

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\
 Data File : BF093210.D
 Acq On : 27 Feb 2017 15:47
 Operator : SJ/MA
 Sample : I1972-02 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2B15-BASE-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-BF013017.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	4.373	198	200	204	rVB	43669	63018	1.92%	0.282%
2	4.978	250	253	262	rBV	239740	355287	10.83%	1.590%
3	5.356	284	286	290	rBV2	86097	145940	4.45%	0.653%
4	5.424	290	292	301	rVB	900995	931608	28.40%	4.169%
5	5.596	305	307	309	rBV	35527	36320	1.11%	0.163%
6	6.476	382	384	389	rBV	1067033	1038886	31.67%	4.650%
7	6.601	392	395	397	rBV	884322	1026997	31.30%	4.596%
8	6.830	413	415	417	rBV	1021855	985179	30.03%	4.409%
9	6.979	426	428	431	rVB	559933	671943	20.48%	3.007%
10	7.264	451	453	461	rVB	85370	107814	3.29%	0.483%
11	7.402	462	465	467	rBV	491977	524455	15.99%	2.347%
12	8.122	526	528	530	rBV	1477993	1361694	41.51%	6.094%
13	8.842	589	591	593	rBV	56423	43957	1.34%	0.197%
14	9.150	614	618	620	rBV2	22837	49567	1.51%	0.222%
15	9.207	620	623	625	rVV	1019823	917615	27.97%	4.107%
16	9.287	628	630	635	rVB2	26138	54985	1.68%	0.246%
17	9.470	643	646	649	rBV	30274	44868	1.37%	0.201%
18	9.539	651	652	654	rVV2	30383	40021	1.22%	0.179%
19	9.608	656	658	659	rVV	85318	70793	2.16%	0.317%
20	9.676	661	664	669	rBV4	29313	53722	1.64%	0.240%
21	9.745	669	670	672	rBV	27419	33621	1.02%	0.150%
22	9.813	674	676	678	rVV2	35463	42095	1.28%	0.188%
23	9.893	680	683	684	rVV	818683	1160790	35.38%	5.195%
24	9.916	684	685	687	rVB	77159	68027	2.07%	0.304%
25	10.088	698	700	703	rBV3	53275	77217	2.35%	0.346%
26	10.293	716	718	719	rBV2	21849	33487	1.02%	0.150%
27	10.316	719	720	722	rVB	62381	46670	1.42%	0.209%
28	10.430	728	730	733	rBV	65242	76548	2.33%	0.343%
29	10.533	737	739	743	rBV	57738	117535	3.58%	0.526%
30	10.682	749	752	754	rBV	327766	394208	12.02%	1.764%
31	10.796	760	762	766	rVB	153214	277802	8.47%	1.243%
