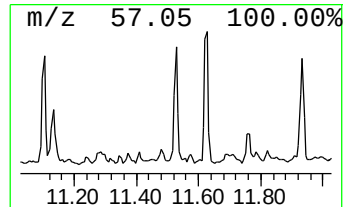
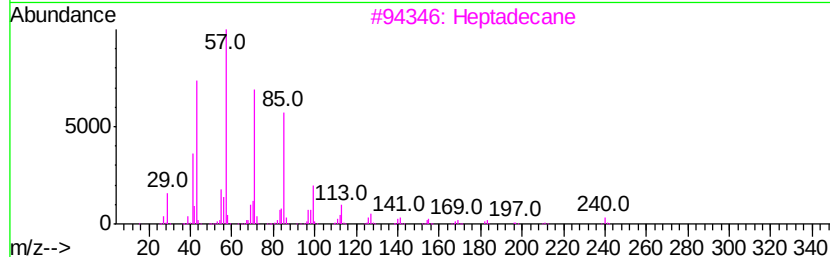
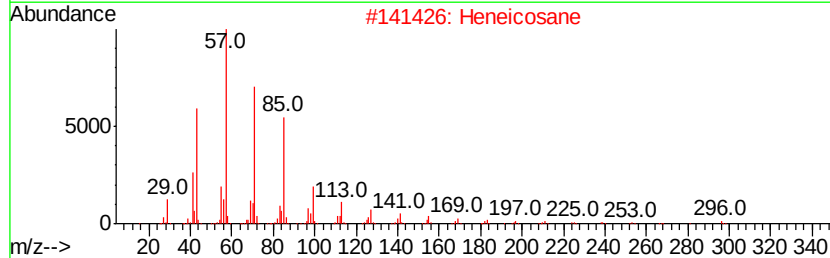
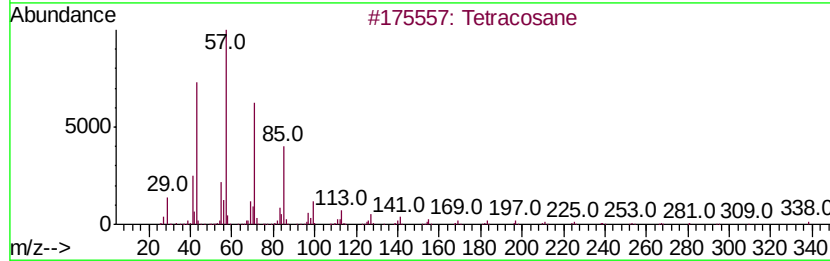
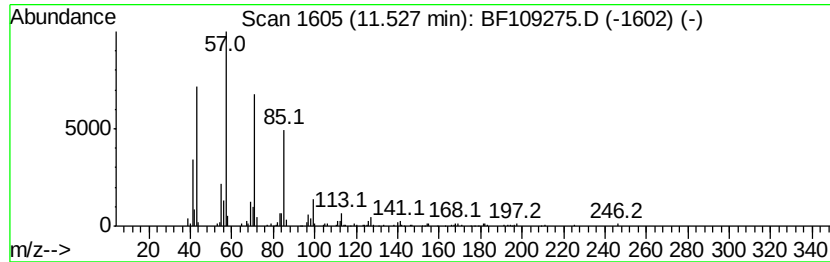
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Tetracosane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.53	2.99 ng	135045	Phenanthrene-d10	11.22	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	91
2	Heneicosane	296	C21H44	000629-94-7	91
3	Heptadecane	240	C17H36	000629-78-7	91
4	Nonadecane	268	C19H40	000629-92-5	90
5	Hexadecane	226	C16H34	000544-76-3	90



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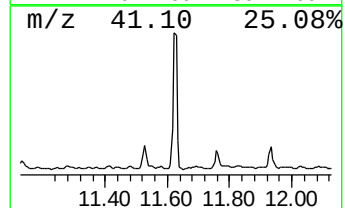
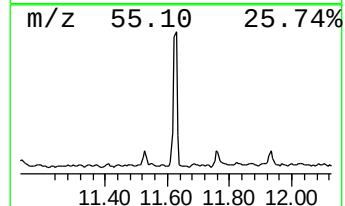
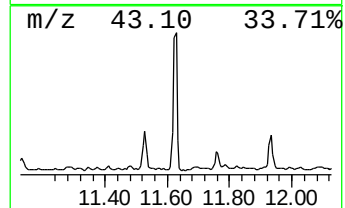
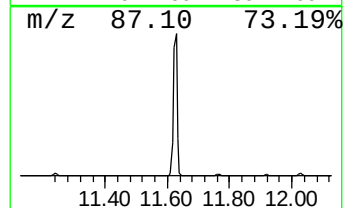
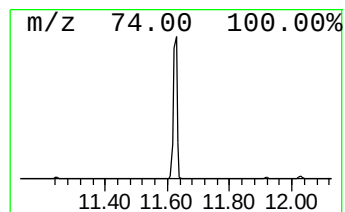
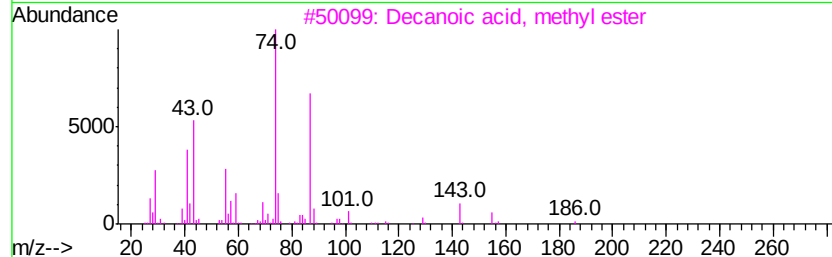
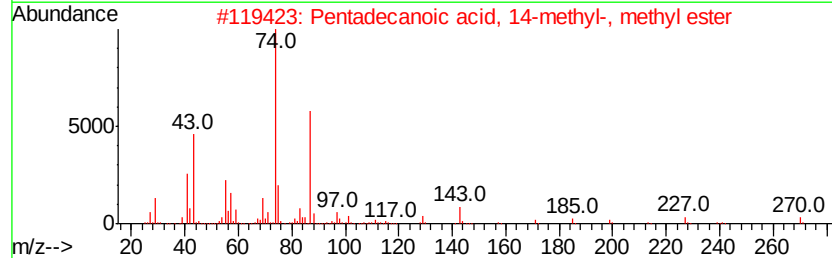
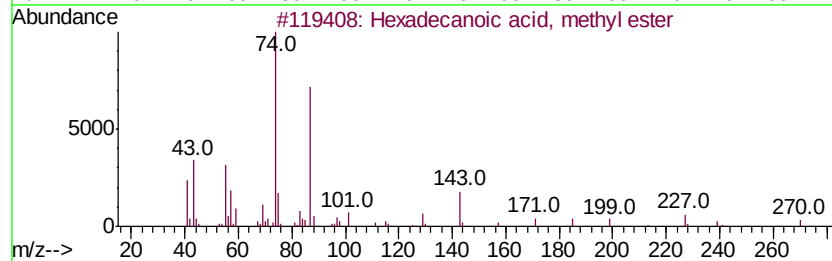
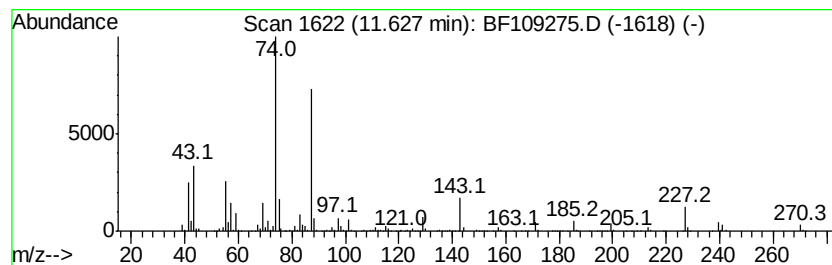
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.63	26.81 ng	1209890	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	92
3		Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	87
4		Dodecanoic acid, 10-methyl-, met...	228	C14H28O2	005129-65-7	86
5		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86



