

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020616\
 Data File : BF084916.D
 Acq On : 6 Feb 2016 17:16
 Operator : SJ/IZ
 Sample : H1338-02
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-2

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-BF020116.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	4.087	172	175	180	rBV	37318	61608	1.48%	0.189%
2	4.796	235	237	243	rBV	323947	484098	11.60%	1.488%
3	5.241	274	276	279	rBV	1747335	1735234	41.58%	5.334%
4	5.653	310	312	315	rBV	49910	47960	1.15%	0.147%
5	6.304	367	369	372	rBV	1398419	1175774	28.17%	3.614%
6	6.430	377	380	382	rBV	3417917	3361434	80.55%	10.333%
7	6.659	398	400	402	rBV	1170353	954761	22.88%	2.935%
8	6.819	411	414	416	rBV	3330299	3354737	80.38%	10.313%
9	7.230	447	450	452	rBV	2732052	2527966	60.57%	7.771%
10	7.950	510	513	515	rBV	1056165	1212752	29.06%	3.728%
11	8.476	556	559	563	rBV3	18868	61871	1.48%	0.190%
12	8.579	564	568	570	rBV2	74322	136584	3.27%	0.420%
13	8.762	580	584	586	rBV4	22282	46222	1.11%	0.142%
14	9.025	604	607	610	rBV	4285196	4173355	100.00%	12.829%
15	9.356	634	636	640	rBV4	25808	66843	1.60%	0.205%
16	9.425	640	642	644	rVV	407124	357230	8.56%	1.098%
17	9.470	644	646	649	rVB3	29697	61594	1.48%	0.189%
18	9.642	659	661	663	rBV	129272	106449	2.55%	0.327%
19	9.699	663	666	668	rBV	1124969	1306042	31.29%	4.015%
20	9.905	677	684	685	rBV5	45730	126451	3.03%	0.389%
21	9.939	685	687	688	rVV2	71455	88898	2.13%	0.273%
22	10.008	691	693	698	rVB4	40085	77278	1.85%	0.238%
23	10.088	698	700	702	rBV2	36523	60393	1.45%	0.186%
24	10.145	702	705	706	rVB2	141839	241832	5.79%	0.743%
