

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050616\
 Data File : BF086737.D
 Acq On : 6 May 2016 21:30
 Operator : UM/SJ
 Sample : H2770-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW27

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:\HPCHEM1\BNA_F\METHODS\8270-BF050416.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.899	244	246	253	rBV	63456	101774	2.15%	0.280%
2	5.310	280	282	297	rBV	1647931	1822082	38.49%	5.015%
3	6.362	372	374	380	rBV	1028540	1184417	25.02%	3.260%
4	6.476	382	384	387	rBV	3266140	3331433	70.37%	9.170%
5	6.704	401	404	406	rBV	1095447	1012532	21.39%	2.787%
6	6.864	415	418	420	rBV	2742712	3308996	69.89%	9.108%
7	7.276	451	454	457	rBV	2595711	2757282	58.24%	7.589%
8	7.402	463	465	469	rVB	47932	69529	1.47%	0.191%
9	7.996	514	517	519	rBV	1036503	1304817	27.56%	3.591%
10	9.070	608	611	613	rBV	4721189	4734401	100.00%	13.031%
11	9.470	644	646	650	rVB	119534	130631	2.76%	0.360%
12	9.676	663	664	667	rBV	44244	51317	1.08%	0.141%
13	9.745	667	670	672	rBV	1335269	1461049	30.86%	4.022%
14	10.156	703	706	707	rBV2	40006	62043	1.31%	0.171%
15	10.179	707	708	710	rVB	165160	125007	2.64%	0.344%
16	10.396	724	727	730	rBV	52464	93687	1.98%	0.258%
17	10.533	736	739	741	rBV	3000867	3316426	70.05%	9.128%
18	10.648	747	749	754	rVB	176207	223489	4.72%	0.615%
19	11.093	787	788	790	rBV	120706	128831	2.72%	0.355%
20	11.219	797	799	802	rBV	1523233	1368792	28.91%	3.768%
21	11.276	802	804	808	rVB	42404	68603	1.45%	0.189%
22	11.528	823	826	829	rBV	44619	70813	1.50%	0.195%
23	11.768	844	847	850	rBV	758178	767713	16.22%	2.113%
24	12.454	905	907	913	rBV2	55437	138949	2.93%	0.382%
25	12.545	913	915	921	rVV	689840	855389	18.07%	2.354%
26	12.636	921	923	926	rVB2	37532	52684	1.11%	0.145%
27	12.808	935	938	940	rBV	4149526	4123468	87.10%	11.350%
28	13.036	957	958	960	rVB	60921	67808	1.43%	0.187%
29	13.379	986	988	992	rVB	116153	130650	2.76%	0.360%
30	13.699	1014	1016	1019	rBV	501201	668595	14.12%	1.840%
31	13.848	1027	1029	1033	rBV	1100099	1078405	22.78%	2.968%
32	14.019	1042	1044	1047	rBV	140075	207400	4.38%	0.571%
33	14.328	1069	1071	1077	rBV	212382	253928	5.36%	0.699%
34	14.625	1094	1097	1099	rBV	142516	170984	3.61%	0.471%

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35	14.922	1121	1123	1128	rVB	115194	137515	2.90%	0.379%
36	15.231	1147	1150	1158	rVB	556199	799315	16.88%	2.200%
37	15.642	1184	1186	1188	rBV	49042	68945	1.46%	0.190%
38	15.688	1188	1190	1194	rVB2	52832	81009	1.71%	0.223%

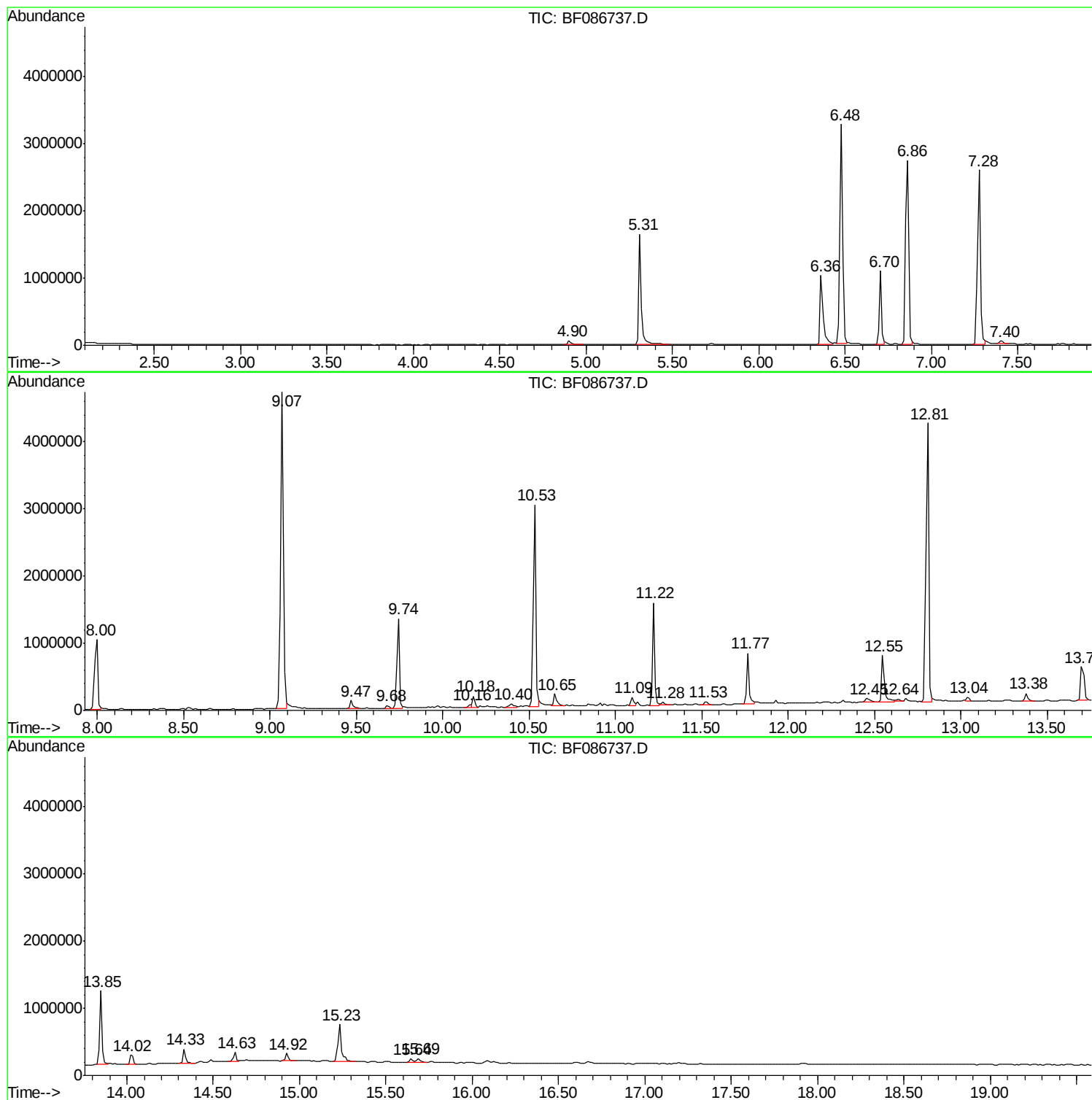
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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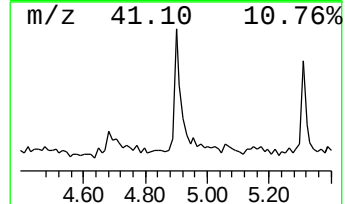
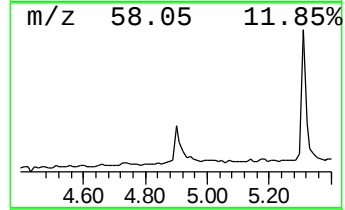
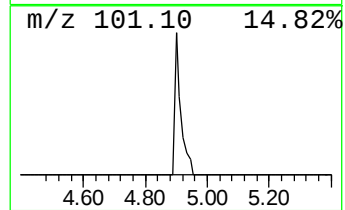
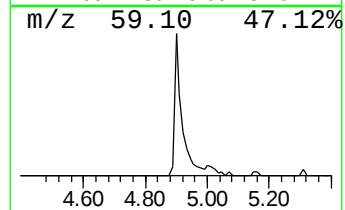
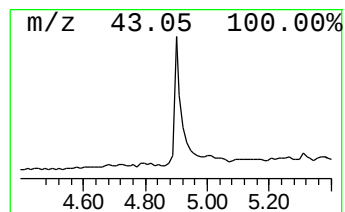
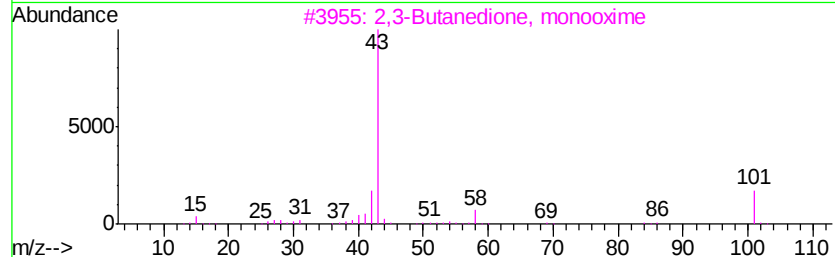
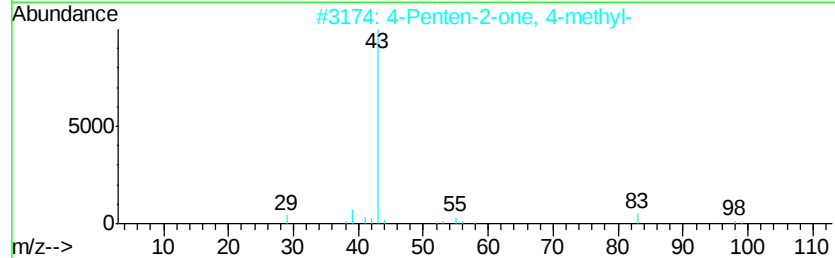
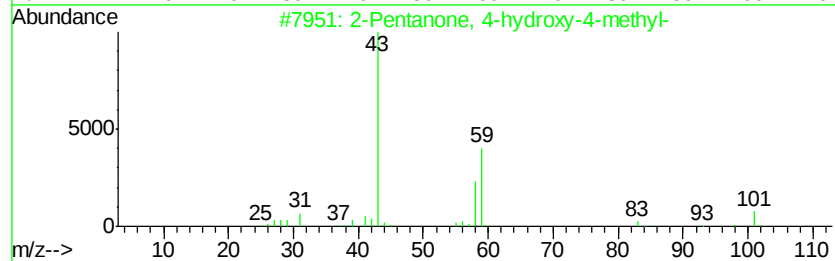
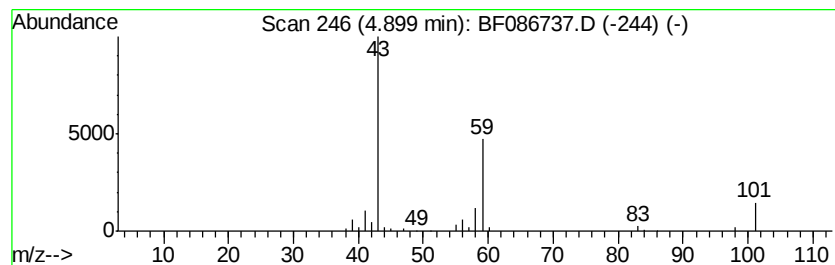
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.90	4.02 ng	101774	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	10
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		Diacetamide	101	C4H7NO2	000625-77-4	9
5		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9