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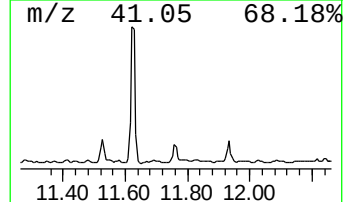
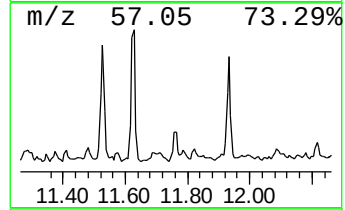
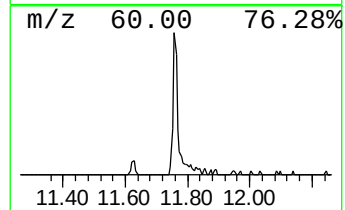
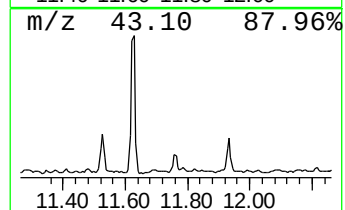
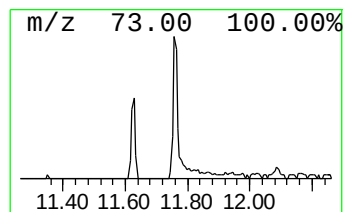
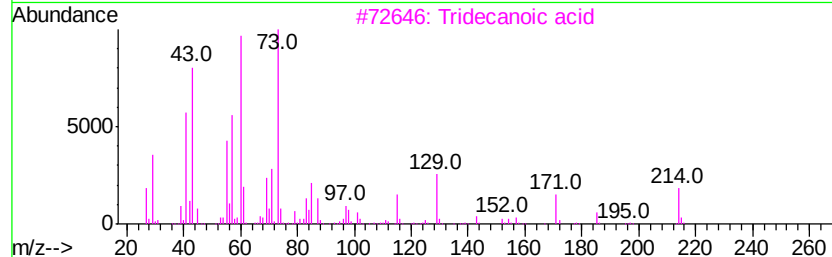
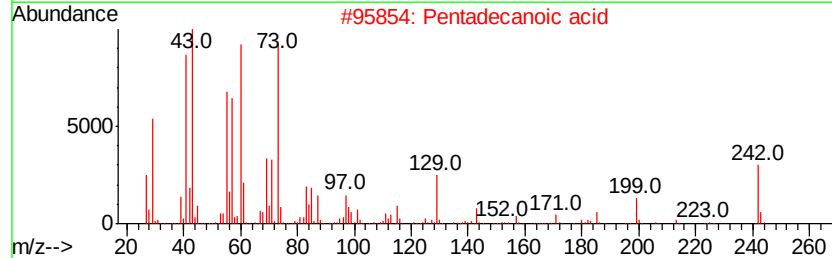
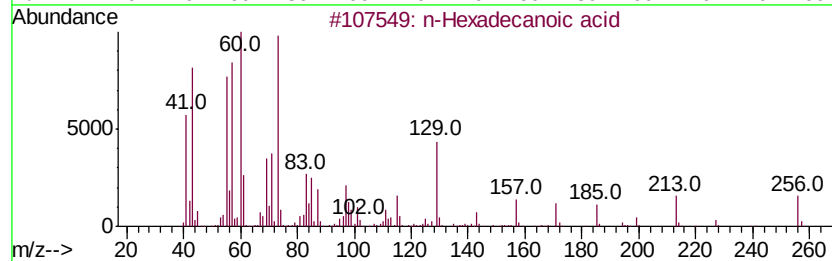
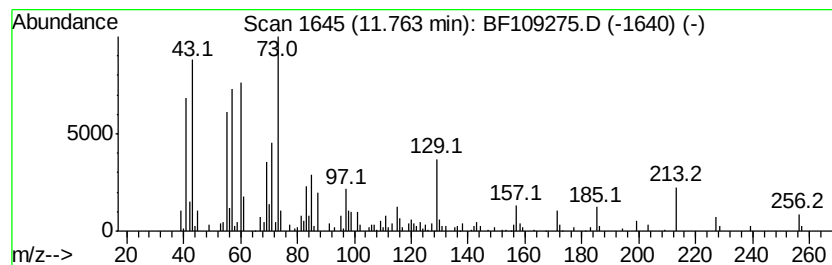
Instrument :
BNA_F
ClientSampleId :
JC-02-092118-G

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 7 n-Hexadecanoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.76	3.32 ng	149771	Phenanthrene-d10	11.22		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3		Tridecanoic acid	214	C13H26O2	000638-53-9	81
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	64
5		n-Decanoic acid	172	C10H20O2	000334-48-5	62



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092418\
Data File : BF109275.D
Acq On : 24 Sep 2018 18:37
Operator : JU/SJ
Sample : J4770-13 2X
Misc :
ALS Vial : 22 Sample Multiplier: 1