25	10.362	722	724	725	rBV	64193	79574	1.91%	0.245%
26	10.488	732	735	737	rBV2	1890790	2477803	59.37%	7.617%
27	10.613	744	746	750	rVB	198617	231491	5.55%	0.712%
28	10.819	762	764	765	rBV2	45943	61493	1.47%	0.189%
29	10.865	765	768	769	rVV3	27455	51891	1.24%	0.160%
30	10.933	772	774	776	rVB2	41397	52448	1.26%	0.161%
31	11.059	783	785	786	rBV	88266	79468	1.90%	0.244%
32	11.093	786	788	791	rVB	33121	48702	1.17%	0.150%
33	11.173	792	795	797	rBV	1414476	1199522	28.74%	3.687%
34	11.288	803	805	811	rVB7	24447	58773	1.41%	0.181%

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

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35	11.722	842	843	846	rVV	129933	125726	3.01%	0.386%
36	11.779	846	848	853	rVB3	41362	89525	2.15%	0.275%
37	12.225	884	887	888	rBV	32971	46150	1.11%	0.142%
38	12.602	918	920	922	rBV	114869	136082	3.26%	0.418%
39	12.671	925	926	928	rVB	49110	42815	1.03%	0.132%
40	12.762	931	934	936	rBV	3363187	3121594	74.80%	9.596%
41	13.299	979	981	983	rBV	96804	88693	2.13%	0.273%
42	13.345	983	985	986	rBV2	46365	43187	1.03%	0.133%
43	13.677	1011	1014	1018	rBV	240512	362295	8.68%	1.114%
44	13.802	1023	1025	1027	rBV	1025490	913348	21.89%	2.808%
45	13.985	1039	1041	1049	rBV	51670	91628	2.20%	0.282%
46	14.294	1066	1068	1075	rBV	75368	118237	2.83%	0.363%
47	14.591	1091	1094	1096	rBV	77730	93577	2.24%	0.288%
48	14.888	1117	1120	1122	rVB	60830	65019	1.56%	0.200%
49	15.174	1142	1145	1147	rBV	727983	829269	19.87%	2.549%
50	15.220	1147	1149	1152	rVB	49065	70634	1.69%	0.217%
51	15.597	1179	1182	1184	rBV	36592	58017	1.39%	0.178%