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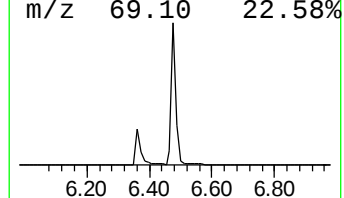
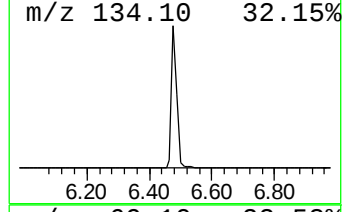
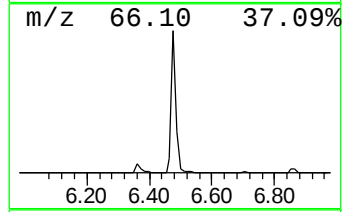
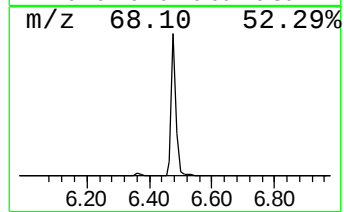
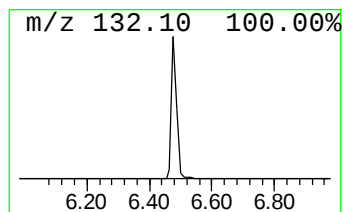
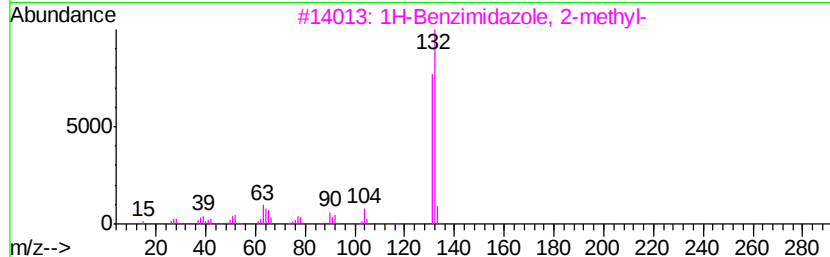
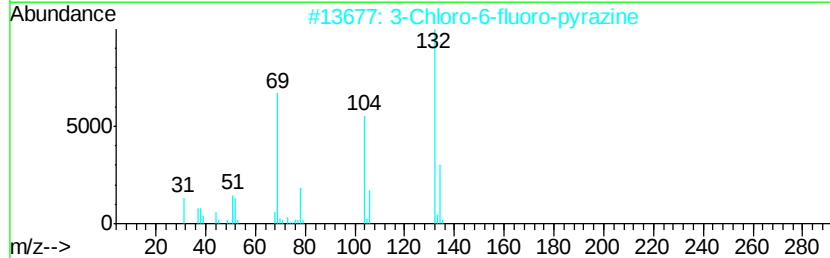
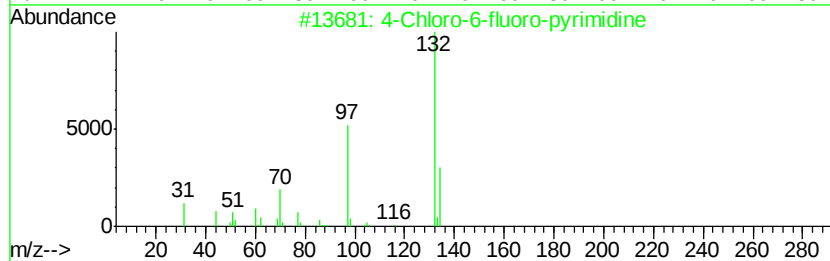
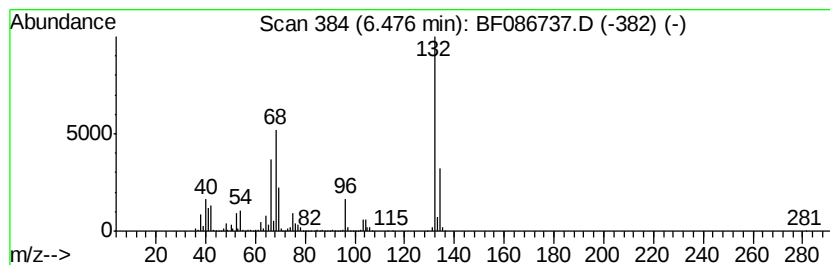
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 unknown6.48 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.48	131.61 ng	3331430	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
5		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	22



