

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061419\
 Data File : BF115050.D
 Acq On : 14 Jun 2019 19:58
 Operator : HP/JU
 Sample : K3351-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RT-3407

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-BF061219.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	5.263	534	540	543	rBV	5563248	6157884	92.06%	10.438%
2	6.292	709	715	718	rBV	5030687	6512937	97.37%	11.040%
3	6.416	731	736	739	rBV	6054308	6688763	100.00%	11.338%
4	6.645	771	775	785	rVB	1806063	1557840	23.29%	2.641%
5	6.804	797	802	805	rBV	5251900	5002733	74.79%	8.480%
6	7.210	865	871	874	rBV	3942572	4252155	63.57%	7.208%
7	7.922	988	992	999	rBV	2016464	2010943	30.06%	3.409%
8	9.004	1170	1176	1179	rBV	6604181	6424687	96.05%	10.890%
9	9.392	1238	1242	1255	rVB	842786	781595	11.69%	1.325%
10	9.669	1284	1289	1293	rBV	2274513	2259246	33.78%	3.830%
11	10.457	1418	1423	1426	rBV	4134296	3893372	58.21%	6.599%
12	11.145	1536	1540	1543	rBV	2572485	2195631	32.83%	3.722%
13	11.169	1543	1544	1548	rVB	174959	94630	1.41%	0.160%
