

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041516\
 Data File : BF086265.D
 Acq On : 16 Apr 2016 14:53
 Operator : UM/SJ
 Sample : H2471-13
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CDMSHDUP

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-BF041516.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	2.155	4	6	24	rVB	252300	590801	9.21%	0.964%
2	5.081	260	262	269	rVB	300300	436137	6.80%	0.711%
3	5.493	295	298	300	rBV	5187583	6252386	97.49%	10.197%
4	6.521	384	388	390	rBV	5448034	6413531	100.00%	10.460%
5	6.659	397	400	402	rBV	5600911	5757528	89.77%	9.390%
6	6.887	418	420	423	rVB	1452734	1640088	25.57%	2.675%
7	7.047	432	434	436	rVB	4666672	4245425	66.19%	6.924%
8	7.459	467	470	472	rBV	3943261	4452321	69.42%	7.261%
9	7.539	475	477	479	rVB	79789	90147	1.41%	0.147%
10	8.179	530	533	535	rBV	2637157	2289139	35.69%	3.733%
11	9.253	624	627	630	rBV	5481985	5677685	88.53%	9.260%
12	9.653	659	662	664	rBV	808905	1116051	17.40%	1.820%
13	9.870	677	681	683	rBV	176839	156613	2.44%	0.255%
14	9.928	684	686	689	rVB2	2165161	2327792	36.30%	3.796%
15	10.110	699	702	705	rBV4	48354	110484	1.72%	0.180%
16	10.373	721	725	726	rBV2	201365	342209	5.34%	0.558%
17	10.590	741	744	747	rBV	121067	195645	3.05%	0.319%
18	10.716	752	755	758	rBV	3358874	3467565	54.07%	5.655%
19	10.842	764	766	773	rVB	377816	458322	7.15%	0.747%
20	11.036	781	783	785	rBV2	49681	77839	1.21%	0.127%
21	11.288	803	805	806	rVV	285389	233829	3.65%	0.381%
22	11.322	806	808	811	rVV	66612	93700	1.46%	0.153%
23	11.413	814	816	818	rVV	2555838	2256040	35.18%	3.679%
24	11.470	818	821	825	rVB	42462	73321	1.14%	0.120%
25	11.642	834	836	841	rVV	130898	148304	2.31%	0.242%
26	11.711	841	842	846	rVV	90615	99967	1.56%	0.163%
27	11.939	861	862	864	rBV	132300	165665	2.58%	0.270%
28	11.973	864	865	867	rVB	81069	72059	1.12%	0.118%
29	12.111	874	877	883	rVV3	75027	145558	2.27%	0.237%
30	12.236	886	888	894	rVB	247427	227778	3.55%	0.371%
31	12.511	910	912	916	rVB2	67524	83925	1.31%	0.137%
32	12.613	919	921	923	rBV	43447	69444	1.08%	0.113%
33	12.831	937	940	943	rBV2	241872	321318	5.01%	0.524%
34	12.876	943	944	950	rVB2	106107	206694	3.22%	0.337%

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

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

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35	13.002	952	955	957	rBV	4572581	4667612	72.78%	7.612%
36	13.242	972	976	978	rBV2	124095	195987	3.06%	0.320%
37	13.528	999	1001	1004	rBV2	133358	177473	2.77%	0.289%
38	13.574	1004	1005	1010	rVB	114324	216922	3.38%	0.354%
39	13.665	1010	1013	1015	rBV2	155533	154233	2.40%	0.252%
40	13.905	1031	1034	1041	rBV	298251	477693	7.45%	0.779%
41	14.042	1044	1046	1049	rBV	1566561	1493273	23.28%	2.435%
42	14.225	1059	1062	1064	rBV	112862	115960	1.81%	0.189%
43	14.534	1086	1089	1095	rBV	104029	164588	2.57%	0.268%
44	14.831	1112	1115	1119	rBV	59130	69324	1.08%	0.113%
45	15.162	1141	1144	1150	rBV2	74436	157034	2.45%	0.256%
46	15.482	1169	1172	1175	rBV	1062901	1374714	21.43%	2.242%
47	15.951	1210	1213	1216	rBV	62545	104860	1.63%	0.171%
48	17.048	1304	1309	1320	rBV7	43390	212656	3.32%	0.347%
49	17.471	1343	1346	1350	rBV4	72830	172928	2.70%	0.282%
50	17.848	1375	1379	1384	rVB4	49234	128529	2.00%	0.210%
51	18.020	1391	1394	1399	rVB5	39217	98894	1.54%	0.161%
52	18.271	1410	1416	1425	rBV2	177281	678976	10.59%	1.107%
53	18.523	1434	1438	1448	rVB6	62259	226125	3.53%	0.369%
54	19.151	1490	1493	1501	rVB8	31593	133881	2.09%	0.218%

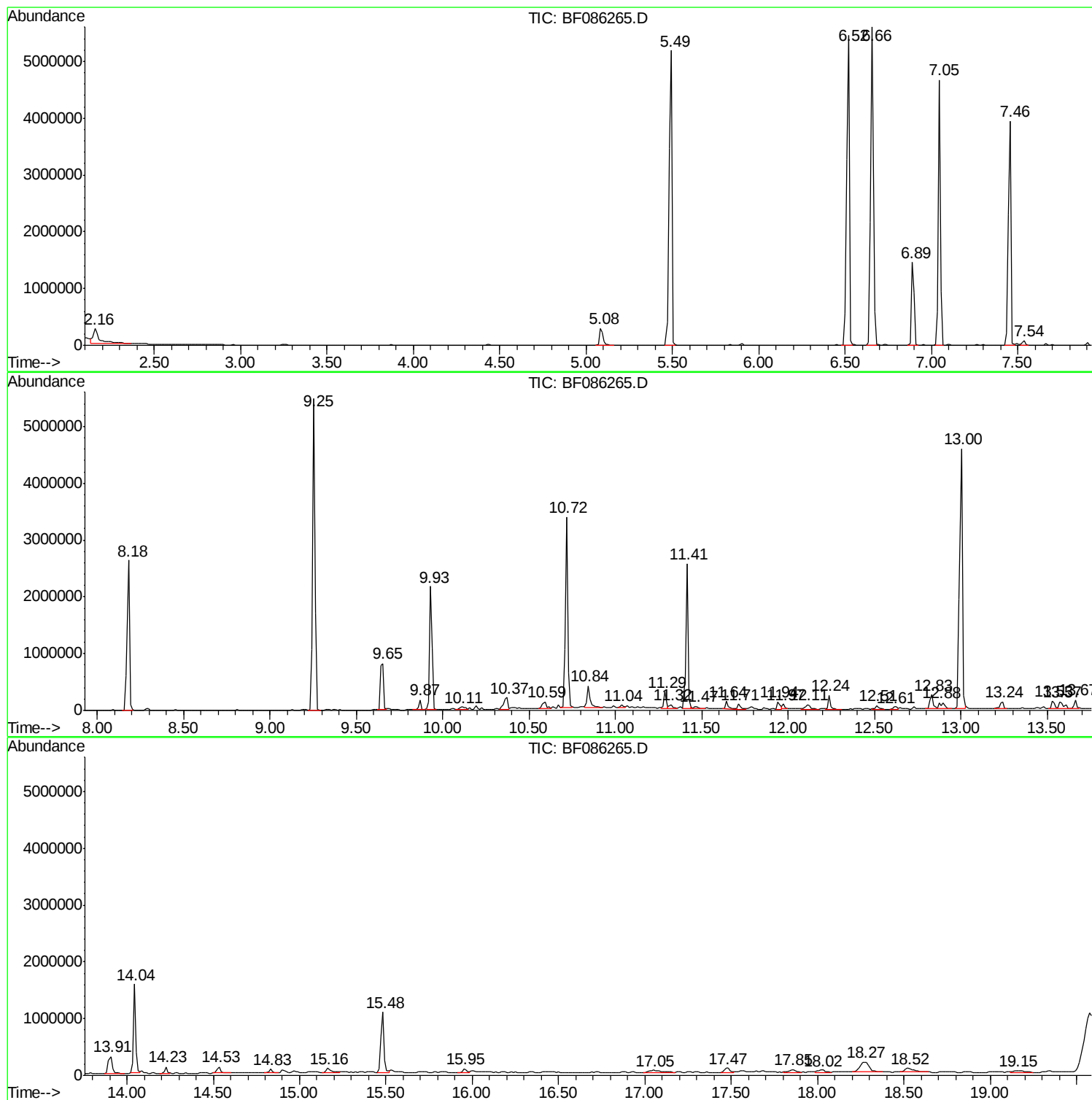
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF041516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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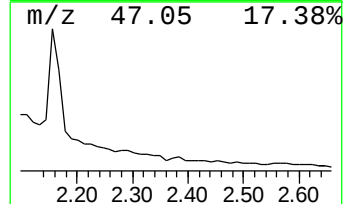
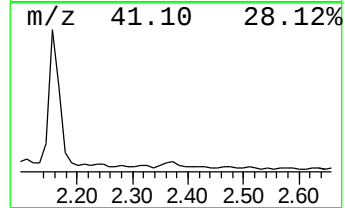
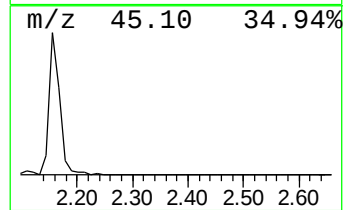
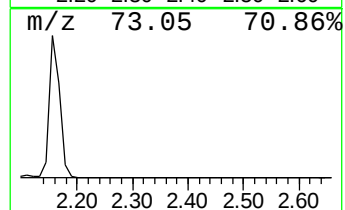
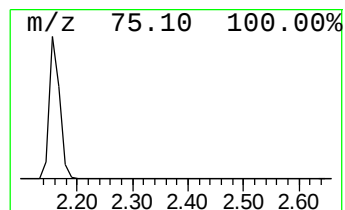
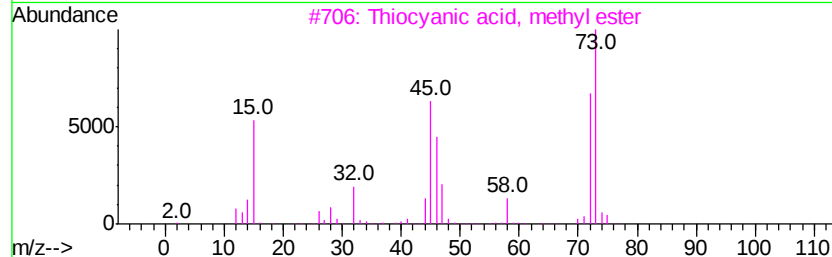
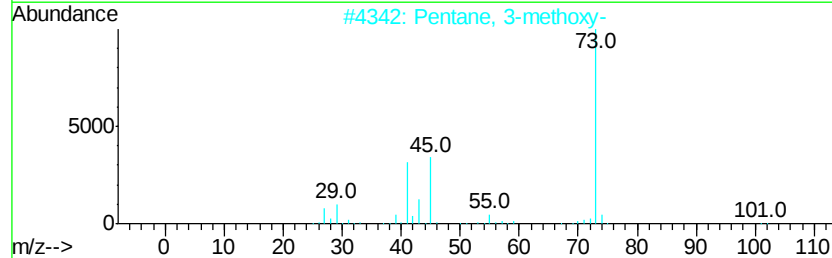
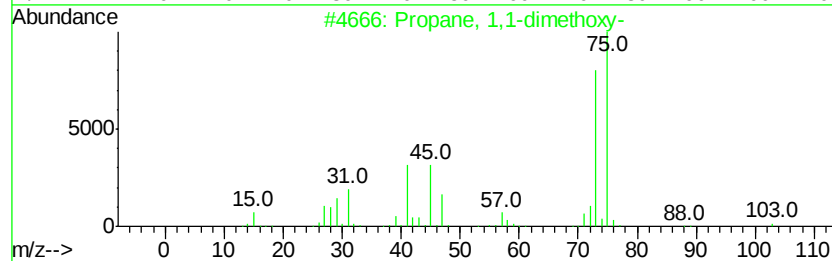
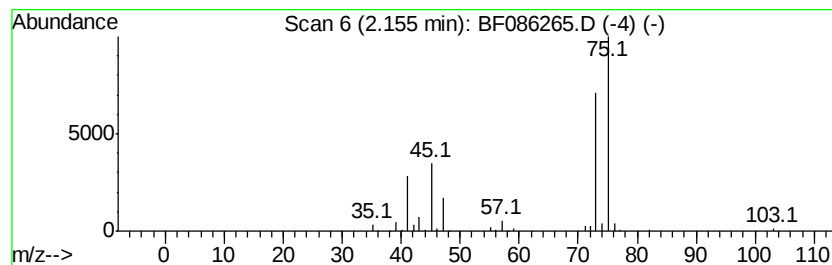
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Propane, 1,1-dimethoxy- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.16	7.20 ng	590801	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	78
2		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	10
3		Thiocyanic acid, methyl ester	73	C2H3NS	000556-64-9	9
4		Silanol, trimethyl-	90	C3H10OSi	001066-40-6	9
5		1,3-Dioxolane, 2-(chloromethyl)-	122	C4H7ClO2	002568-30-1	9