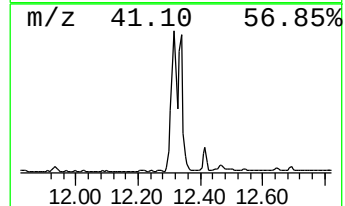
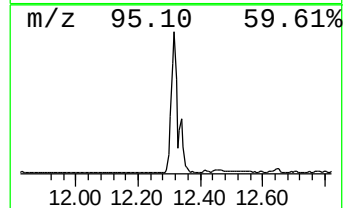
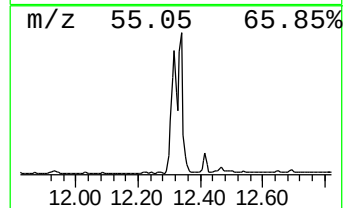
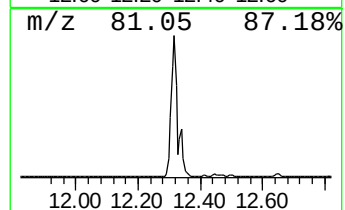
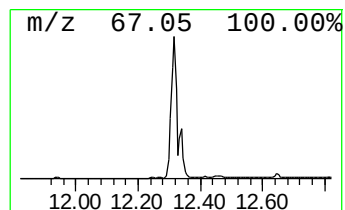
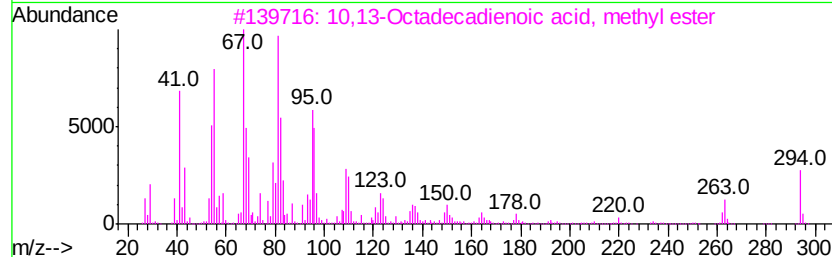
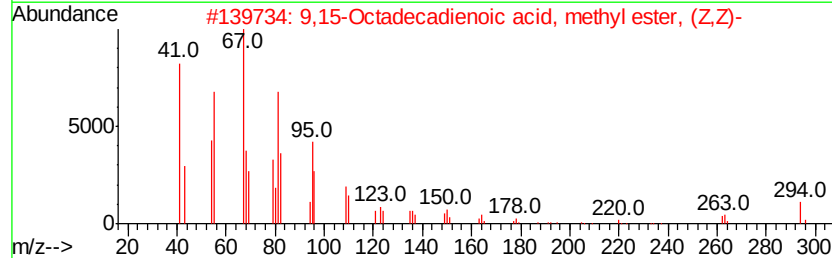
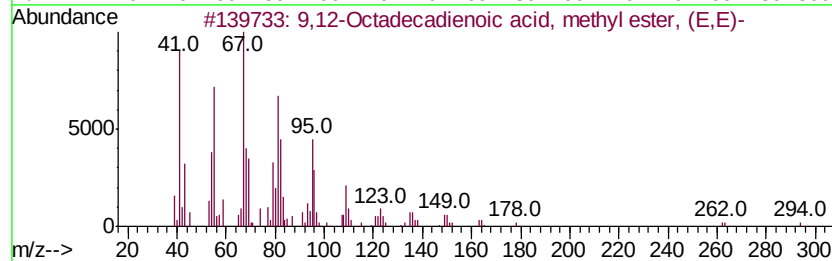
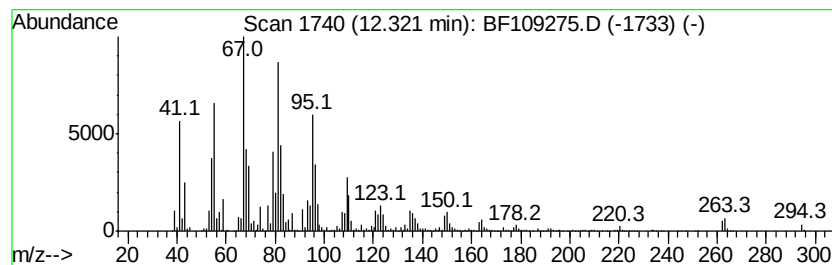
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.32	156.45 ng	7059330	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	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99	
3	10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	98	
4	Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	1000336-44-0	97	
5	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	96	