14	11.692	1630	1633	1642	rBV2	262350	336284	5.03%	0.570%
15	12.345	1741	1744	1749	rBV	179088	169321	2.53%	0.287%
16	12.474	1763	1766	1774	rBV6	37690	69265	1.04%	0.117%
17	12.574	1778	1783	1786	rBV	132742	142523	2.13%	0.242%
18	12.727	1804	1809	1812	rBV	5722770	5584819	83.50%	9.467%
19	13.645	1962	1965	1971	rVB	89087	91226	1.36%	0.155%
20	13.768	1981	1986	1989	rBV	1746205	1768714	26.44%	2.998%
21	13.957	2015	2018	2027	rVB	135631	153756	2.30%	0.261%
22	14.263	2066	2070	2074	rBV	157216	162914	2.44%	0.276%
23	14.415	2092	2096	2100	rBV2	61975	91418	1.37%	0.155%
24	14.551	2116	2119	2124	rBV	168533	161671	2.42%	0.274%
25	14.763	2152	2155	2163	rVB	29178	68230	1.02%	0.116%
26	14.857	2167	2171	2179	rVV	205730	218677	3.27%	0.371%
27	15.145	2215	2220	2225	rBV	1195528	1466560	21.93%	2.486%
28	15.198	2225	2229	2237	rVB	207247	268637	4.02%	0.455%
29	15.586	2289	2295	2302	rBV	165245	250911	3.75%	0.425%
30	16.027	2366	2370	2381	rVB3	92064	157729	2.36%	0.267%

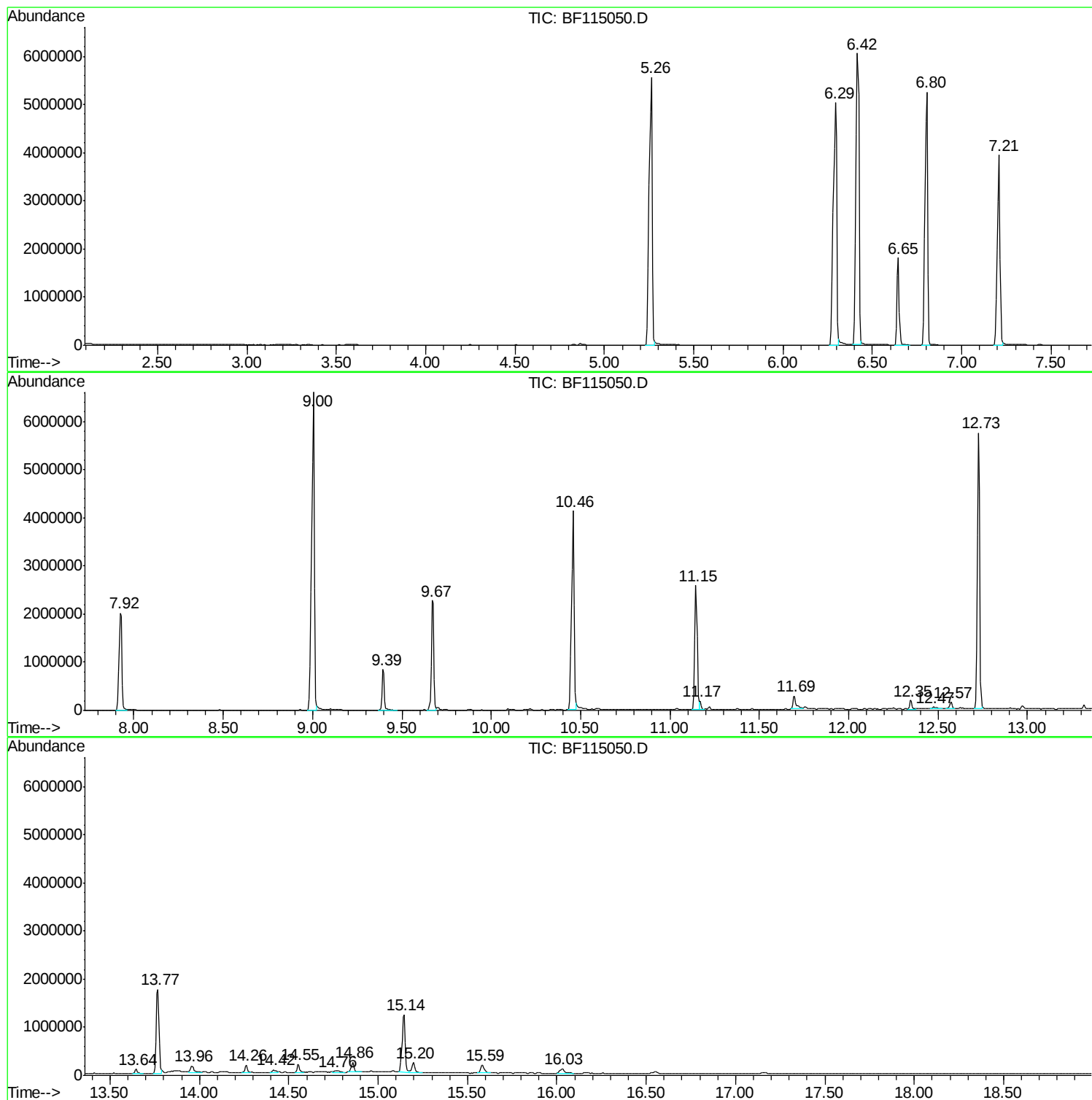
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ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