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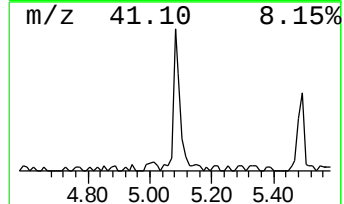
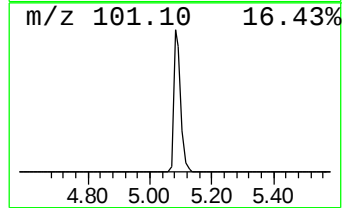
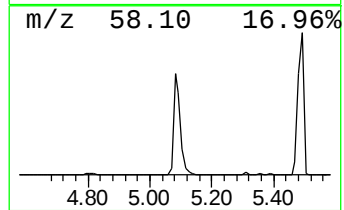
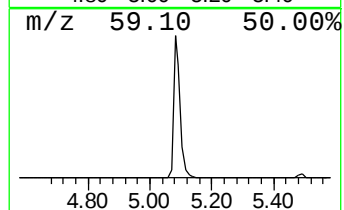
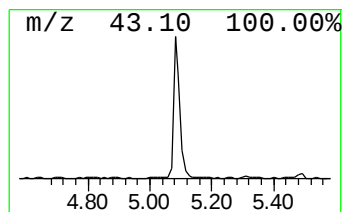
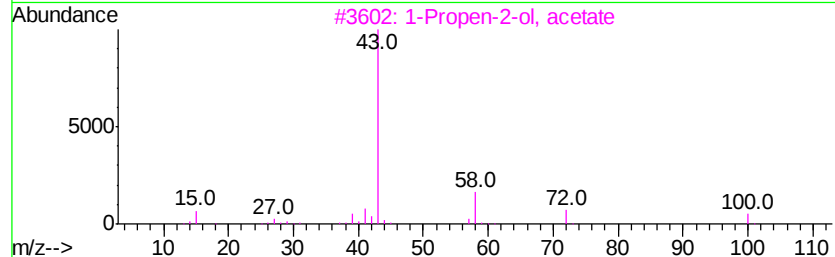
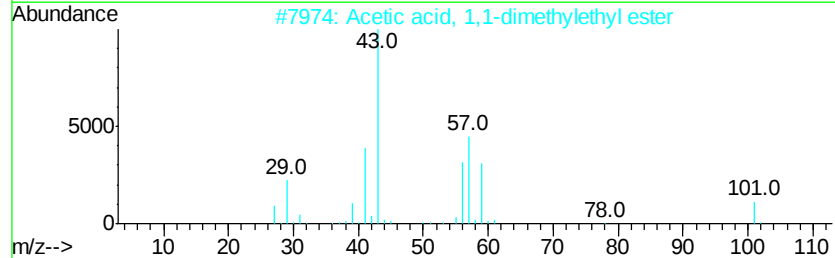
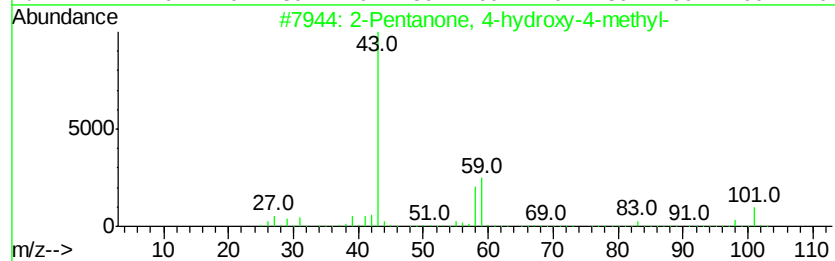
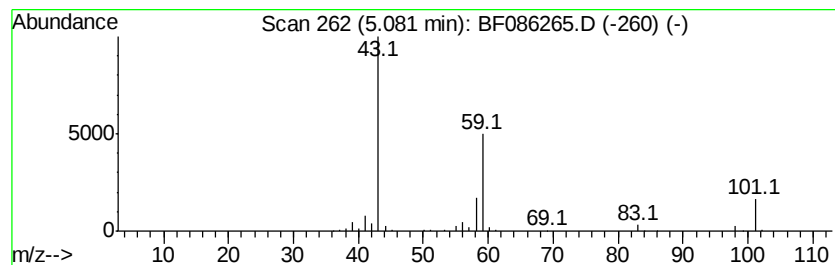
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.08	5.32 ng	436137	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10
5		Acetone	58	C3H6O	000067-64-1	9