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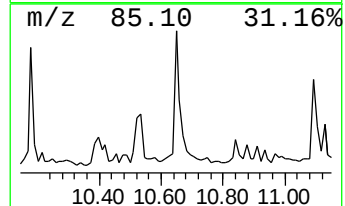
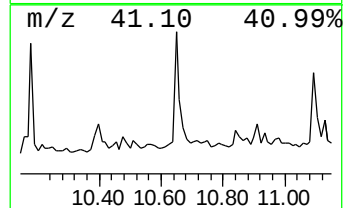
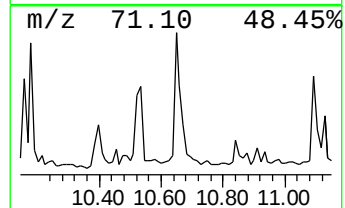
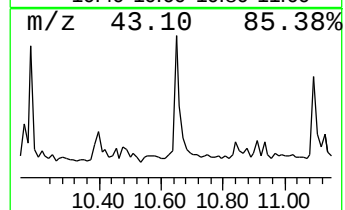
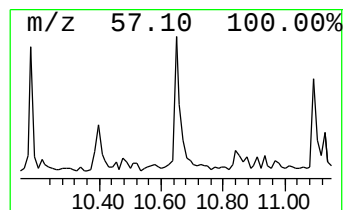
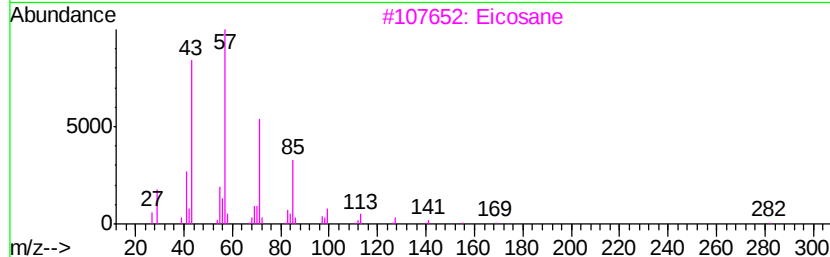
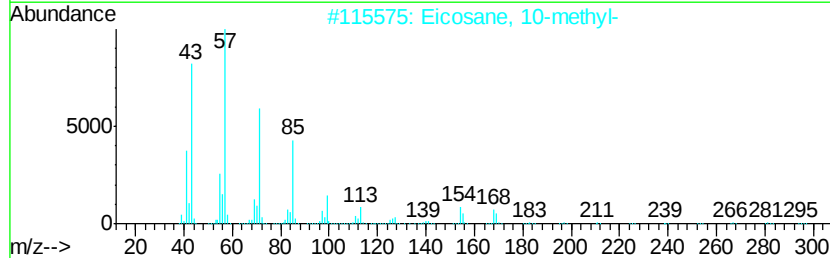
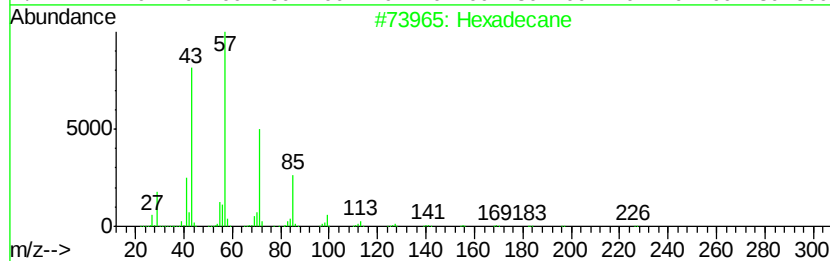
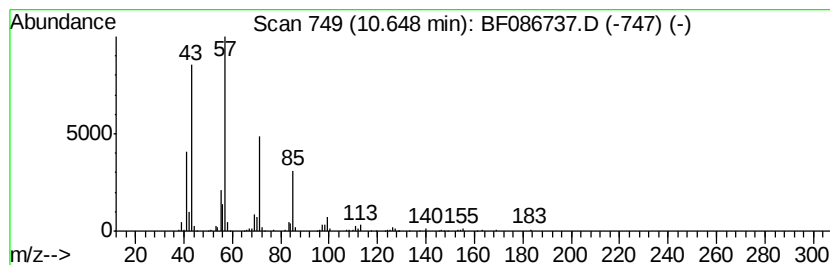
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 Hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.65	6.53 ng	223489	Phenanthrene-d10	11.22

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	91
2	Eicosane, 10-methyl-		296	C21H44	054833-23-7	90
3	Eicosane		282	C20H42	000112-95-8	80
4	Heptadecane, 2,6-dimethyl-		268	C19H40	054105-67-8	72
5	1-Iodo-2-methylnonane		268	C10H21I	1000101-47-9	72



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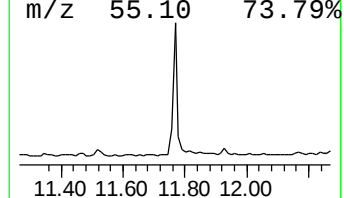
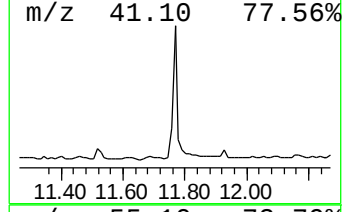
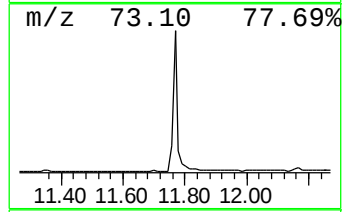
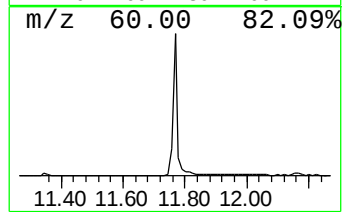
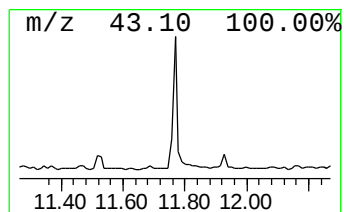
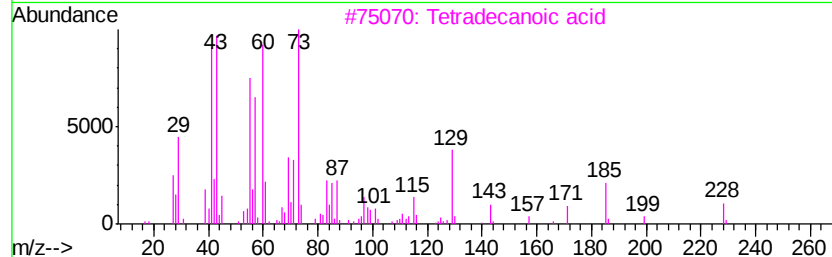
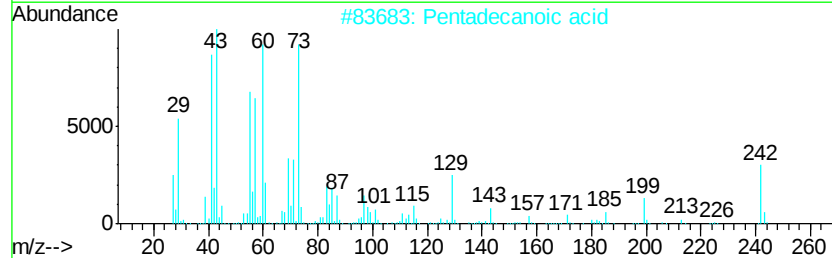
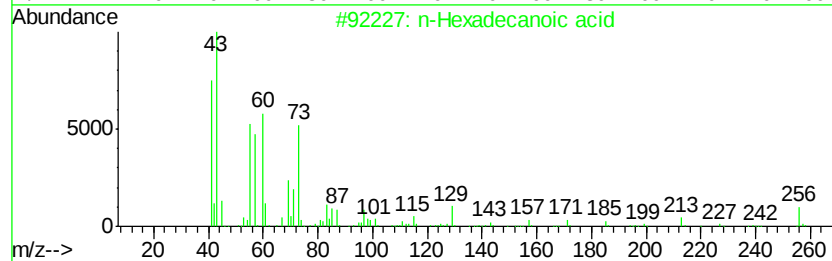
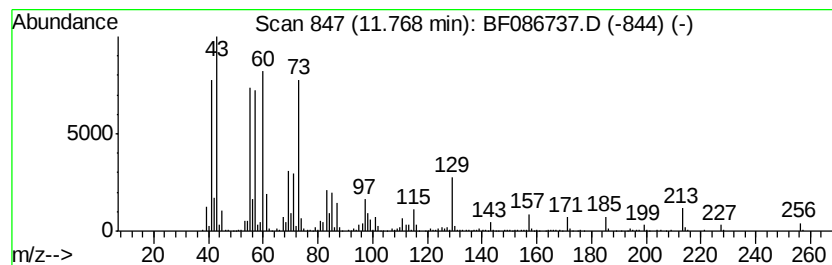
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	22.43 ng	767713	Phenanthrene-d10	11.22

