

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
 Data File : BF086393.D
 Acq On : 23 Apr 2016 16:31
 Operator : UM/SJ
 Sample : H2640-05
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WPG-B13(0-5)-C

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-BF042216.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.024	254	257	265	rBV	592178	877221	19.89%	2.008%
2	5.436	290	293	296	rBV	2692703	3793494	86.01%	8.686%
3	5.847	327	329	331	rBV	71053	66600	1.51%	0.152%
4	6.476	381	384	393	rBV	3293144	3891580	88.23%	8.910%
5	6.613	393	396	398	rVV	3794995	4327543	98.12%	9.908%
6	6.682	400	402	404	rVB2	35227	50495	1.14%	0.116%
7	6.842	414	416	420	rBV	1131974	1084184	24.58%	2.482%
8	7.002	427	430	432	rBV	3179999	3433583	77.85%	7.861%
9	7.402	463	465	468	rVV	1752611	2446794	55.48%	5.602%
10	7.447	468	469	471	rVV	103836	132041	2.99%	0.302%
11	7.493	471	473	481	rVB2	60491	148017	3.36%	0.339%
12	7.870	502	506	510	rBV3	24055	58519	1.33%	0.134%
13	8.122	526	528	533	rBV	1144846	1498039	33.97%	3.430%
14	8.727	579	581	582	rBV	51056	62451	1.42%	0.143%
15	8.842	587	591	593	rVV2	48633	63080	1.43%	0.144%
16	8.933	597	599	601	rBV	48611	54093	1.23%	0.124%
17	9.207	620	623	625	rBV	3736091	4410476	100.00%	10.098%
18	9.287	628	630	632	rVB2	102313	110058	2.50%	0.252%
19	9.470	643	646	647	rBV	29754	46925	1.06%	0.107%
20	9.539	649	652	655	rBV2	64827	104334	2.37%	0.239%
21	9.596	655	657	659	rBV	635580	582164	13.20%	1.333%
22	9.813	675	676	679	rVB	151789	118666	2.69%	0.272%
23	9.882	679	682	683	rBV	1088484	1339323	30.37%	3.066%
24	9.905	683	684	686	rVB	102808	101733	2.31%	0.233%
25	10.042	693	696	698	rBV4	29175	66484	1.51%	0.152%
26	10.076	698	699	702	rBV2	62816	69055	1.57%	0.158%
27	10.133	702	704	706	rVB3	38853	55147	1.25%	0.126%
28	10.248	710	714	715	rBV4	25487	50866	1.15%	0.116%
29	10.282	715	717	718	rVV	76707	104506	2.37%	0.239%
30	10.316	718	720	721	rVV	233911	223559	5.07%	0.512%
31	10.339	721	722	728	rVB5	28096	72136	1.64%	0.165%
32	10.419	728	729	733	rVB	81276	92961	2.11%	0.213%
33	10.533	736	739	742	rVB2	105152	157835	3.58%	0.361%
34	10.579	742	743	745	rVB2	41774	49156	1.11%	0.113%

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35	10.659	748	750	753	rBV	1466890	1753691	39.76%	4.015%
36	10.785	759	761	765	rVB	314447	392805	8.91%	0.899%
37	10.979	775	778	780	rBV3	37065	75693	1.72%	0.173%
38	11.231	796	800	801	rBV	206782	219409	4.97%	0.502%
39	11.265	801	803	806	rVB2	107160	146702	3.33%	0.336%
40	11.356	809	811	815	rBV2	1028968	1625523	36.86%	3.722%
41	11.436	815	818	823	rVB	124017	262651	5.96%	0.601%
42	11.608	830	833	836	rBV2	36916	70215	1.59%	0.161%
43	11.665	836	838	839	rVV	68322	92891	2.11%	0.213%
44	11.859	853	855	856	rBV	70585	72961	1.65%	0.167%
45	11.882	856	857	859	rVV	58250	82303	1.87%	0.188%
46	11.973	863	865	869	rVB2	88751	167206	3.79%	0.383%
47	12.065	872	873	875	rVV	73368	77060	1.75%	0.176%
48	12.156	878	881	886	rBV3	52027	133078	3.02%	0.305%
49	12.419	902	904	905	rBV	48528	83267	1.89%	0.191%
50	12.454	905	907	909	rVV2	113940	122685	2.78%	0.281%
51	12.579	916	918	920	rBV	544756	498793	11.31%	1.142%
52	12.796	934	937	938	rBV	453292	478015	10.84%	1.094%
53	12.945	947	950	955	rBV	2902308	3048621	69.12%	6.980%
54	13.185	968	971	973	rBV2	143080	212043	4.81%	0.485%
55	13.528	998	1001	1002	rBV	133663	207980	4.72%	0.476%
56	13.848	1027	1029	1034	rVB	298546	373953	8.48%	0.856%
57	13.997	1039	1042	1050	rVV2	702329	1498472	33.98%	3.431%
58	14.168	1055	1057	1059	rBV	157520	161491	3.66%	0.370%
59	14.465	1082	1083	1085	rBV	119857	185635	4.21%	0.425%
60	15.002	1128	1130	1134	rBV	130133	281799	6.39%	0.645%
61	15.414	1163	1166	1168	rBV	386025	583751	13.24%	1.337%
62	16.077	1222	1224	1232	rVB3	176390	467085	10.59%	1.069%
63	16.488	1258	1260	1264	rVB	116083	211333	4.79%	0.484%
64	16.957	1299	1301	1305	rVB2	174516	345908	7.84%	0.792%