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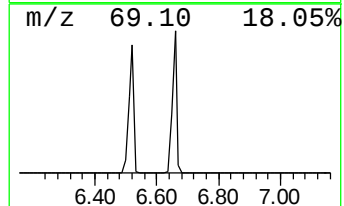
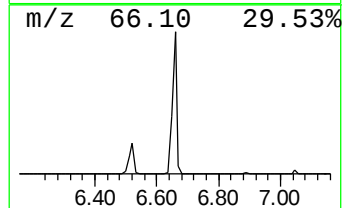
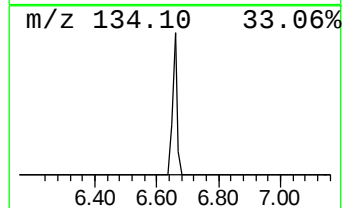
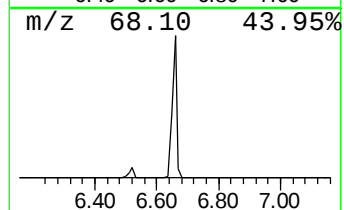
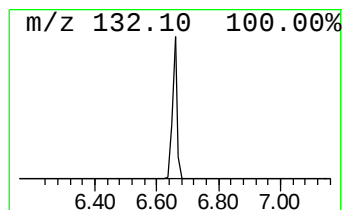
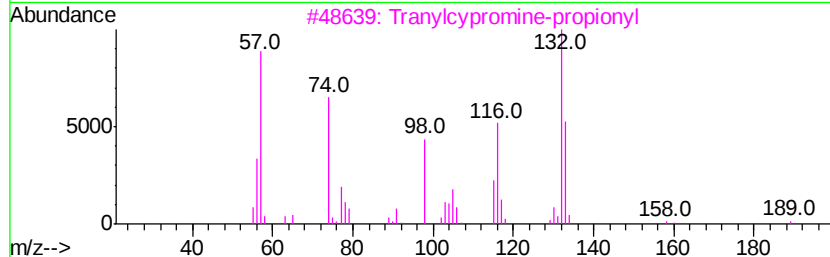
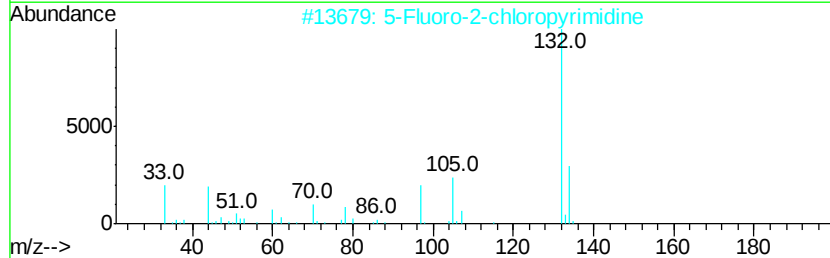
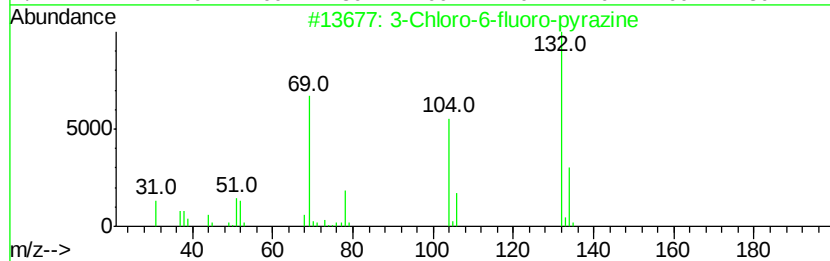
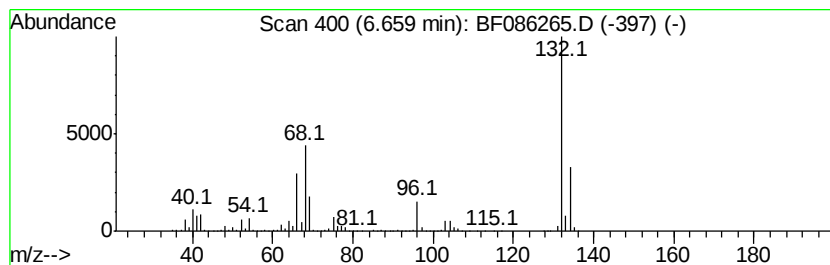
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.66 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.66	70.21 ng	5757530	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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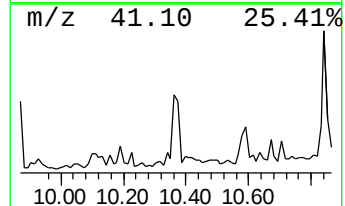
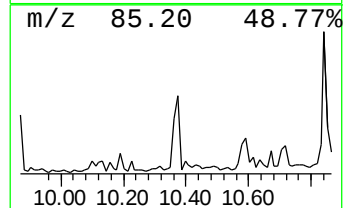
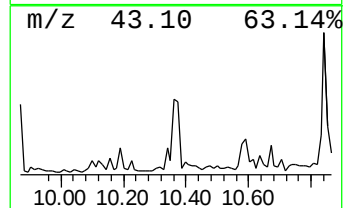
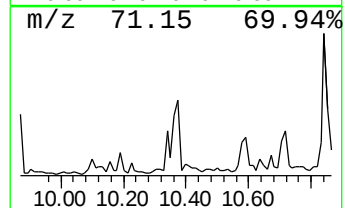
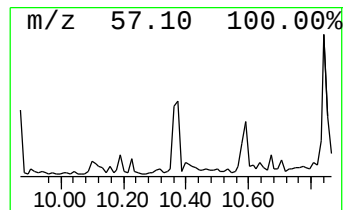
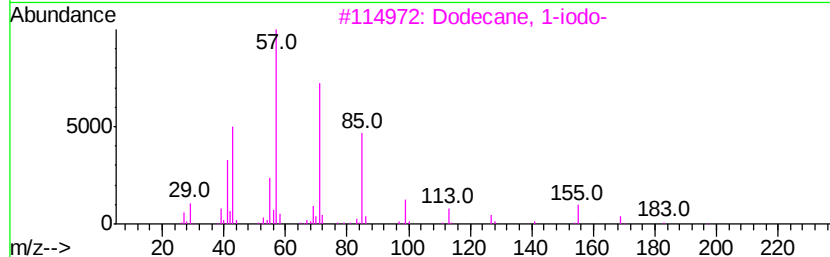
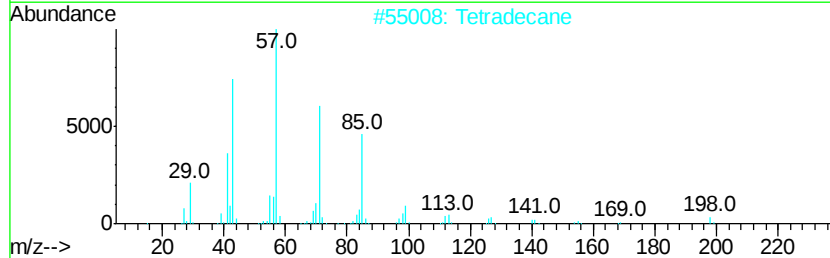
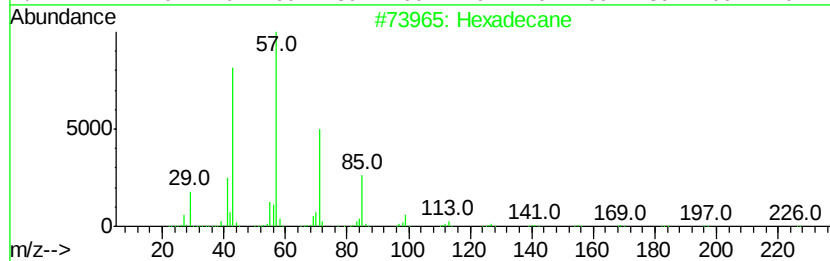
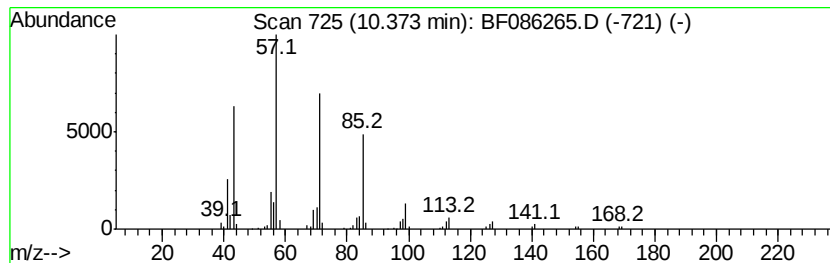
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Peak Number 5 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.37	2.94 ng	342209	Acenaphthene-d10	9.93	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	90
2	Tetradecane	198	C14H30	000629-59-4	87
3	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	86
4	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	78
5	Eicosane	282	C20H42	000112-95-8	72



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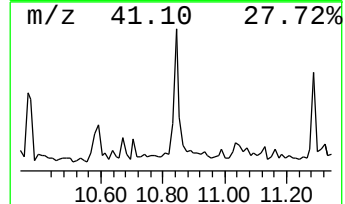
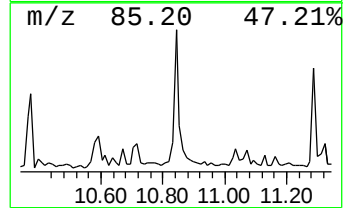
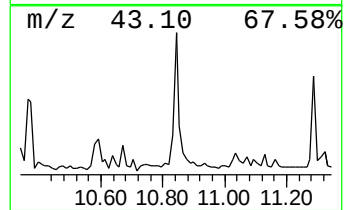
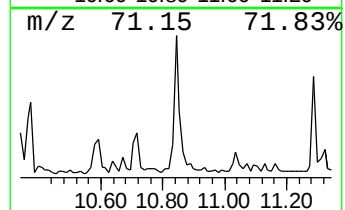
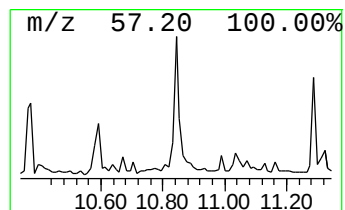
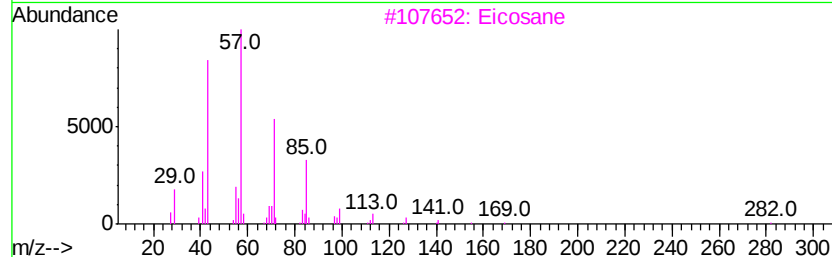
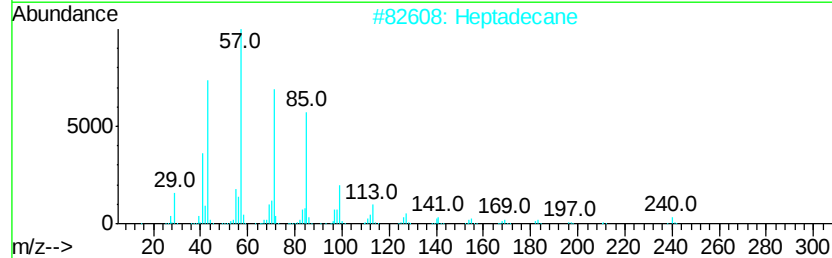
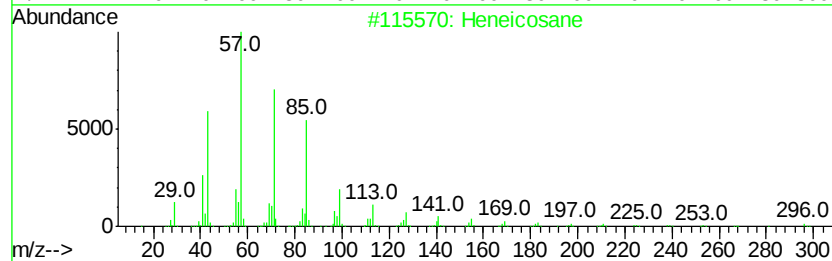
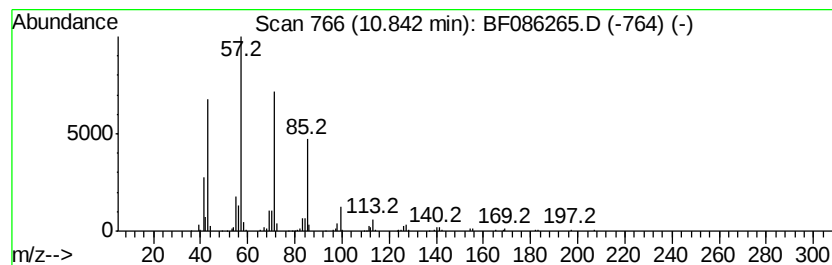
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Heneicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.84	4.06 ng	458322	Phenanthrene-d10	11.41