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4		Tridecanoic acid	214	C13H26O2	000638-53-9	80
5		Undecanoic acid	186	C11H22O2	000112-37-8	72



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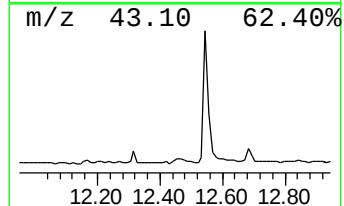
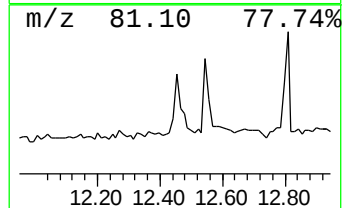
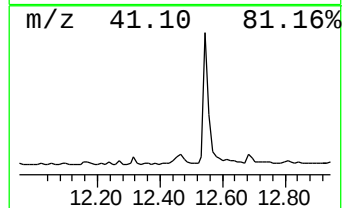
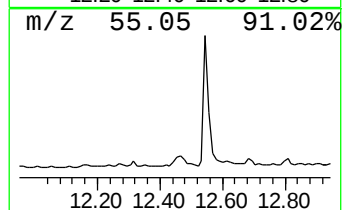
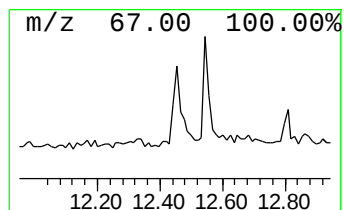
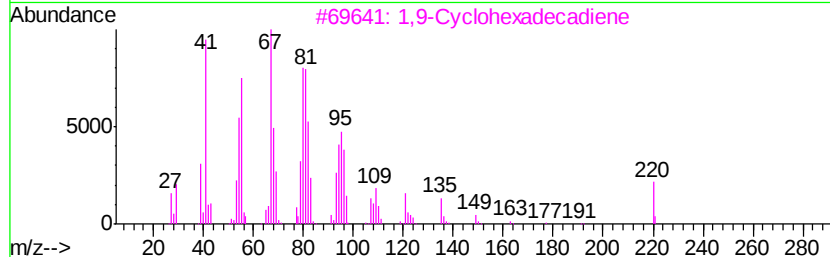
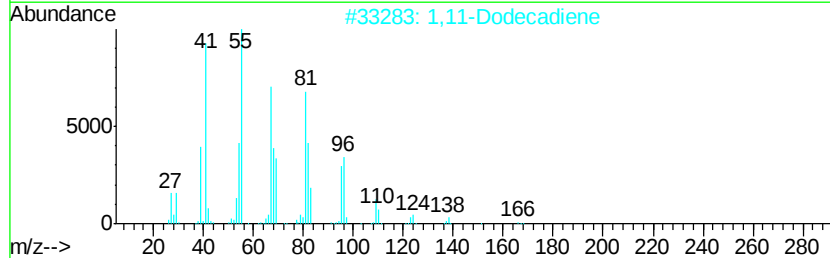
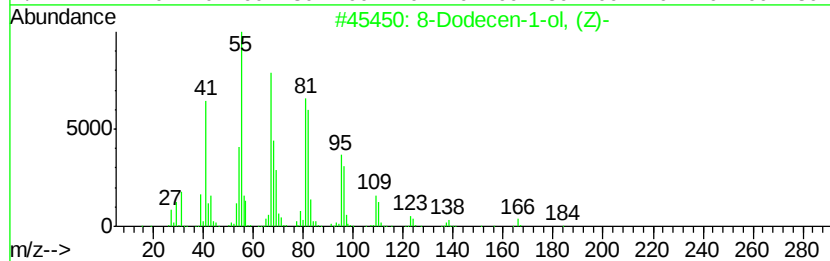
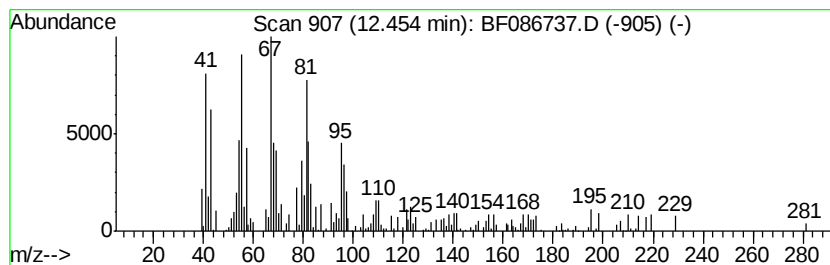
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Peak Number 6 8-Dodecen-1-ol, (Z)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.45	4.06 ng	138949	Phenanthrene-d10	11.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	8-Dodecen-1-ol, (Z)-	184	C12H24O	040642-40-8	90
2		1,11-Dodecadiene	166	C12H22	005876-87-9	89
3		1,9-Cyclohexadecadiene	220	C16H28	004277-06-9	87
4		9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	86
5		2-Chloroethyl linoleate	342	C20H35ClO2	025525-76-2	74



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ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW27

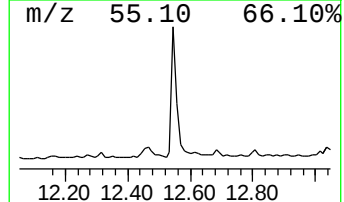
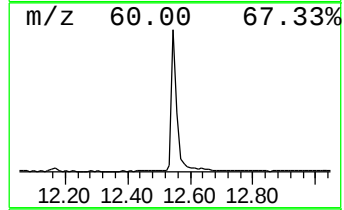
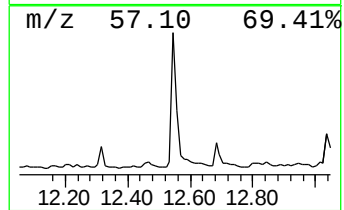
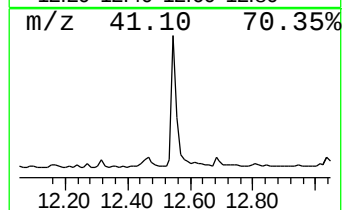
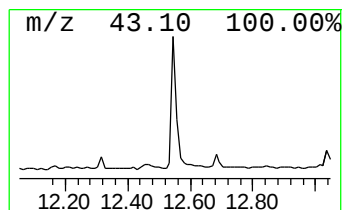
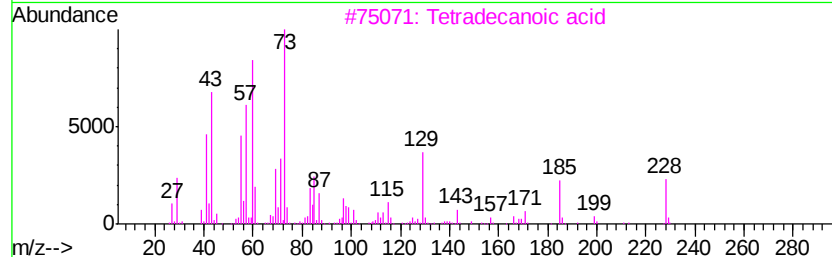
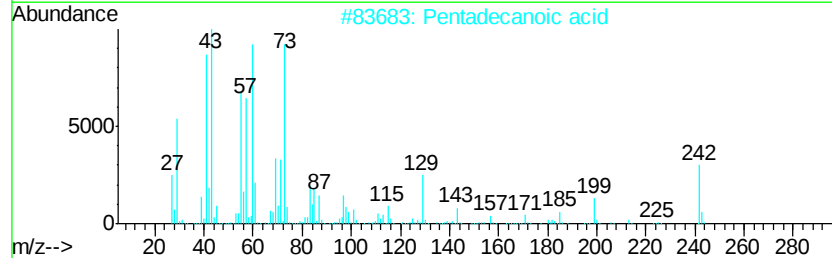
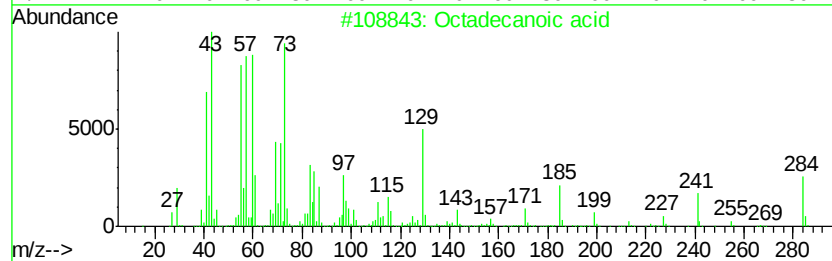
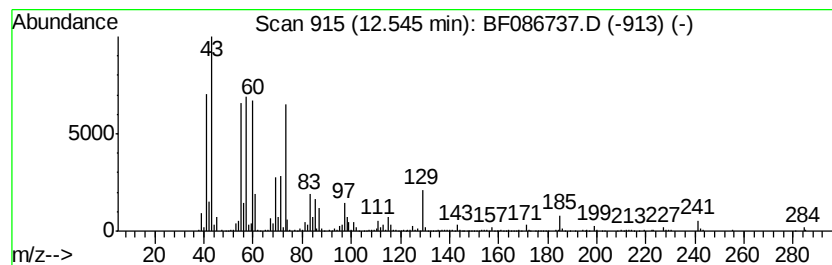
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Octadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.55	31.73 ng	855389	Chrysene-d12	13.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	98
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	91
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
5		n-Decanoic acid	172	C10H20O2	000334-48-5	76