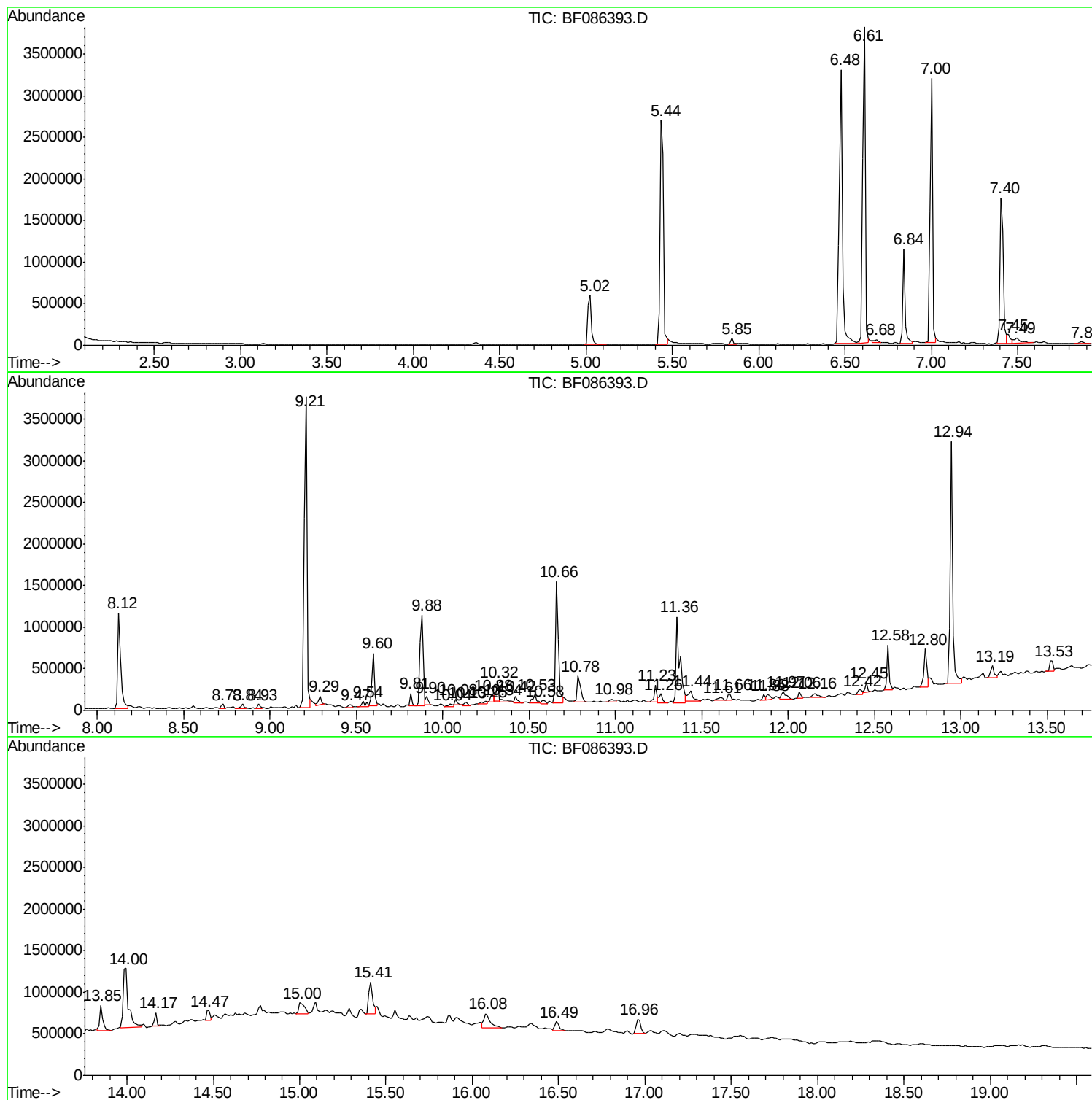
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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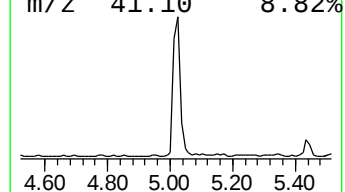
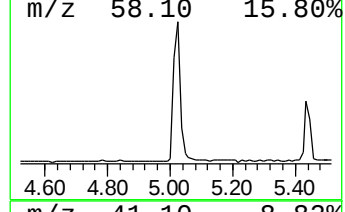
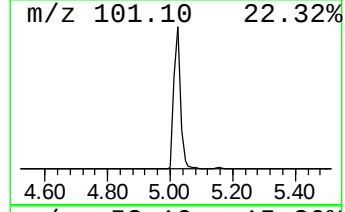
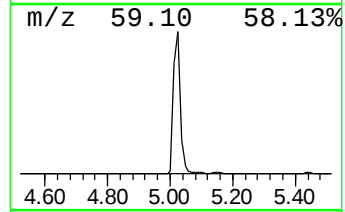
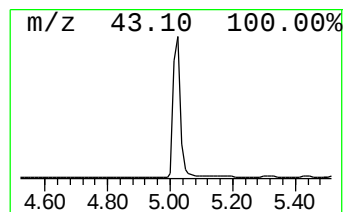
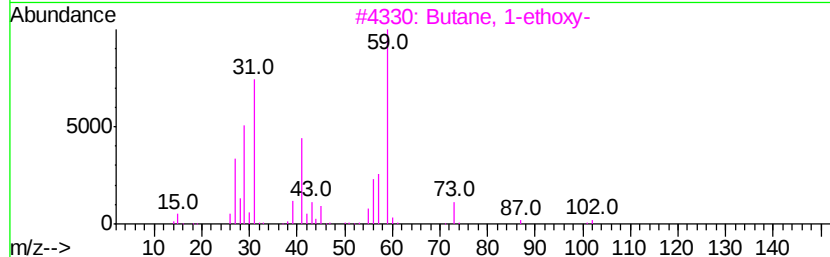
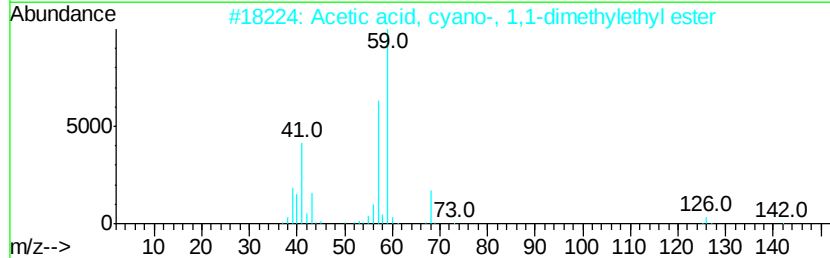
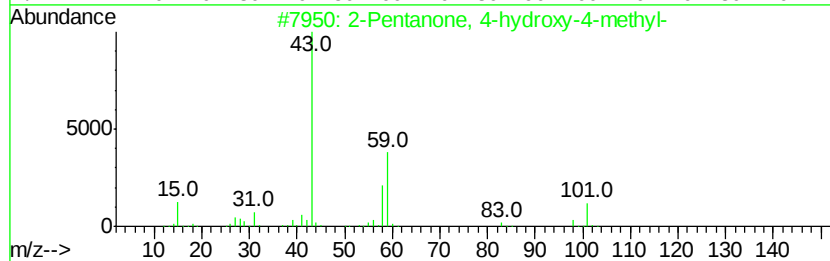
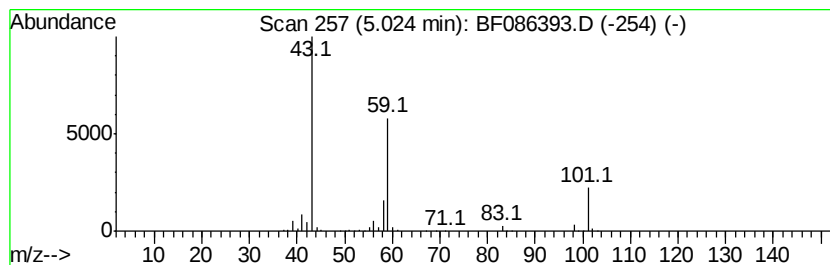
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
5.02	16.18 ng	877221	1,4-Dichlorobenzene-d4	6.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4			5-Hexen-2-one	98	C6H10O	000109-49-9	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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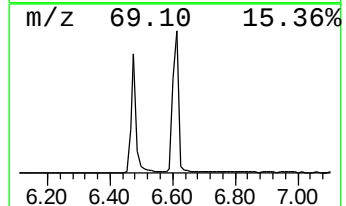
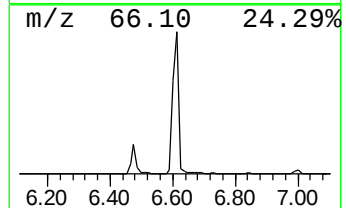
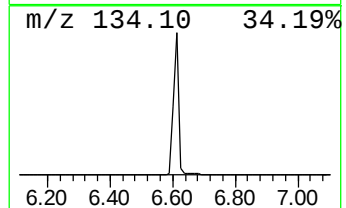
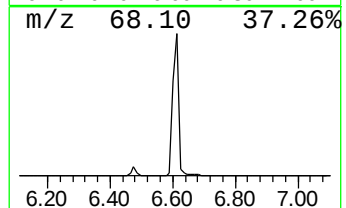
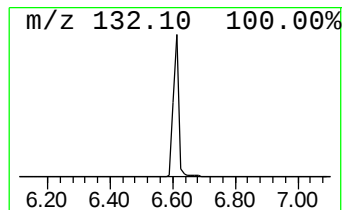
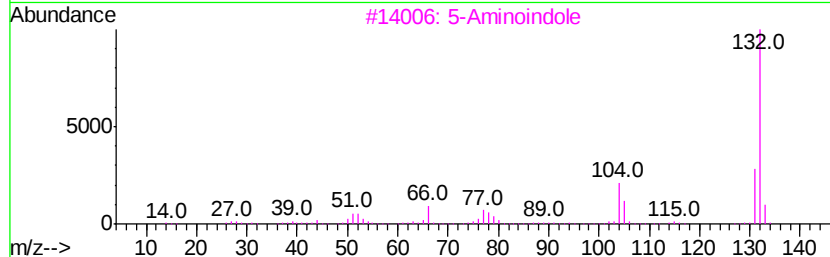
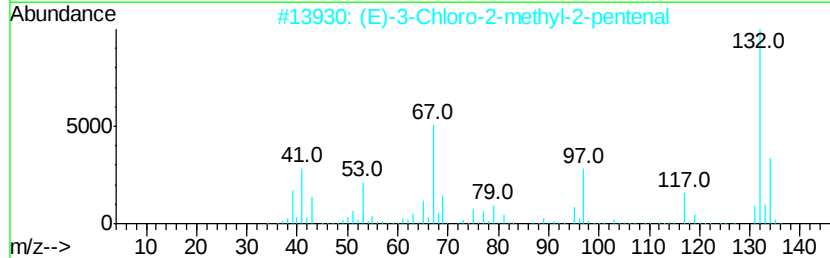
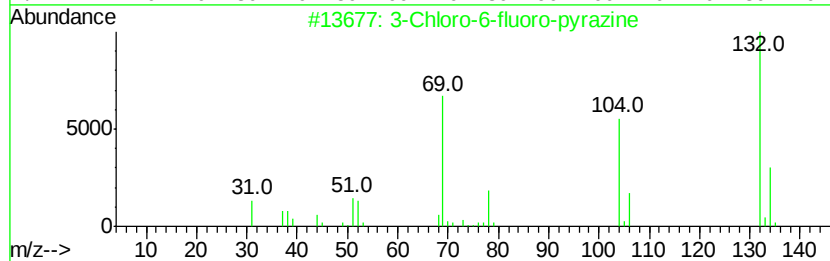
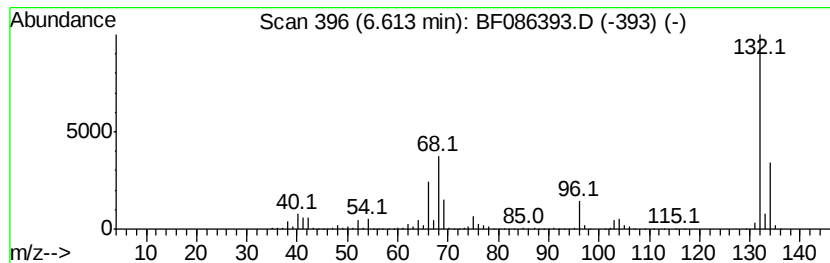
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.61	79.83 ng	4327540	1,4-Dichlorobenzene-d4	6.84