52	16.020	1217	1219	1225	rVB2	25830	65630	1.57%	0.202%

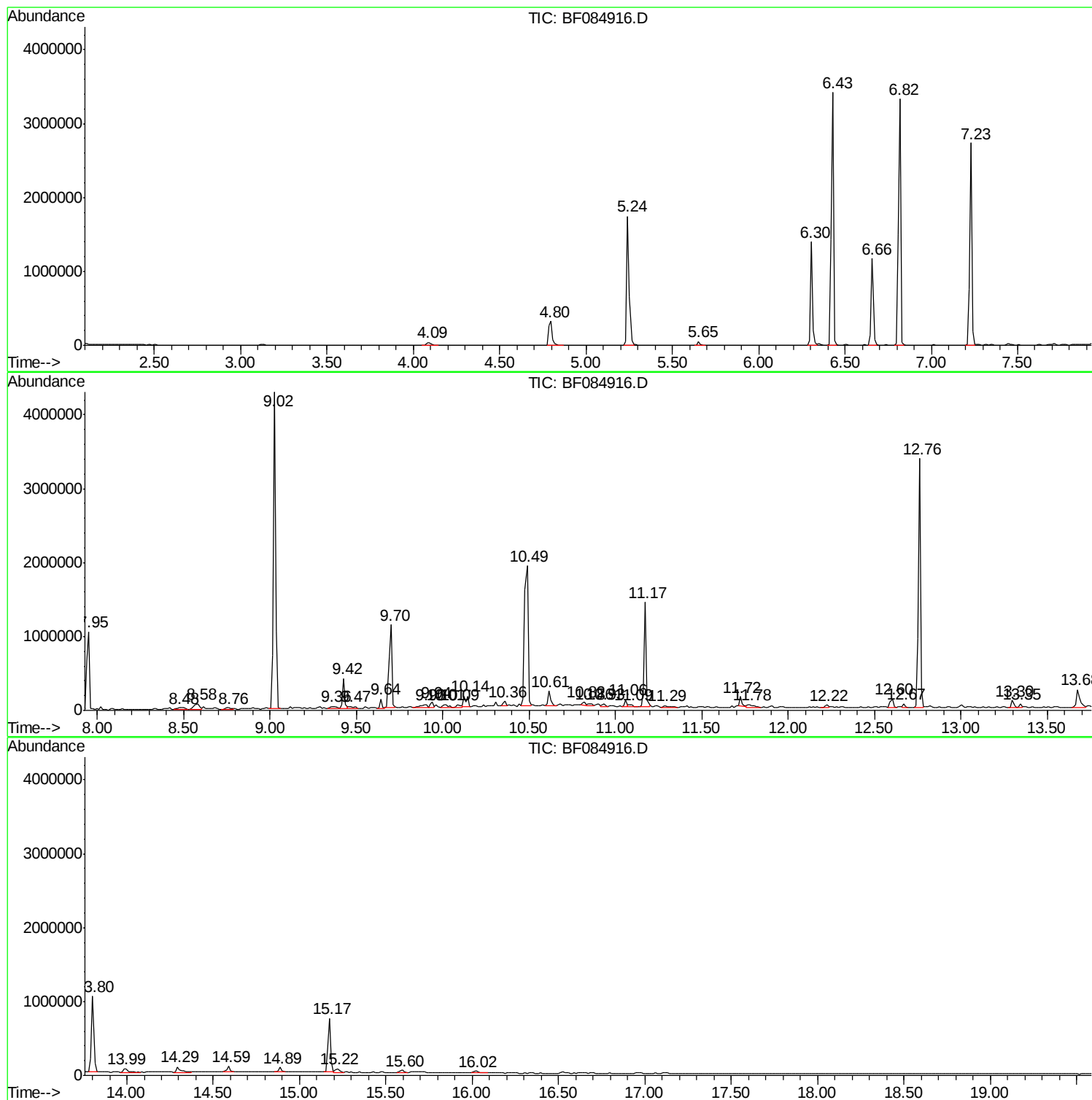
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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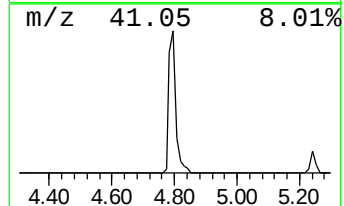
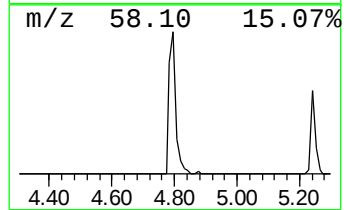
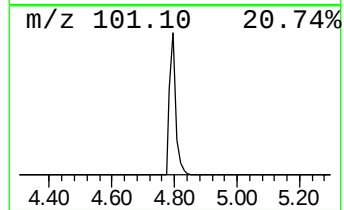
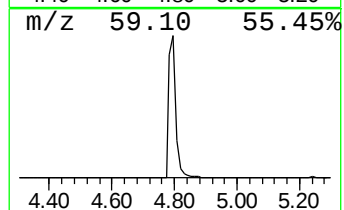
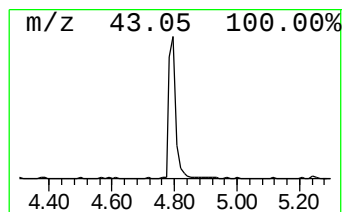
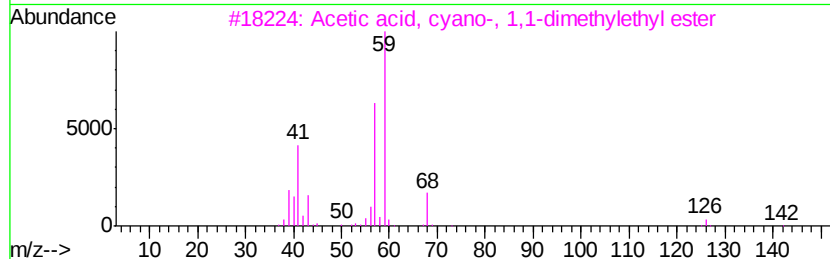
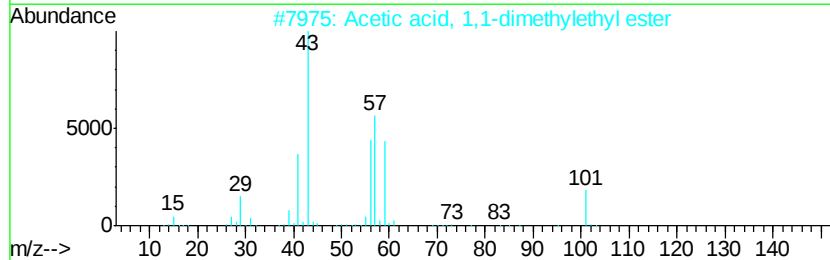
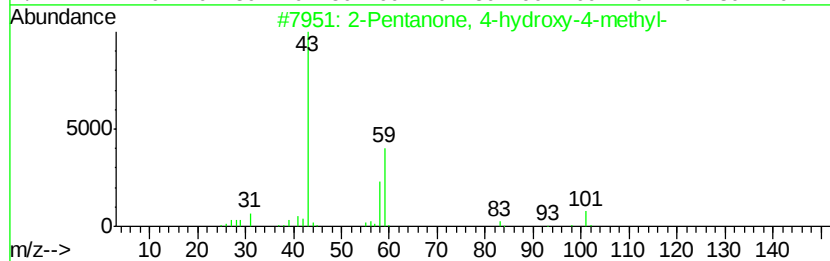
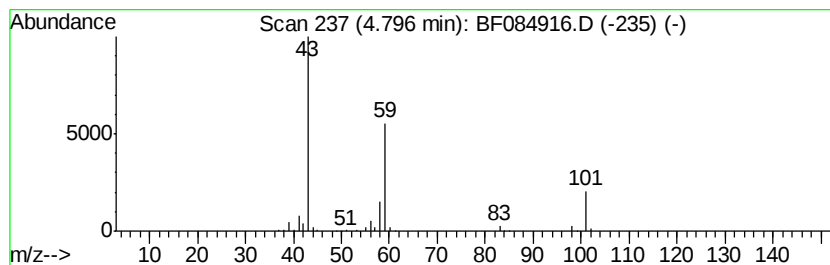
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.80	10.14 ng	484098	1,4-Dichlorobenzene-d4	6.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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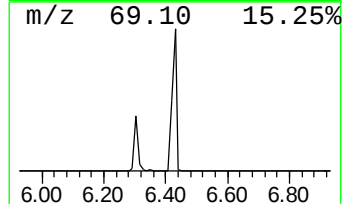
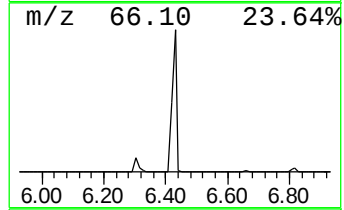
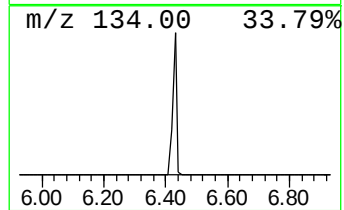
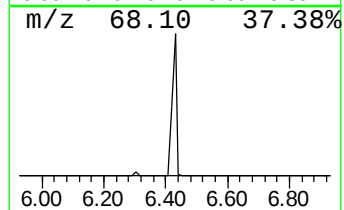
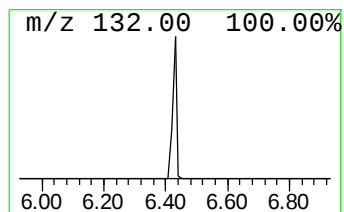
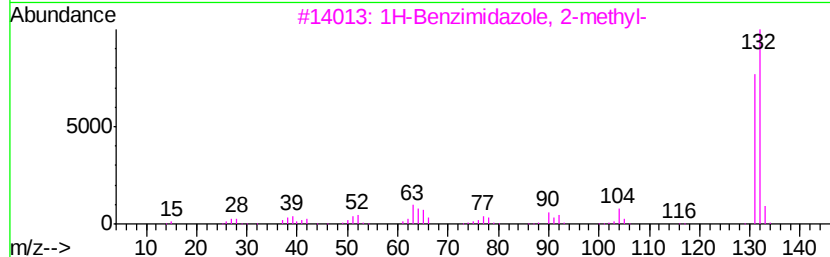
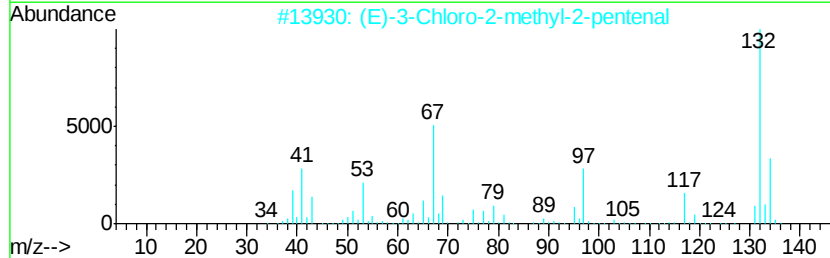
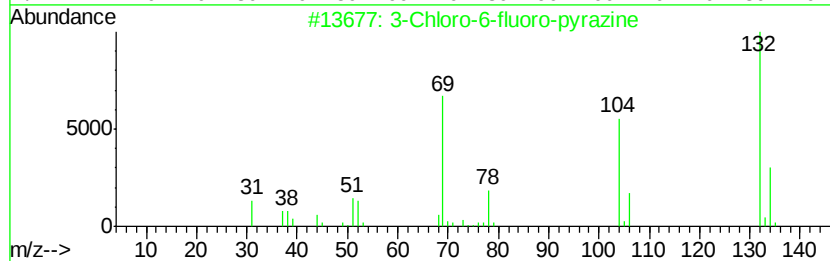
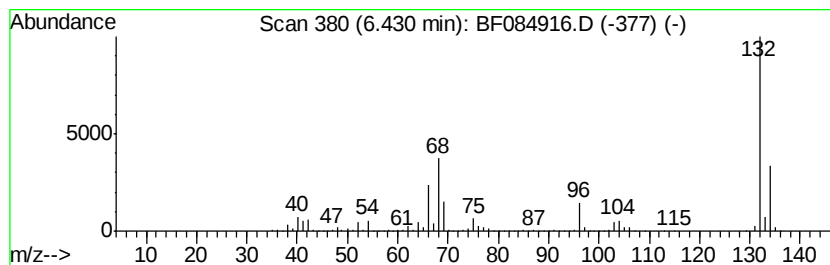
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.43 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.43	70.41 ng	3361430	1,4-Dichlorobenzene-d4	6.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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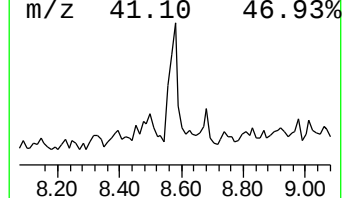
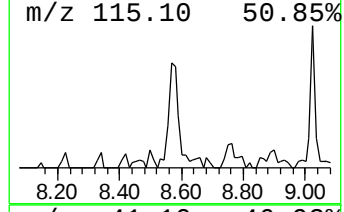
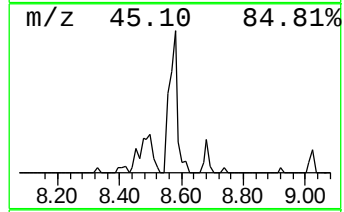
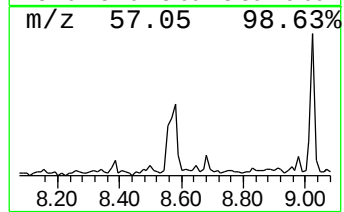
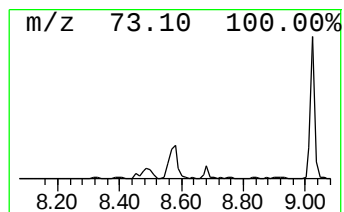