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050616\
Data File : BF086737.D
Acq On : 6 May 2016 21:30
Operator : UM/SJ
Sample : H2770-02
Misc :
ALS Vial : 12 Sample Multiplier: 1

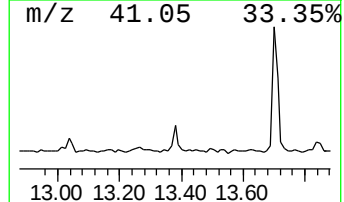
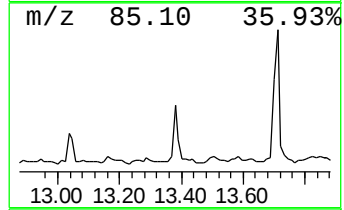
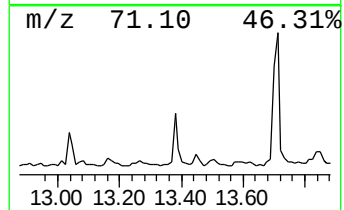
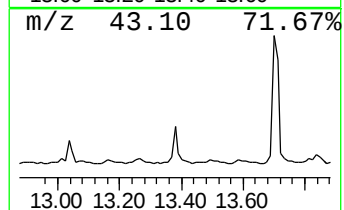
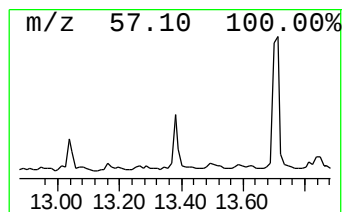
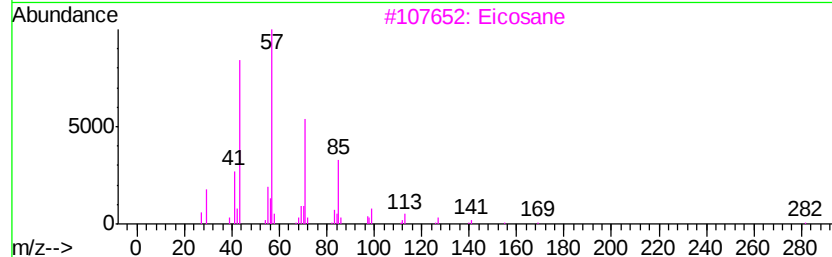
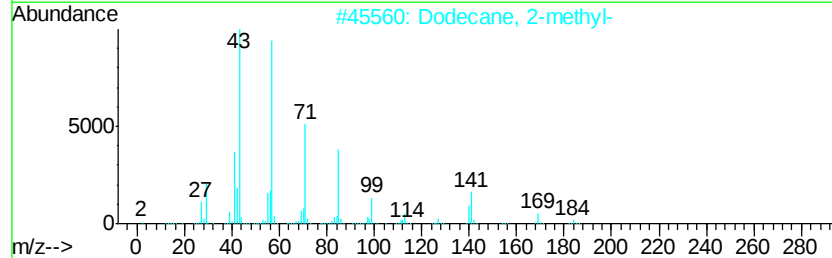
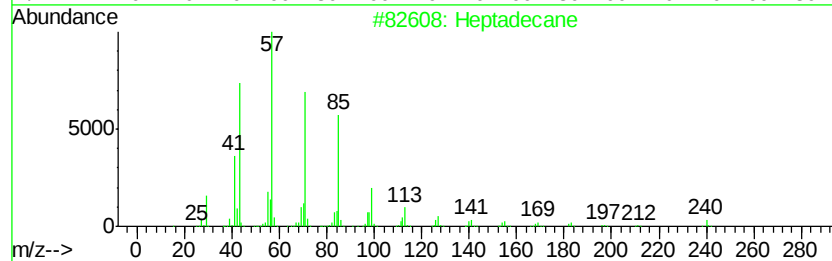
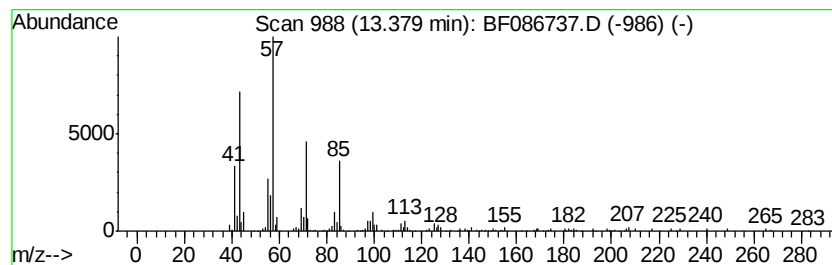
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Heptadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.38	4.85 ng	130650	Chrysene-d12	13.85	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	93
2	Dodecane, 2-methyl-	184	C13H28	001560-97-0	87
3	Eicosane	282	C20H42	000112-95-8	83
4	Tridecane, 2-methyl-	198	C14H30	001560-96-9	83
5	Tridecane	184	C13H28	000629-50-5	81



