

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042319\
 Data File : BF113930.D
 Acq On : 23 Apr 2019 19:38
 Operator : JU/SJ
 Sample : K2529-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P3P4042219

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-BF042219.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.216	17	22	30	rVB	47280	69779	2.09%	0.238%
2	5.528	580	585	588	rBV	3180535	3059281	91.76%	10.441%
3	6.528	749	755	758	rBV	2938865	3074113	92.20%	10.491%
4	6.669	774	779	782	rBV	3310065	3226163	96.76%	11.010%
5	6.898	814	818	822	rBV	948259	744628	22.33%	2.541%
6	7.057	841	845	848	rBV	3066224	2582443	77.46%	8.813%
7	7.457	907	913	916	rBV	2025105	1889110	56.66%	6.447%
8	8.180	1032	1036	1039	rBV	938956	830065	24.90%	2.833%
9	9.257	1214	1219	1222	rBV	3503754	3334043	100.00%	11.378%
10	9.639	1281	1284	1289	rBV	315631	281334	8.44%	0.960%
11	9.933	1330	1334	1345	rVB	1012953	850906	25.52%	2.904%
12	10.721	1463	1468	1471	rBV	1983748	1869318	56.07%	6.380%
13	11.416	1582	1586	1590	rBV	936410	824510	24.73%	2.814%
14	11.939	1672	1675	1680	rBV	151389	129350	3.88%	0.441%
15	12.721	1805	1808	1813	rBV	34872	37922	1.14%	0.129%
16	13.004	1851	1856	1859	rBV	3527359	3315015	99.43%	11.313%
17	13.233	1892	1895	1898	rBV	68926	58486	1.75%	0.200%
18	13.574	1950	1953	1956	rBV	136347	119728	3.59%	0.409%