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Heptadecane		240	C17H36	000629-78-7	83
3	Eicosane		282	C20H42	000112-95-8	80
4	Undecane, 5-ethyl-5-propyl-		226	C16H34	002755-07-9	78
5	Nonadecane		268	C19H40	000629-92-5	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041516\
Data File : BF086265.D
Acq On : 16 Apr 2016 14:53
Operator : UM/SJ
Sample : H2471-13
Misc :
ALS Vial : 46 Sample Multiplier: 1

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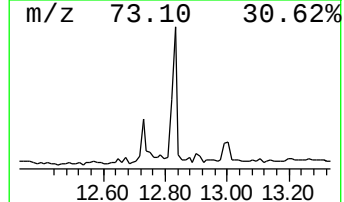
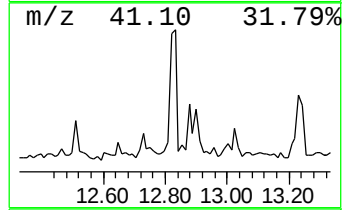
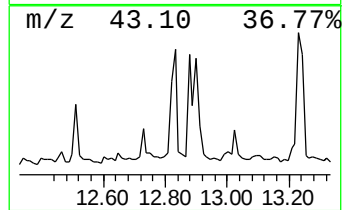
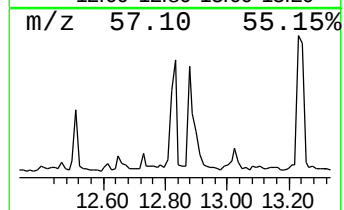
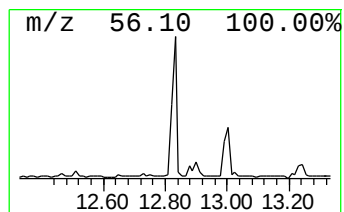
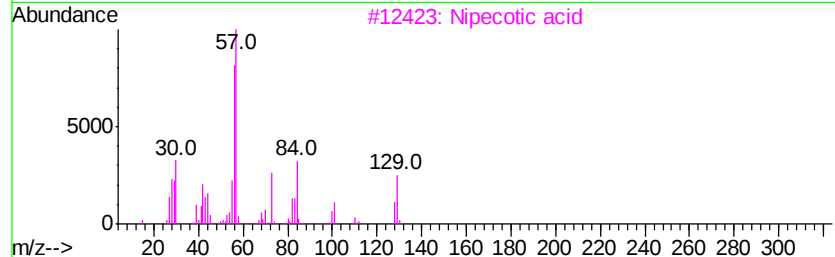
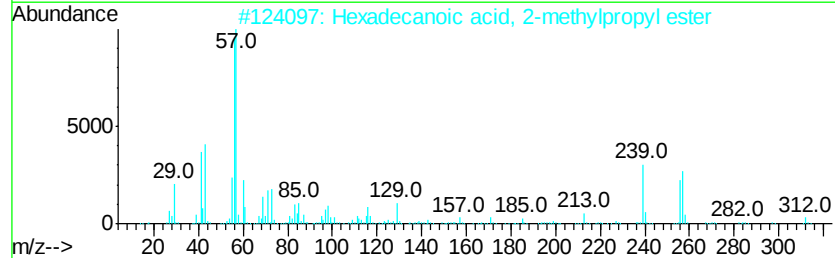
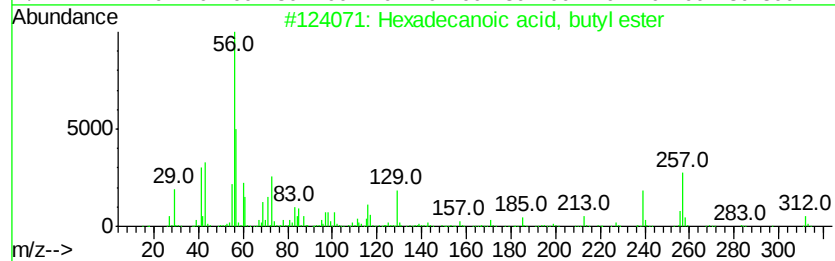
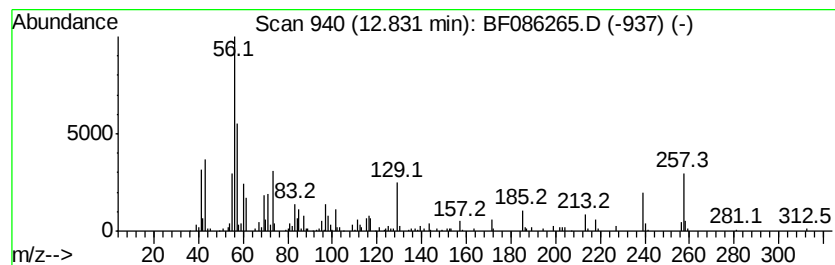
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexadecanoic acid, butyl ester Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.83	4.30 ng	321318	Chrysene-d12	14.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	91
2		Hexadecanoic acid, 2-methylpropyl...	312	C20H40O2	000110-34-9	50
3		Nipecotic acid	129	C6H11NO2	000498-95-3	47
4		2,2-Dimethyl-1,3-butanediol	118	C6H14O2	000076-35-7	35
5		Hexadecanoic acid, 1,1-dimethyl...	312	C20H40O2	031158-91-5	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041516\
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Acq On : 16 Apr 2016 14:53
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ClientSampleId :
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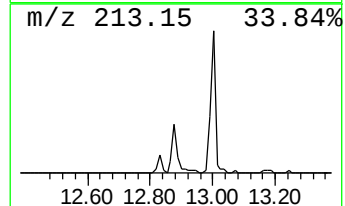
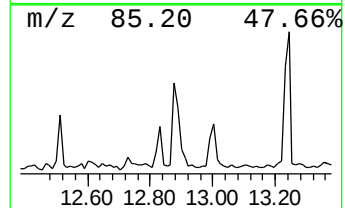
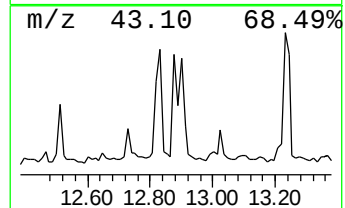
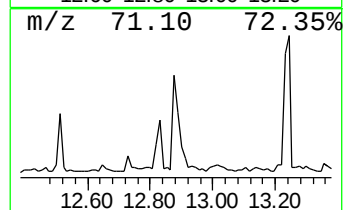
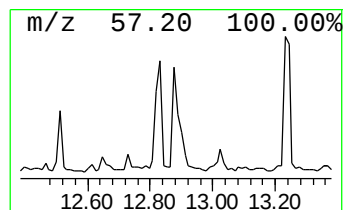
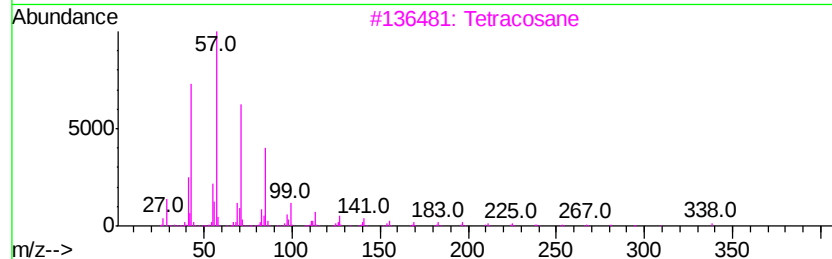
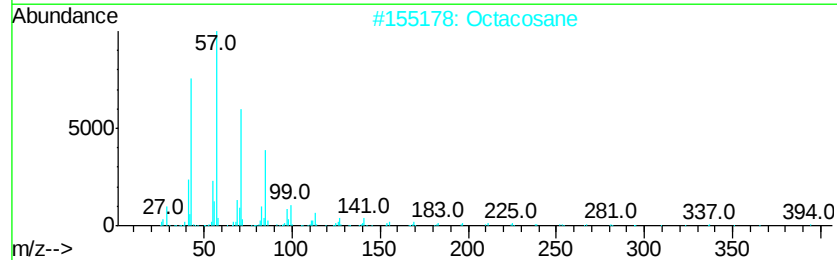
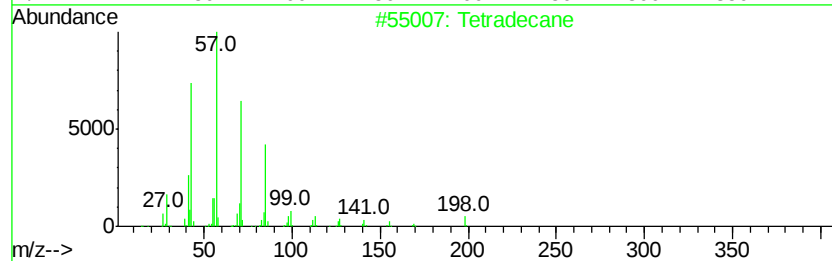
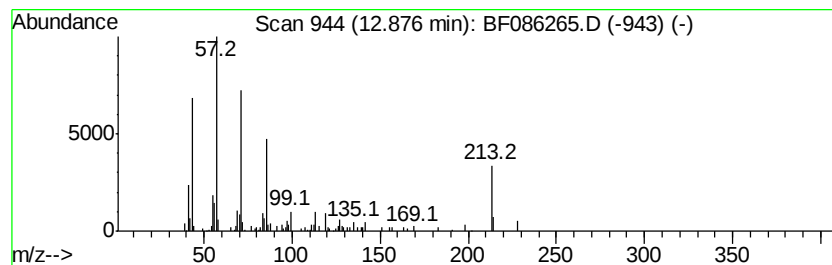
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetradecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.88	2.77 ng	206694	Chrysene-d12	14.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	76