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	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	16
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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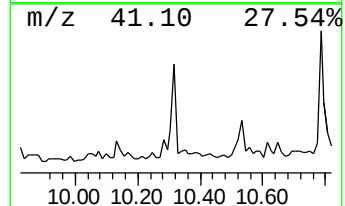
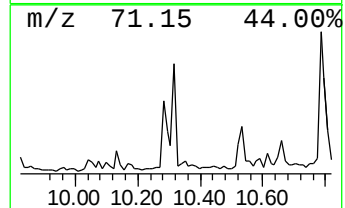
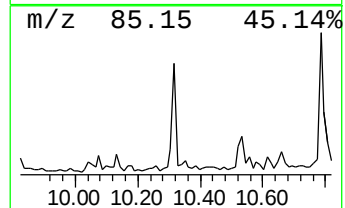
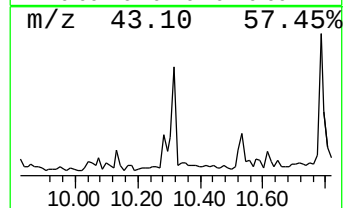
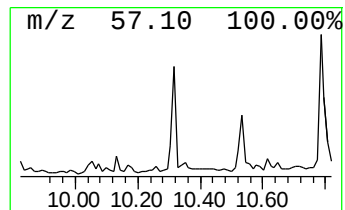
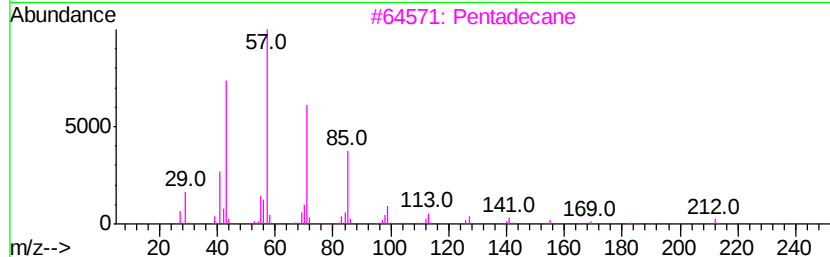
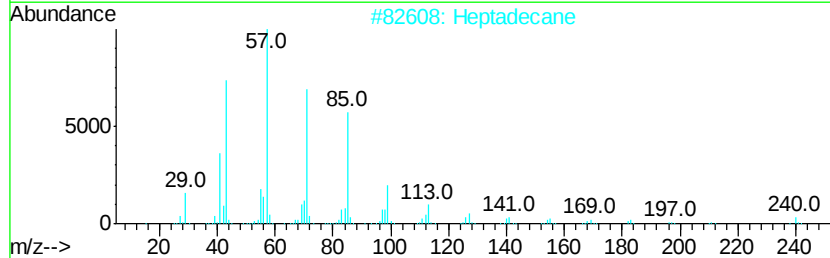
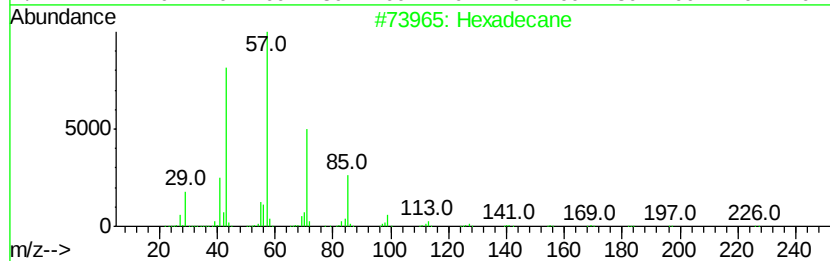
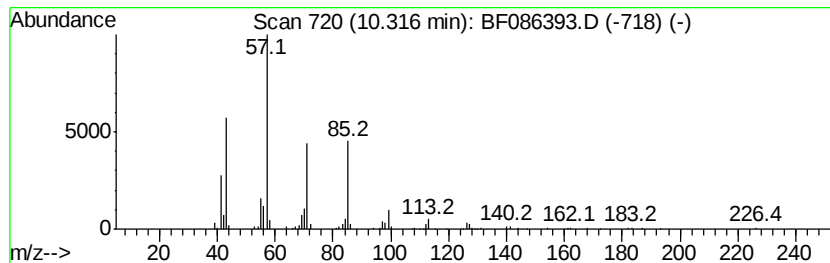
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Hexadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.32	3.34 ng	223559	Acenaphthene-d10	9.88	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	94
2	Heptadecane	240	C17H36	000629-78-7	83
3	Pentadecane	212	C15H32	000629-62-9	72
4	Eicosane, 10-methyl-	296	C21H44	054833-23-7	59
5	Eicosane	282	C20H42	000112-95-8	58



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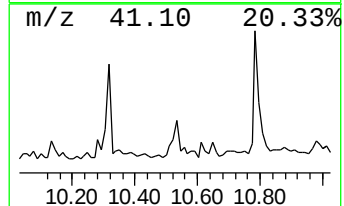
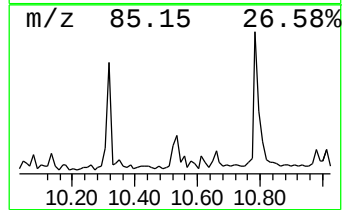
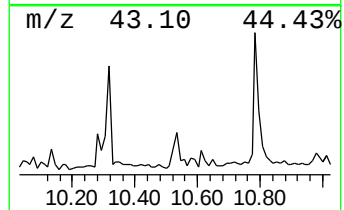
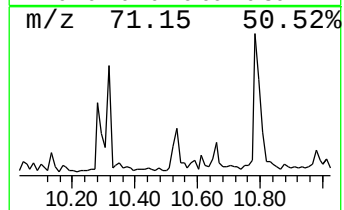
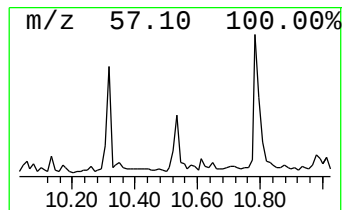
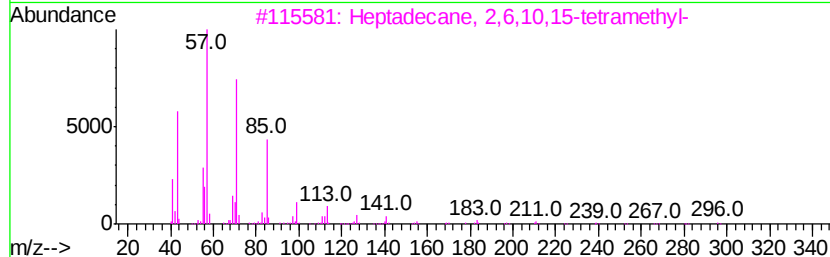
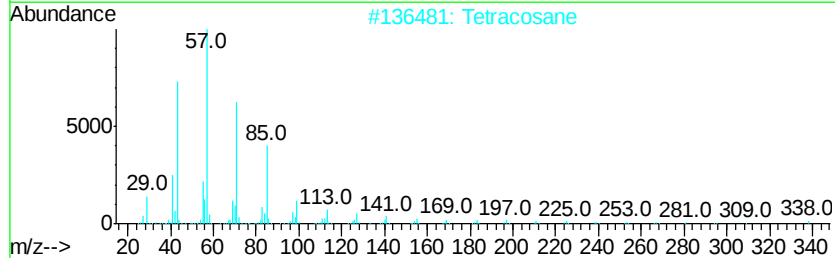
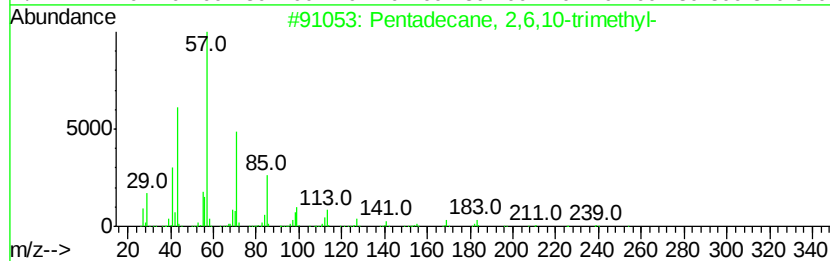
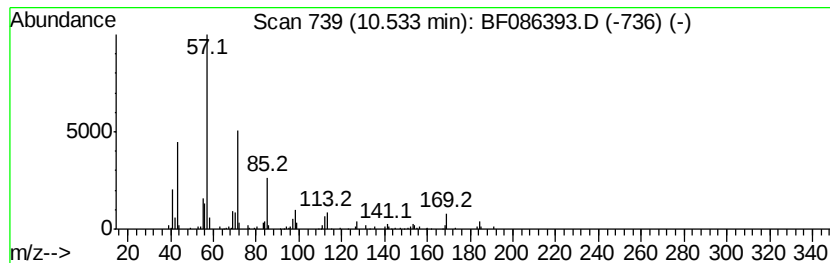
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Pentadecane, 2,6,10-trimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.53	2.36 ng	157835	Acenaphthene-d10	9.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	78
2		Tetracosane	338	C24H50	000646-31-1	58
3		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	58
4		Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	58
5		Octadecane	254	C18H38	000593-45-3	53