32	10.979	776	778	779	rBV2	35935	50942	1.55%	0.228%
33	11.242	799	801	802	rBV	78872	89397	2.72%	0.400%
34	11.265	802	803	806	rVB	170100	167505	5.11%	0.750%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.379	810	813	817	rBV2	716358	1373350	41.86%	6.146%
36	11.448	817	819	821	rVB	103248	84256	2.57%	0.377%
37	11.482	821	822	823	rBV	39484	39477	1.20%	0.177%
38	11.619	832	834	836	rBV2	101381	144448	4.40%	0.646%
39	11.665	836	838	840	rVV	120717	118273	3.61%	0.529%
40	12.008	864	868	870	rVB3	300684	441660	13.46%	1.977%
41	12.111	873	877	880	rVB2	1695723	3280689	100.00%	14.683%
42	12.202	883	885	886	rBV	922730	832487	25.38%	3.726%
43	12.351	896	898	900	rBV	95406	101498	3.09%	0.454%
44	12.396	900	902	904	rVV	168574	200702	6.12%	0.898%
45	12.465	906	908	911	rVB3	106554	139185	4.24%	0.623%
46	12.591	917	919	921	rVB	501142	434164	13.23%	1.943%
47	12.819	936	939	942	rBV	446090	638236	19.45%	2.856%
48	12.911	945	947	948	rVV	140441	160865	4.90%	0.720%