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092418\
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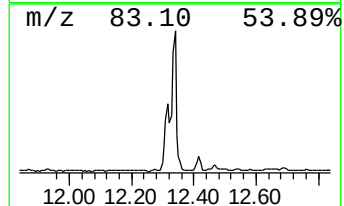
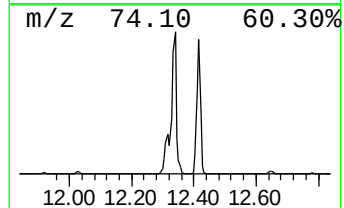
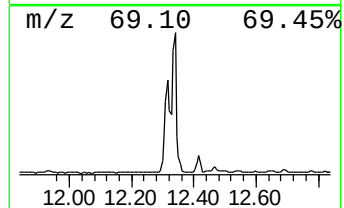
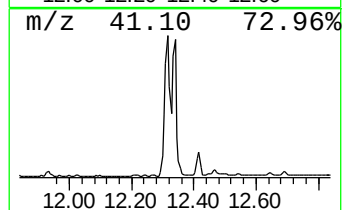
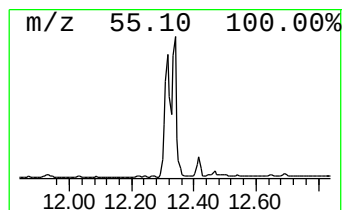
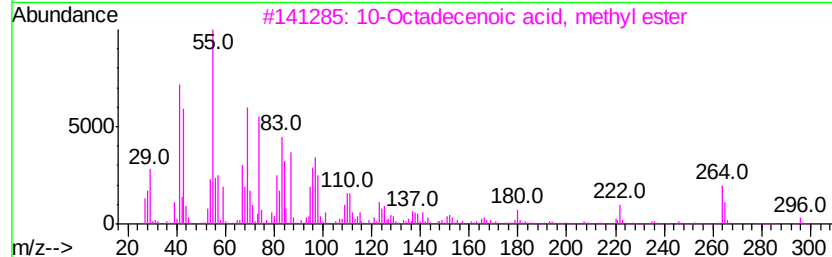
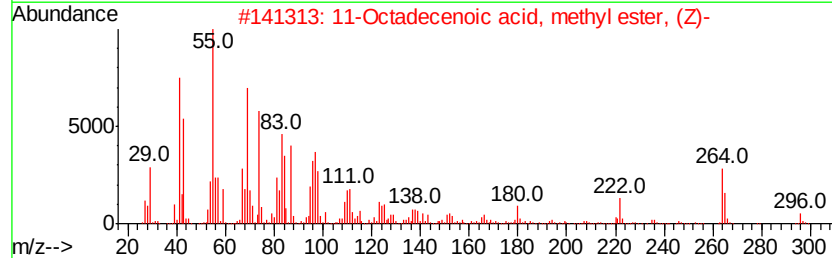
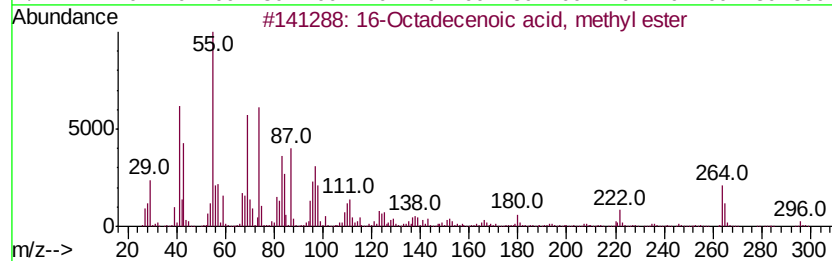
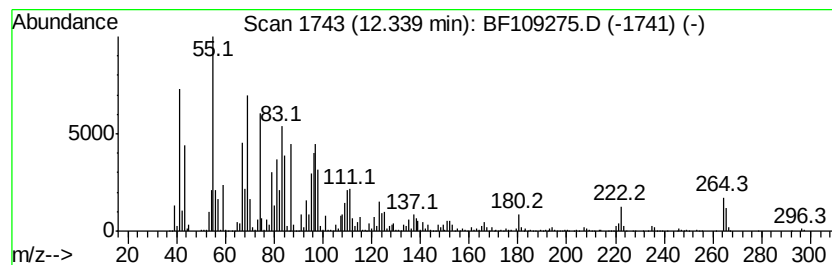
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 16-Octadecenoic acid, methyl ester Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.34	98.02 ng	4422580	Phenanthrene-d10	11.22	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	99
2	11-Octadecenoic acid, methyl ester	296	C19H36O2	001937-63-9	99
3	10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	99
4	12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	99
5	15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	99



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Sample : J4770-13 2X
Misc :
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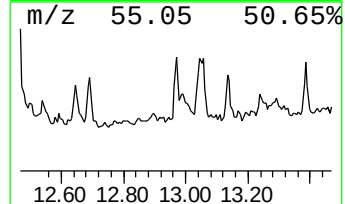
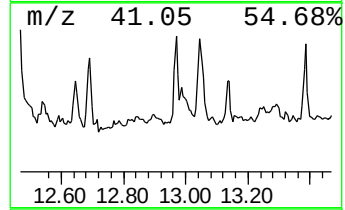
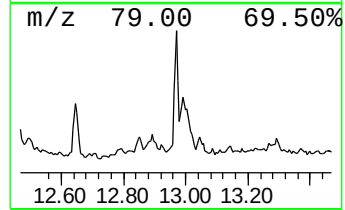
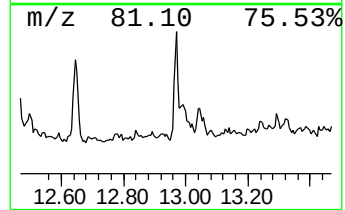
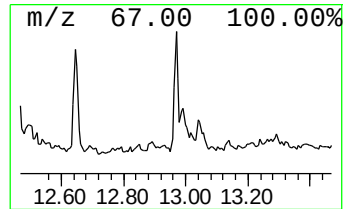
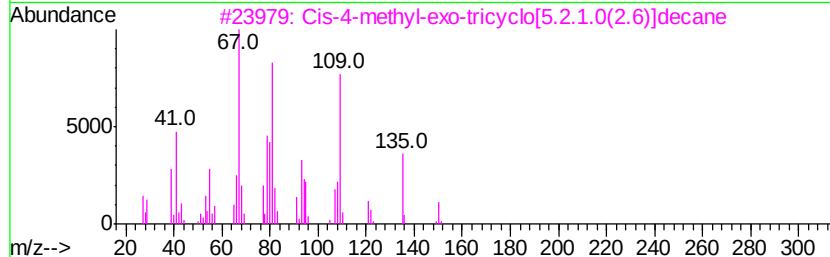
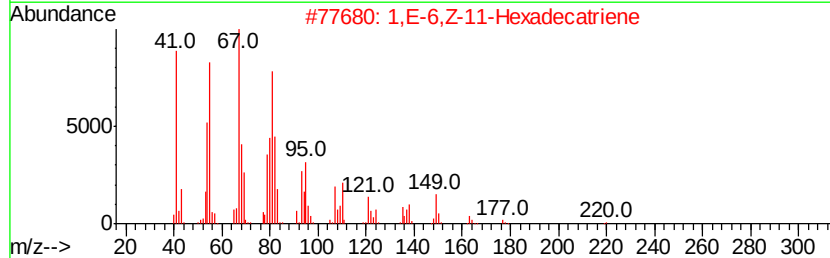
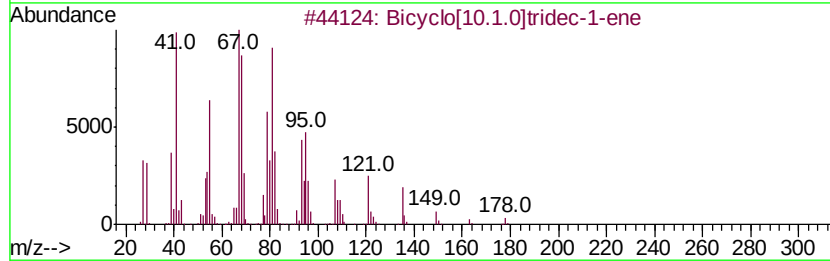
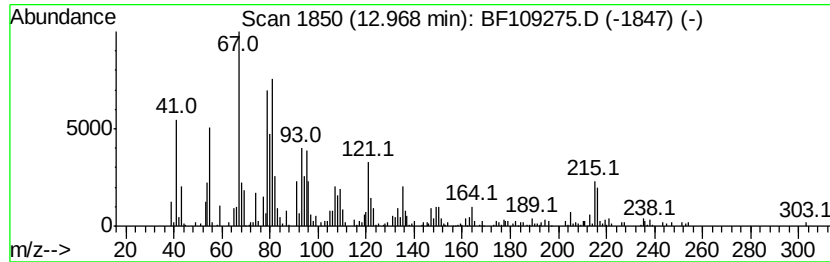
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Bicyclo[10.1.0]tridec-1-ene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.97	7.62 ng	379849	Chrysene-d12	13.84	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[10.1.0]tridec-1-ene	178	C13H22	054766-91-5	76
2	1,E-6,Z-11-Hexadecatriene	220	C16H28	083483-55-0	64
3	Cis-4-methyl-exo-tricyclo[5.2.1....	150	C11H18	1000215-29-2	58
4	Cyclooctene, 4-ethenyl-	136	C10H16	001124-45-4	55
5	Cis-8-methyl-exo-tricyclo[5.2.1....	150	C11H18	1000215-29-3	49