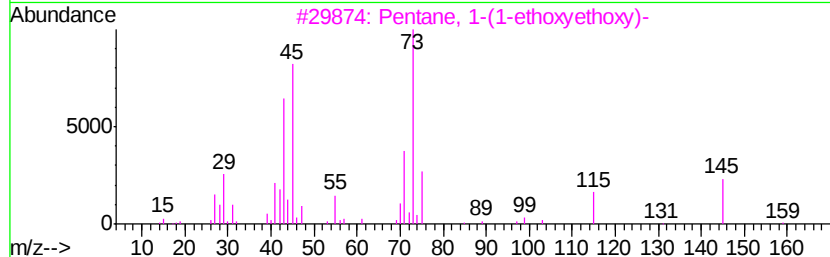
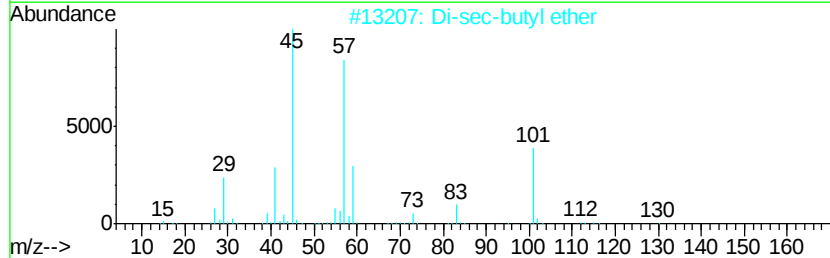
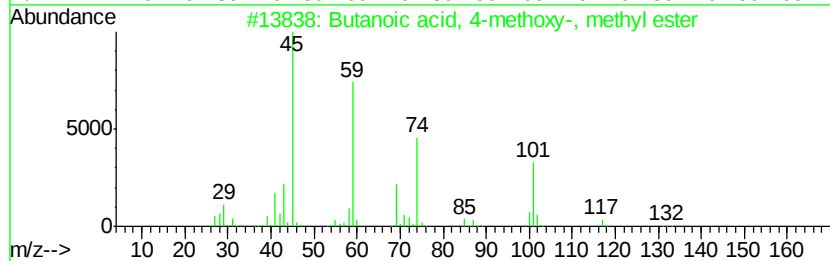
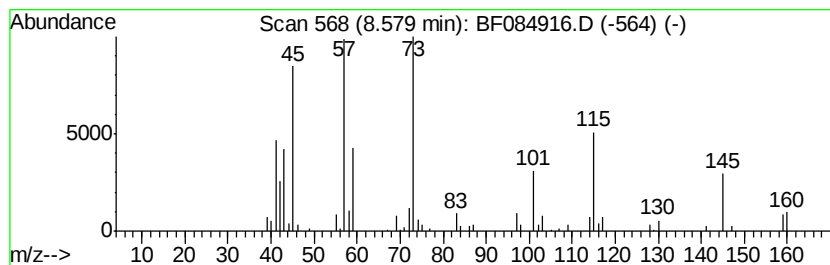
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown8.58 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.58	2.25 ng	136584	Napthalene-d8	7.95

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butanoic acid, 4-methoxy-, methy...	132	C6H12O3	029006-01-7	43
2		Di-sec-butyl ether	130	C8H18O	006863-58-7	38
3		Pentane, 1-(1-ethoxyethoxy)-	160	C9H20O2	013442-89-2	37
4		3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	35
5		1,3-Dioxan-5-ol	104	C4H8O3	004740-78-7	27



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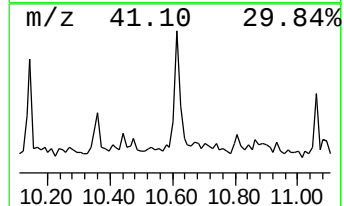
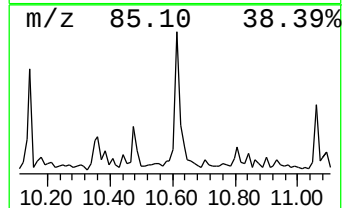
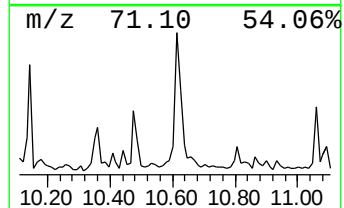
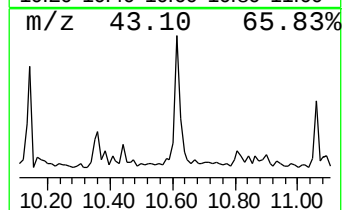
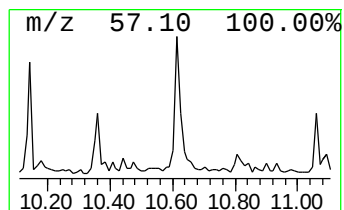
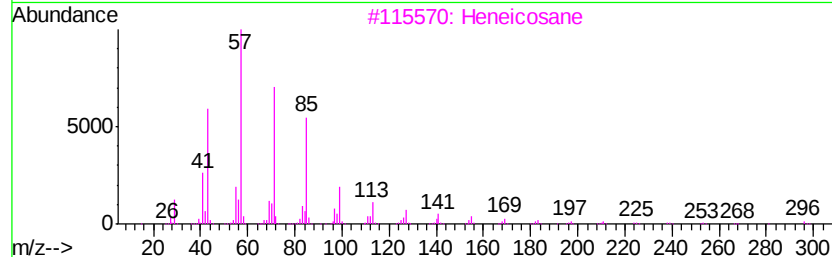
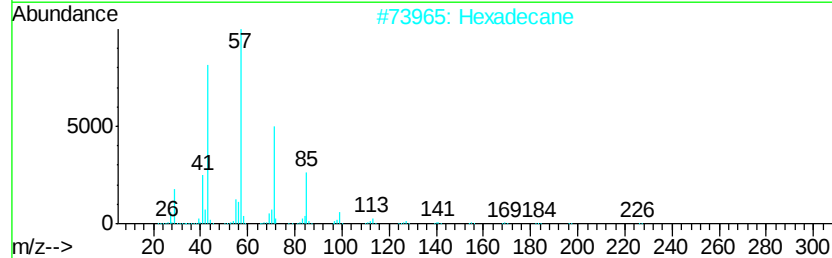
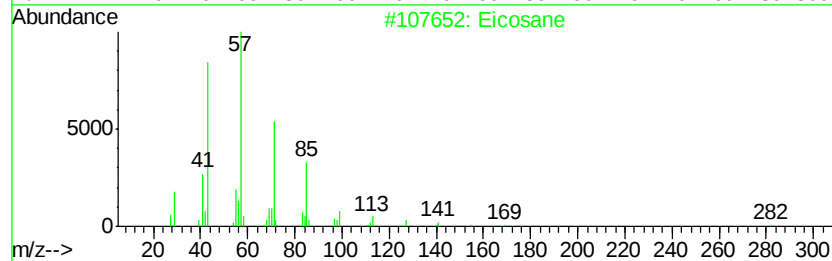
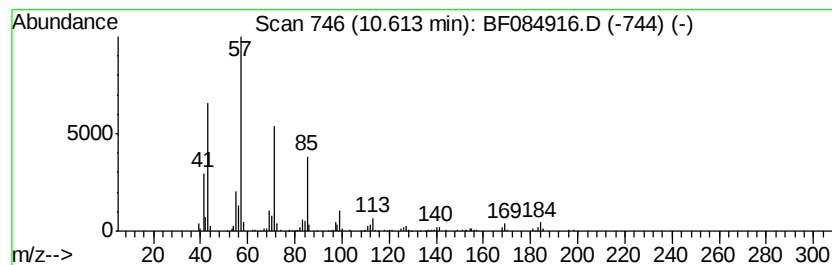
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 Eicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.61	3.86 ng	231491	Phenanthrene-d10	11.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	90
2	Hexadecane		226	C16H34	000544-76-3	86
3	Heneicosane		296	C21H44	000629-94-7	86
4	Nonadecane		268	C19H40	000629-92-5	83
5	Tetradecane		198	C14H30	000629-59-4	81



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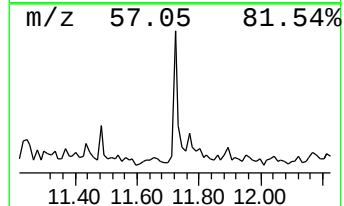
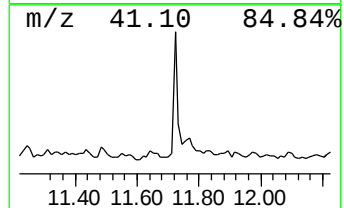