19	13.909	2006	2010	2014	rVB	199726	172030	5.16%	0.587%
20	14.021	2021	2029	2032	rBV5	45595	120400	3.61%	0.411%
21	14.062	2032	2036	2040	rVB	1166864	1054606	31.63%	3.599%
22	14.233	2062	2065	2068	rBV	185334	159105	4.77%	0.543%
23	14.551	2116	2119	2123	rVB	165226	148360	4.45%	0.506%
24	14.868	2170	2173	2177	rBV	146034	139852	4.19%	0.477%
25	15.215	2228	2232	2236	rBV	123427	149846	4.49%	0.511%
26	15.568	2287	2292	2296	rBV	683830	852582	25.57%	2.910%
27	15.604	2296	2298	2303	rVB	89768	114774	3.44%	0.392%
28	16.051	2371	2374	2381	rBV	59410	93993	2.82%	0.321%

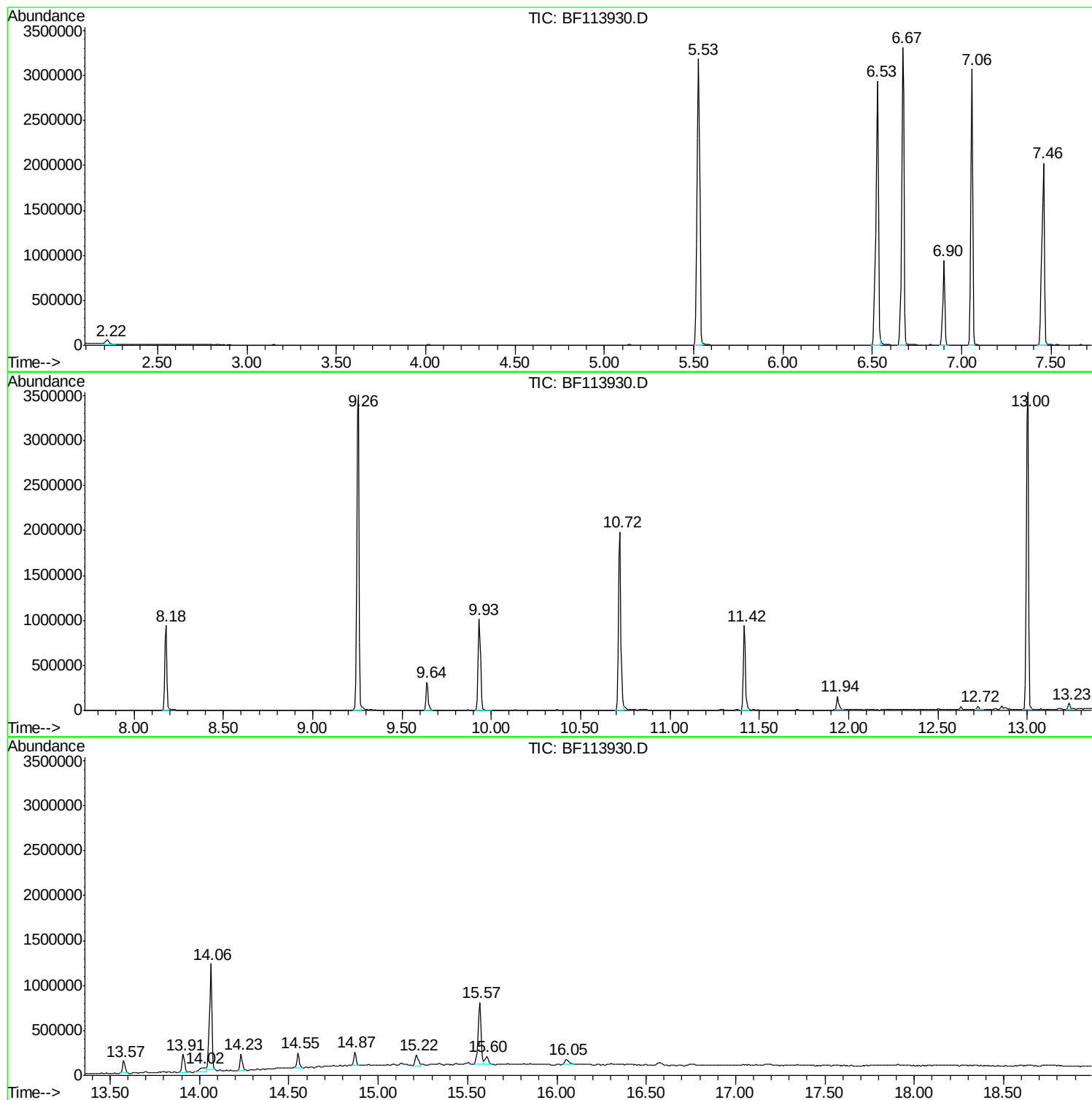
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF042219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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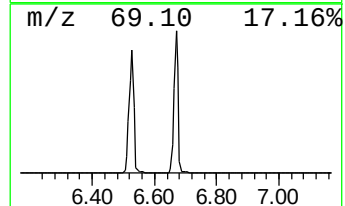
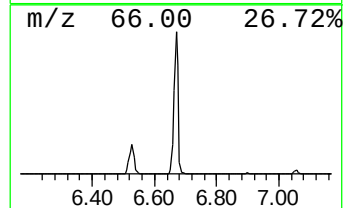
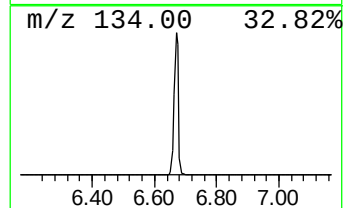
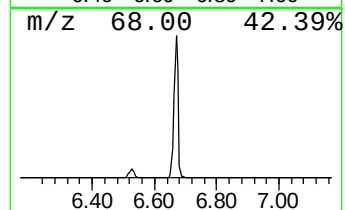
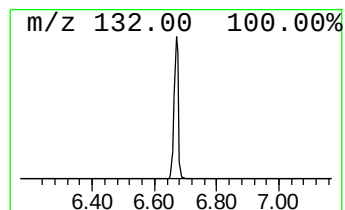
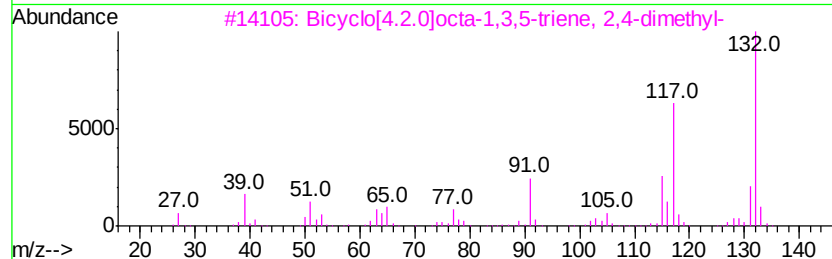
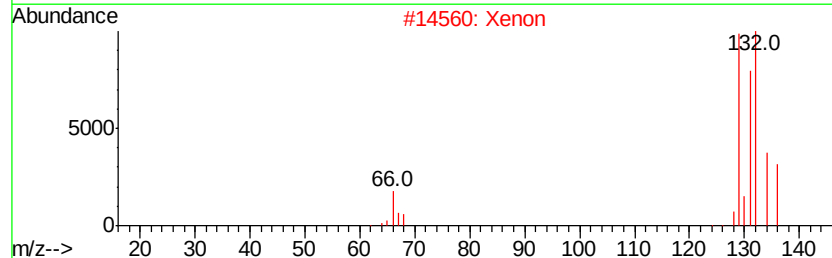
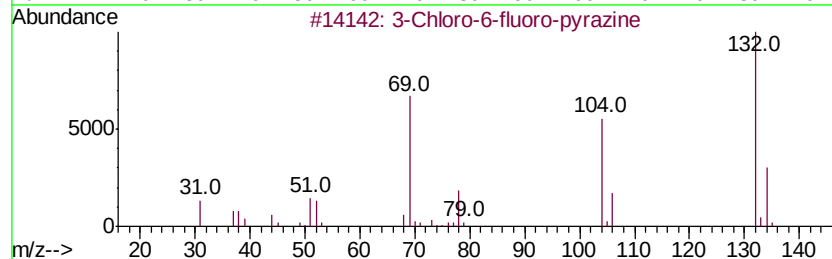
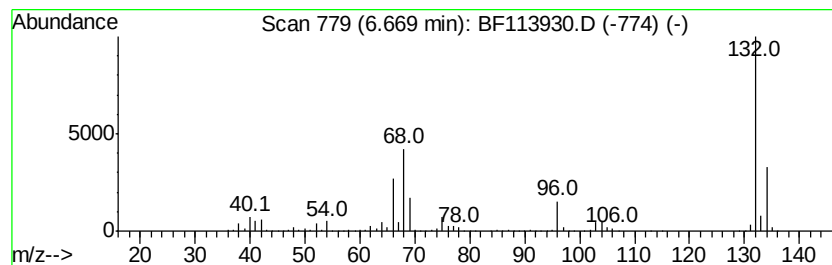
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF042219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.67 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.67	86.65 ng	3226160	1,4-Dichlorobenzene-d4	6.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	27