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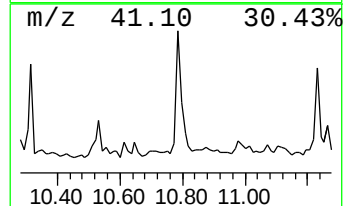
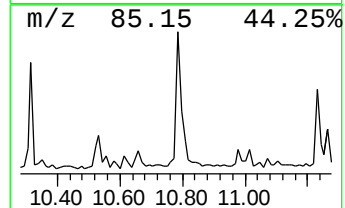
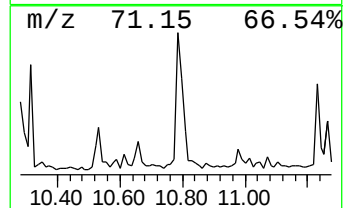
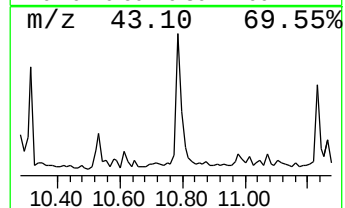
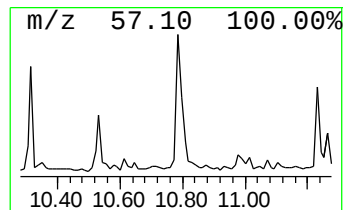
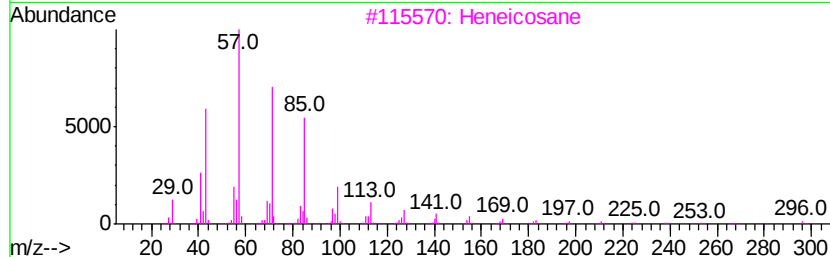
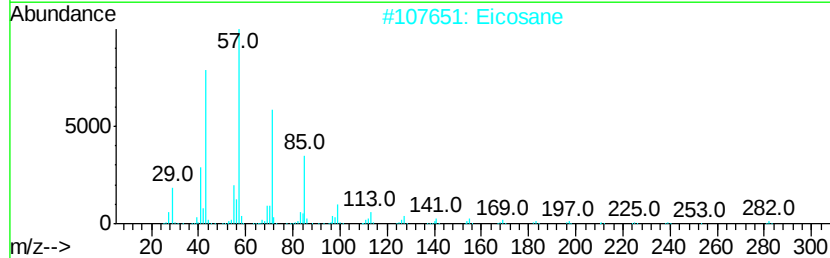
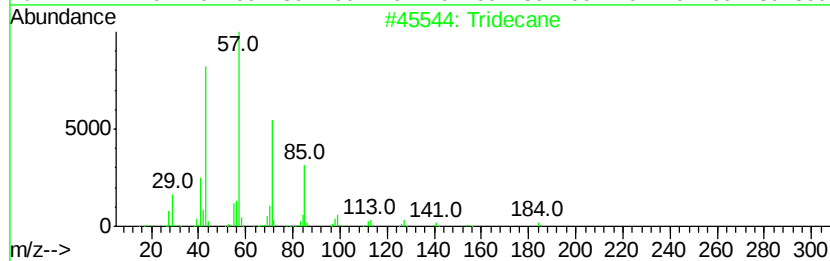
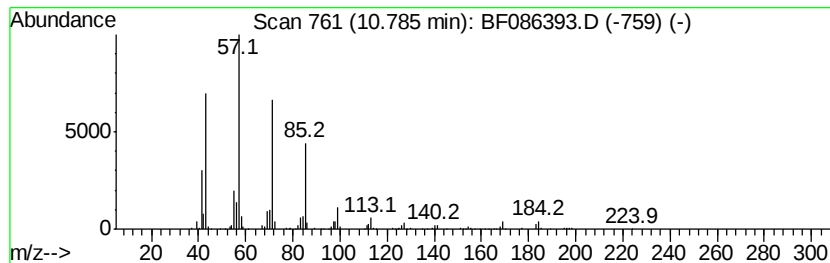
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Tridecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.78	4.83 ng	392805	Phenanthrene-d10	11.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane	184	C13H28	000629-50-5	87
2		Eicosane	282	C20H42	000112-95-8	86
3		Heneicosane	296	C21H44	000629-94-7	86
4		Tetradecane	198	C14H30	000629-59-4	86
5		Undecane, 5-ethyl-5-propyl-	226	C16H34	002755-07-9	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
Data File : BF086393.D
Acq On : 23 Apr 2016 16:31
Operator : UM/SJ
Sample : H2640-05
Misc :
ALS Vial : 37 Sample Multiplier: 1