2		Octacosane	394	C28H58	000630-02-4	62
3		Tetracosane	338	C24H50	000646-31-1	62
4		Eicosane	282	C20H42	000112-95-8	58
5		Octane, 2-methyl-	128	C9H20	003221-61-2	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041516\
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Operator : UM/SJ
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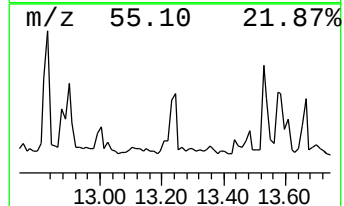
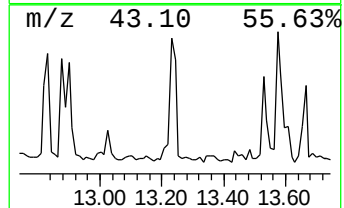
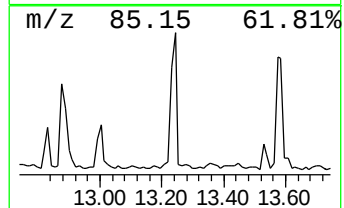
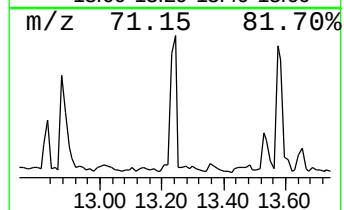
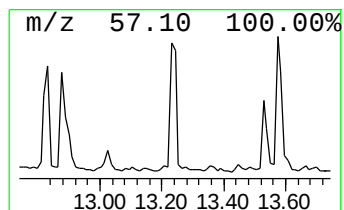
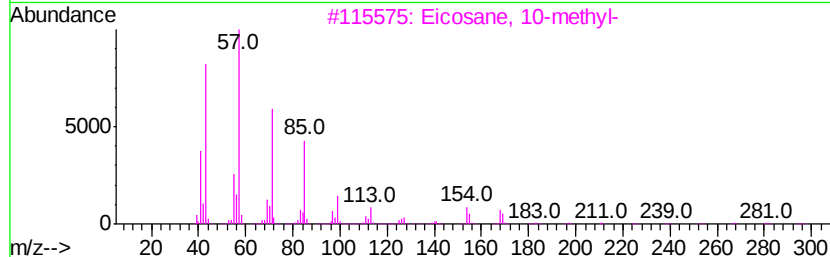
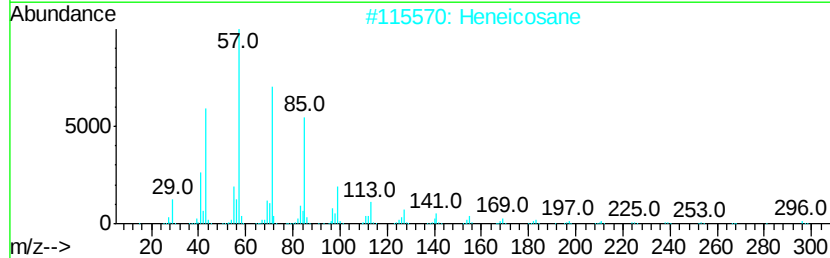
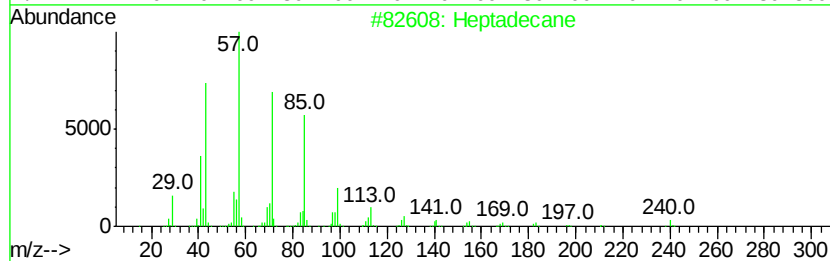
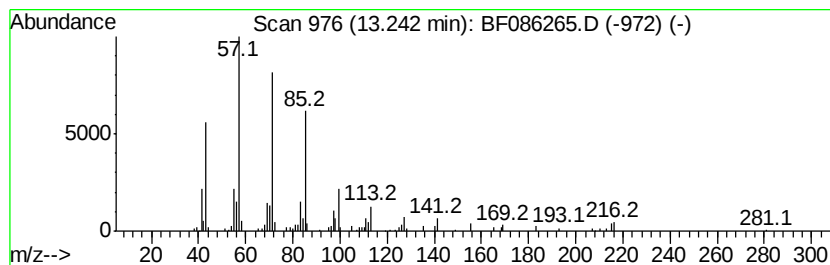
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.24	2.62 ng	195987	Chrysene-d12	14.04
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptadecane	240 C17H36	000629-78-7	90
2	Heneicosane	296 C21H44	000629-94-7	90
3	Eicosane, 10-methyl-	296 C21H44	054833-23-7	87
4	Nonacosane	408 C29H60	000630-03-5	80
5	Heptadecane, 2,6,10,15-tetramethyl-	296 C21H44	054833-48-6	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041516\
Data File : BF086265.D
Acq On : 16 Apr 2016 14:53
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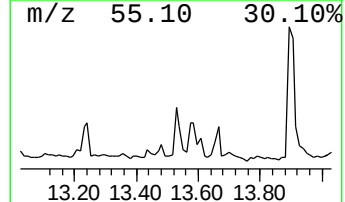
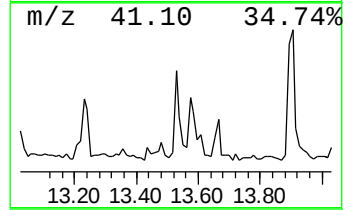
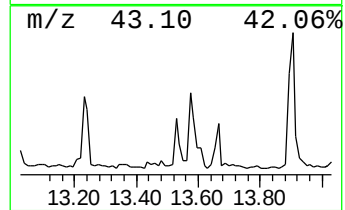
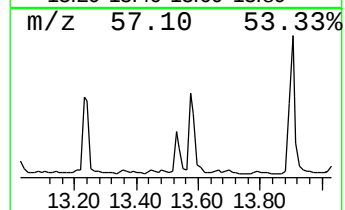
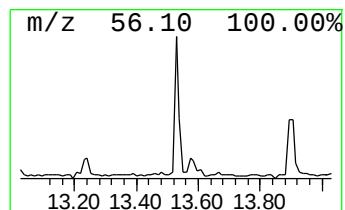
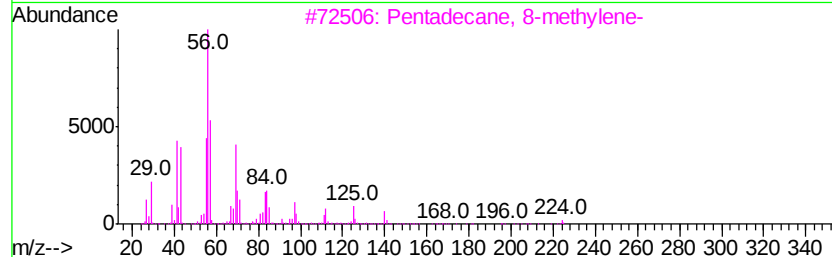
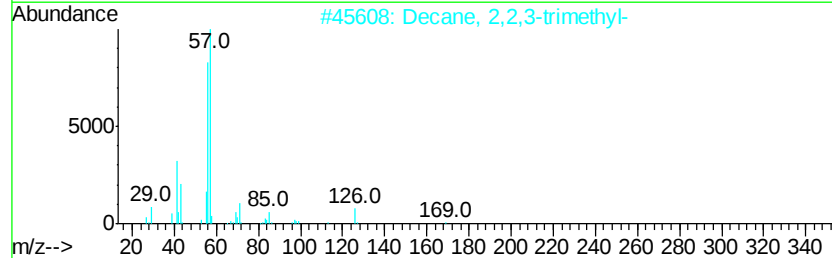
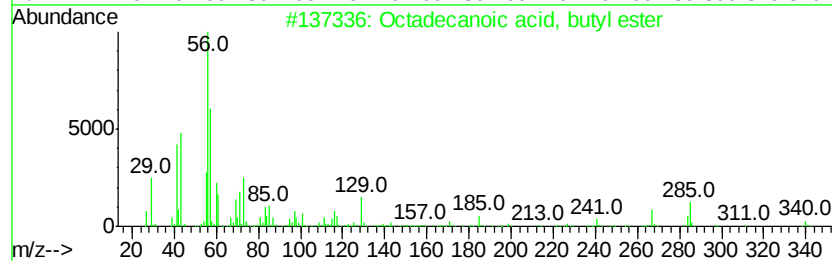
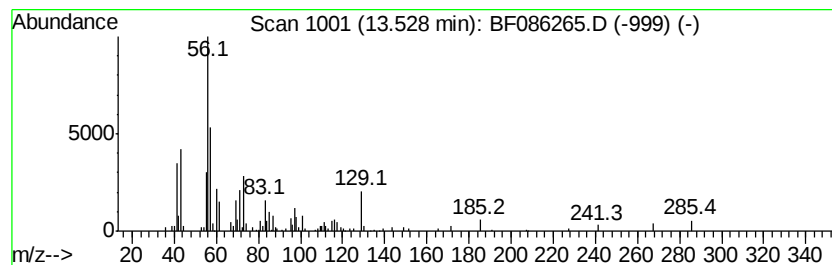
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadecanoic acid, butyl ester Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.53	2.38 ng	177473	Chrysene-d12	14.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	64
2			Decane, 2,2,3-trimethyl-	184	C13H28	062338-09-4	43
3			Pentadecane, 8-methyl-	224	C16H32	055668-09-2	37
4			2-Methyl-n-1-tridecene	196	C14H28	018094-01-4	35
5			1-Hexene, 4-methyl-	98	C7H14	003769-23-1	30