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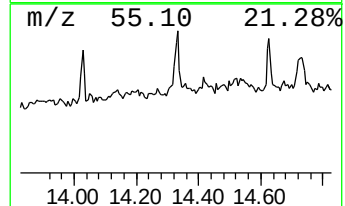
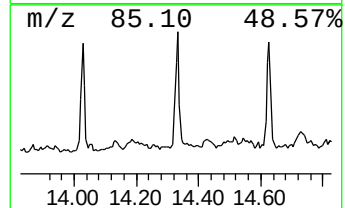
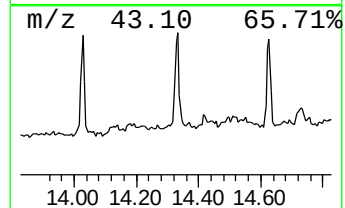
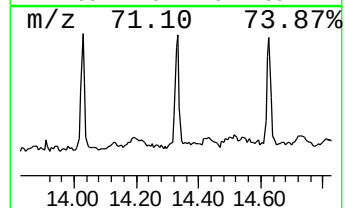
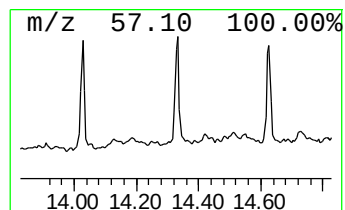
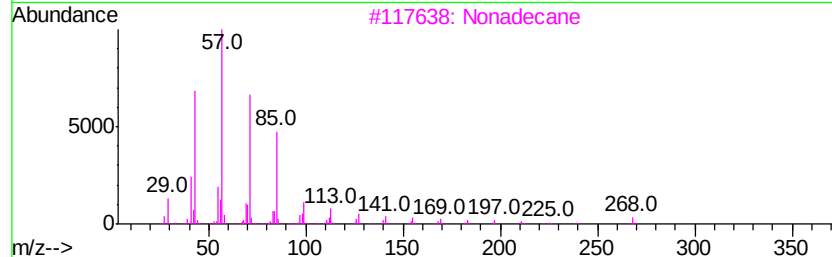
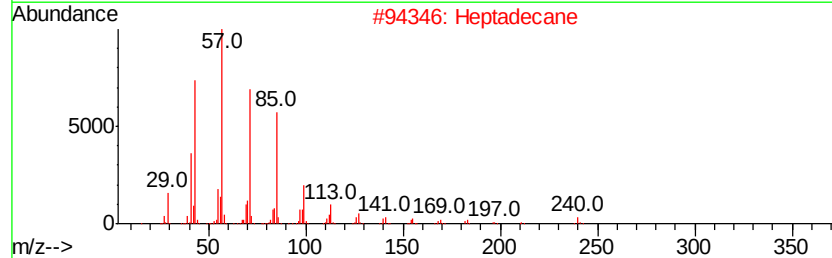
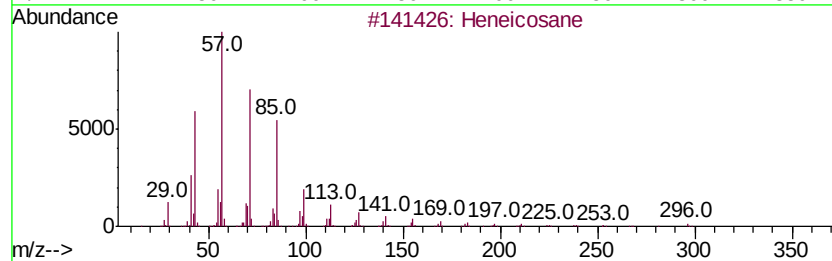
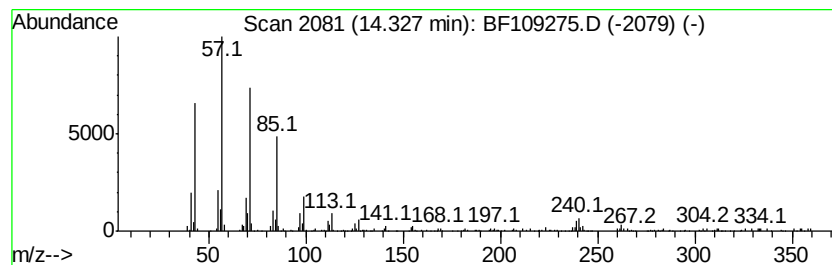
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Heneicosane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.33	3.22 ng	160240	Chrysene-d12	13.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Heptadecane	240	C17H36	000629-78-7	90
3		Nonadecane	268	C19H40	000629-92-5	86
4		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	86
5		Heptacosane	380	C27H56	000593-49-7	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092418\
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Sample : J4770-13 2X
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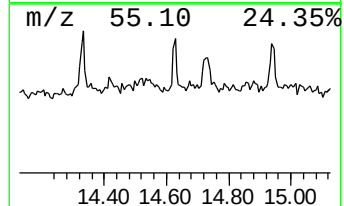
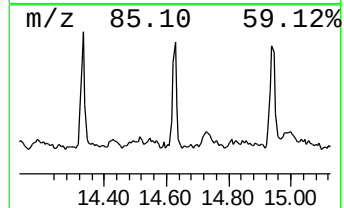
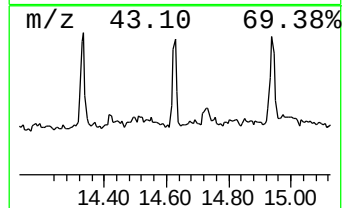
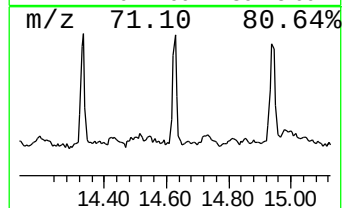
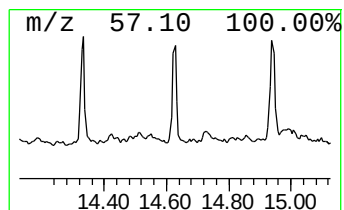
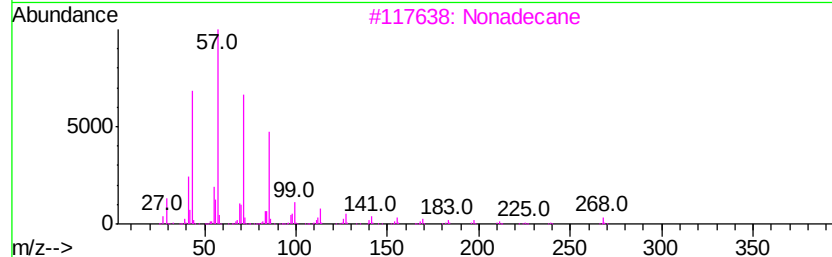
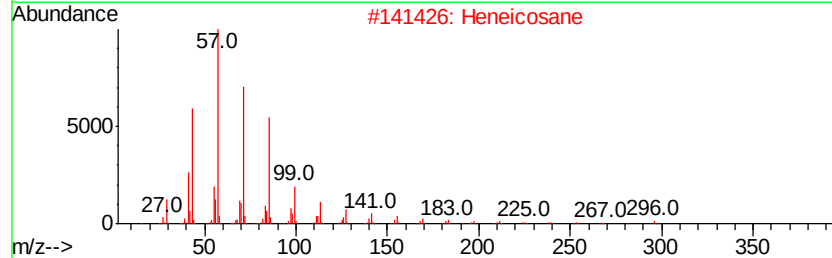
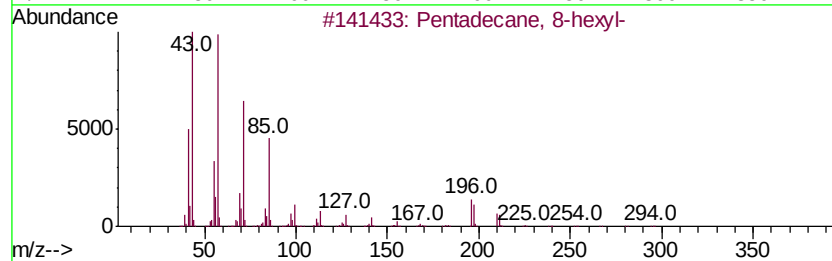
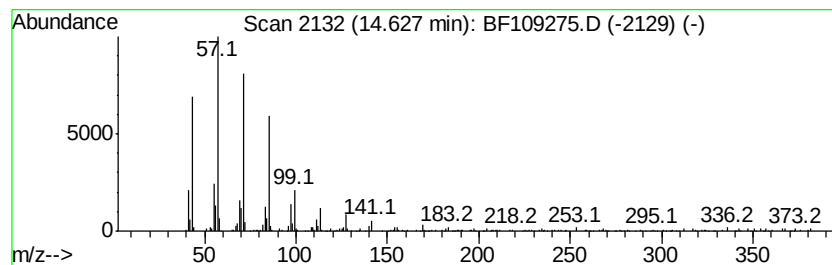
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Pentadecane, 8-hexyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.63	6.61 ng	163656	Perylene-d12	15.25	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93
2	Heneicosane	296	C21H44	000629-94-7	91
3	Nonadecane	268	C19H40	000629-92-5	90
4	2-methyloctacosane	408	C29H60	1000376-72-8	90
5	Triacontane	422	C30H62	000638-68-6	90