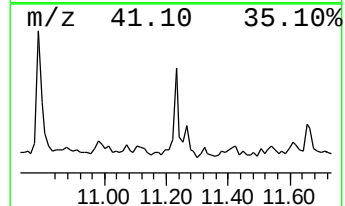
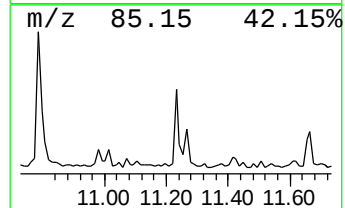
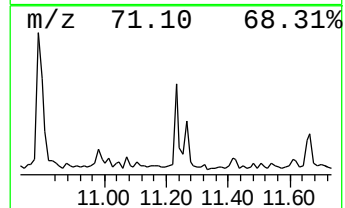
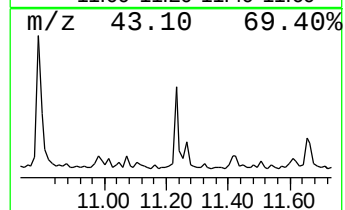
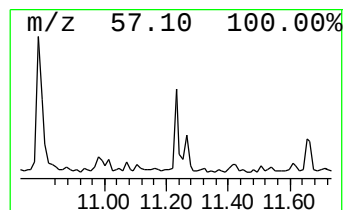
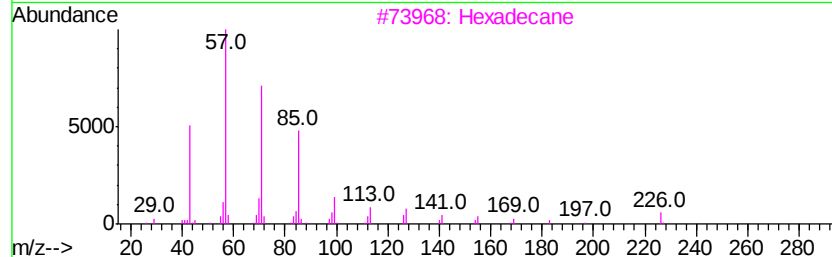
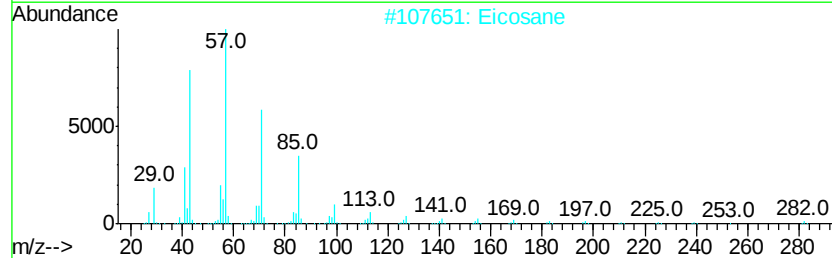
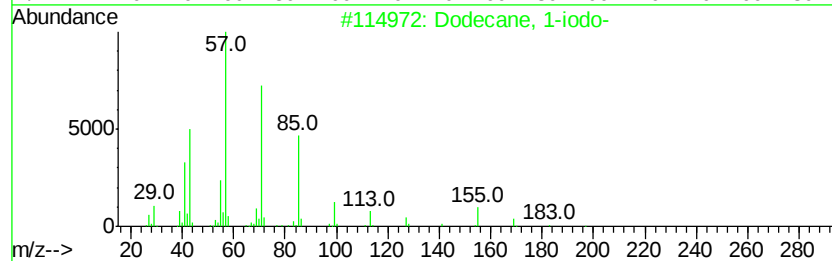
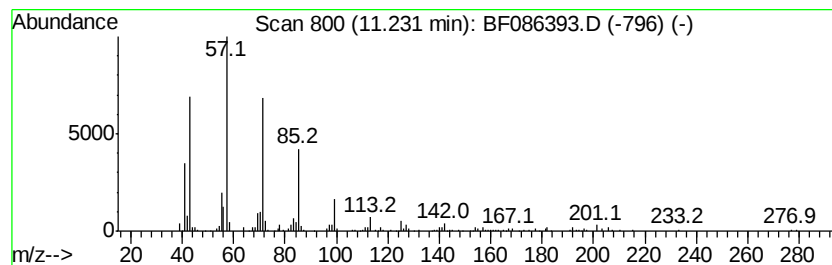
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Dodecane, 1-iodo- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.23	2.70 ng	219409	Phenanthrene-d10		11.36
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	80
2	Eicosane	282	C20H42	000112-95-8	80
3	Hexadecane	226	C16H34	000544-76-3	72
4	Heneicosane	296	C21H44	000629-94-7	72
5	Decane, 6-ethyl-2-methyl-	184	C13H28	062108-21-8	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
Data File : BF086393.D
Acq On : 23 Apr 2016 16:31
Operator : UM/SJ
Sample : H2640-05
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Instrument :
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ClientSampleId :
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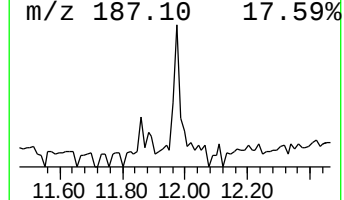
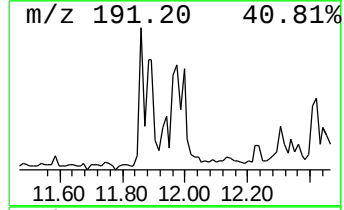
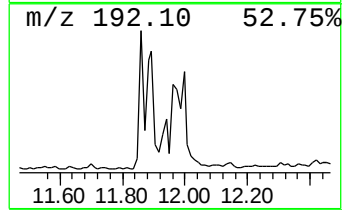
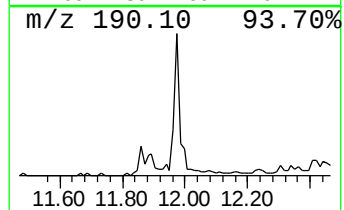
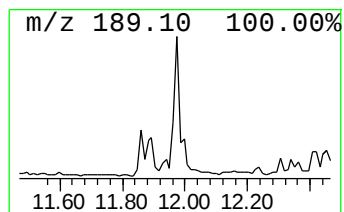
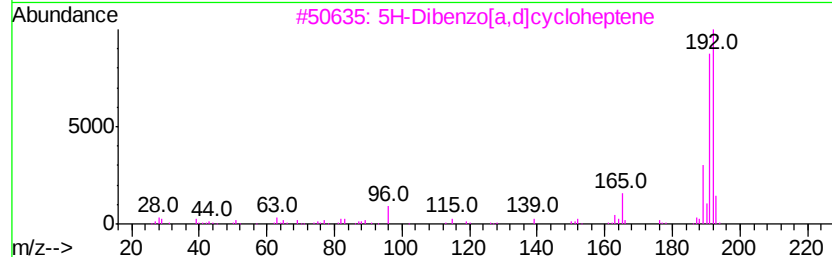
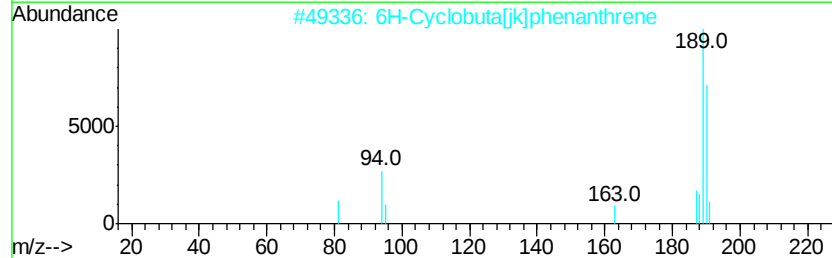
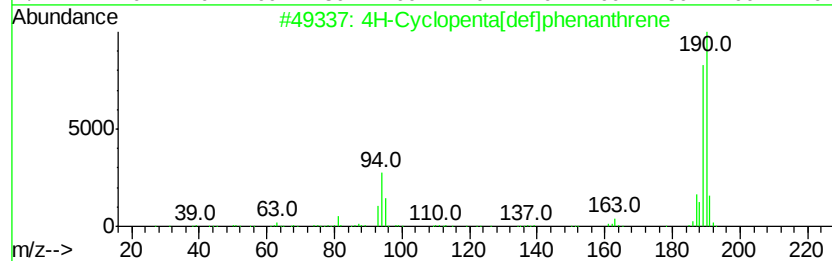
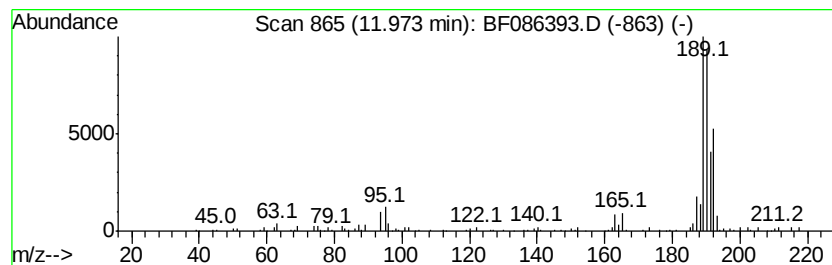
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 4H-Cyclopenta[def]phenanthrene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.97	2.06 ng	167206	Phenanthrene-d10	11.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2		6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	58
3		5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	25
4		Benzo[1,2-b:5,4-b']dithiophene	190	C10H6S2	000267-61-8	22
5		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
Data File : BF086393.D
Acq On : 23 Apr 2016 16:31
Operator : UM/SJ
Sample : H2640-05
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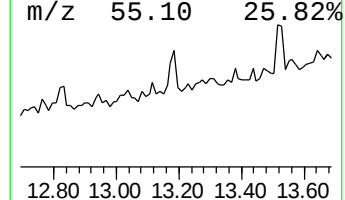
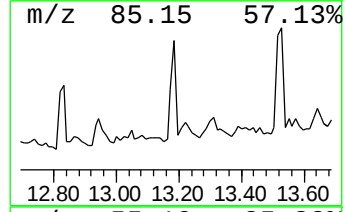
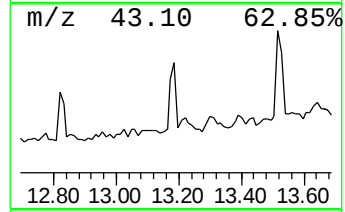
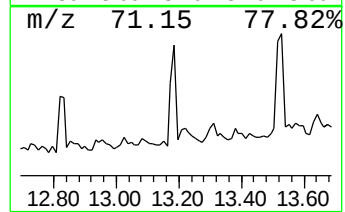
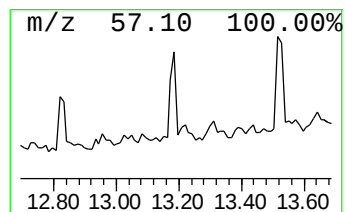
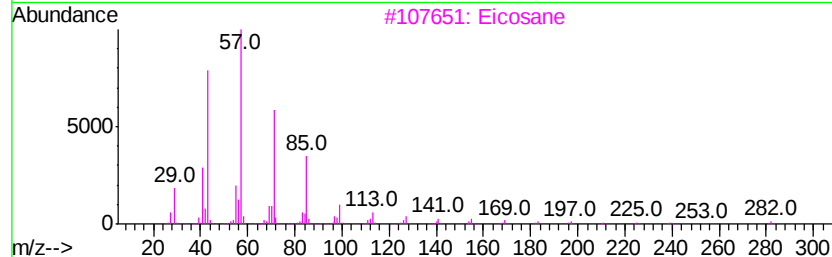
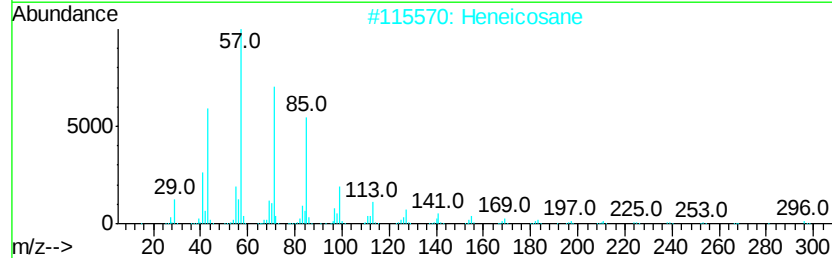
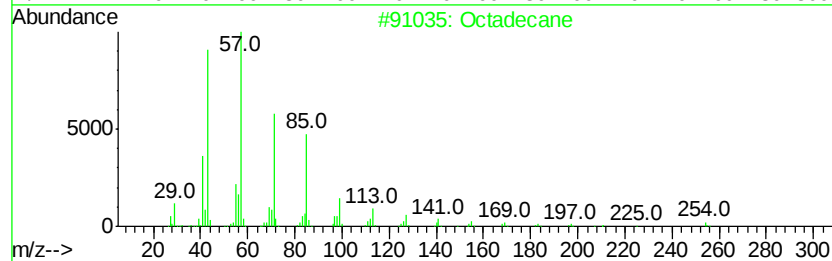
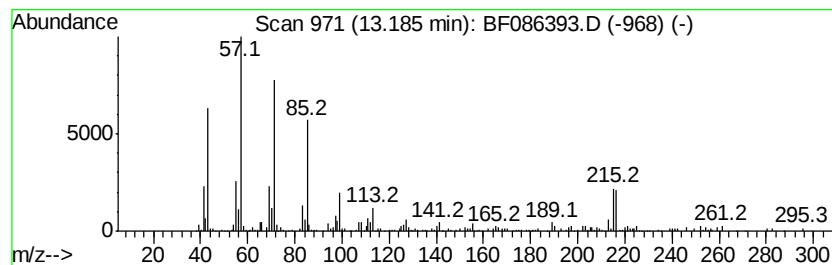
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.19	2.83 ng	212043	Chrysene-d12		14.00
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	86
2	Heneicosane	296	C21H44	000629-94-7	76
3	Eicosane	282	C20H42	000112-95-8	76
4	Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	70
5	Nonacosane	408	C29H60	000630-03-5	62