Data Path : Z:\HPCHEM1\BNA F\DATA\BF041516\
Data File : BF086265.D
Acq On : 16 Apr 2016 14:53
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Sample : H2471-13
Misc :
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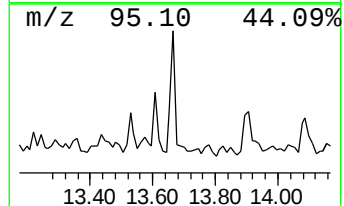
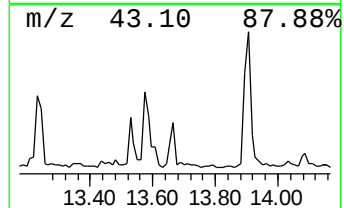
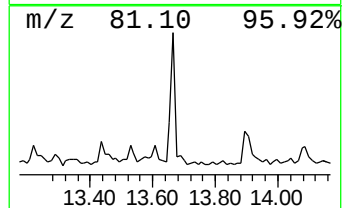
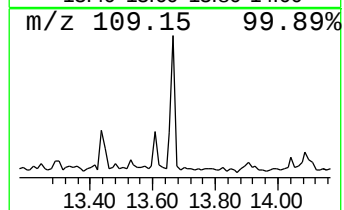
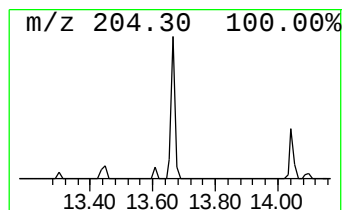
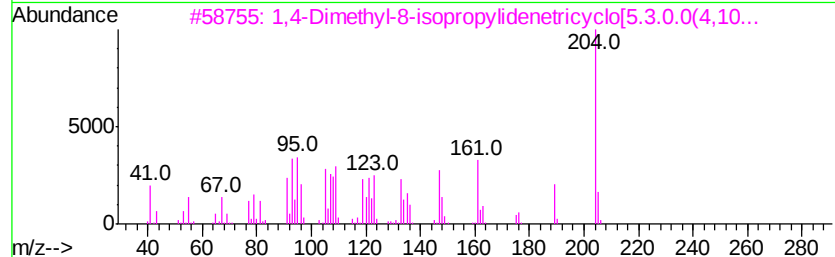
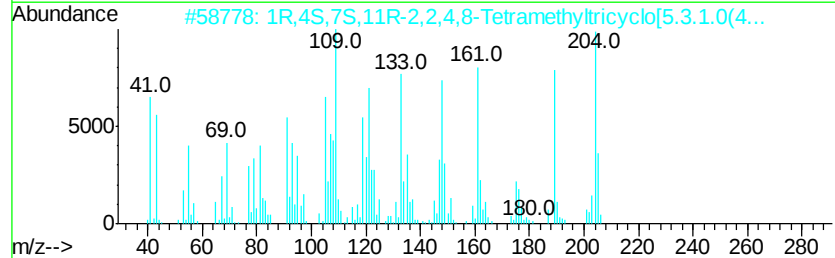
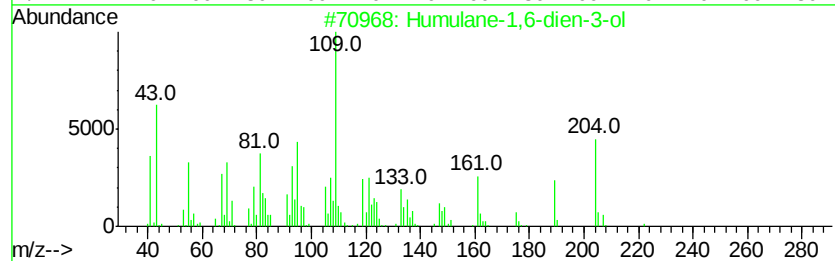
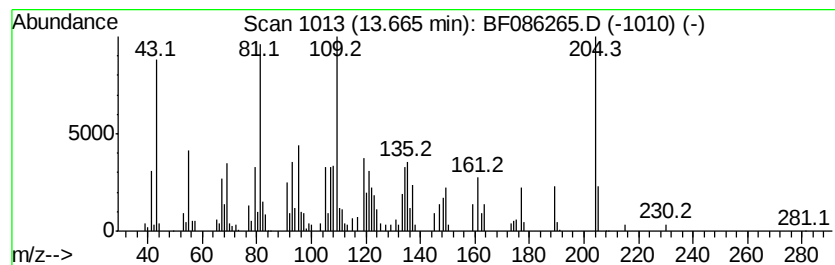
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown13.67 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.67	2.07 ng	154233	Chrysene-d12	14.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Humulane-1,6-dien-3-ol	222	C15H26O	1000140-23-1	43
2		1R,4S,7S,11R-2,2,4,8-Tetramethyl...	204	C15H24	1000140-07-6	41
3		1,4-Dimethyl-8-isopropylidenetri...	204	C15H24	1000140-07-7	38
4		1H-Cyclopropylazulene, 1a,2,3,4...	204	C15H24	000489-40-7	38
5		1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	30