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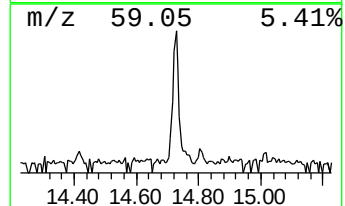
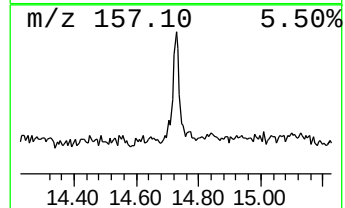
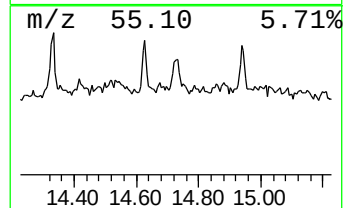
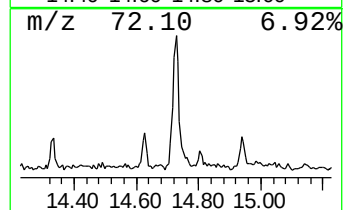
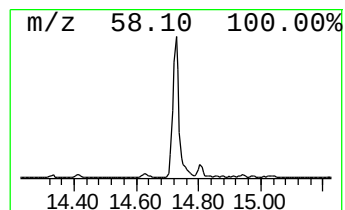
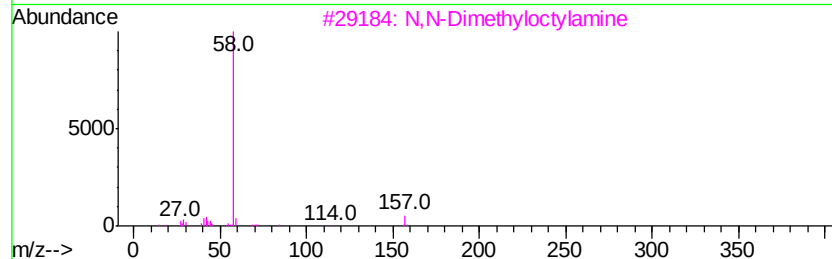
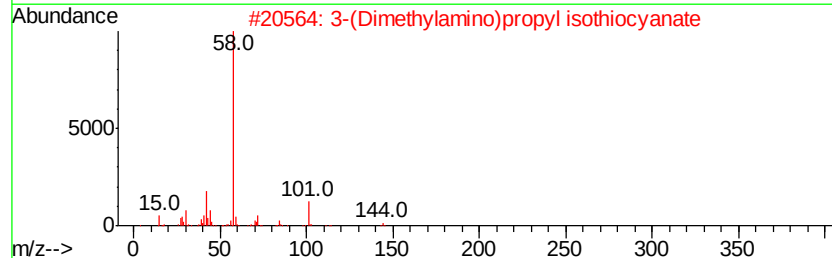
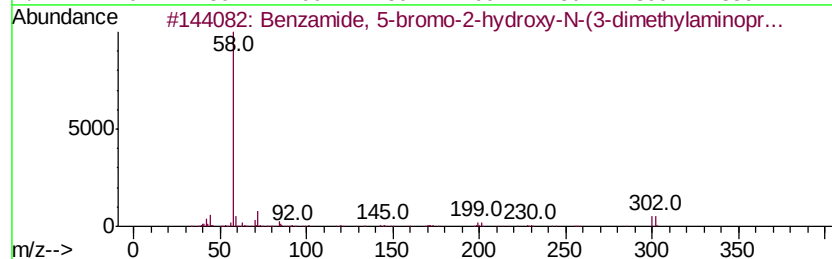
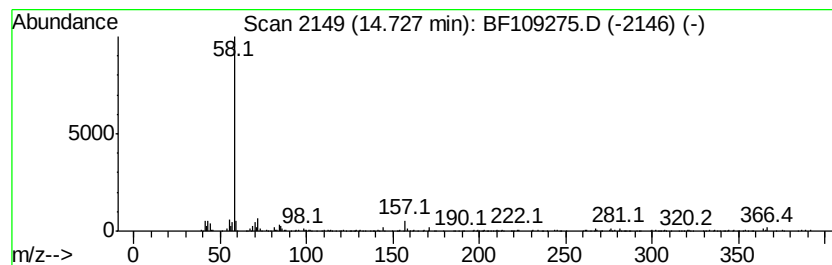
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Benzamide, 5-bromo-2-hydrox... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.73	11.25 ng	278614	Perylene-d12	15.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzamide, 5-bromo-2-hydroxy-N-(...	300	C12H17BrN2O2	289660-53-3	72
2		3-(Dimethylamino)propyl isothiocy...	144	C6H12N2S	027421-70-1	72
3		N,N-Dimethyloctylamine	157	C10H23N	007378-99-6	72
4		9-Azabicyclo[6.1.0]non-4-en-9-am...	138	C8H14N2	066387-78-8	68
5		3(N,N-Dimethylmyristylammonio)pr...	363	C19H41N03S	014933-09-6	64