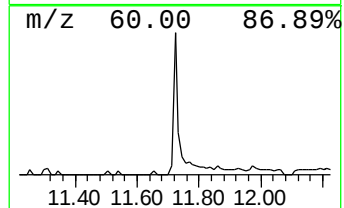
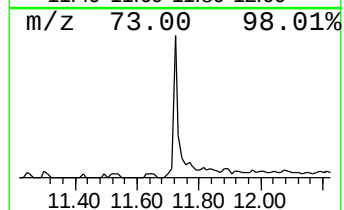
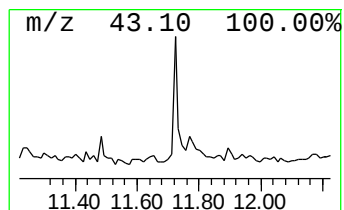
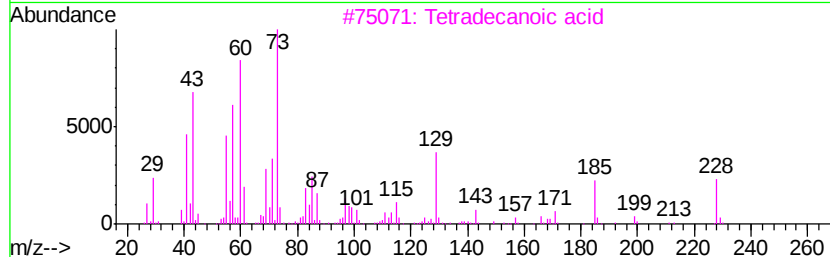
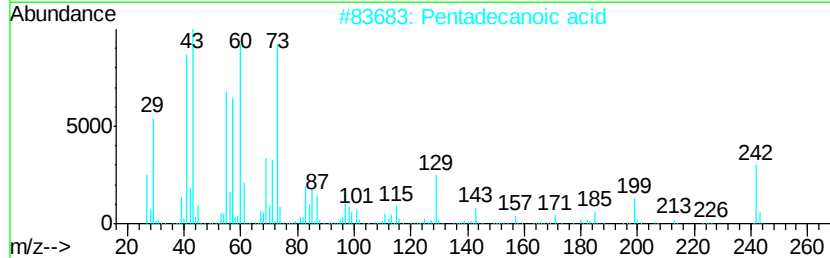
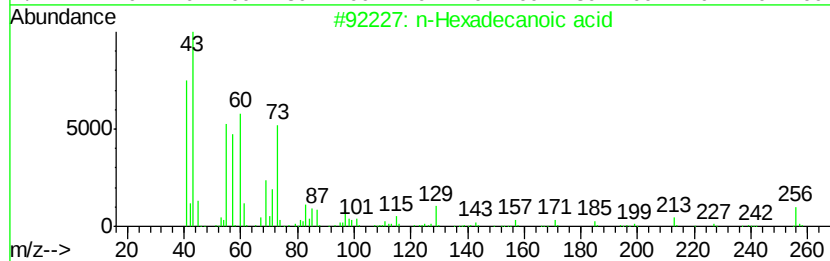
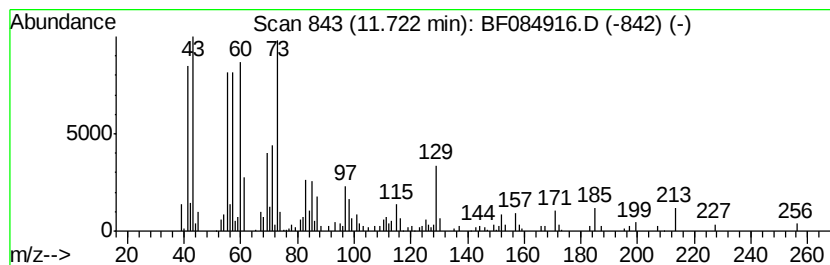
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.72	2.10 ng	125726	Phenanthrene-d10	11.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	86
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	72
4		n-Decanoic acid	172	C10H20O2	000334-48-5	53
5		N-Cyclooct-4-enylacetamide	167	C10H17NO	170952-69-9	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020616\
Data File : BF084916.D
Acq On : 6 Feb 2016 17:16
Operator : SJ/IZ
Sample : H1338-02
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-2

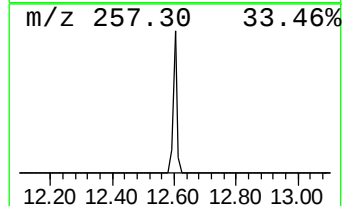
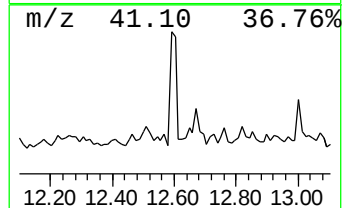
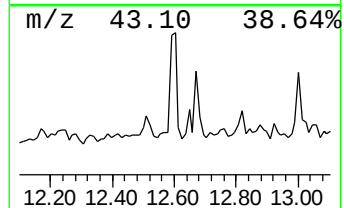
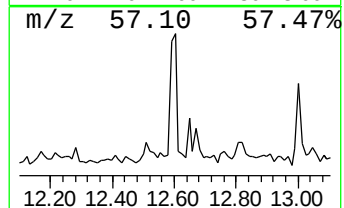
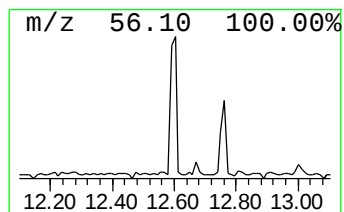
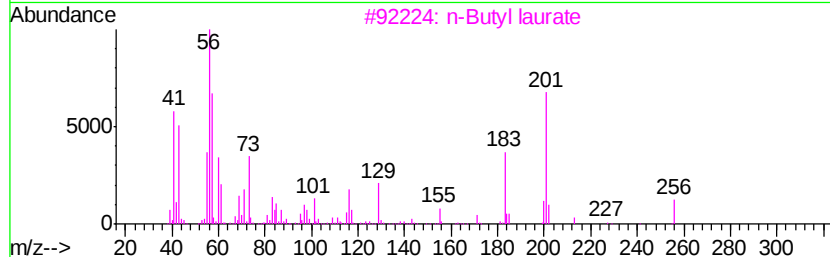
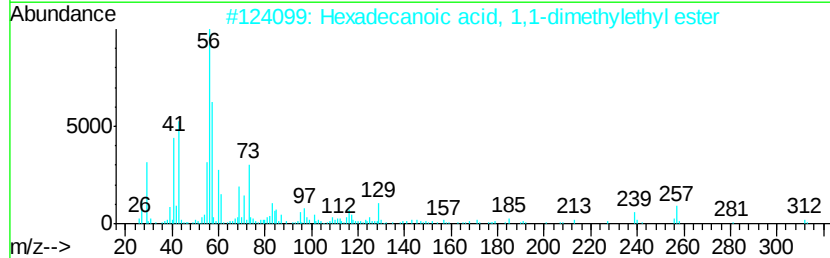
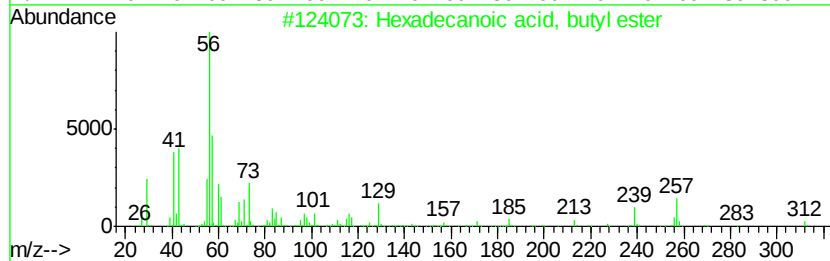
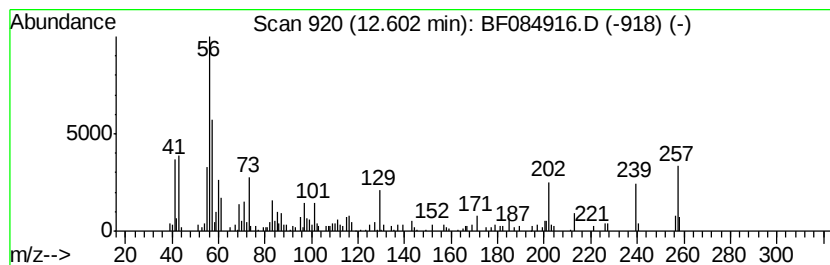
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecanoic acid, butyl ester Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.60	2.98 ng	136082	Chrysene-d12	13.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	62
2		Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	49
3		n-Butyl laurate	256	C16H32O2	000106-18-3	37
4		1-Hexene, 2-methyl-	98	C7H14	006094-02-6	25