Data Path : Z:\HPCHEM1\BNA F\DATA\BF041516\
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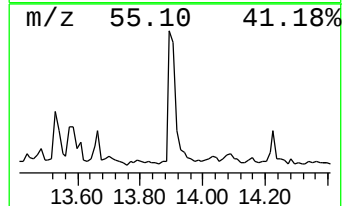
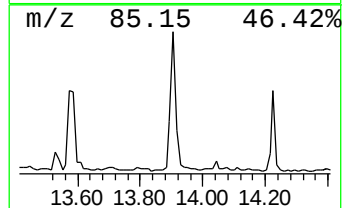
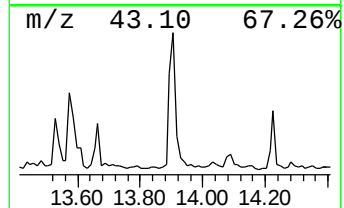
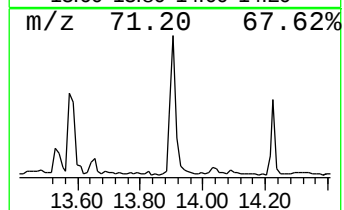
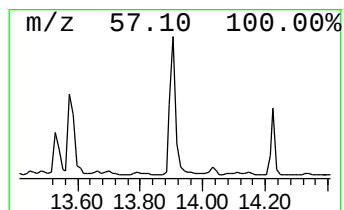
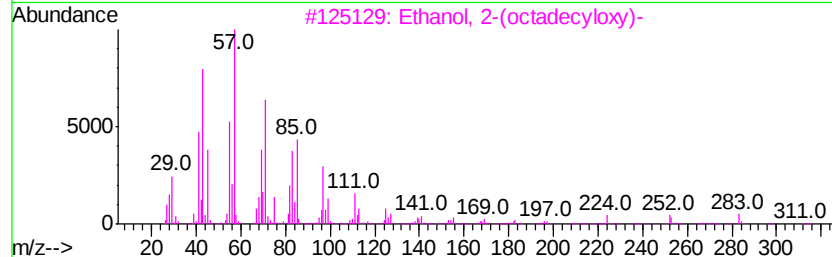
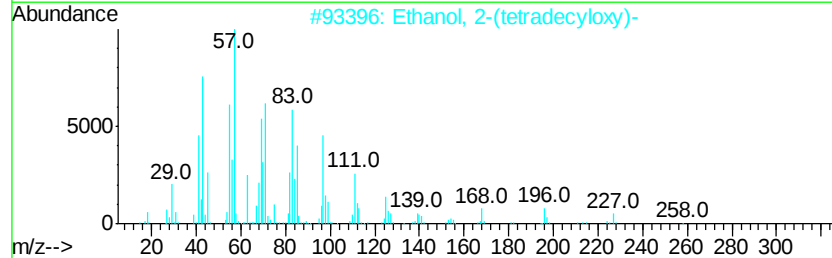
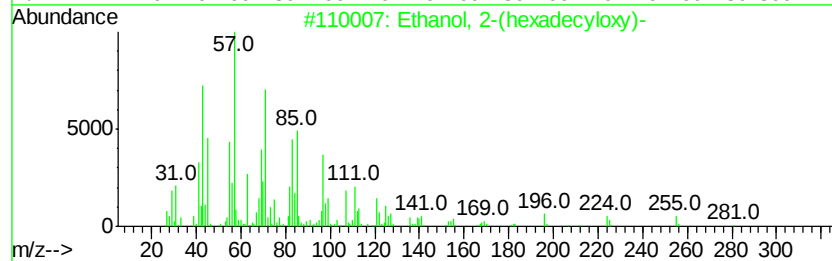
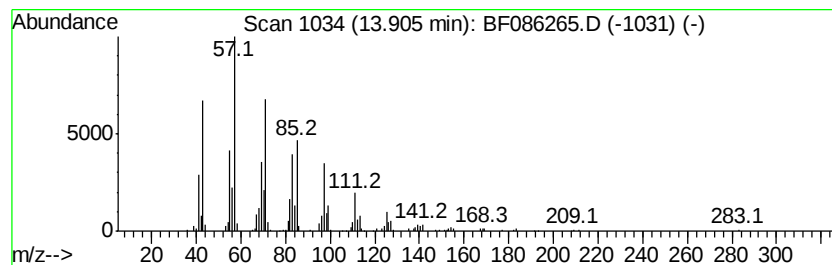
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Ethanol, 2-(hexadecyloxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.91	6.40 ng	477693	Chrysene-d12	14.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(hexadecyloxy)-	286	C18H38O2	002136-71-2	91
2		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	74
3		Ethanol, 2-(octadecyloxy)-	314	C20H42O2	002136-72-3	74
4		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	70
5		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	68



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041516\
Data File : BF086265.D
Acq On : 16 Apr 2016 14:53
Operator : UM/SJ
Sample : H2471-13
Misc :
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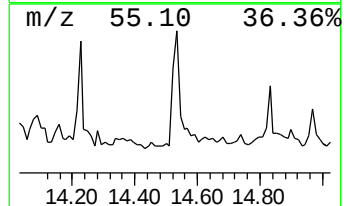
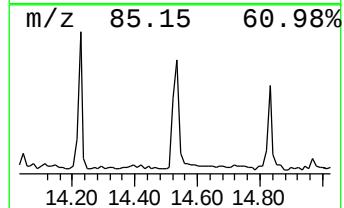
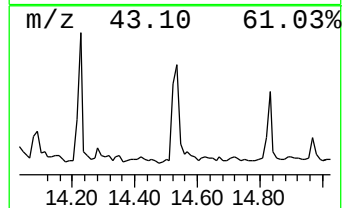
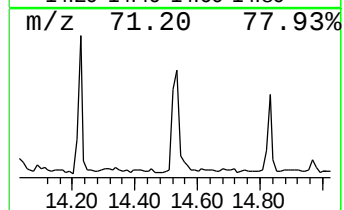
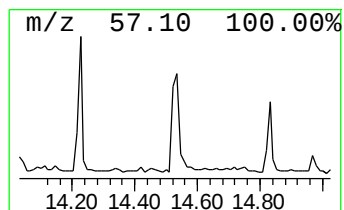
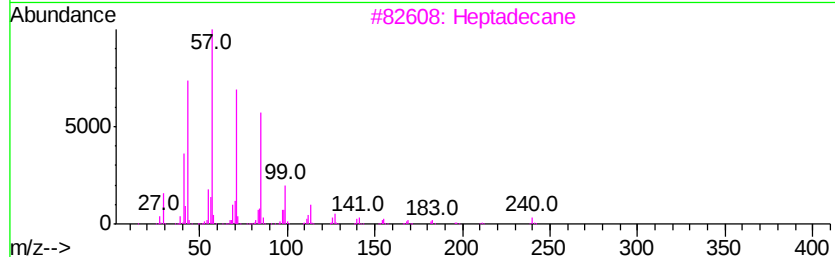
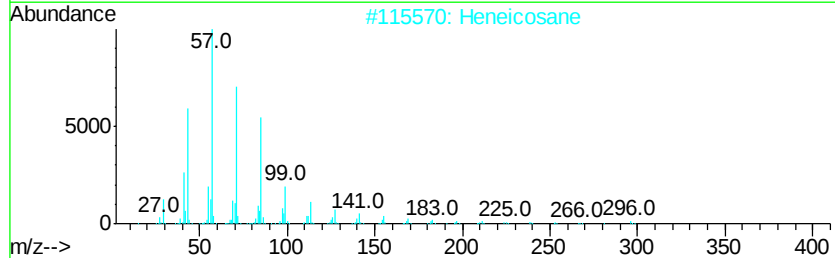
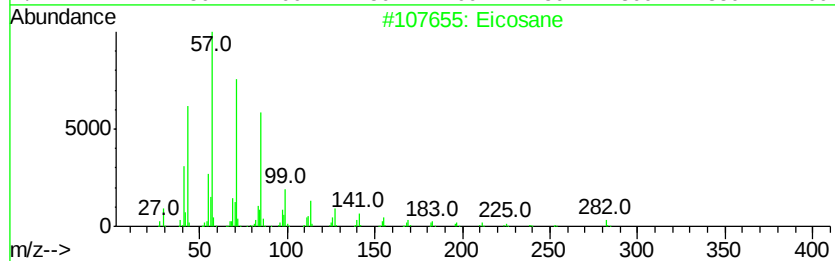
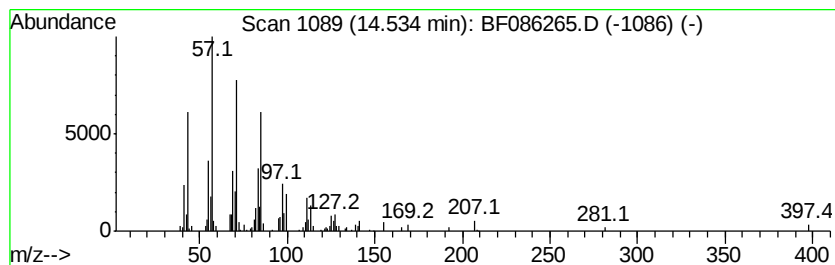
Instrument :
BNA_F
ClientSampleId :
CDMSHDUP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF041516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Eicosane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.53	2.20 ng	164588	Chrysene-d12	14.04
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane	282 C20H42	000112-95-8	74
2	Heneicosane	296 C21H44	000629-94-7	72
3	Heptadecane	240 C17H36	000629-78-7	68
4	Hexadecane	226 C16H34	000544-76-3	68
5	Octadecane	254 C18H38	000593-45-3	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF041516\
Data File : BF086265.D
Acq On : 16 Apr 2016 14:53
Operator : UM/SJ
Sample : H2471-13
Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CDMSHDUP