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
Data File : BF086393.D
Acq On : 23 Apr 2016 16:31
Operator : UM/SJ
Sample : H2640-05
Misc :
ALS Vial : 37 Sample Multiplier: 1

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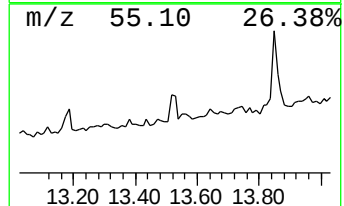
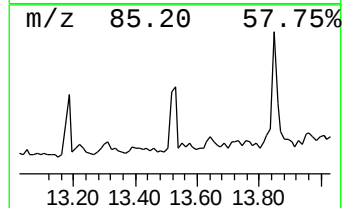
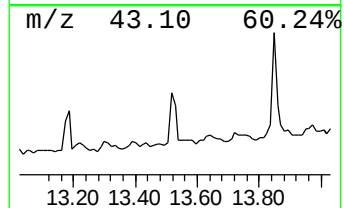
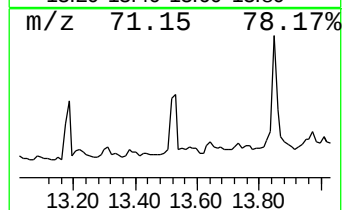
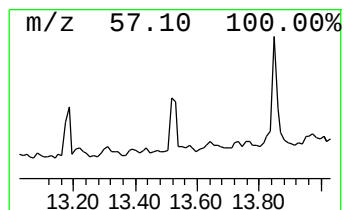
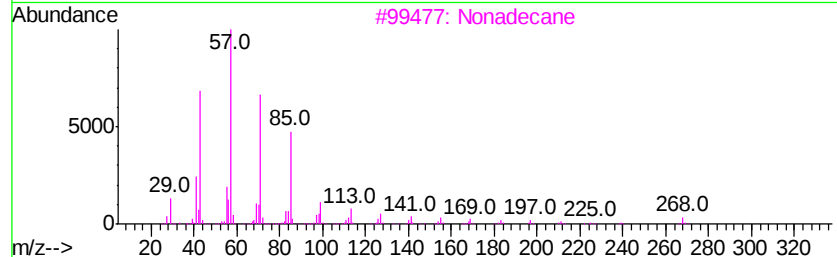
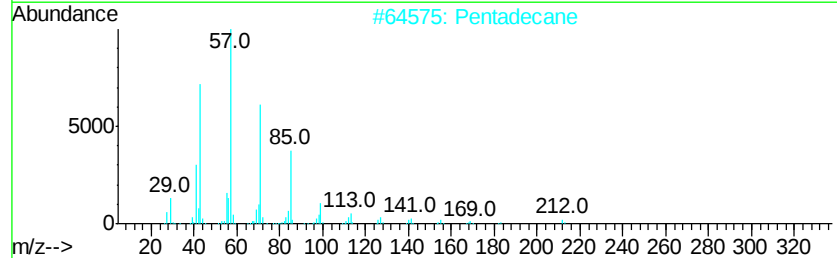
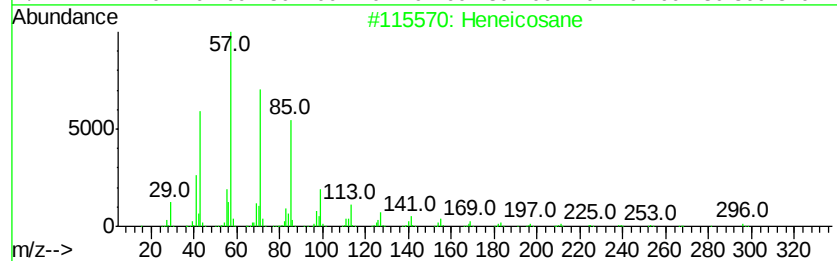
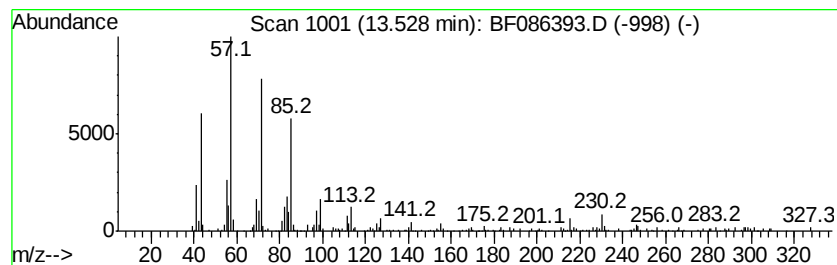
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heneicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.53	2.78 ng	207980	Chrysene-d12	14.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	96
2			Pentadecane	212	C15H32	000629-62-9	93
3			Nonadecane	268	C19H40	000629-92-5	91
4			Tetracosane	338	C24H50	000646-31-1	91
5			Heneicosane, 11-decyl-	437	C31H64	055320-06-4	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
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Acq On : 23 Apr 2016 16:31
Operator : UM/SJ
Sample : H2640-05
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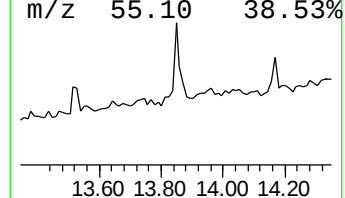
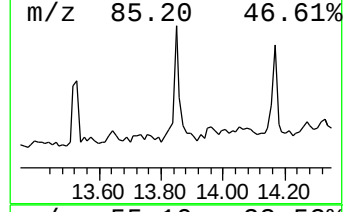
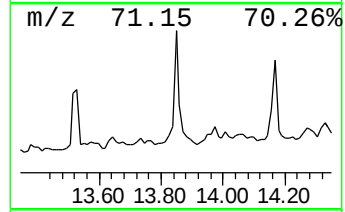
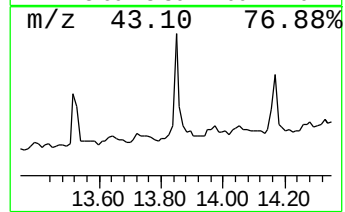
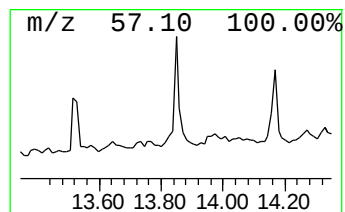
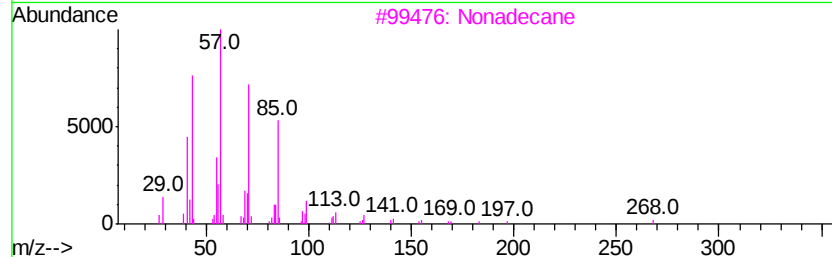
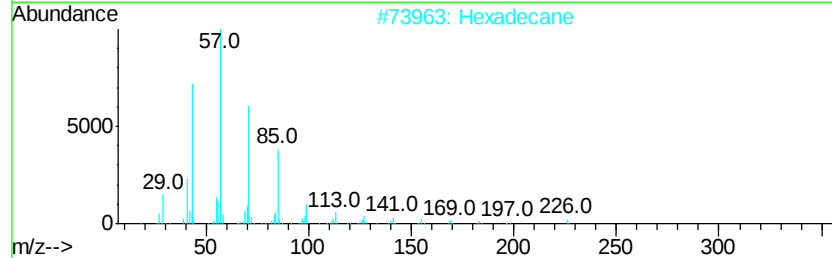
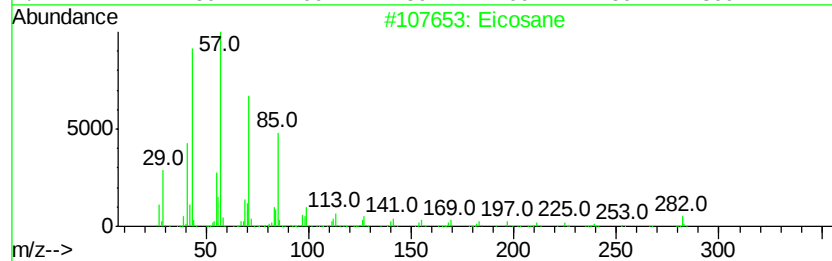
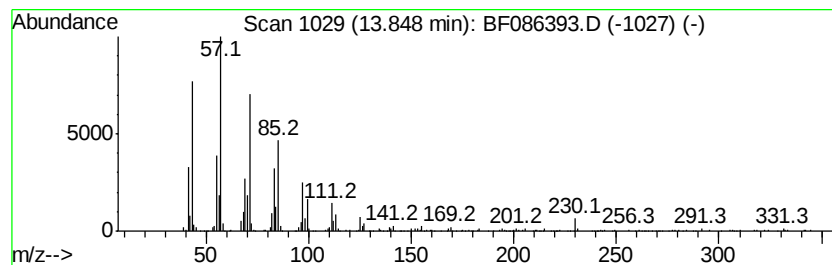
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.85	4.99 ng	373953	Chrysene-d12	14.00
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane	282 C20H42	000112-95-8	68
2	Hexadecane	226 C16H34	000544-76-3	60
3	Nonadecane	268 C19H40	000629-92-5	58
4	1-Iodoundecane	282 C11H23I	004282-44-4	53
5	Pentadecane	212 C15H32	000629-62-9	52