RT-3407

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061219.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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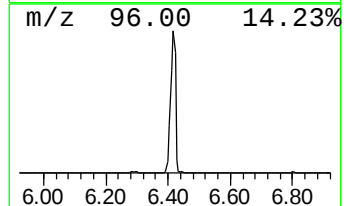
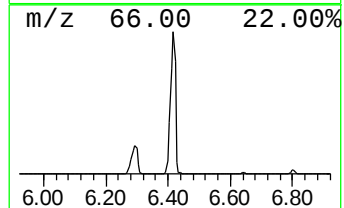
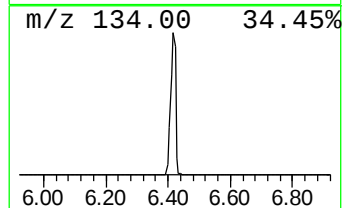
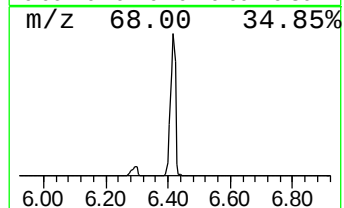
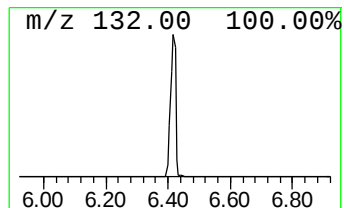
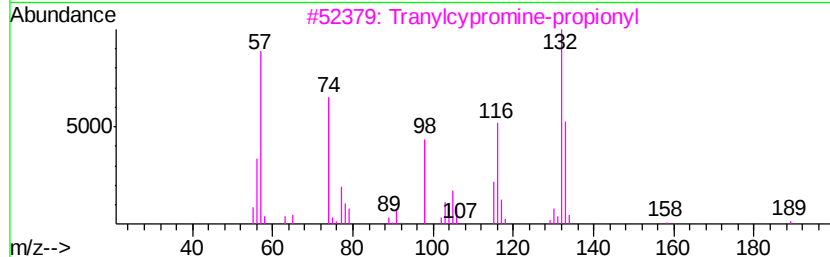
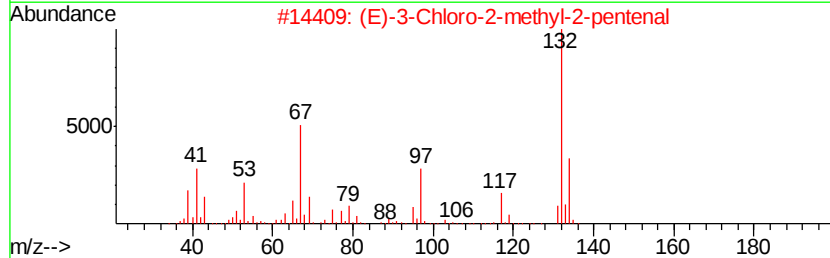
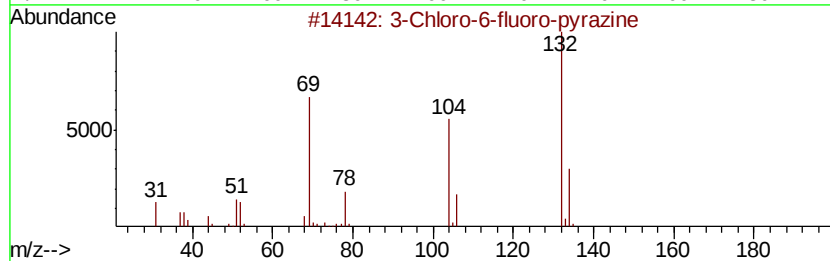
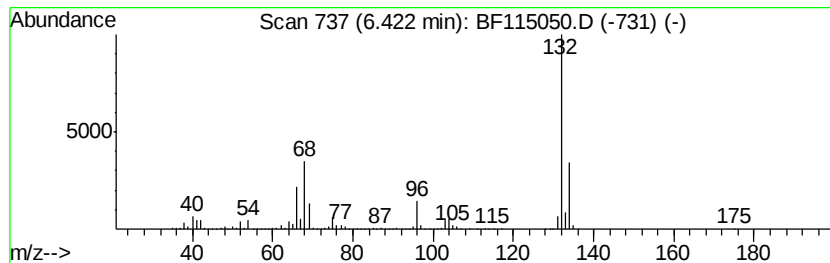
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.42 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.42	85.87 ng	6688760	1,4-Dichlorobenzene-d4	6.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Tranvlcvpromine-propionyl	189	C12H15NO	1000123-86-3	12
4		Xenon	132	Xe	007440-63-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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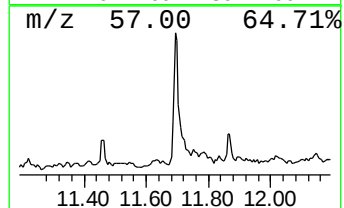
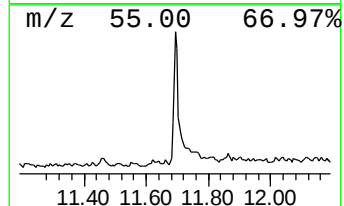
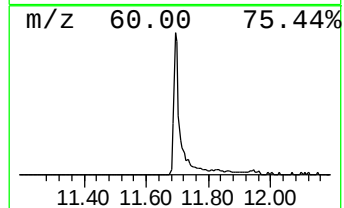