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ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
JC-02-092118-G

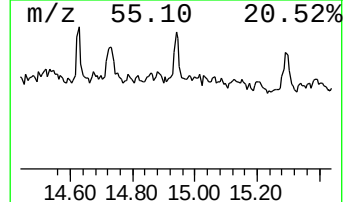
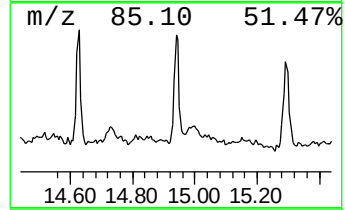
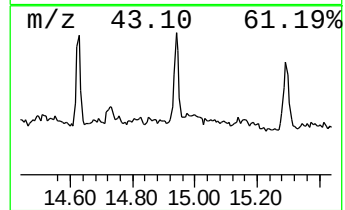
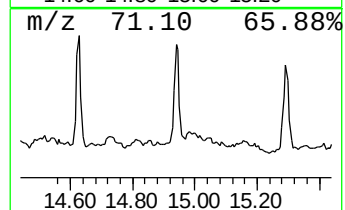
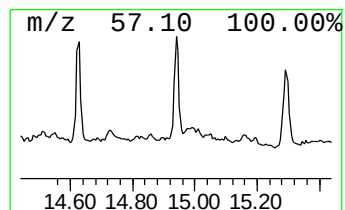
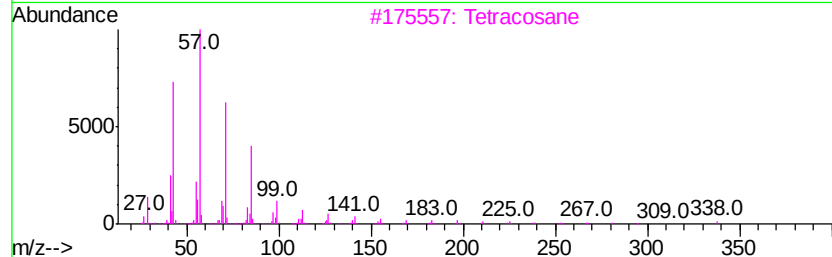
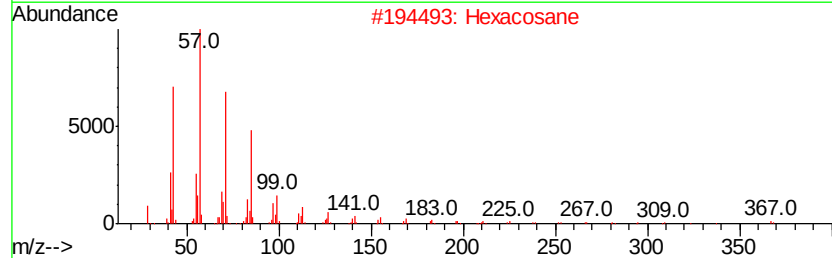
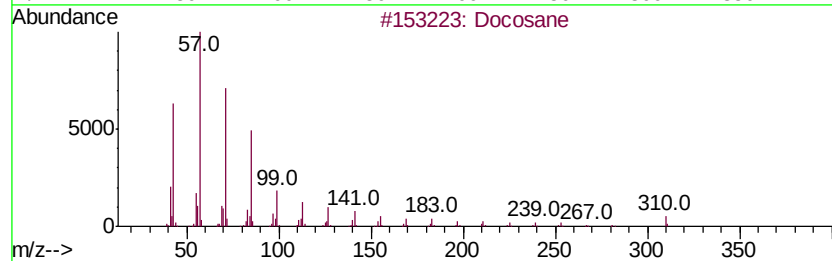
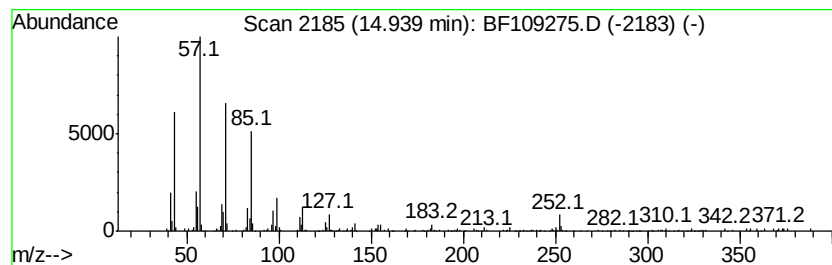
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Docosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.94	7.82 ng	193777	Perylene-d12	15.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosane	310	C22H46	000629-97-0	93
2		Hexacosane	366	C26H54	000630-01-3	90
3		Tetracosane	338	C24H50	000646-31-1	87
4		Heneicosane	296	C21H44	000629-94-7	87
5		Octacosane	394	C28H58	000630-02-4	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092418\
Data File : BF109275.D
Acq On : 24 Sep 2018 18:37
Operator : JU/SJ
Sample : J4770-13 2X
Misc :
ALS Vial : 22 Sample Multiplier: 1