2	Xenon			132	Xe	007440-63-3	9
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	1-Methylpyrrolo[1,2-a]pyrazine			132	C8H8N2	064608-59-9	9



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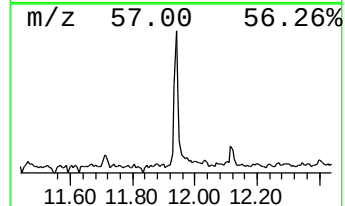
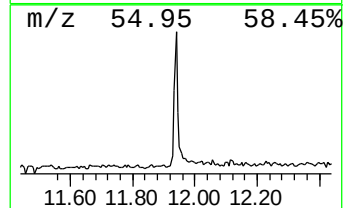
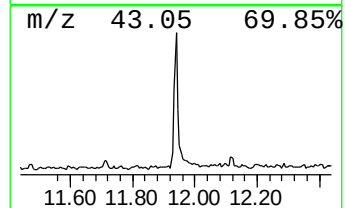
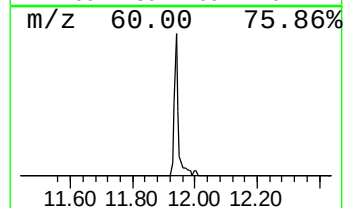
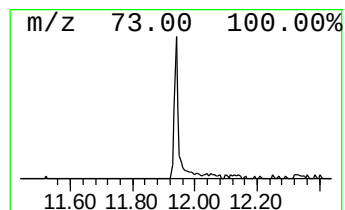
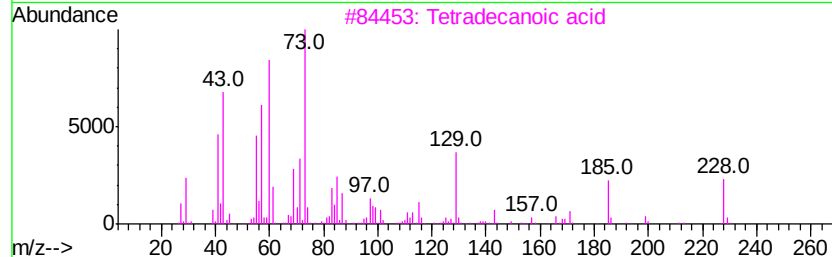
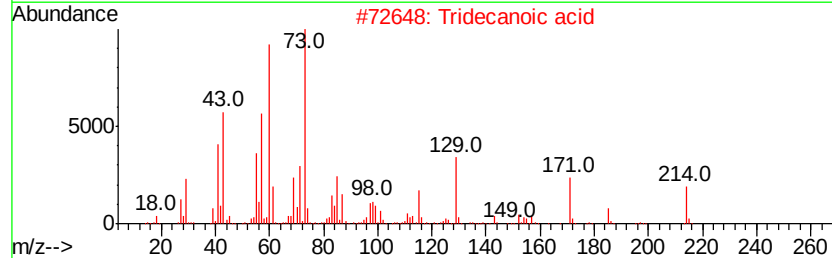
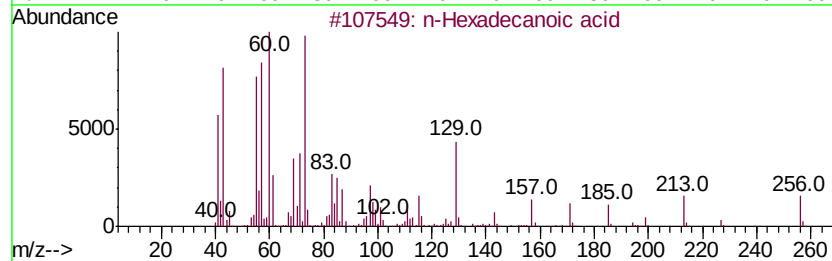
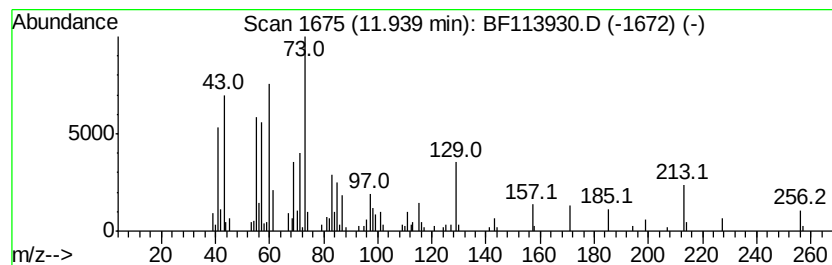
Instrument :
BNA_F
ClientSampleId :
P3P4042219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF042219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.94	3.14 ng	129350	Phenanthrene-d10	11.42	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tridecanoic acid	214	C13H26O2	000638-53-9	83
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	74
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	68
5	Sulfurous acid, pentadecyl 2-pro...	334	C18H38O3S	1000309-12-6	18



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Instrument :
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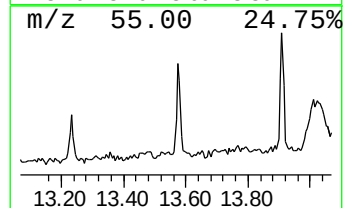
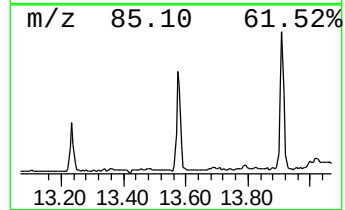
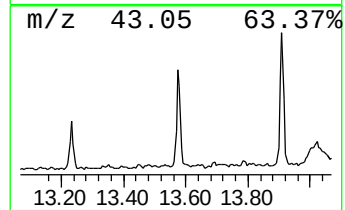