49	12.956	949	951	955	rVB	595633	609192	18.57%	2.726%
50	13.151	966	968	970	rVB	285561	277896	8.47%	1.244%
51	14.008	1041	1043	1047	rBV2	748319	1219187	37.16%	5.456%
52	15.037	1131	1133	1137	rBV	175604	364146	11.10%	1.630%
53	15.448	1167	1169	1172	rBV	535455	723589	22.06%	3.238%

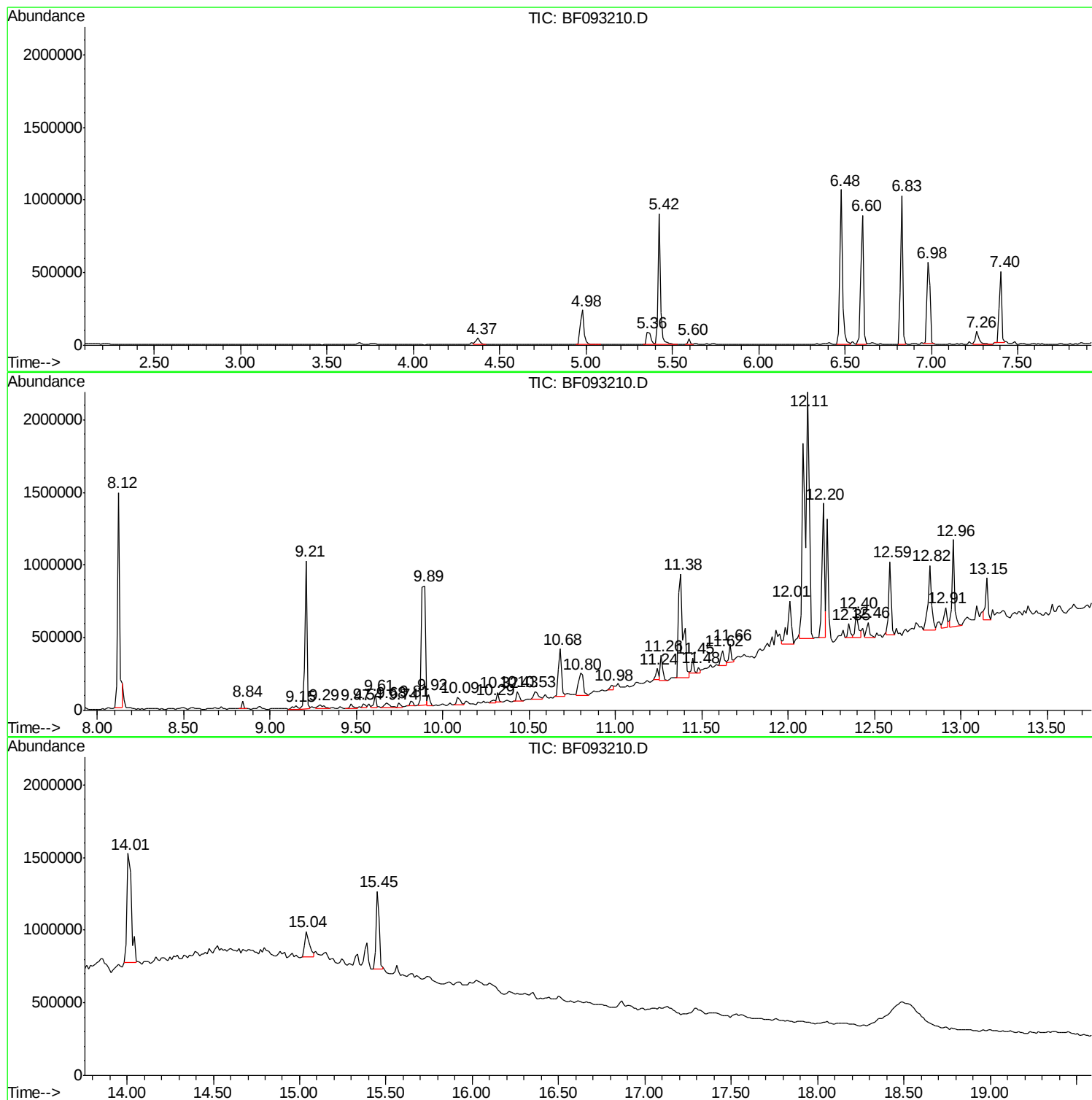
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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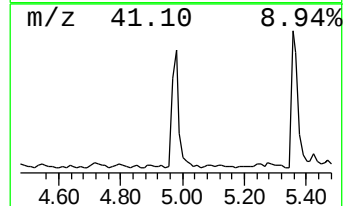
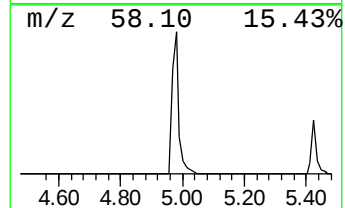
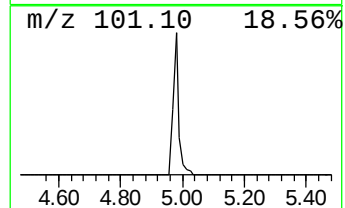
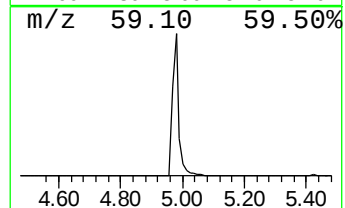
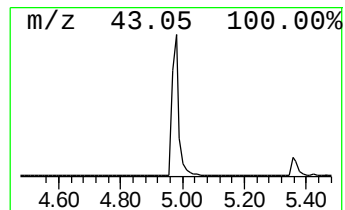
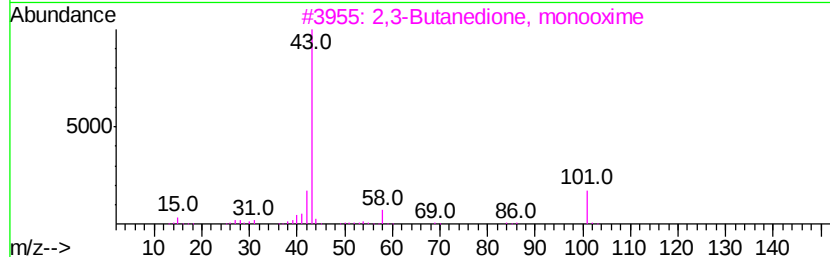
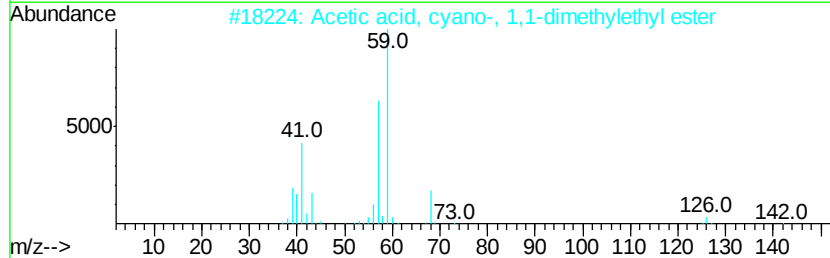
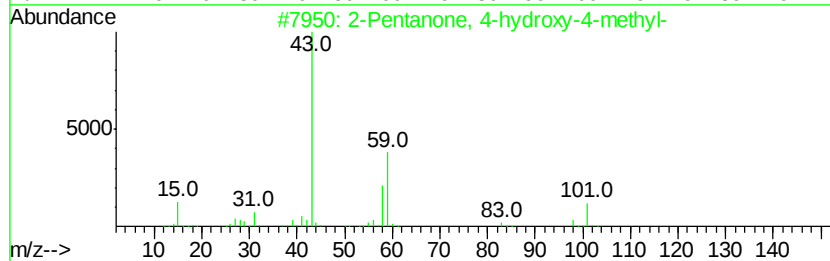
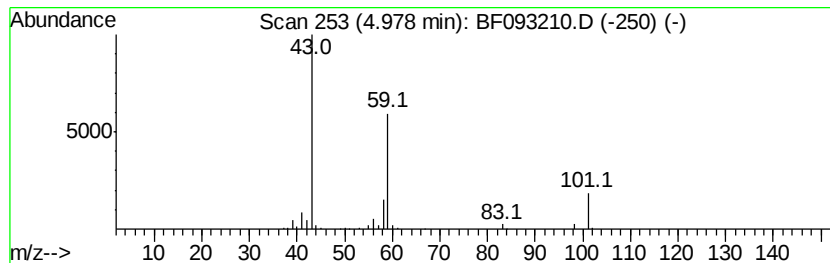
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.98	7.21 ng	355287	1,4-Dichlorobenzene-d4	6.83