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050616\
Data File : BF086737.D
Acq On : 6 May 2016 21:30
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Instrument :
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ClientSampleId :
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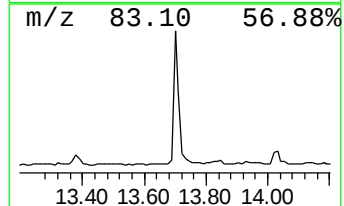
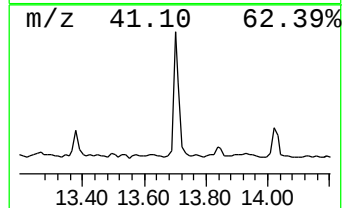
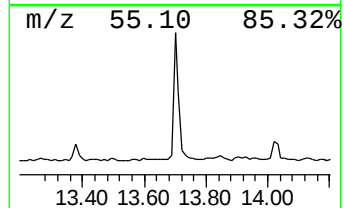
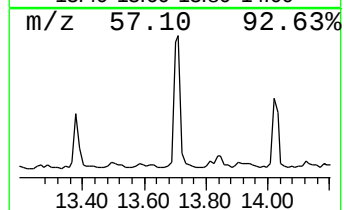
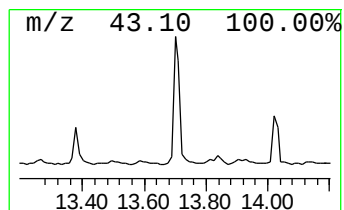
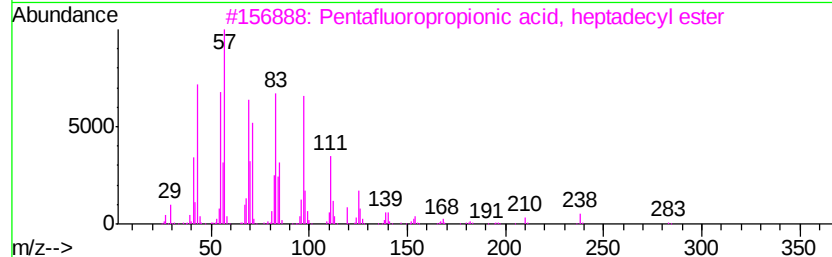
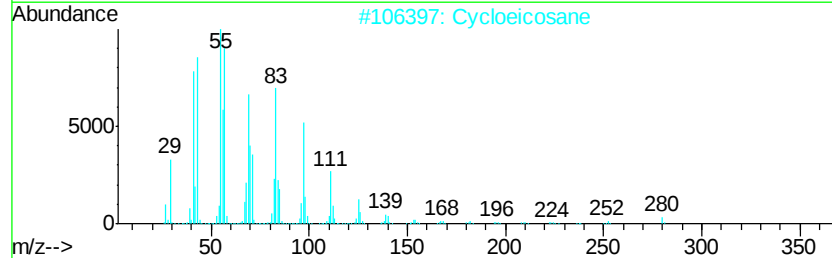
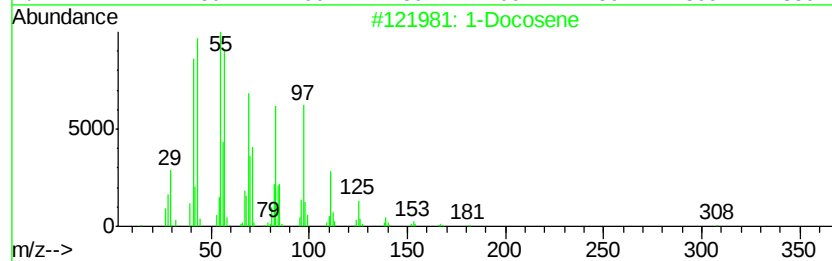
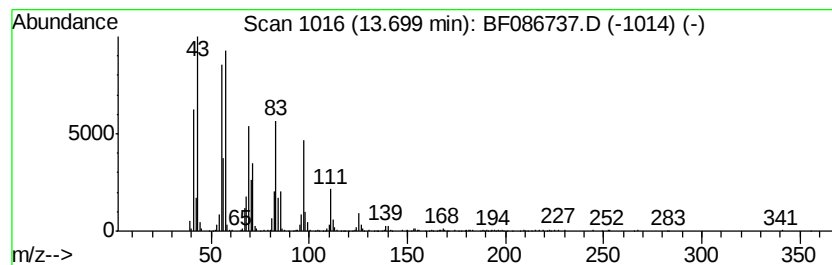
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 1-Docosene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.70	24.80 ng	668595	Chrysene-d12	13.85

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene		308	C22H44	001599-67-3	94
2	Cycloeicosane		280	C20H40	000296-56-0	91
3	Pentafluoropropionic acid, hepta...		402	C20H35F5O2	1000283-04-2	91
4	3-Eicosene, (E)-		280	C20H40	074685-33-9	91
5	Cyclotetracosane		336	C24H48	000297-03-0	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050616\
Data File : BF086737.D
Acq On : 6 May 2016 21:30
Operator : UM/SJ
Sample : H2770-02
Misc :
ALS Vial : 12 Sample Multiplier: 1