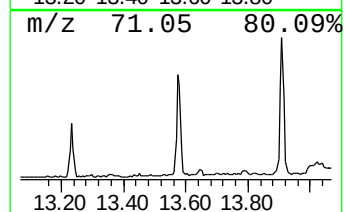
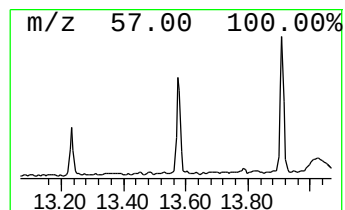
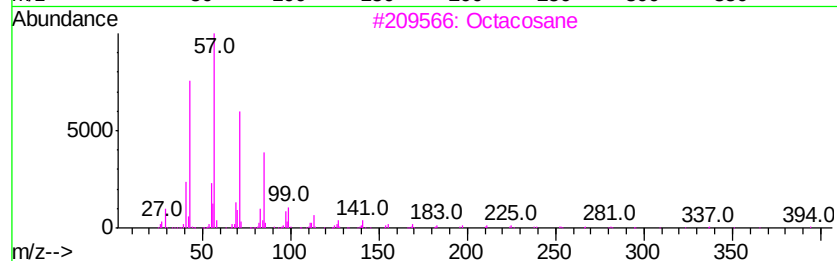
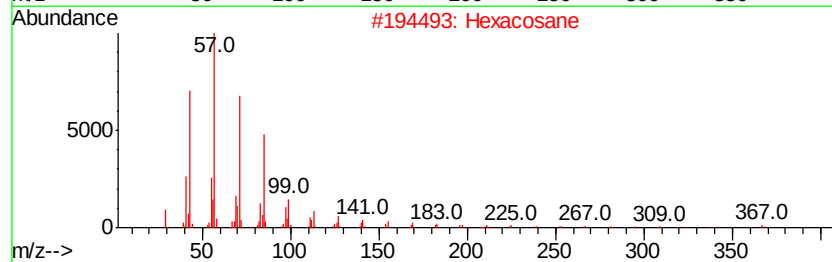
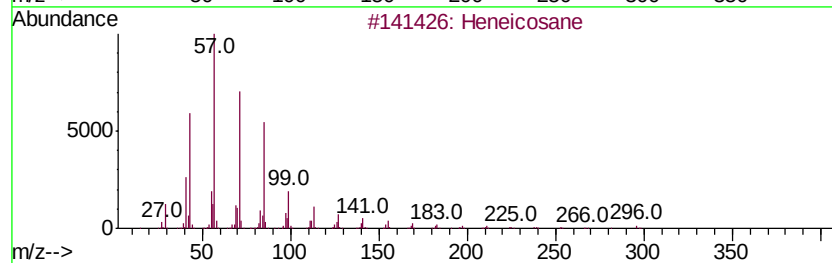
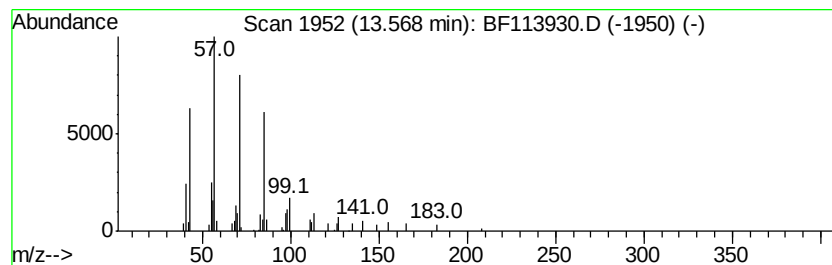
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Heneicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.57	2.27 ng	119728	Chrysene-d12	14.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Hexacosane		366	C26H54	000630-01-3	91
3	Octacosane		394	C28H58	000630-02-4	91
4	Octadecane		254	C18H38	000593-45-3	90
5	Docosane		310	C22H46	000629-97-0	90



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ALS Vial : 16 Sample Multiplier: 1

Instrument :
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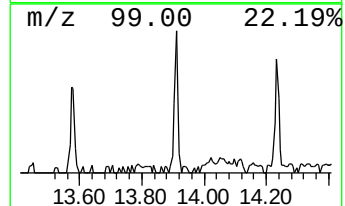
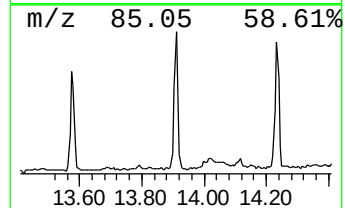
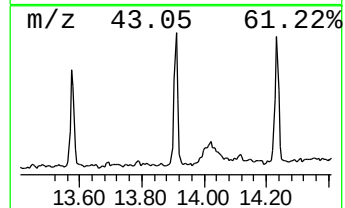
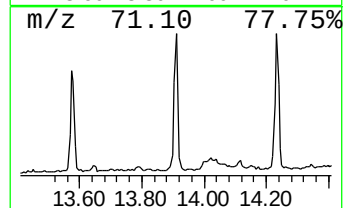
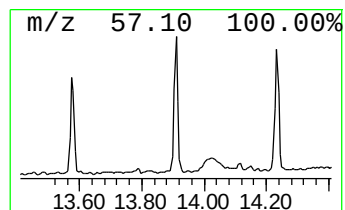
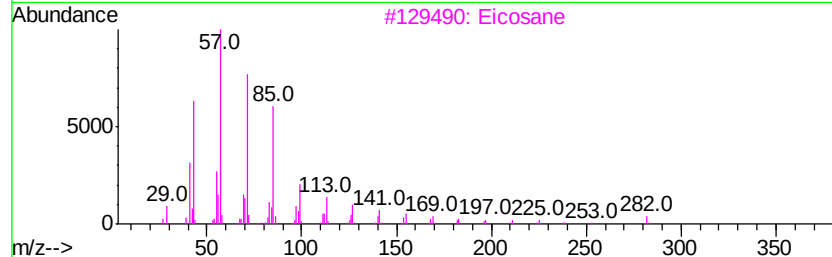
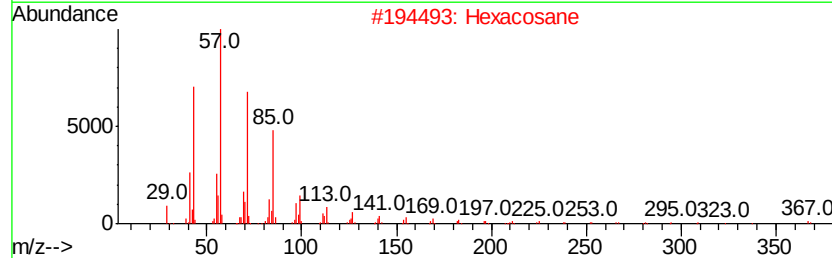
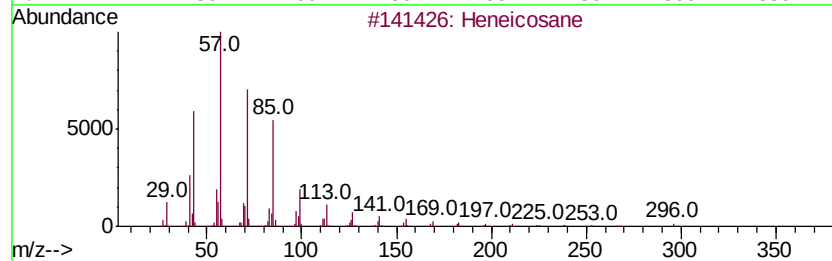
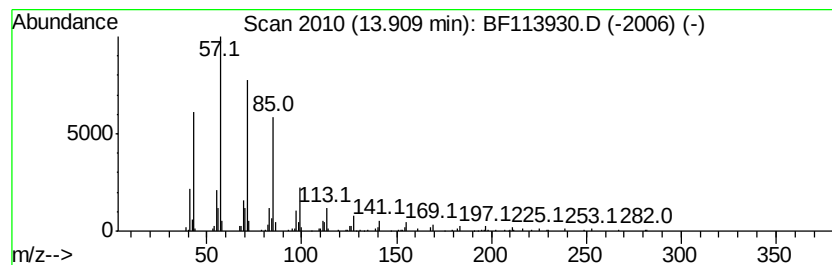
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Hexacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.91	3.26 ng	172030	Chrysene-d12	14.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Hexacosane	366	C26H54	000630-01-3	90