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, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	37
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	9



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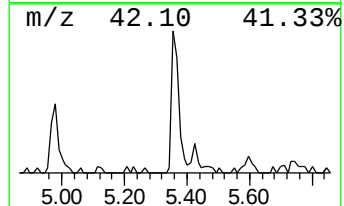
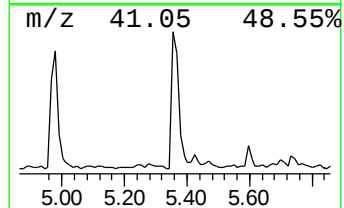
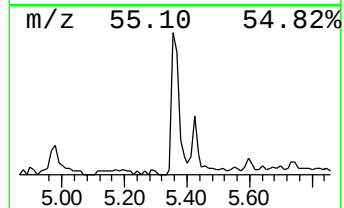
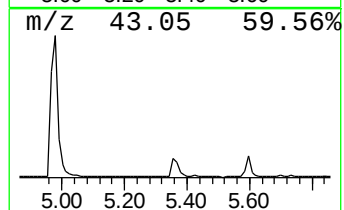
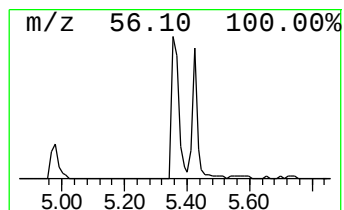
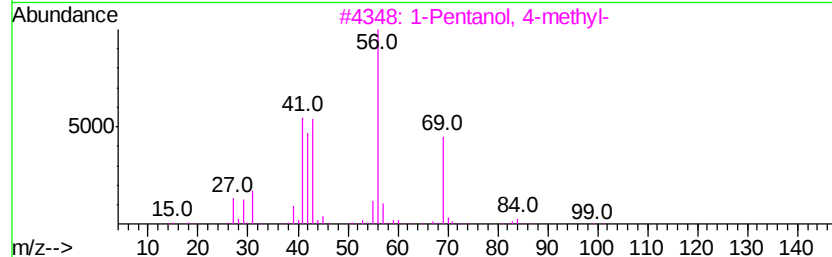
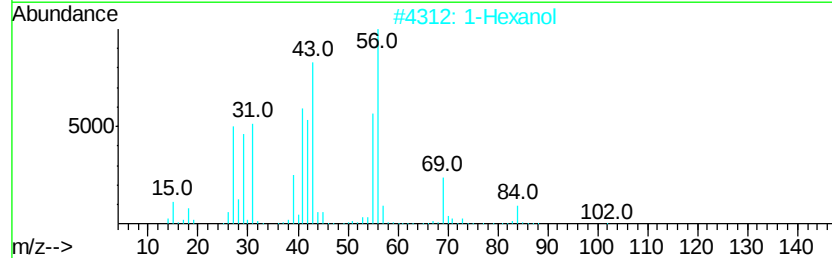
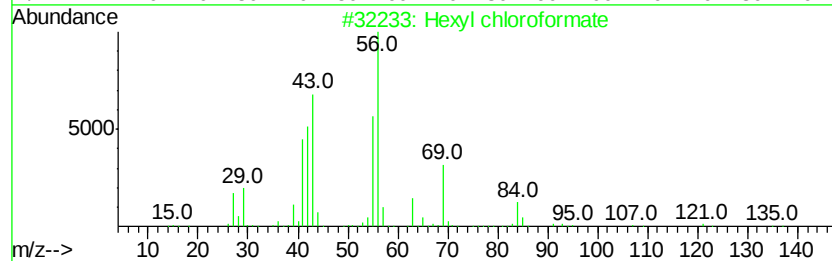
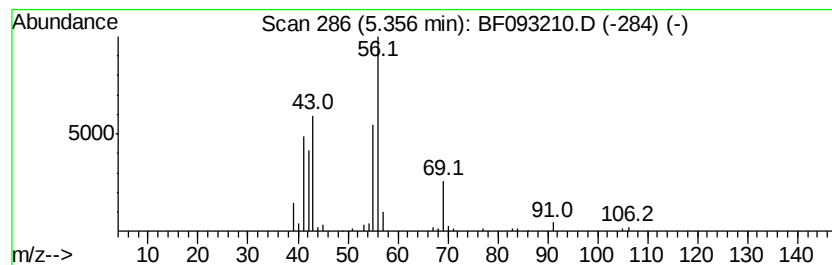
Instrument :
BNA_F
ClientSampled :
E2B15-BASE-2

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Hexyl chloroformate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
5.36	2.96 ng	145940	1,4-Dichlorobenzene-d4	6.83		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexyl chloroformate	164	C7H13ClO2	006092-54-2	78
2		1-Hexanol	102	C6H14O	000111-27-3	78
3		1-Pentanol, 4-methyl-	102	C6H14O	000626-89-1	72