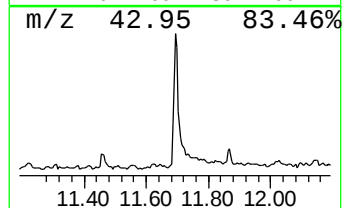
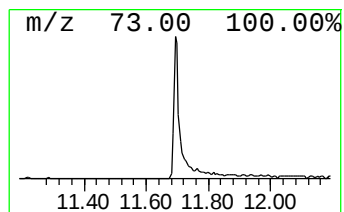
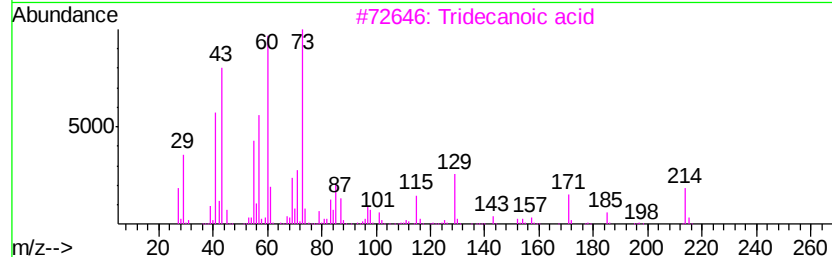
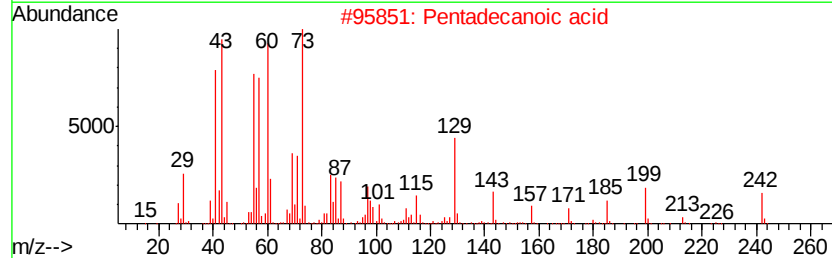
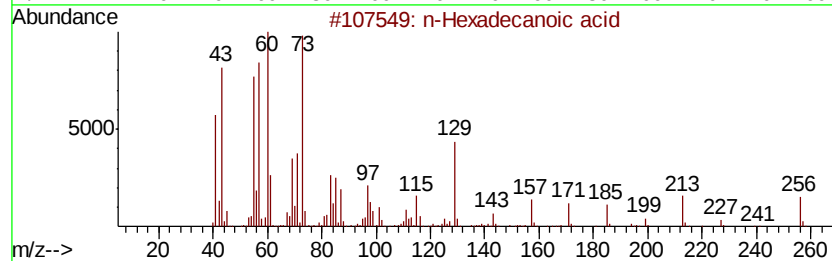
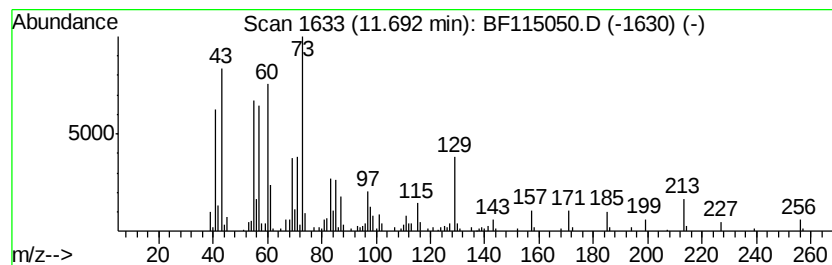
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Peak Number 3 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.69	3.06 ng	336284	Phenanthrene-d10	11.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	93
3	Tridecanoic acid	214	C13H26O2	000638-53-9	89
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5	n-Decanoic acid	172	C10H20O2	000334-48-5	60



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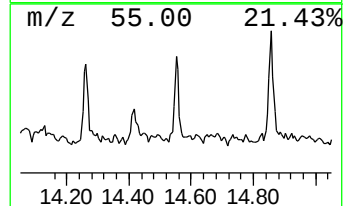
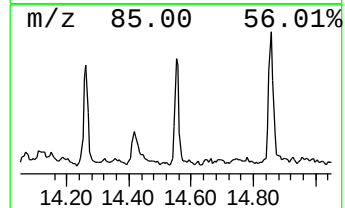
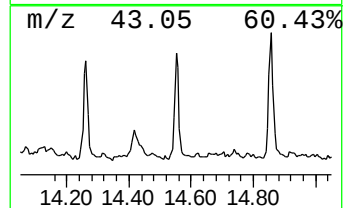
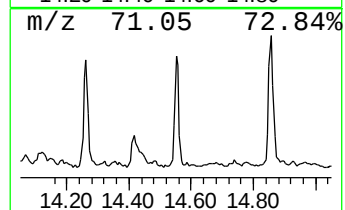
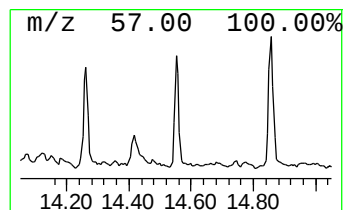
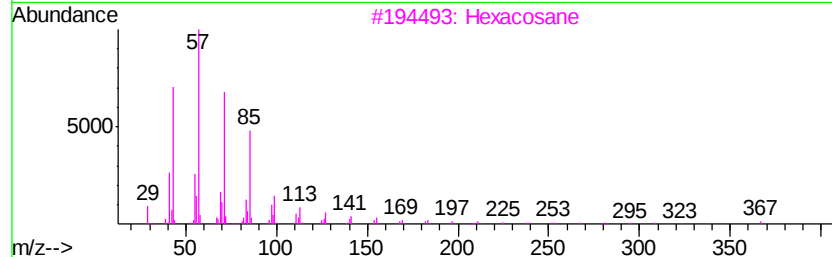
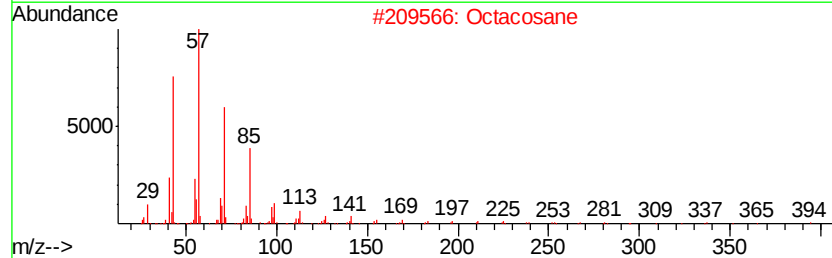
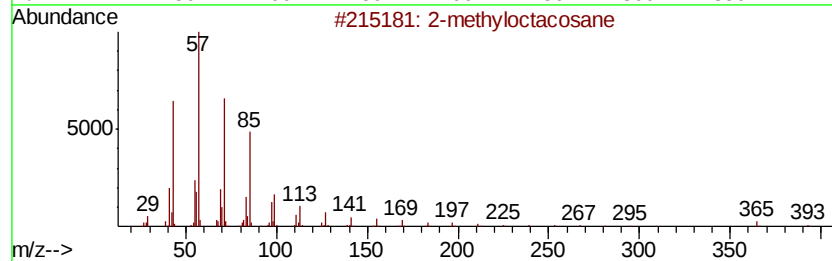
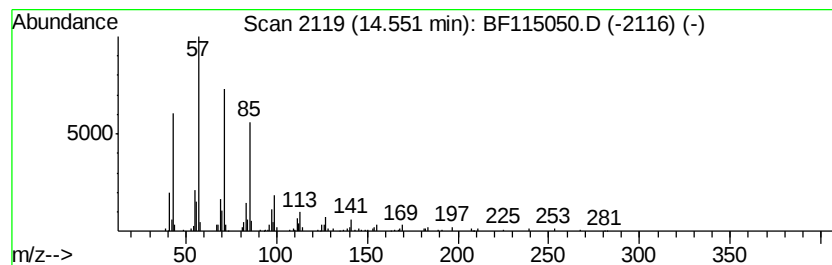