3		Eicosane	282	C20H42	000112-95-8	90
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
5		Tetracosane	338	C24H50	000646-31-1	87



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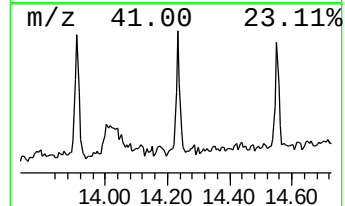
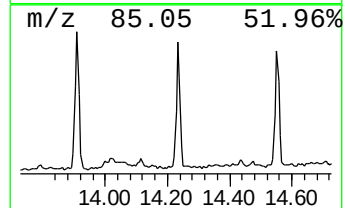
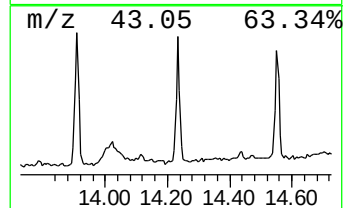
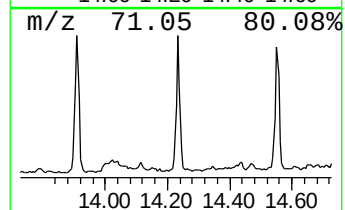
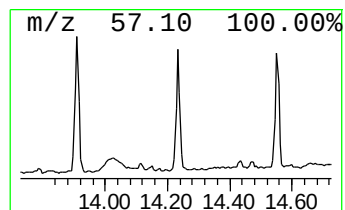
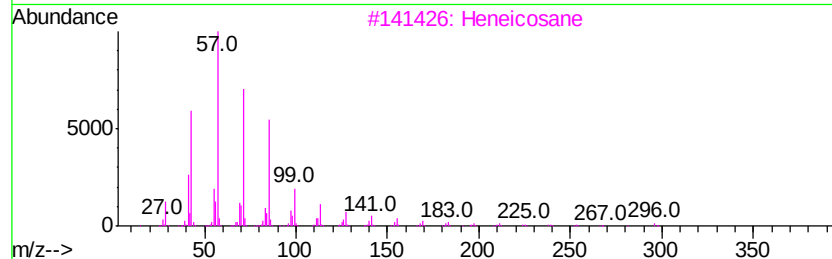
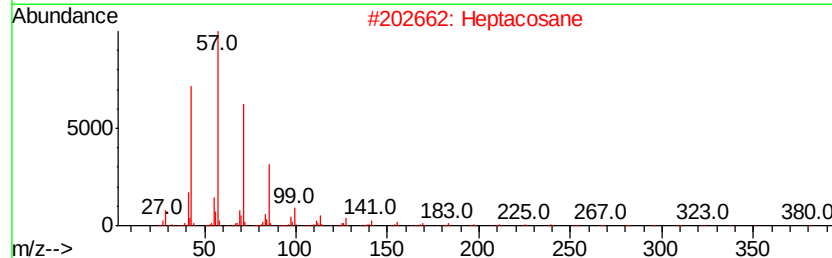
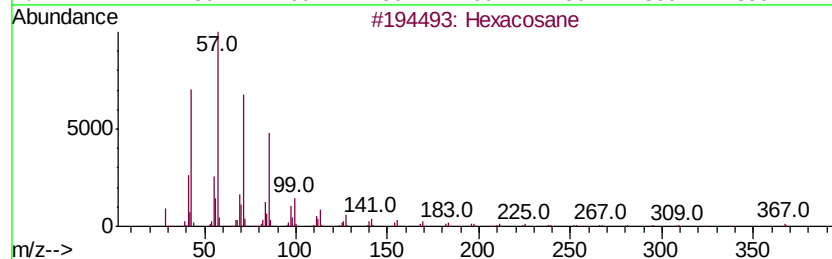
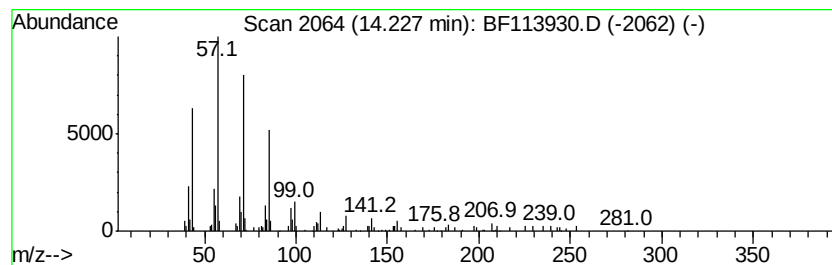
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Heptacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.23	3.02 ng	159105	Chrysene-d12	14.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Heptacosane	380	C27H56	000593-49-7	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Triacontane	422	C30H62	000638-68-6	91



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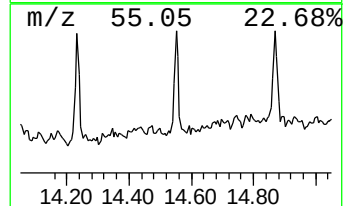
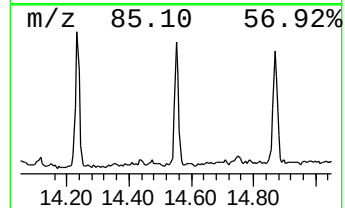
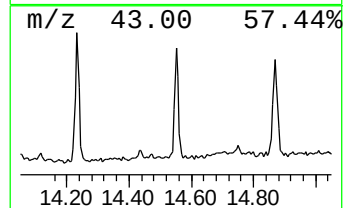
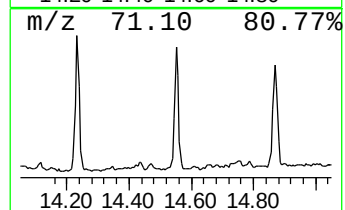
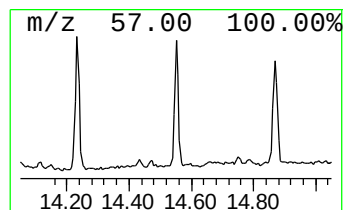
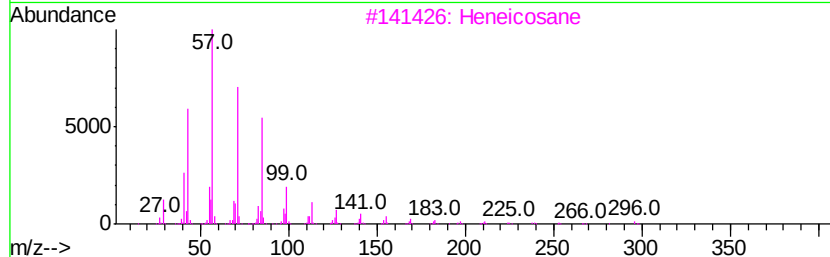
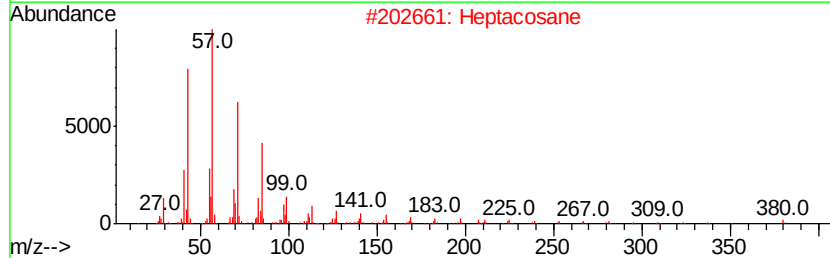
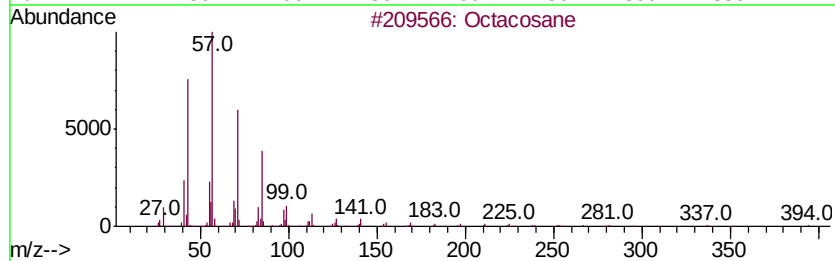
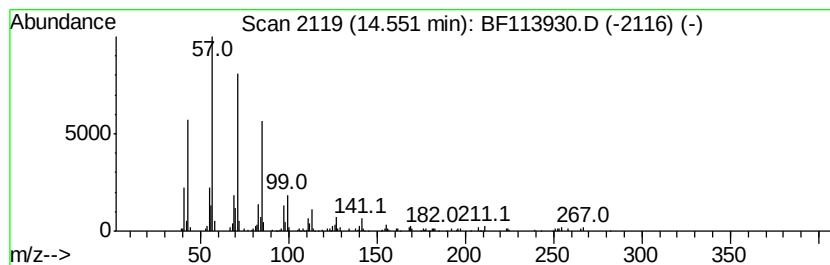