5		3-Bromopropyl isocyanate	163	C4H6BrNO	056017-72-2	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020616\
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Operator : SJ/IZ
Sample : H1338-02
Misc :
ALS Vial : 58 Sample Multiplier: 1

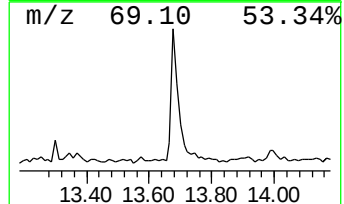
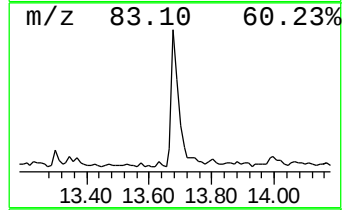
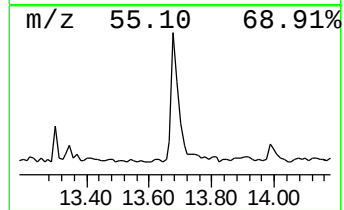
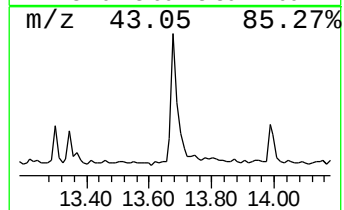
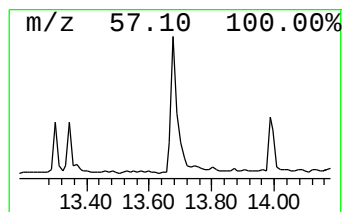
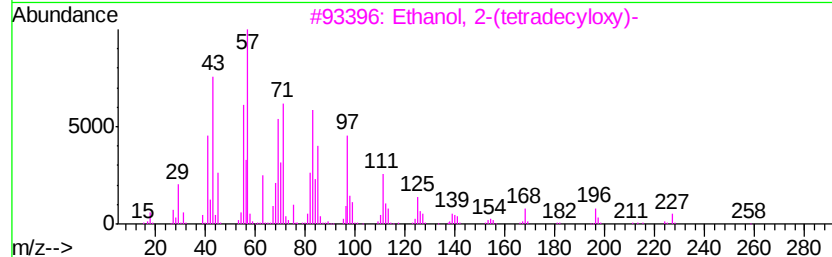
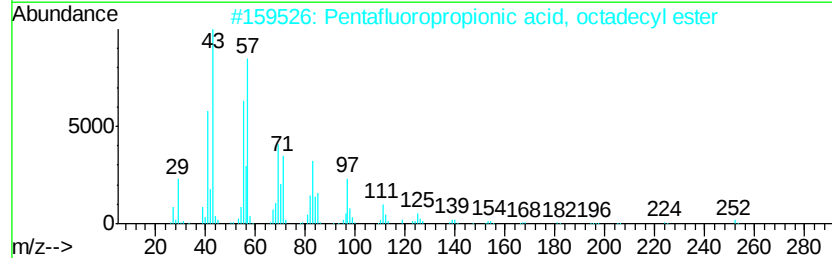
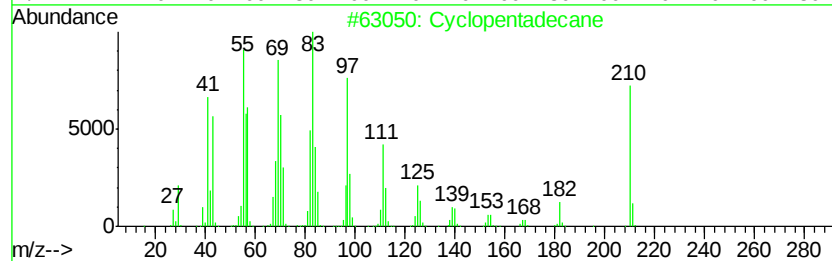
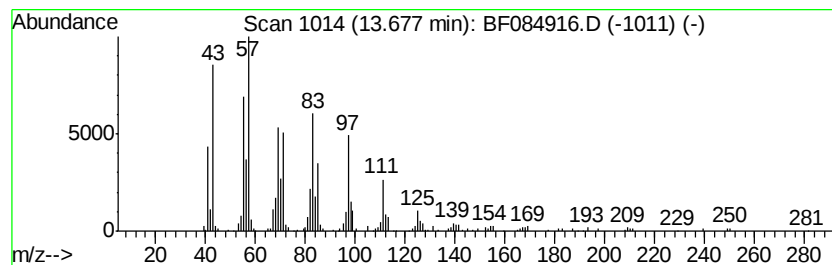
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Cyclopentadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.68	7.93 ng	362295	Chrysene-d12	13.80		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cvclopentadecane	210	C15H30	000295-48-7	91
2		Pentafluoropropionic acid, octad...	416	C21H37F5O2	1000280-07-7	91
3		Ethanol, 2-(tetradecvloxy)-	258	C16H34O2	002136-70-1	91
4		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	91
5		17-Pentatriacontene	491	C35H70	006971-40-0	90



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020616\
Data File : BF084916.D
Acq On : 6 Feb 2016 17:16
Operator : SJ/IZ
Sample : H1338-02
Misc :
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Instrument :
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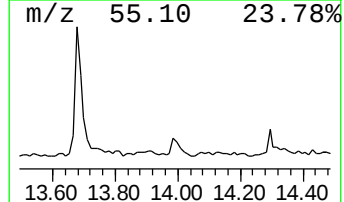
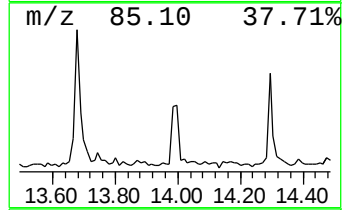
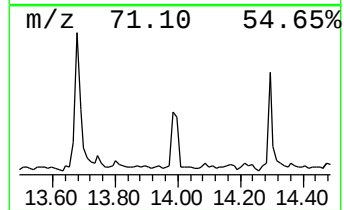
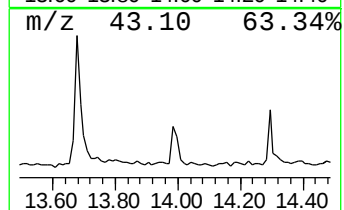
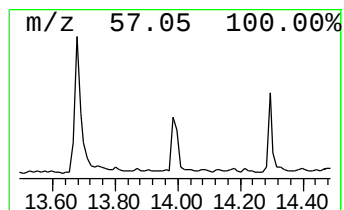
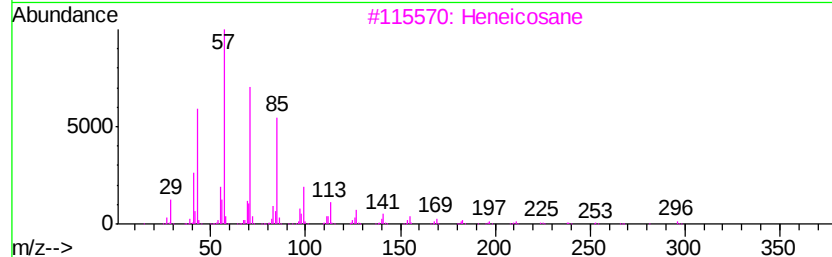