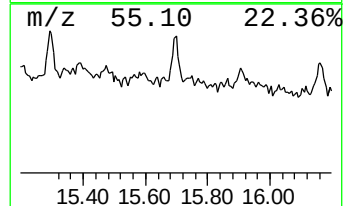
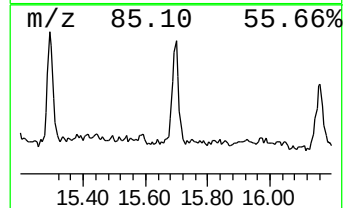
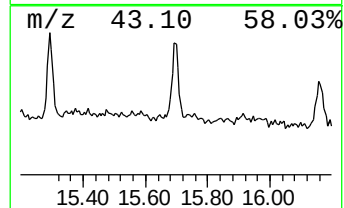
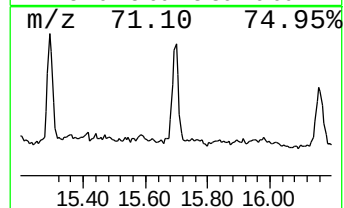
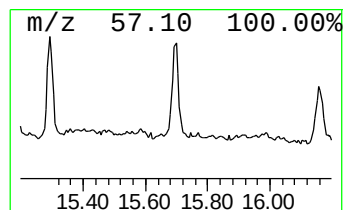
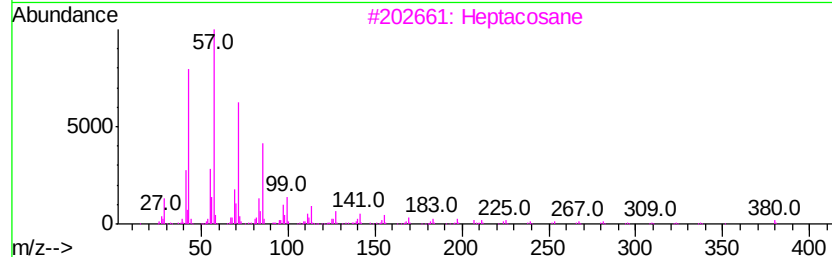
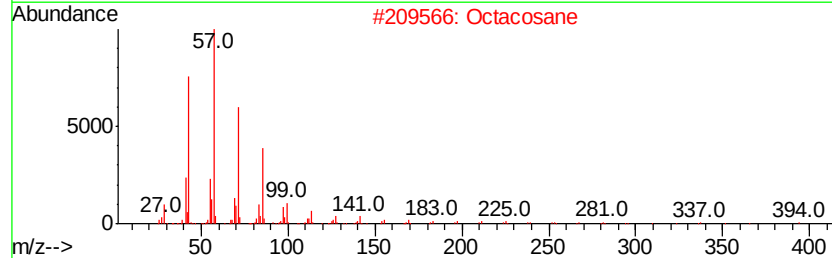
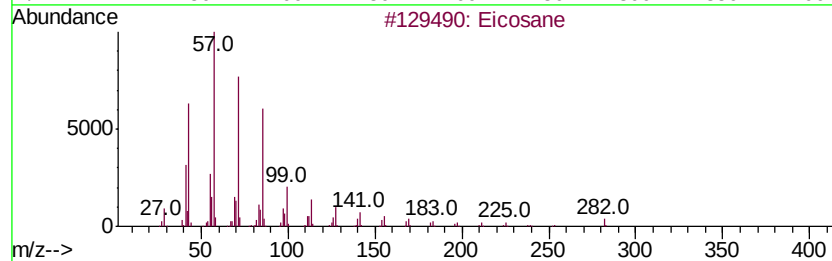
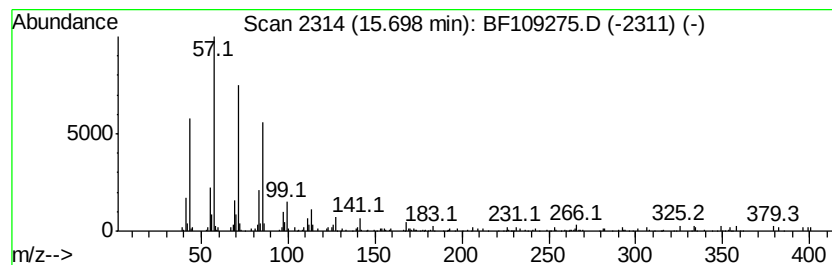
Instrument :
BNA_F
ClientSampleId :
JC-02-092118-G

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 20 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.70	5.98 ng	148185	Perylene-d12	15.25
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane	282 C20H42	000112-95-8	91
2	Octacosane	394 C28H58	000630-02-4	86
3	Heptacosane	380 C27H56	000593-49-7	86
4	Decane, 3,8-dimethyl-	170 C12H26	017312-55-9	81
5	2-Bromotetradecane	276 C14H29Br	074036-95-6	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092418\
 Data File : BF109275.D
 Acq On : 24 Sep 2018 18:37
 Operator : JU/SJ
 Sample : J4770-13 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

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

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	26.0	ng	906515	1	6.71	696936	20.0
Tridecane	10.66	2.9	ng	132829	4	11.22	902421	20.0
Tetracosane	11.53	3.0	ng	135045	4	11.22	902421	20.0
Hexadecanoic acid...	11.63	26.8	ng	1209890	4	11.22	902421	20.0
n-Hexadecanoic acid	11.76	3.3	ng	149771	4	11.22	902421	20.0
9,12-Octadecadien...	12.32	156.4	ng	7059330	4	11.22	902421	20.0
16-Octadecenoic a...	12.34	98.0	ng	4422580	4	11.22	902421	20.0
Bicyclo[10.1.0]tr...	12.97	7.6	ng	379849	5	13.84	996639	20.0
Heneicosane	14.33	3.2	ng	160240	5	13.84	996639	20.0
Pentadecane, 8-he...	14.63	6.6	ng	163656	6	15.25	495343	20.0
Benzamide, 5-brom...	14.73	11.3	ng	278614	6	15.25	495343	20.0
Docosane	14.94	7.8	ng	193777	6	15.25	495343	20.0
Eicosane	15.70	6.0	ng	148185	6	15.25	495343	20.0