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Peak Number 4 2-methyloctacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.55	2.20 ng	161671	Perylene-d12	15.14

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-methyloctacosane		408	C29H60	1000376-72-8	91
2	Octacosane		394	C28H58	000630-02-4	90
3	Hexacosane		366	C26H54	000630-01-3	90
4	Heneicosane		296	C21H44	000629-94-7	90
5	Tetracosane		338	C24H50	000646-31-1	87



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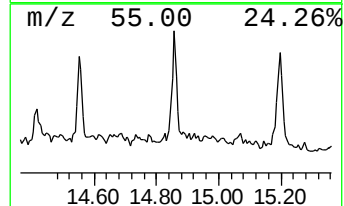
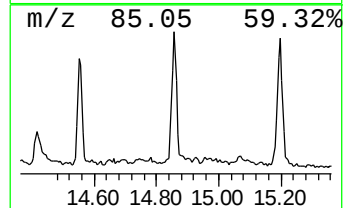
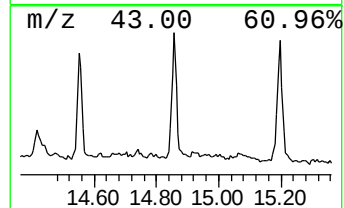
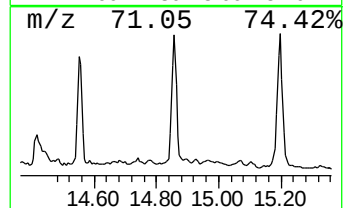
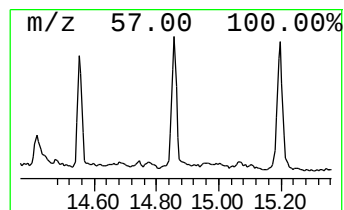
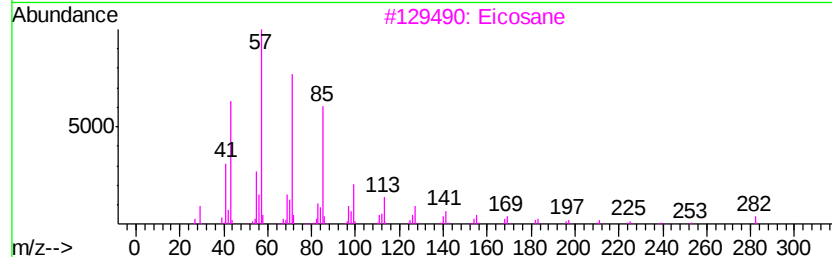
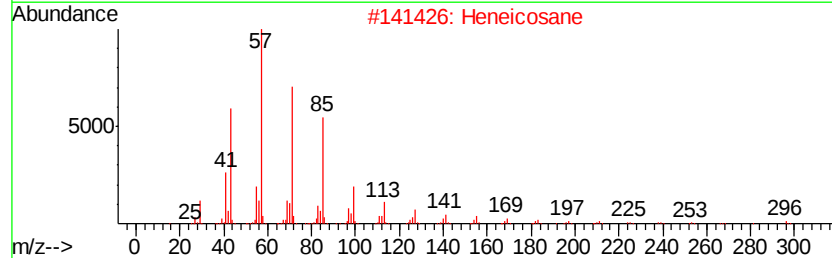
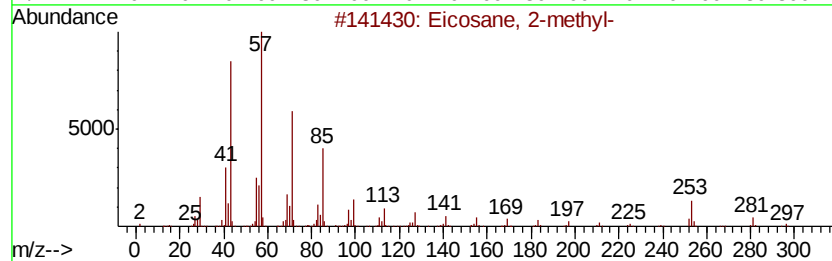
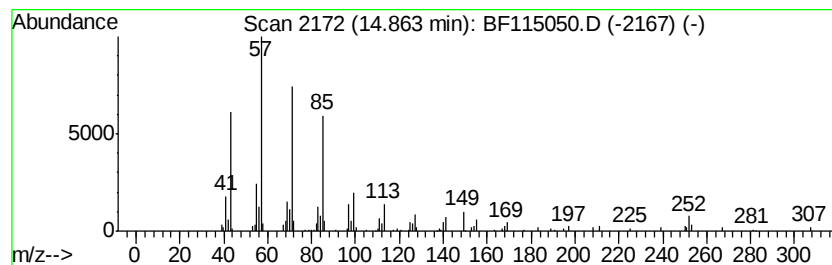
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Eicosane, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
14.86	2.98 ng	218677	Perylene-d12		15.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane, 2-methyl-			296	C21H44	001560-84-5	91