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.55	2.81 ng	148360	Chrysene-d12	14.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octacosane	394 C28H58	000630-02-4	91
2	Heptacosane	380 C27H56	000593-49-7	91
3	Heneicosane	296 C21H44	000629-94-7	91
4	Hexacosane	366 C26H54	000630-01-3	91
5	Hentriacontane	437 C31H64	000630-04-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042319\
Data File : BF113930.D
Acq On : 23 Apr 2019 19:38
Operator : JU/SJ
Sample : K2529-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

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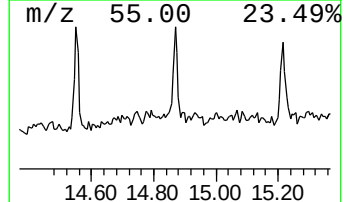
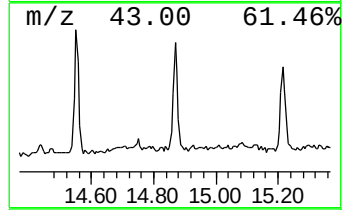
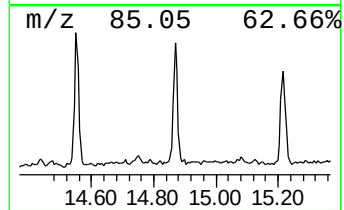
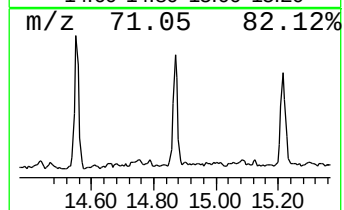
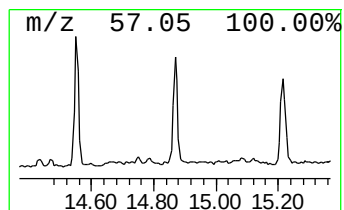
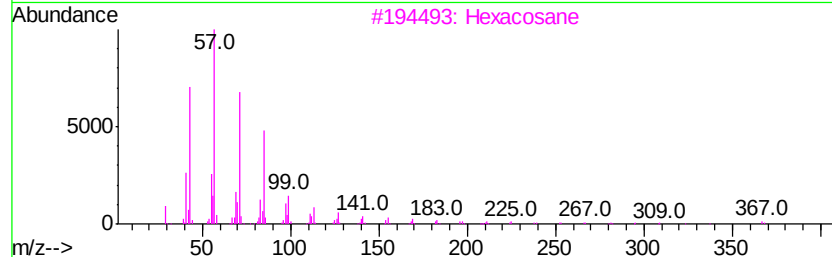
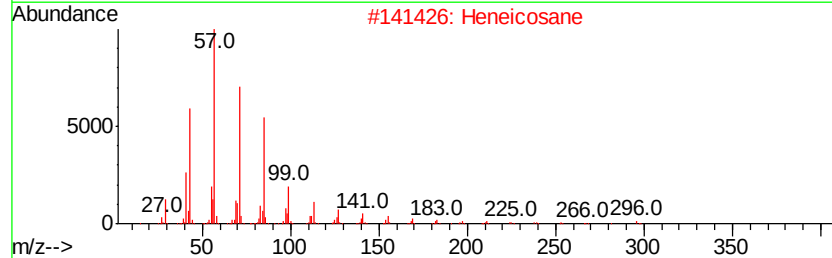
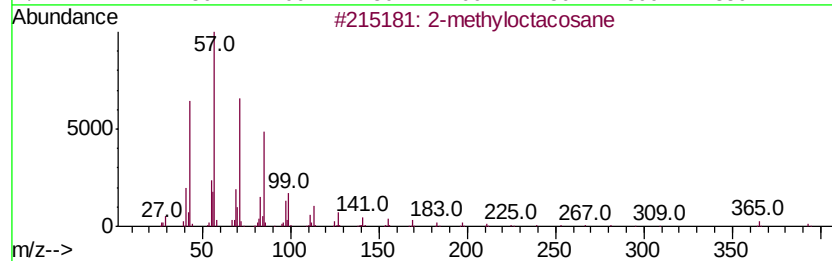
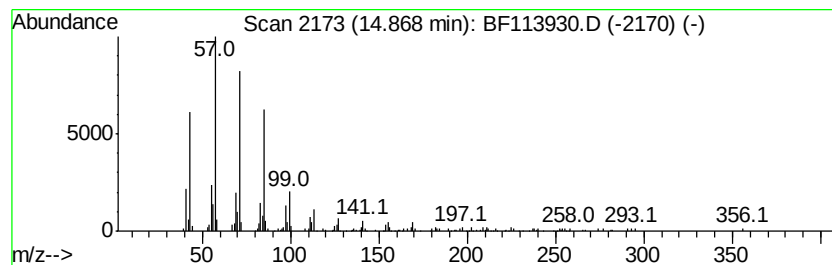
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 2-methyloctacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.87	3.28 ng	139852	Perylene-d12	15.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		Heptadecane	240	C17H36	000629-78-7	91
5		Docosane, 11-butyl-	366	C26H54	013475-76-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042319\
Data File : BF113930.D
Acq On : 23 Apr 2019 19:38