4		1-Hexene	84	C6H12	000592-41-6	64
5		Formic acid, hexyl ester	130	C7H14O2	000629-33-4	56



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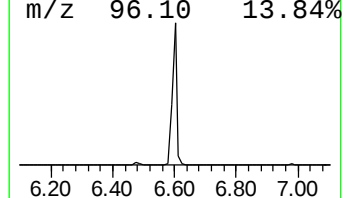
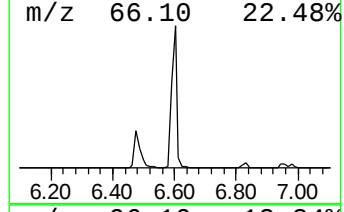
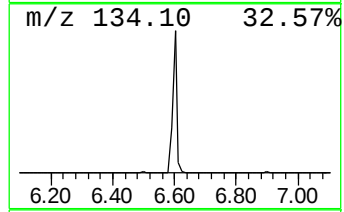
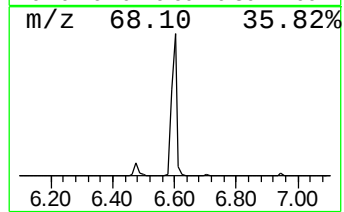
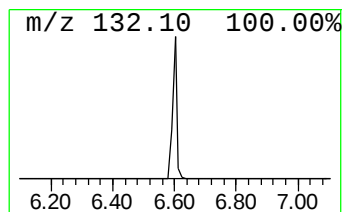
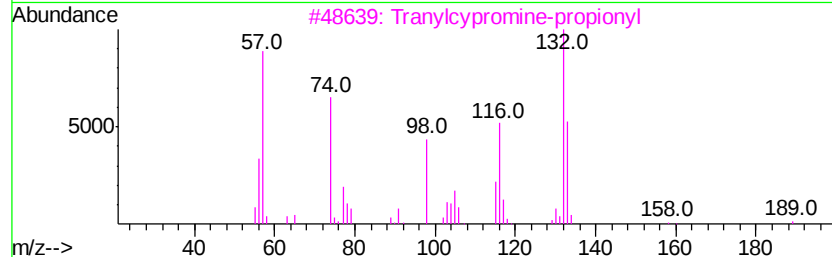
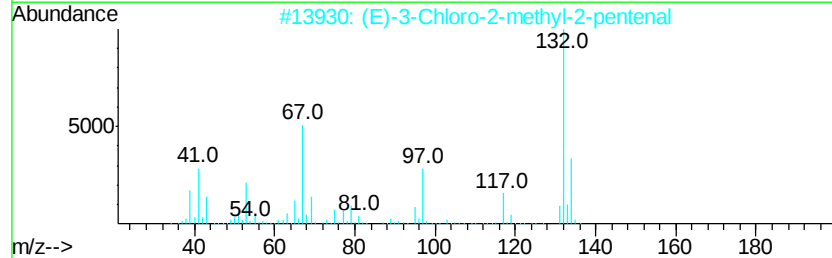
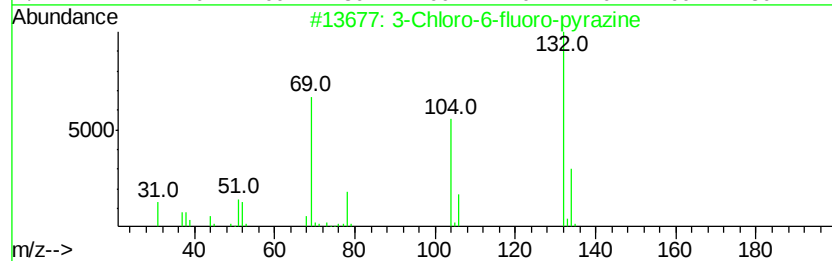
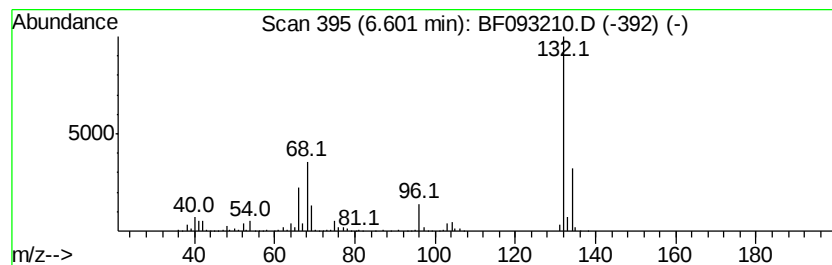
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.60 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.60	20.85 ng	1027000	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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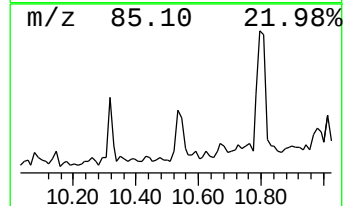
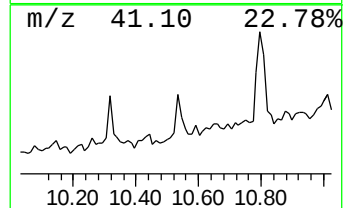
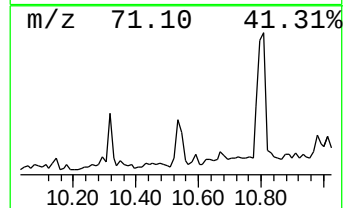
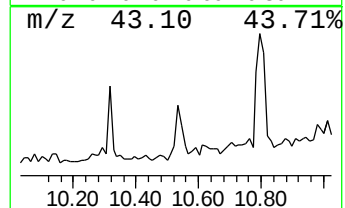
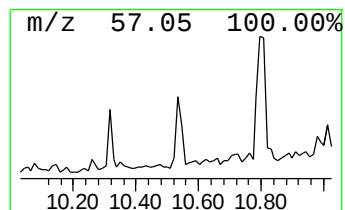