2	Heneicosane			296	C21H44	000629-94-7	90
3	Eicosane			282	C20H42	000112-95-8	90
4	Octacosane			394	C28H58	000630-02-4	87
5	Heptadecane, 9-octyl-			352	C25H52	007225-64-1	87



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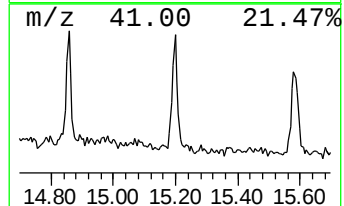
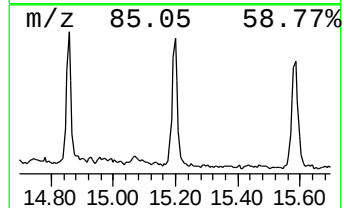
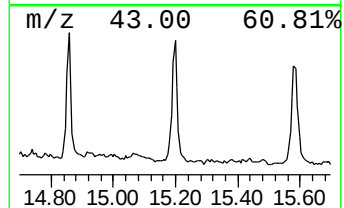
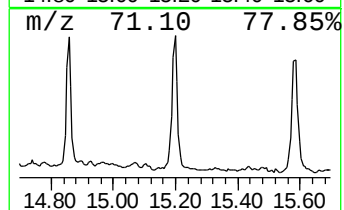
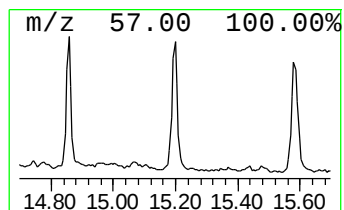
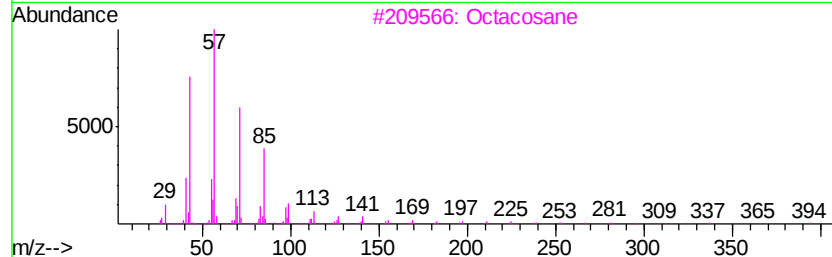
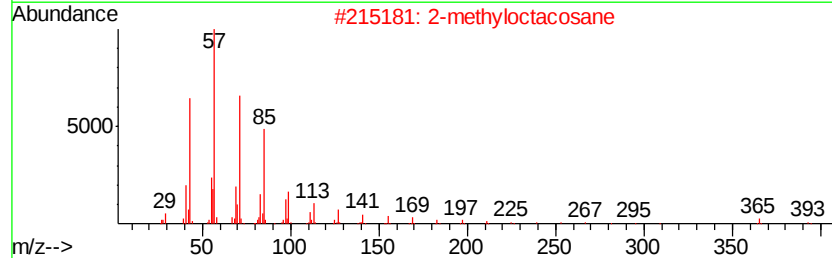
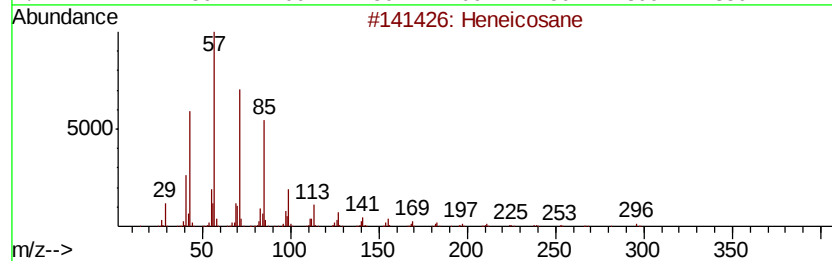
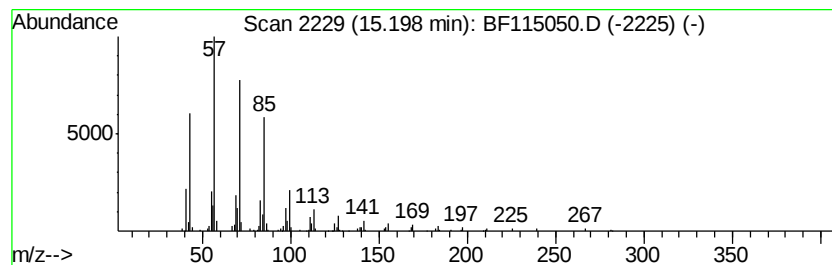
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Heneicosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.20	3.66 ng	268637	Perylene-d12	15.14

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	2-methyloctacosane		408	C29H60	1000376-72-8	91
3	Octacosane		394	C28H58	000630-02-4	91
4	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
5	Hentriacontane		437	C31H64	000630-04-6	91



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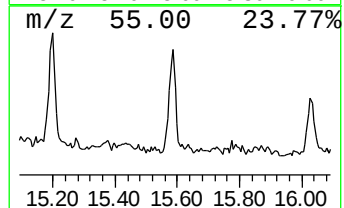