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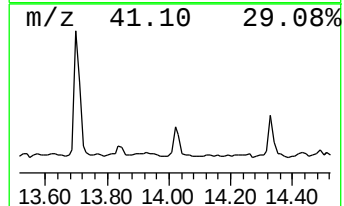
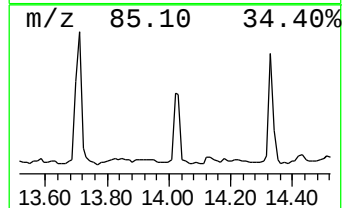
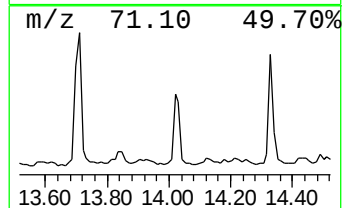
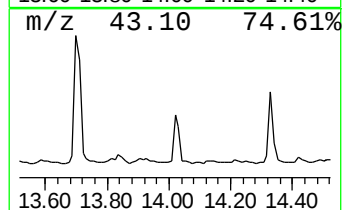
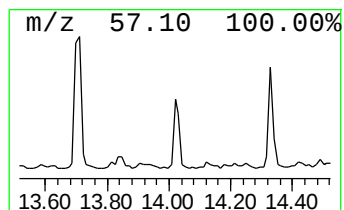
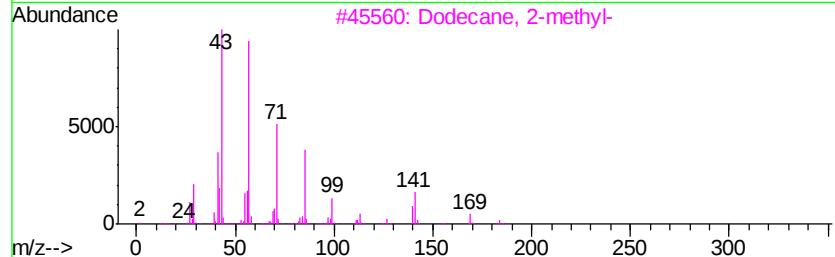
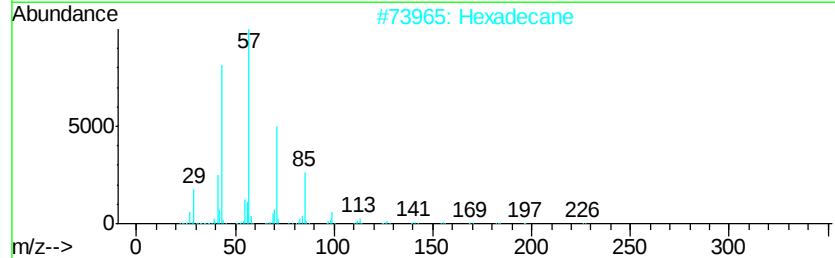
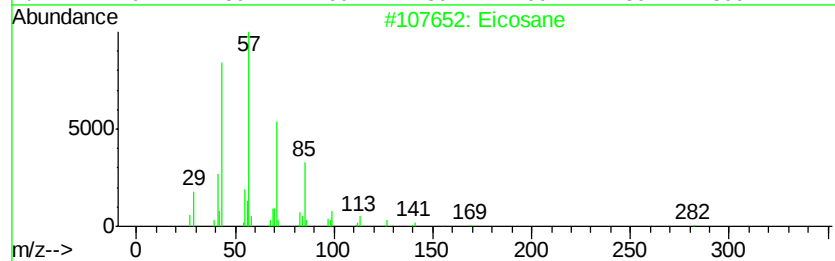
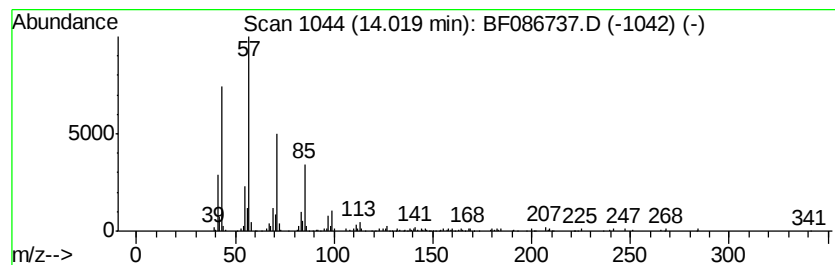
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Eicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.02	7.69 ng	207400	Chrysene-d12	13.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	87
2		Hexadecane	226	C16H34	000544-76-3	83
3		Dodecane, 2-methyl-	184	C13H28	001560-97-0	81
4		2-Bromo dodecane	248	C12H25Br	013187-99-0	80
5		Tetradecane	198	C14H30	000629-59-4	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050616\
Data File : BF086737.D
Acq On : 6 May 2016 21:30
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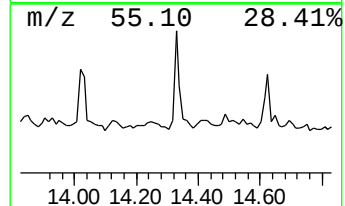
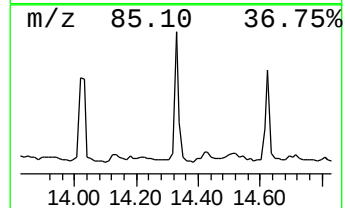
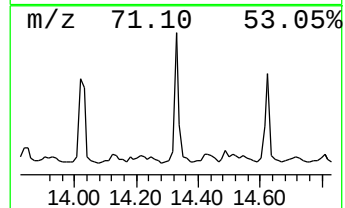
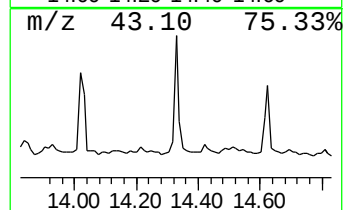
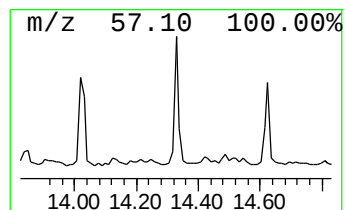
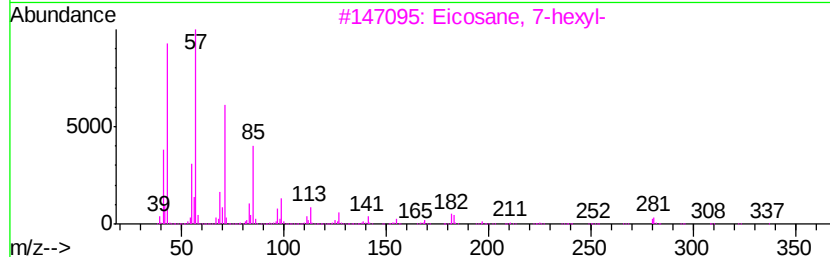
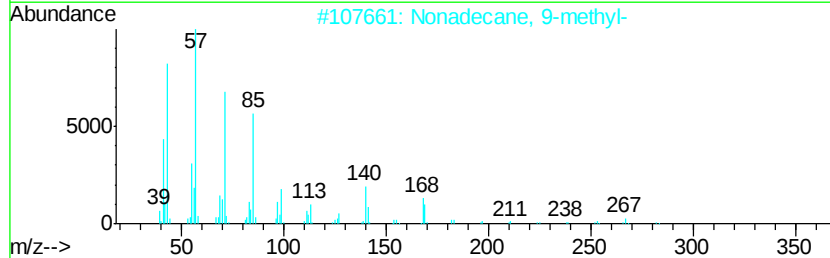
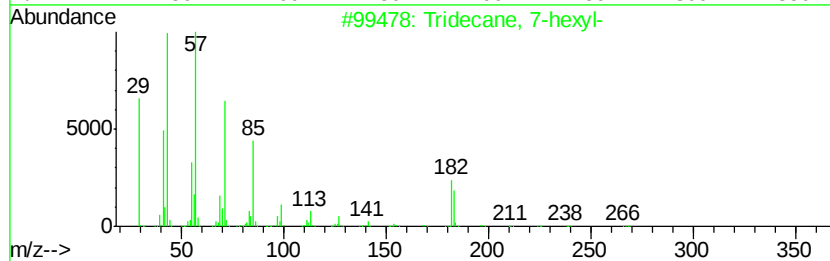
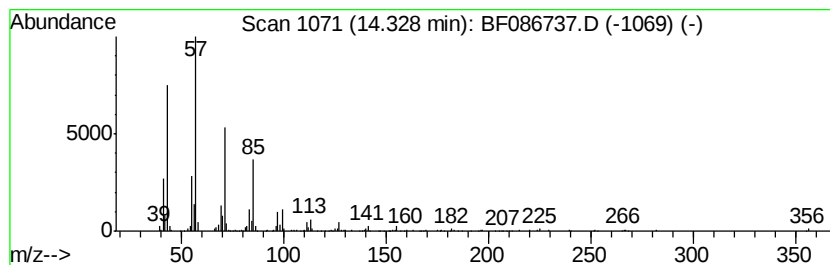
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Tridecane, 7-hexyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.33	9.42 ng	253928	Chrysene-d12	13.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 7-hexyl-	268	C19H40	007225-66-3	95
2		Nonadecane, 9-methyl-	282	C20H42	013287-24-6	93
3		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
4		Heptacosane	380	C27H56	000593-49-7	91
5		Docosane, 7-butyl-	366	C26H54	055282-15-0	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050616\
Data File : BF086737.D
Acq On : 6 May 2016 21:30
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Sample : H2770-02
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ALS Vial : 12 Sample Multiplier: 1