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042216\
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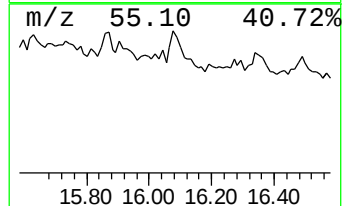
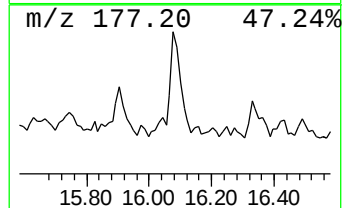
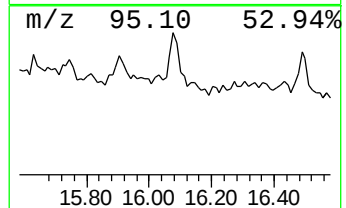
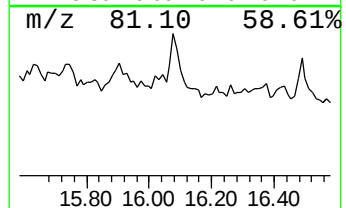
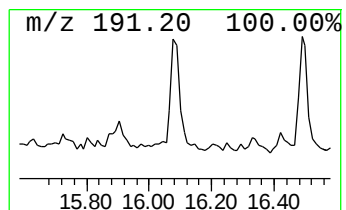
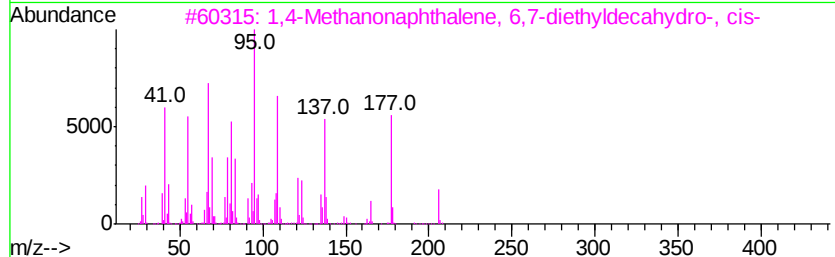
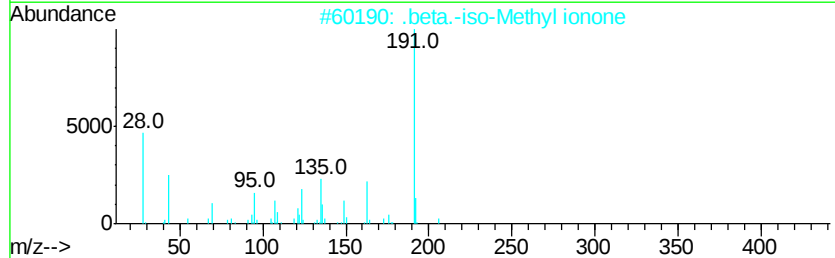
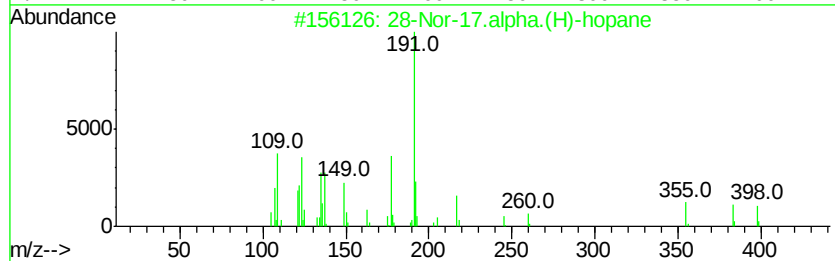
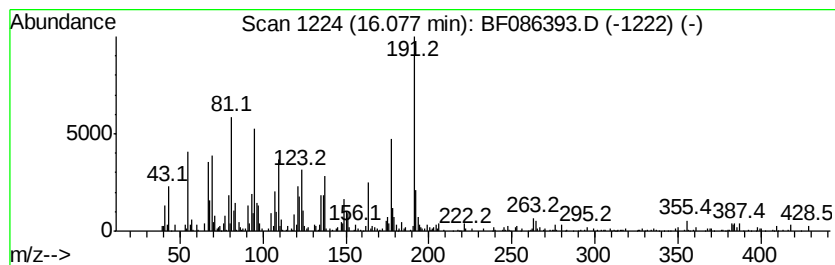
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 28-Nor-17.alpha.(H)-hopane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.08	16.00 ng	467085	Perylene-d12	15.41
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	28-Nor-17.alpha.(H)-hopane	398 C29H50	053584-60-4	72
2	.beta.-iso-Methyl ionone	206 C14H22O	1000285-40-2	43
3	1,4-Methanonaphthalene, 6,7-diet...	206 C15H26	016539-02-9	27
4	1-Penten-3-one, 1-(2,6,6-trimeth...	206 C14H22O	000127-43-5	25
5	2'-Fluoro-6'-(trifluoromethyl)-a...	206 C9H6F4O	174013-29-7	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
Data File : BF086393.D
Acq On : 23 Apr 2016 16:31
Operator : UM/SJ
Sample : H2640-05
Misc :
ALS Vial : 37 Sample Multiplier: 1