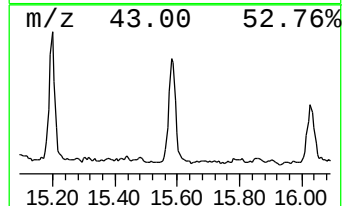
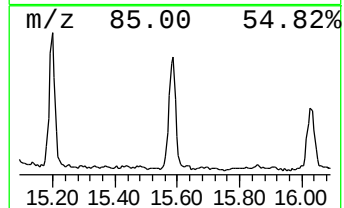
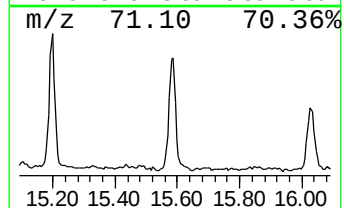
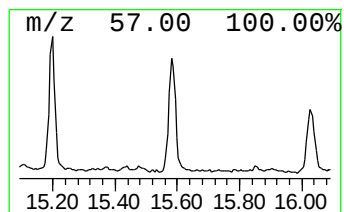
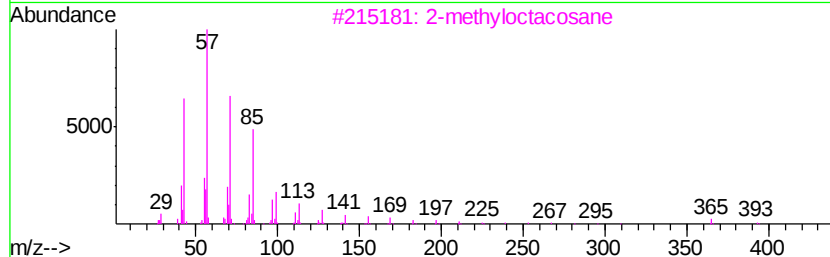
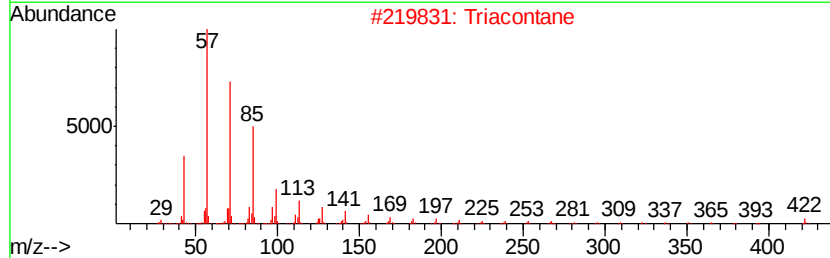
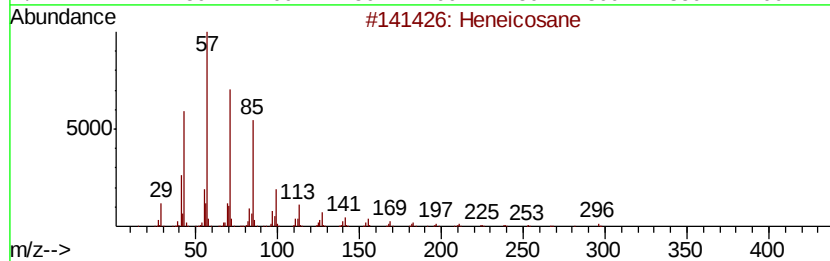
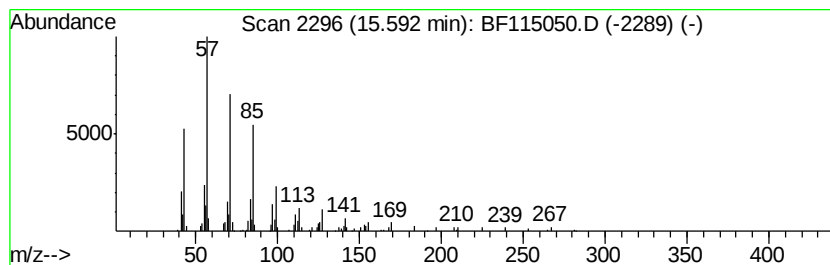
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Triacontane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.59	3.42 ng	250911	Perylene-d12	15.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Triacontane	422	C30H62	000638-68-6	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061419\
Data File : BF115050.D
Acq On : 14 Jun 2019 19:58
Operator : HP/JU
Sample : K3351-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RT-3407

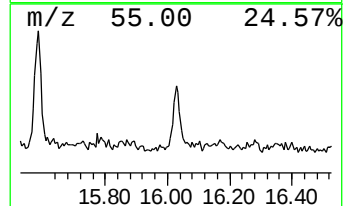
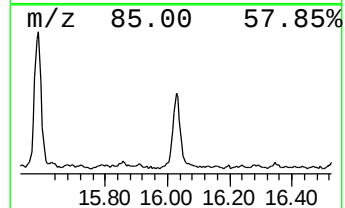
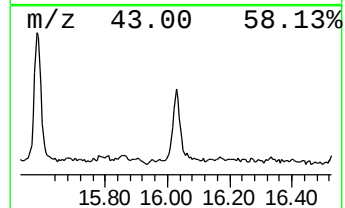
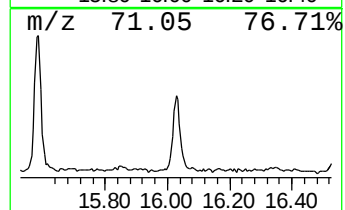
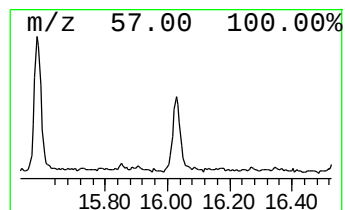
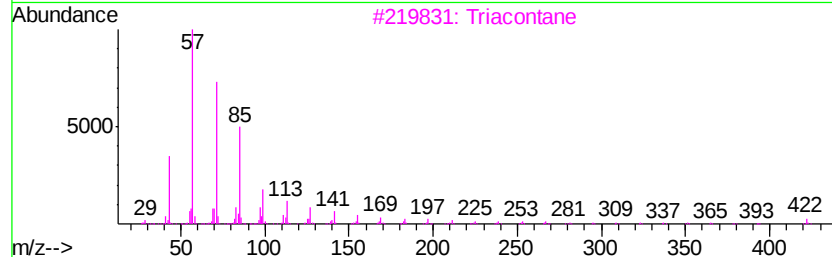
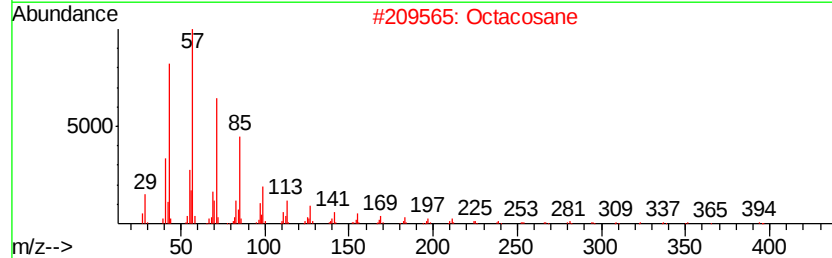