Operator : JU/SJ
Sample : K2529-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

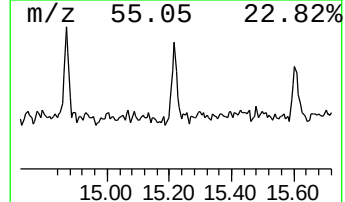
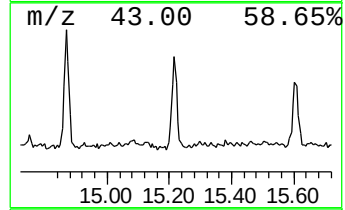
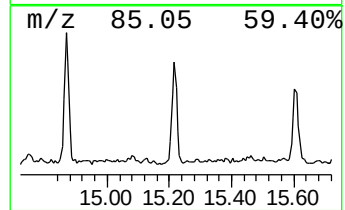
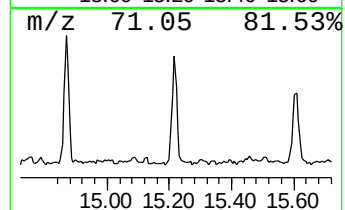
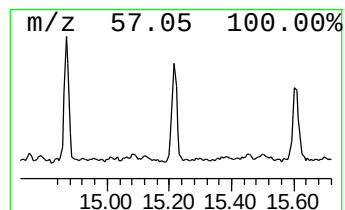
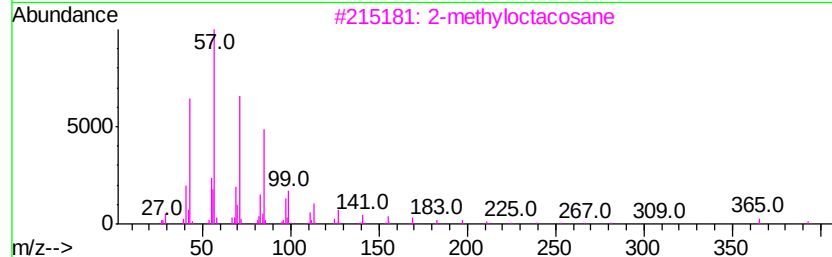
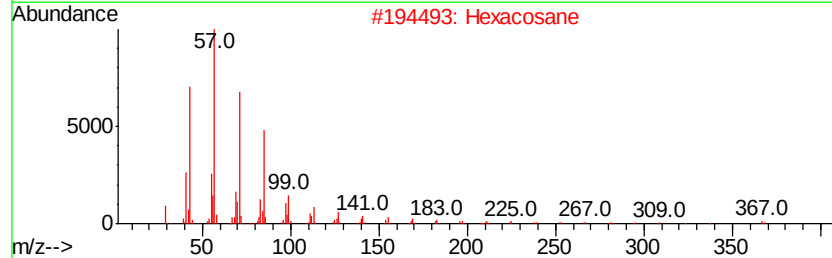
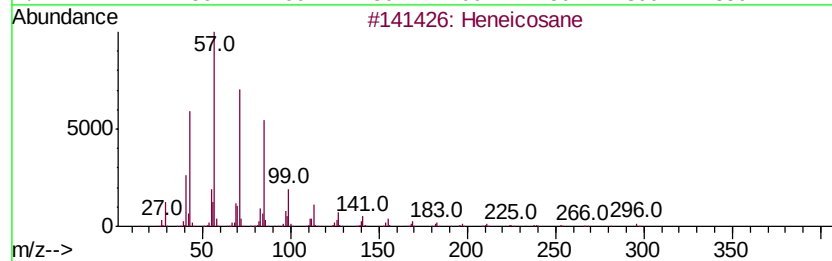
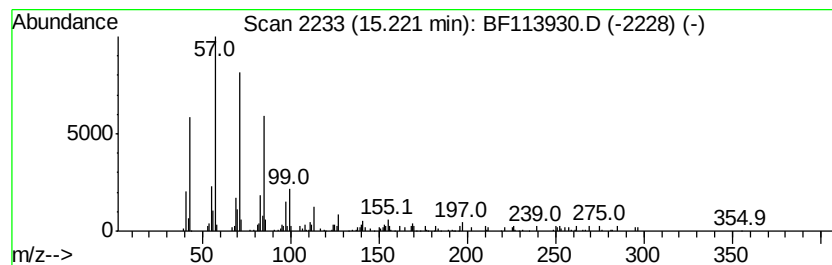
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF042219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecane, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.22	3.52 ng	149846	Perylene-d12	15.57
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heneicosane	296 C21H44	000629-94-7	91
2	Hexacosane	366 C26H54	000630-01-3	91
3	2-methyloctacosane	408 C29H60	1000376-72-8	91
4	Hexadecane	226 C16H34	000544-76-3	90
5	Octadecane, 2-methyl-	268 C19H40	001560-88-9	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042319\
Data File : BF113930.D
Acq On : 23 Apr 2019 19:38
Operator : JU/SJ
Sample : K2529-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
P3P4042219

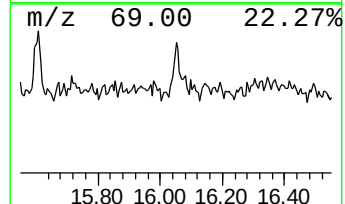
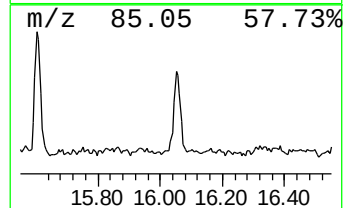
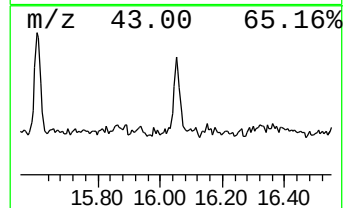
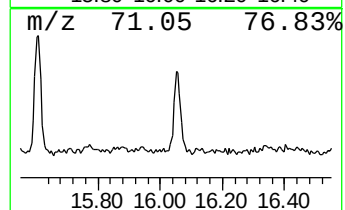
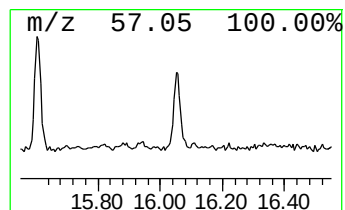