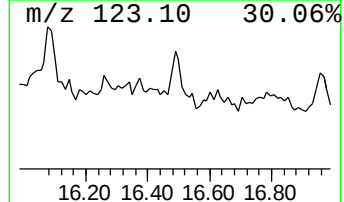
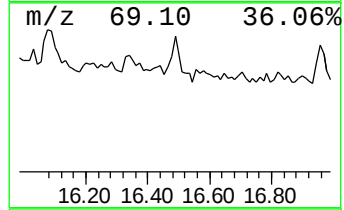
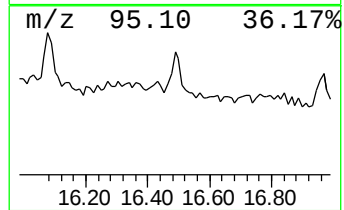
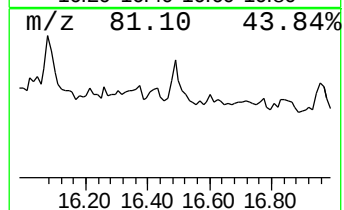
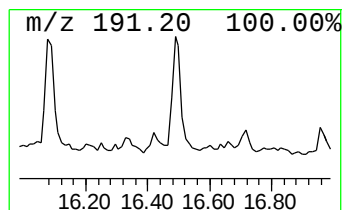
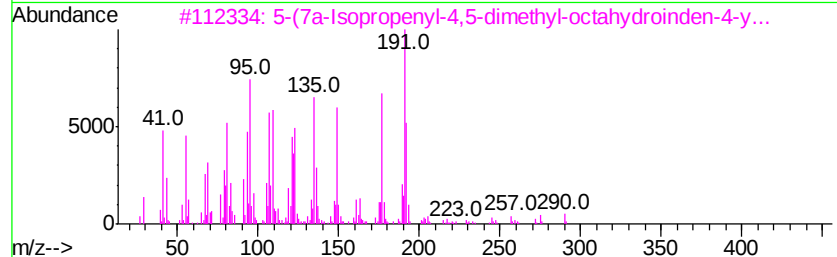
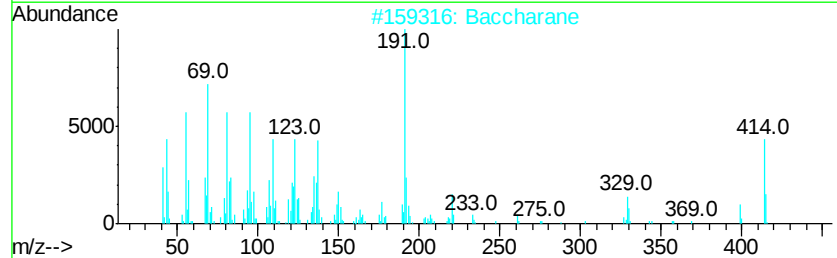
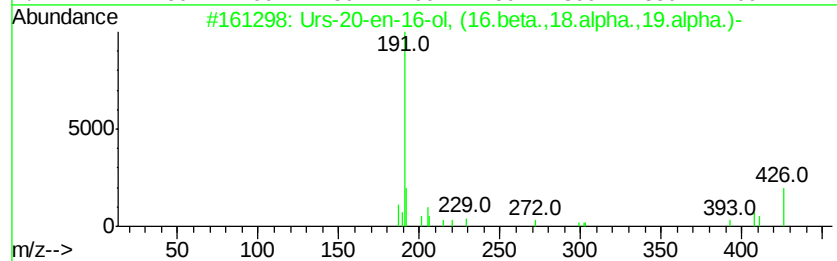
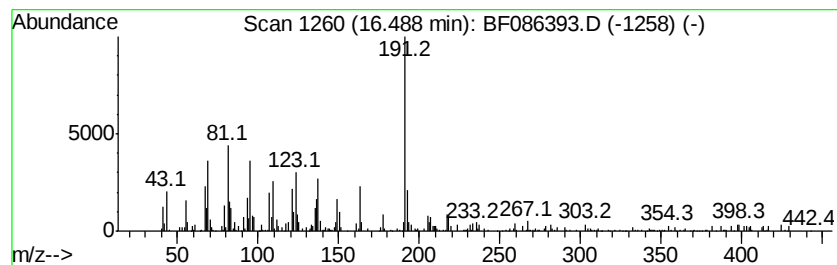
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Urs-20-en-16-ol, (16.beta.,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.49	7.24 ng	211333	Perylene-d12	15.41
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Urs-20-en-16-ol, (16.beta.,18.alpha.,19.alpha.)...	426 C30H50O	066394-74-9 83
2		Baccharane	414 C30H54	036441-74-4 47
3		5-(7a-Isopropenyl-4,5-dimethyl-octahydroinden-4-yl)-...	290 C20H34O	1000193-54-0 46
4		1-Penten-3-one, 1-(2,6,6-trimethyl-2-cyclohexenyl)-...	206 C14H22O	000127-43-5 46
5		1-Cyclohexene, 1,3,3-trimethyl-2-methyl-...	206 C14H22O	1000197-08-4 40



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
Data File : BF086393.D
Acq On : 23 Apr 2016 16:31
Operator : UM/SJ
Sample : H2640-05
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WPG-B13(0-5)-C

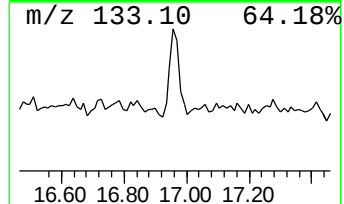
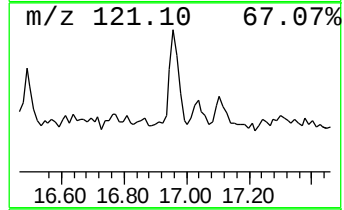
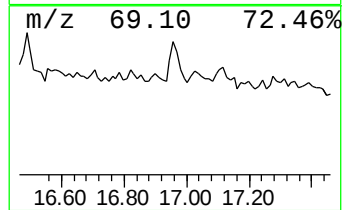
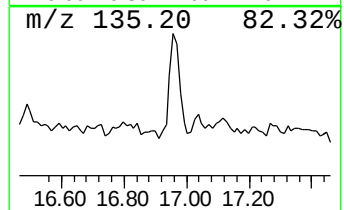
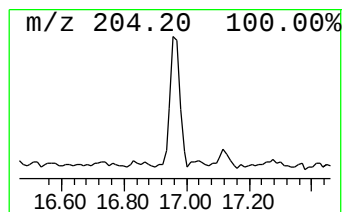
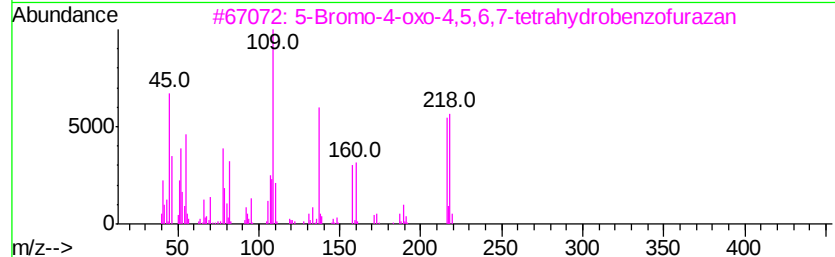
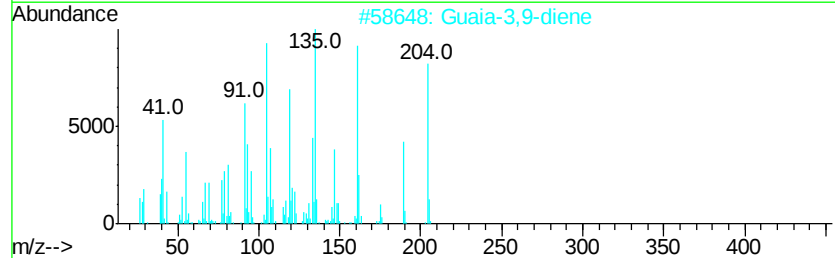
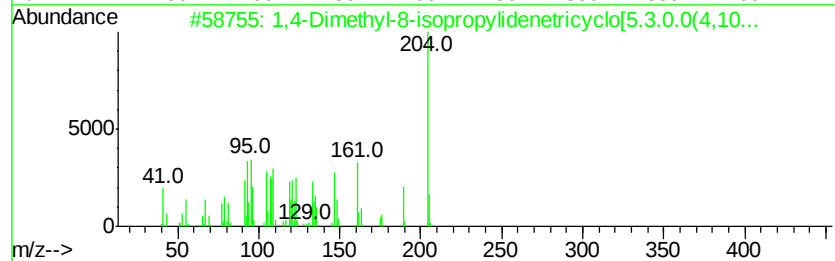
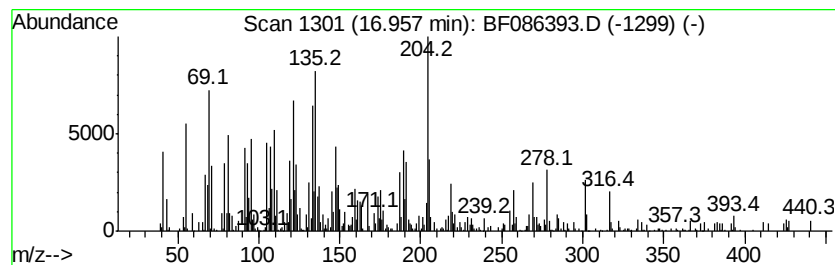
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 1,4-Dimethyl-8-isopropylide... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.96	11.85 ng	345908	Perylene-d12	15.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,4-Dimethyl-8-isopropylidenetri...	204	C15H24	1000140-07-7	78
2		Guaia-3,9-diene	204	C15H24	000489-83-8	58
3		5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	53
4		1,2,4,8-Tetramethylbicyclo[6.3.0]...	204	C15H24	137235-51-9	49
5		Cholest-8(14)-en-3-ol, 4,4-dimet...	414	C29H50O	014772-51-1	46



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
Data File : BF086393.D
Acq On : 23 Apr 2016 16:31
Operator : UM/SJ
Sample : H2640-05
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WPG-B13(0-5)-C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042216.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...	5.02	16.2	ng	877221	1	6.84	1084180	20.0
unknown6.61	6.61	79.8	ng	4327540	1	6.84	1084180	20.0
Hexadecane	10.32	3.3	ng	223559	3	9.88	1339320	20.0
Pentadecane, 2,6,...	10.53	2.4	ng	157835	3	9.88	1339320	20.0
Tridecane	10.78	4.8	ng	392805	4	11.36	1625520	20.0
Dodecane, 1-iodo-	11.23	2.7	ng	219409	4	11.36	1625520	20.0
4H-Cyclopenta[def...	11.97	2.1	ng	167206	4	11.36	1625520	20.0
Octadecane	13.19	2.8	ng	212043	5	14.00	1498470	20.0
Heneicosane	13.53	2.8	ng	207980	5	14.00	1498470	20.0
Eicosane	13.85	5.0	ng	373953	5	14.00	1498470	20.0
28-Nor-17.alpha.(...	16.08	16.0	ng	467085	6	15.41	583751	20.0
Urs-20-en-16-ol, ...	16.49	7.2	ng	211333	6	15.41	583751	20.0
1,4-Dimethyl-8-is...	16.96	11.9	ng	345908	6	15.41	583751	20.0