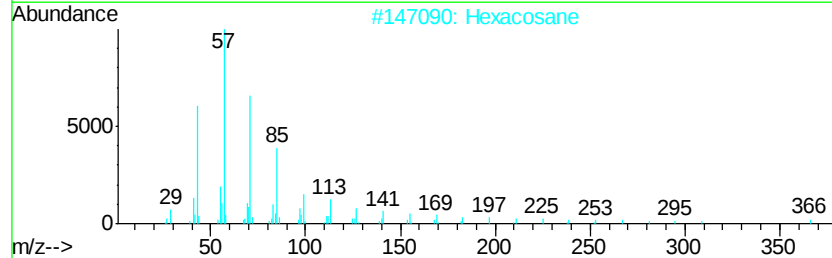
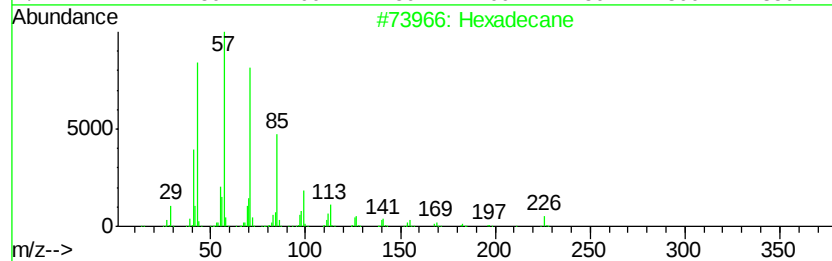
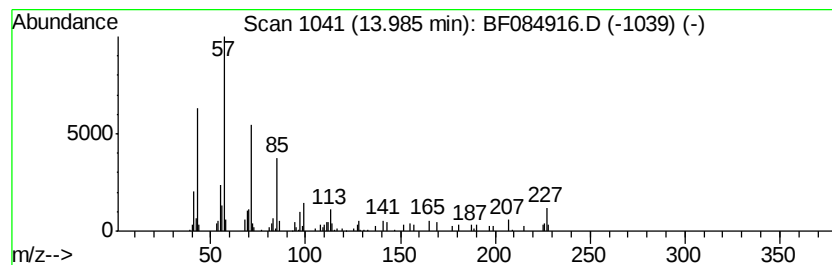
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	2.01 ng	91628	Chrysene-d12	13.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	76
2		Hexacosane	366	C26H54	000630-01-3	74
3		Heneicosane	296	C21H44	000629-94-7	74
4		Hentriacontane	437	C31H64	000630-04-6	72
5		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020616\
Data File : BF084916.D
Acq On : 6 Feb 2016 17:16
Operator : SJ/IZ
Sample : H1338-02
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-2

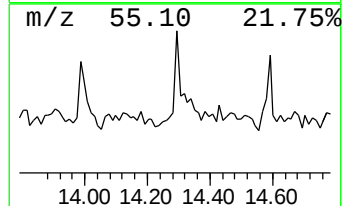
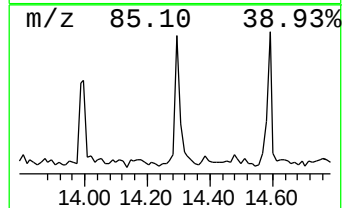
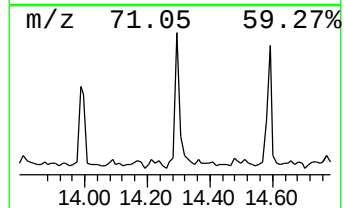
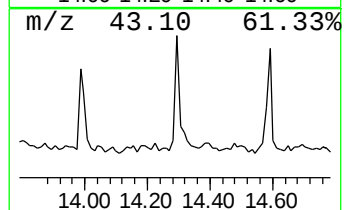
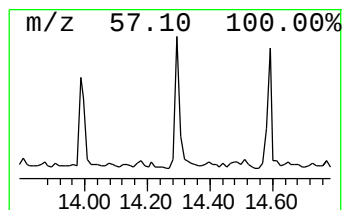
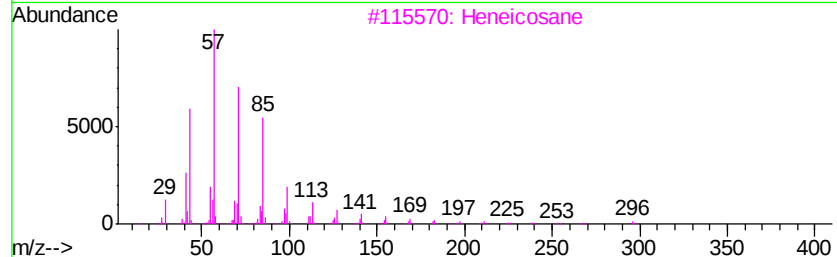
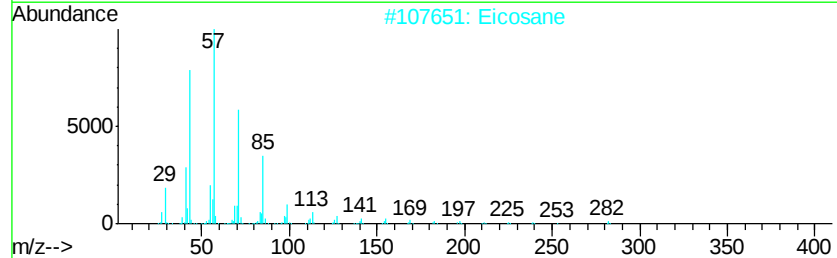
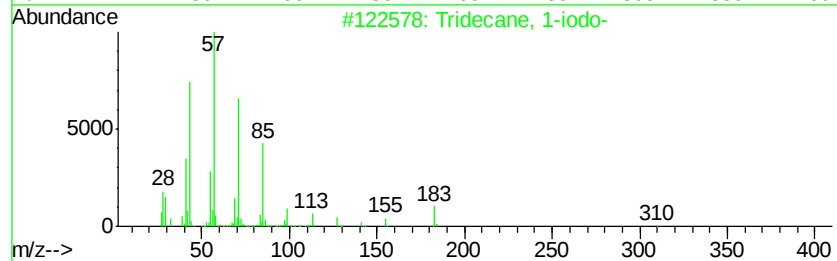
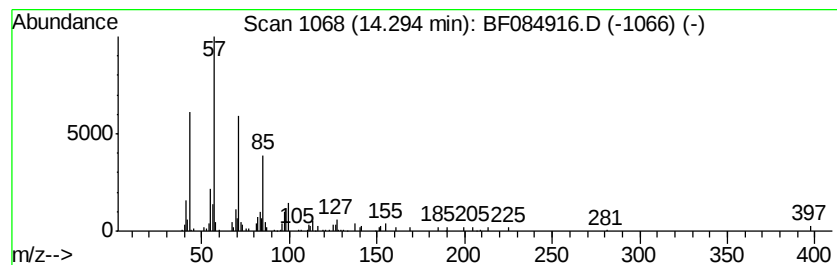
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tridecane, 1-iodo- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.29	2.59 ng	118237	Chrysene-d12	13.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	87
2		Eicosane	282	C20H42	000112-95-8	86
3		Heneicosane	296	C21H44	000629-94-7	86
4		Heptacosane	380	C27H56	000593-49-7	86
5		Pentadecane	212	C15H32	000629-62-9	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020616\
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Sample : H1338-02
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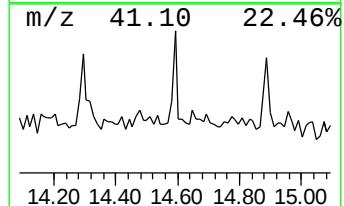