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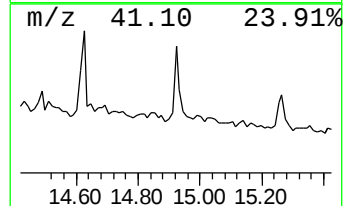
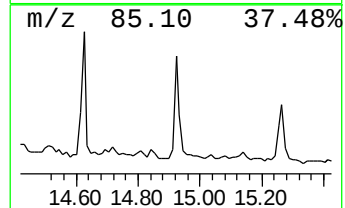
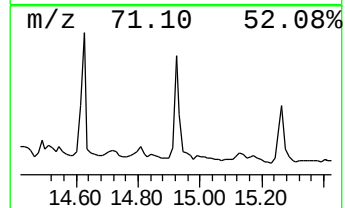
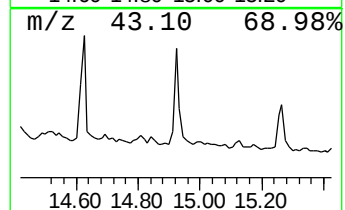
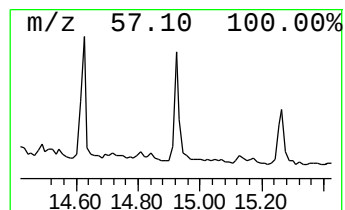
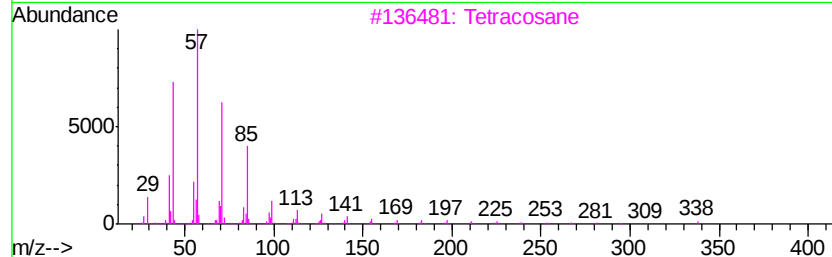
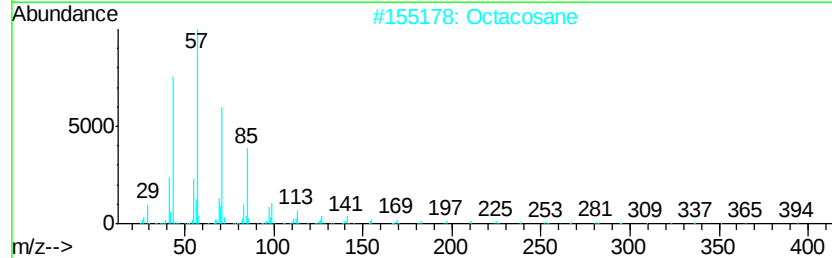
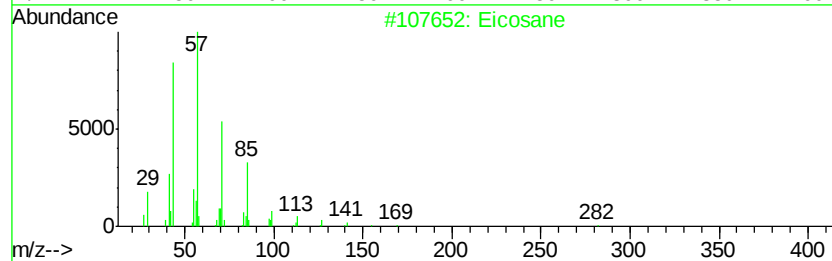
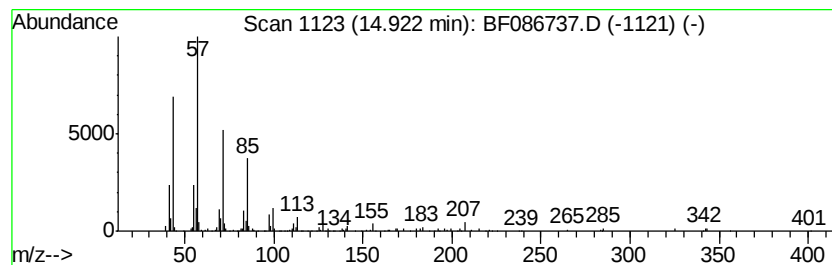
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Octacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.92	6.88 ng	137515	Perylene-d12	15.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane			282	C20H42	000112-95-8	81
2	Octacosane			394	C28H58	000630-02-4	74
3	Tetracosane			338	C24H50	000646-31-1	74
4	Tridecane, 1-iodo-			310	C13H27I	035599-77-0	74
5	Pentadecane			212	C15H32	000629-62-9	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050616\
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Acq On : 6 May 2016 21:30
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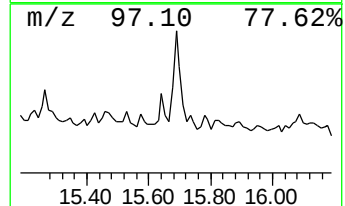
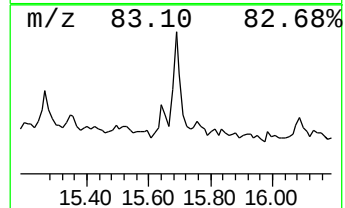
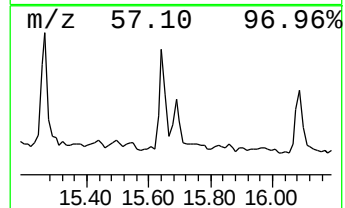
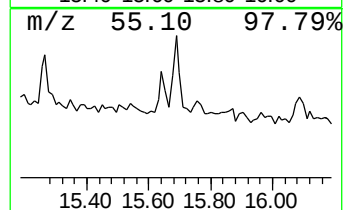
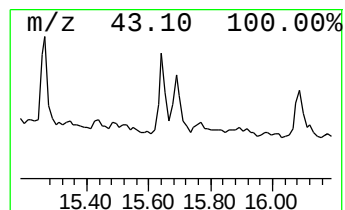
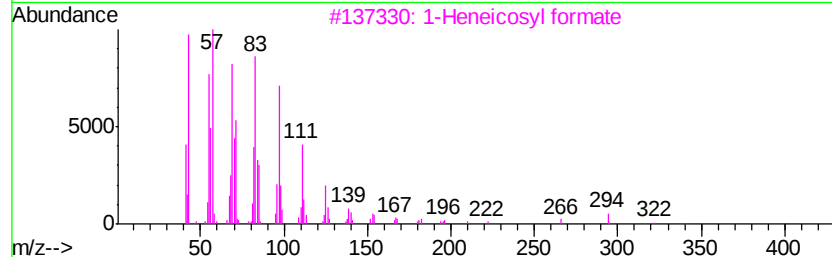
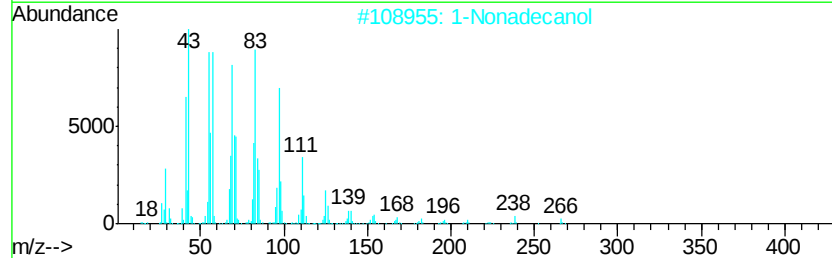
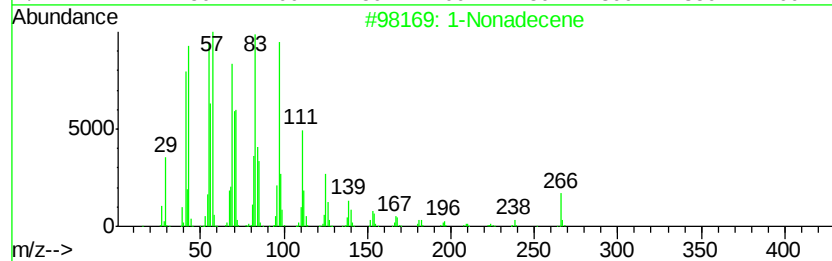
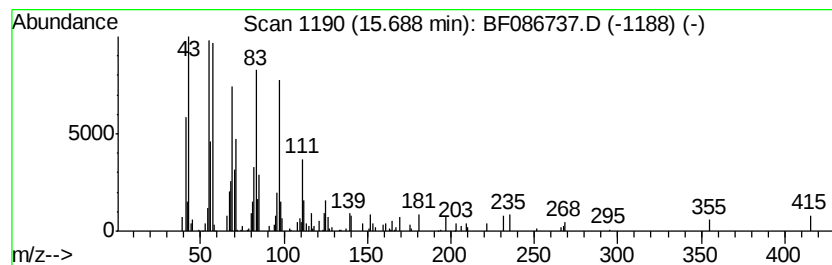
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 1-Nonadecene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.69	4.05 ng	81009	Perylene-d12	15.23	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene	266	C19H38	018435-45-5	94
2	1-Nonadecanol	284	C19H40O	001454-84-8	91
3	1-Heneicosyl formate	340	C22H44O2	077899-03-7	91
4	Chloroacetic acid, dodecyl ester	262	C14H27ClO2	006316-04-7	86
5	Z-5-Nonadecene	266	C19H38	1000131-11-8	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050616\
 Data File : BF086737.D
 Acq On : 6 May 2016 21:30
 Operator : UM/SJ
 Sample : H2770-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW27

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.90	4.0	ng	101774	1	6.70	1012530	40.0
unknown6.48	6.48	131.6	ng	3331430	1	6.70	1012530	40.0
Hexadecane	10.65	6.5	ng	223489	4	11.22	1368790	40.0
n-Hexadecanoic acid	11.77	22.4	ng	767713	4	11.22	1368790	40.0
8-Dodecen-1-ol, (Z)-	12.45	4.1	ng	138949	4	11.22	1368790	40.0
Octadecanoic acid	12.55	31.7	ng	855389	5	13.85	1078410	40.0
Heptadecane	13.38	4.8	ng	130650	5	13.85	1078410	40.0
1-Docosene	13.70	24.8	ng	668595	5	13.85	1078410	40.0
Eicosane	14.02	7.7	ng	207400	5	13.85	1078410	40.0
Tridecane, 7-hexyl-	14.33	9.4	ng	253928	5	13.85	1078410	40.0
Octacosane	14.92	6.9	ng	137515	6	15.23	799315	40.0
1-Nonadecene	15.69	4.0	ng	81009	6	15.23	799315	40.0