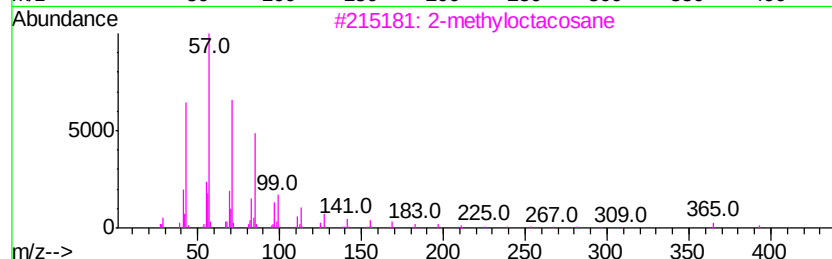
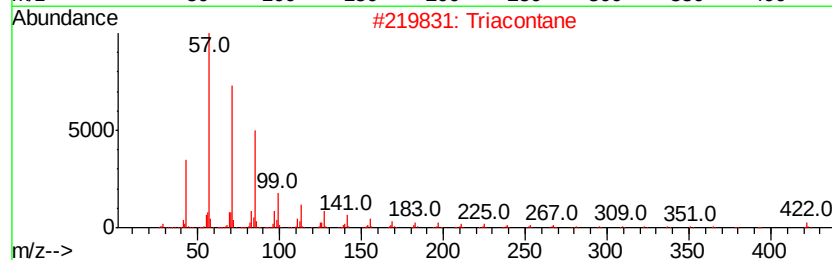
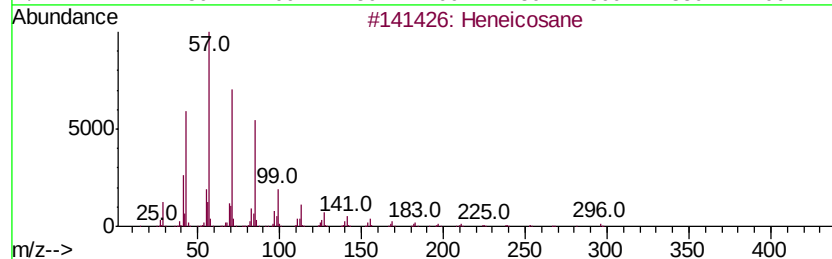
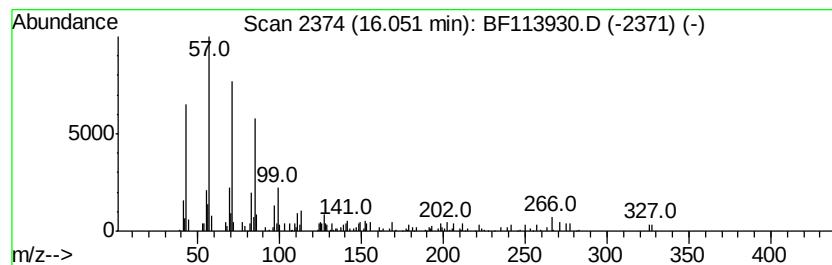
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Docosane, 11-decyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.05	2.20 ng	93993	Perylene-d12	15.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	74
2		Triacontane	422	C30H62	000638-68-6	74
3		2-methyloctacosane	408	C29H60	1000376-72-8	74
4		Tridecane, 2-methyl-	198	C14H30	001560-96-9	72
5		Docosane, 11-decyl-	451	C32H66	055401-55-3	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042319\
 Data File : BF113930.D
 Acq On : 23 Apr 2019 19:38
 Operator : JU/SJ
 Sample : K2529-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF042219.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
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n-Hexadecanoic acid	11.94	3.1	ng	129350	4	11.42	824510	20.0
Heneicosane	13.57	2.3	ng	119728	5	14.06	1054610	20.0
Hexacosane	13.91	3.3	ng	172030	5	14.06	1054610	20.0
Heptacosane	14.23	3.0	ng	159105	5	14.06	1054610	20.0
Octacosane	14.55	2.8	ng	148360	5	14.06	1054610	20.0
2-methyloctacosane	14.87	3.3	ng	139852	6	15.57	852582	20.0
Octadecane, 2-met...	15.22	3.5	ng	149846	6	15.57	852582	20.0
Docosane, 11-decyl-	16.05	2.2	ng	93993	6	15.57	852582	20.0