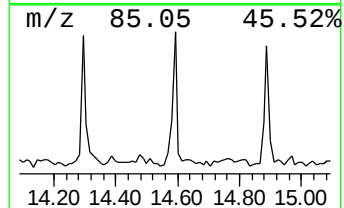
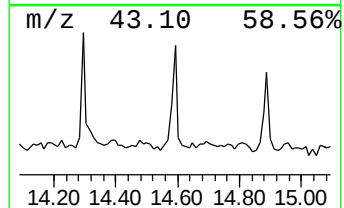
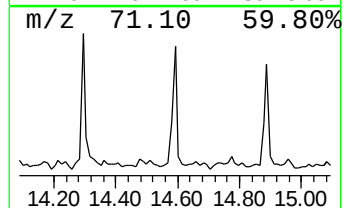
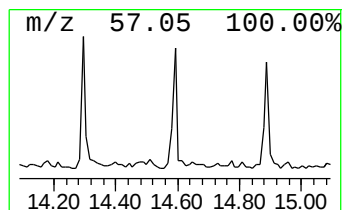
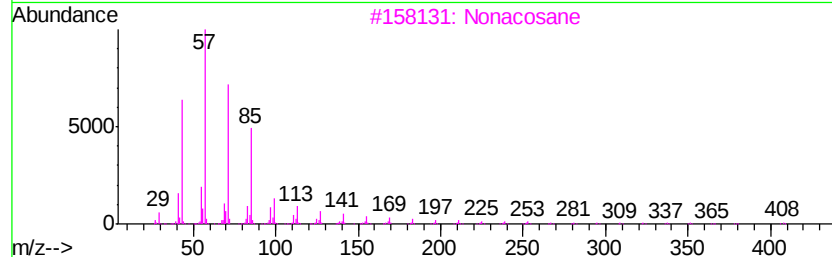
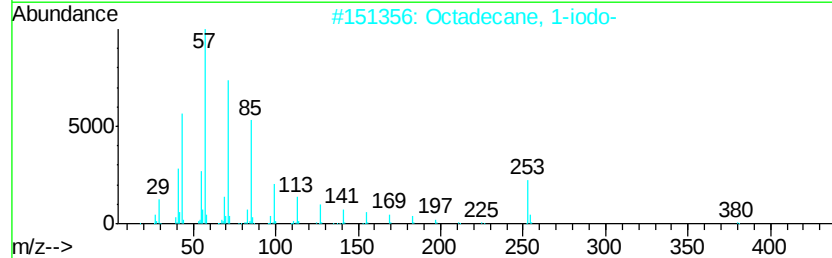
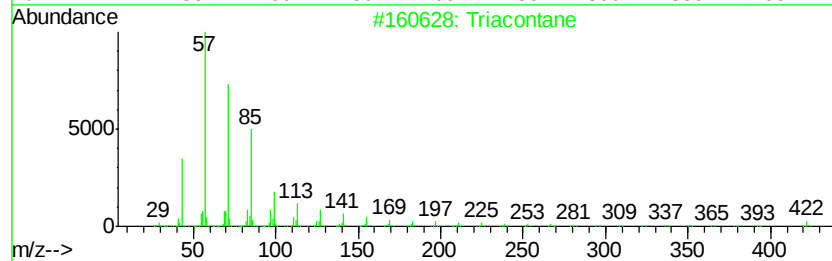
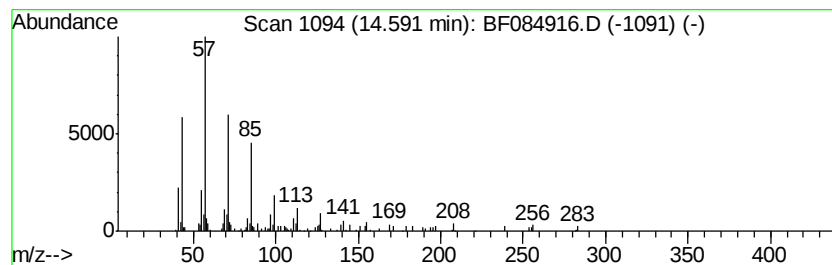
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Triacontane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.59	2.26 ng	93577	Perylene-d12	15.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Triacontane	422	C30H62	000638-68-6	91
2		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	91
3		Nonacosane	408	C29H60	000630-03-5	90
4		Hentriacontane	437	C31H64	000630-04-6	90
5		Tetracosane	338	C24H50	000646-31-1	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020616\
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Acq On : 6 Feb 2016 17:16
Operator : SJ/IZ
Sample : H1338-02
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.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
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unknown6.43	6.43	70.4	ng	3361430	1	6.66	954761	20.0
unknown8.58	8.58	2.3	ng	136584	2	7.95	1212750	20.0
Eicosane	10.61	3.9	ng	231491	4	11.17	1199520	20.0
n-Hexadecanoic acid	11.72	2.1	ng	125726	4	11.17	1199520	20.0
Hexadecanoic acid...	12.60	3.0	ng	136082	5	13.80	913348	20.0
Cyclopentadecane	13.68	7.9	ng	362295	5	13.80	913348	20.0
Hexadecane	13.99	2.0	ng	91628	5	13.80	913348	20.0
Tridecane, 1-iodo-	14.29	2.6	ng	118237	5	13.80	913348	20.0
Triacontane	14.59	2.3	ng	93577	6	15.17	829269	20.0