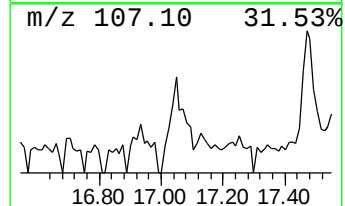
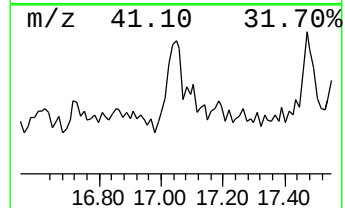
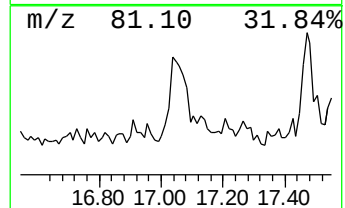
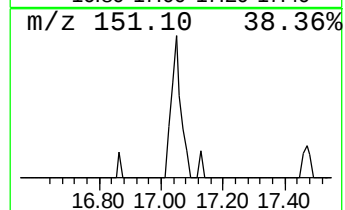
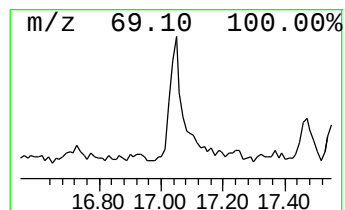
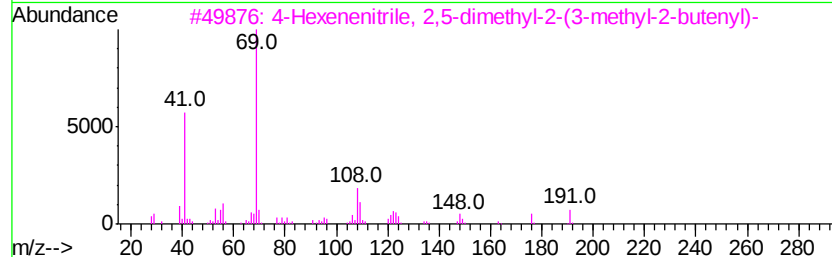
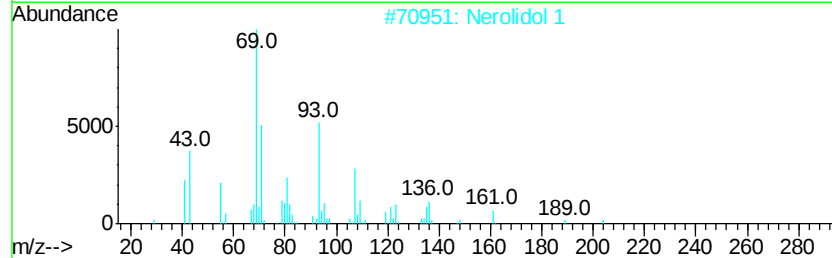
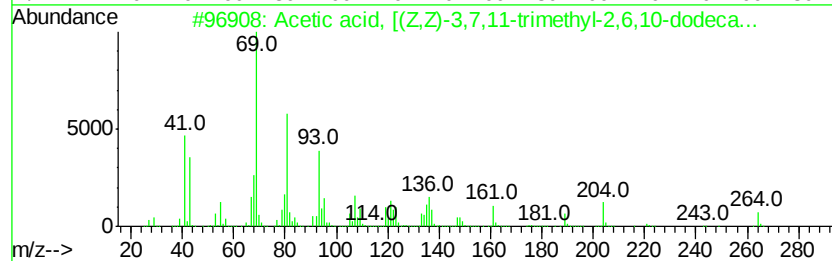
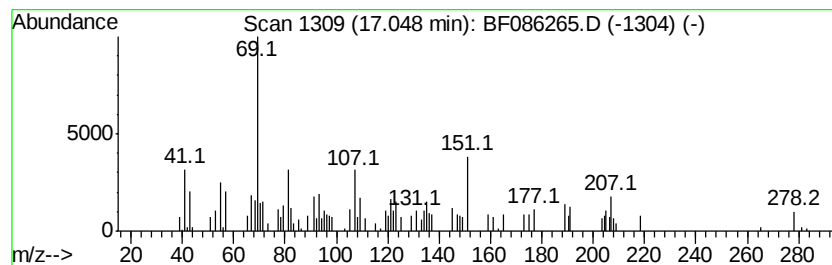
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown17.05 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.05	3.09 ng	212656	Perylene-d12	15.48

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetic acid, [(Z,Z)-3,7,11-trime...	264	C17H28O2	1000164-38-6	38
2			Nerolidol 1	222	C15H26O	1000285-43-5	38
3			4-Hexenenitrile, 2,5-dimethyl-2-...	191	C13H21N	069340-56-3	35
4			2,6,10-Dodecatrien-1-ol, 3,7,11-...	264	C17H28O2	004128-17-0	35
5			Crotonic acid, o-formylphenyl ester	190	C11H10O3	114636-69-0	30



Data Path : Z:\HPCHEM1\BNA F\DATA\BF041516\
Data File : BF086265.D
Acq On : 16 Apr 2016 14:53
Operator : UM/SJ
Sample : H2471-13
Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CDMSHDUP

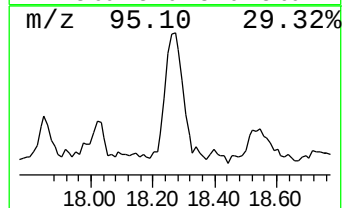
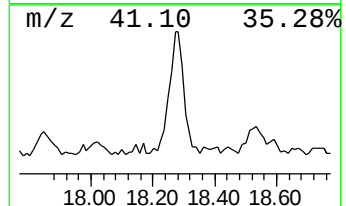
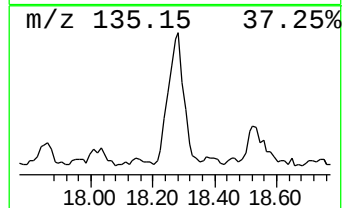
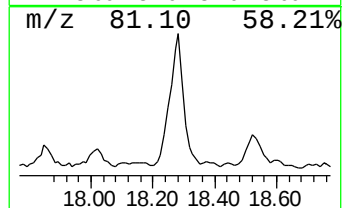
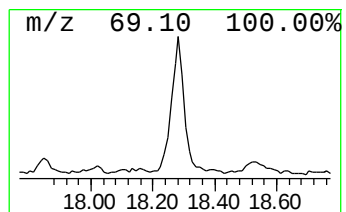
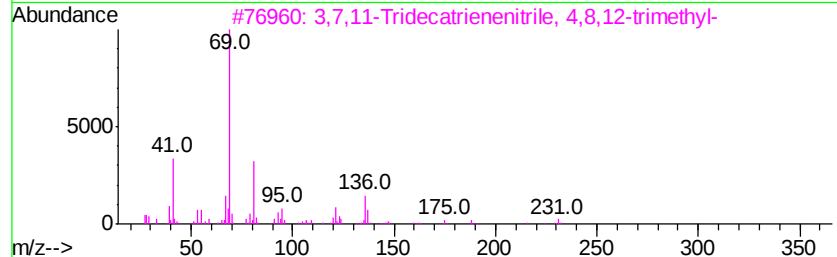
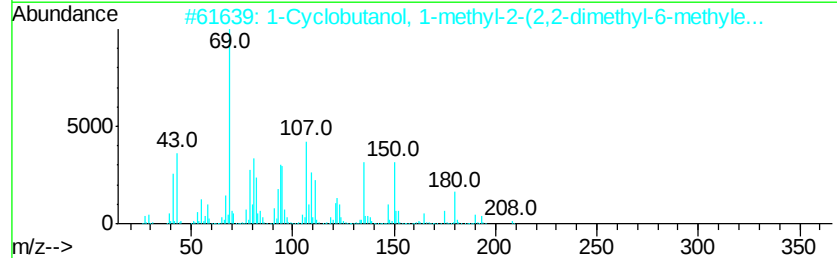
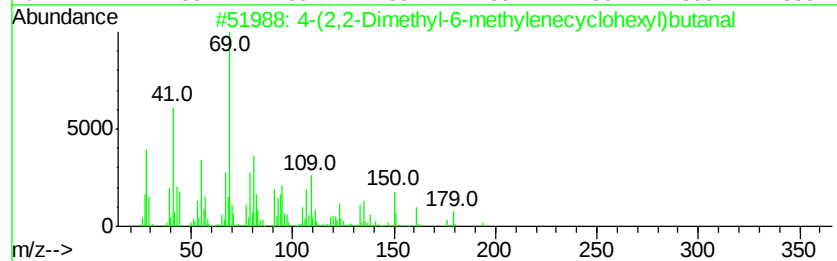
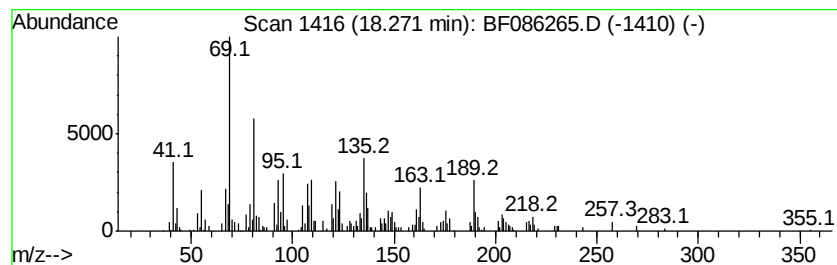
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 4-(2,2-Dimethyl-6-methylene... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.27	9.88 ng	678976	Perylene-d12	15.48

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-(2,2-Dimethyl-6-methylenecyclo...	194	C13H22O	095452-13-4	52
2			1-Cyclobutanol, 1-methyl-2-(2,2-...	208	C14H24O	1000197-21-8	47
3			3,7,11-Tridecatrienitrile, 4,8...	231	C16H25N	006006-01-5	38
4			6,10-Dodecadien-1-yn-3-ol, 3,7,1...	220	C15H24O	002387-68-0	35
5			1-Naphthalenepropanol, .alpha.-e...	306	C20H34O2	004549-12-6	30



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041516\
 Data File : BF086265.D
 Acq On : 16 Apr 2016 14:53
 Operator : UM/SJ
 Sample : H2471-13
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CDMSHDUP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF041516.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
Propane, 1,1-dime...	2.16	7.2	ng	590801	1	6.89	1640090	20.0
2-Pentanone, 4-hy...	5.08	5.3	ng	436137	1	6.89	1640090	20.0
unknown6.66	6.66	70.2	ng	5757530	1	6.89	1640090	20.0
Hexadecane	10.37	2.9	ng	342209	3	9.93	2327790	20.0
Heneicosane	10.84	4.1	ng	458322	4	11.41	2256040	20.0
Hexadecanoic acid...	12.83	4.3	ng	321318	5	14.04	1493270	20.0
Tetradecane	12.88	2.8	ng	206694	5	14.04	1493270	20.0
Heptadecane	13.24	2.6	ng	195987	5	14.04	1493270	20.0
Octadecanoic acid...	13.53	2.4	ng	177473	5	14.04	1493270	20.0
unknown13.67	13.67	2.1	ng	154233	5	14.04	1493270	20.0
Ethanol, 2-(hexad...	13.91	6.4	ng	477693	5	14.04	1493270	20.0
Eicosane	14.53	2.2	ng	164588	5	14.04	1493270	20.0
unknown17.05	17.05	3.1	ng	212656	6	15.48	1374710	20.0
4-(2,2-Dimethyl-6...	18.27	9.9	ng	678976	6	15.48	1374710	20.0