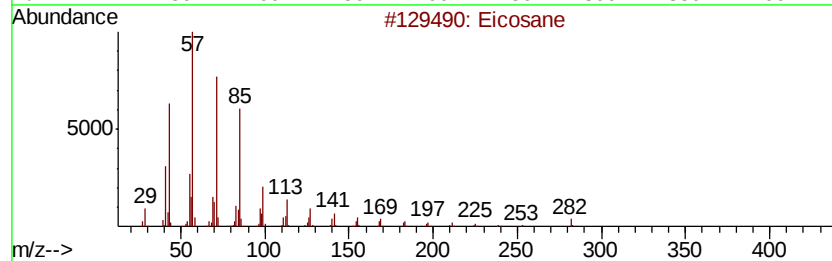
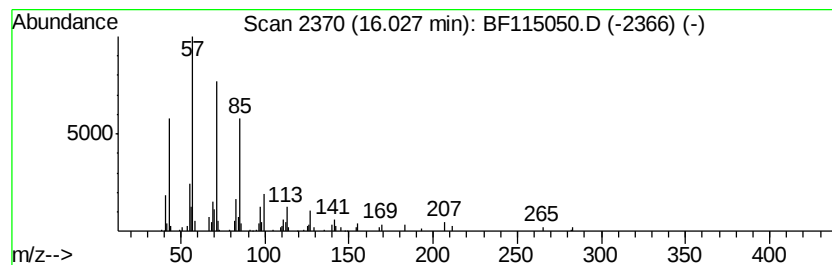
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.03	2.15 ng	157729	Perylene-d12	15.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	97
2		Octacosane	394	C28H58	000630-02-4	91
3		Triacontane	422	C30H62	000638-68-6	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		2-methyloctacosane	408	C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF061419\
Data File : BF115050.D
Acq On : 14 Jun 2019 19:58
Operator : HP/JU
Sample : K3351-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RT-3407

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061219.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
unknown6.42	6.42	85.9	ng	6688760	1	6.65	1557840	20.0
n-Hexadecanoic acid	11.69	3.1	ng	336284	4	11.15	2195630	20.0
2-methyloctacosane	14.55	2.2	ng	161671	6	15.14	1466560	20.0
Eicosane, 2-methyl-	14.86	3.0	ng	218677	6	15.14	1466560	20.0
Heneicosane	15.20	3.7	ng	268637	6	15.14	1466560	20.0
Triacontane	15.59	3.4	ng	250911	6	15.14	1466560	20.0
Octacosane	16.03	2.1	ng	157729	6	15.14	1466560	20.0