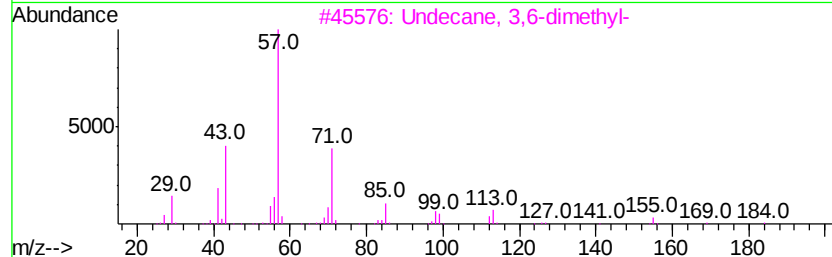
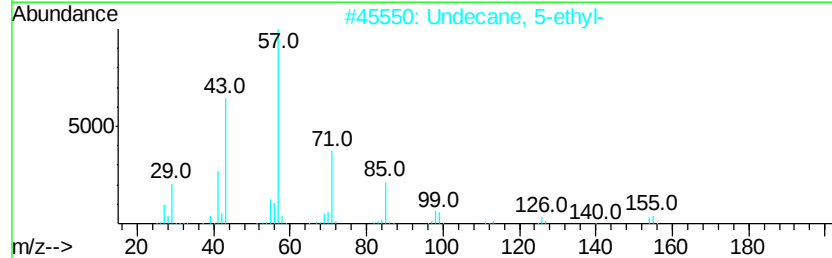
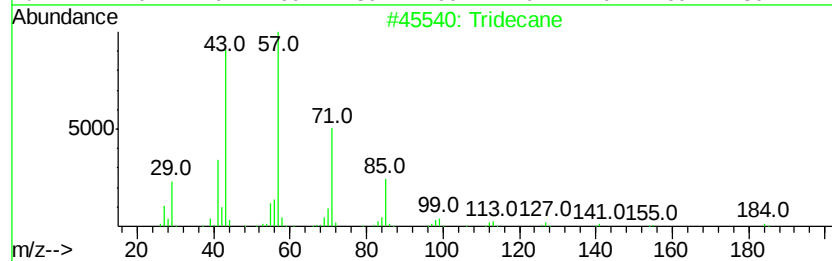
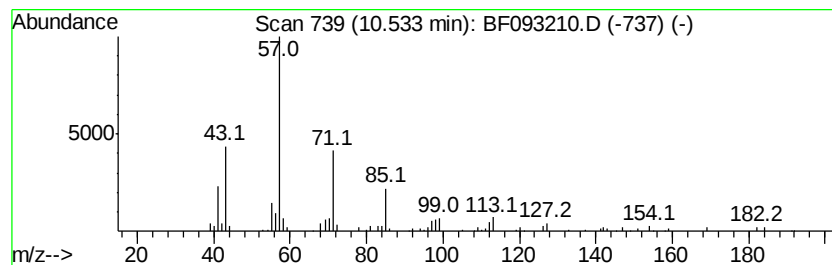
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Tridecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.53	2.03 ng	117535	Acenaphthene-d10		9.89
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	80
2	Undecane, 5-ethyl-	184	C13H28	017453-94-0	72
3	Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	64
4	Dodecane, 3-methyl-	184	C13H28	017312-57-1	64
5	Heptacosane	380	C27H56	000593-49-7	59



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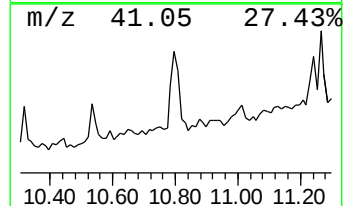
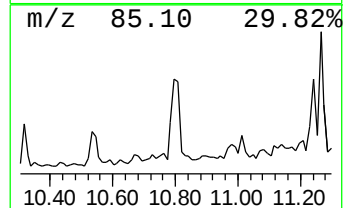
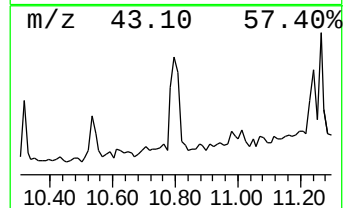
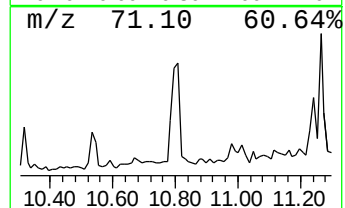
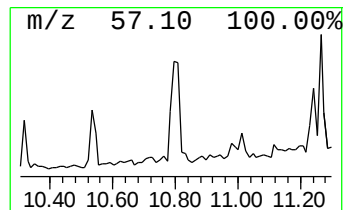
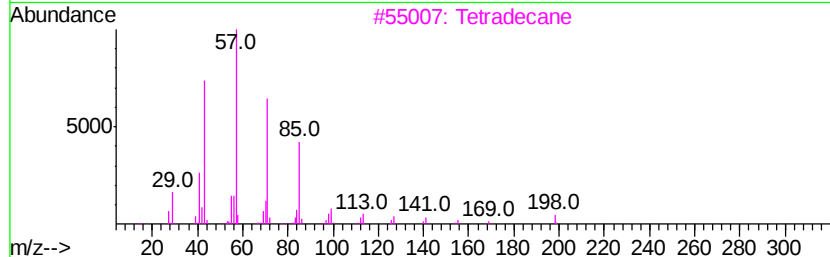
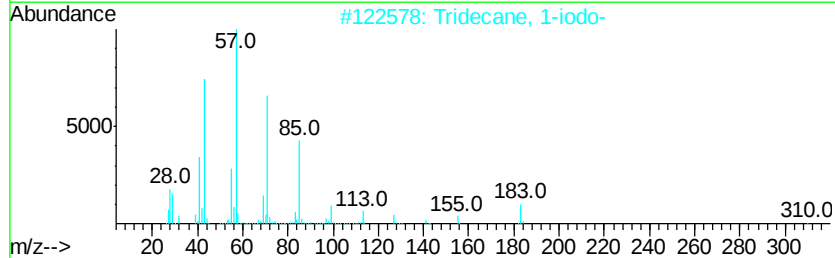
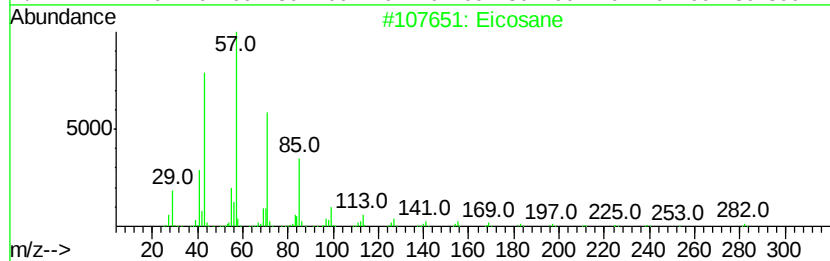
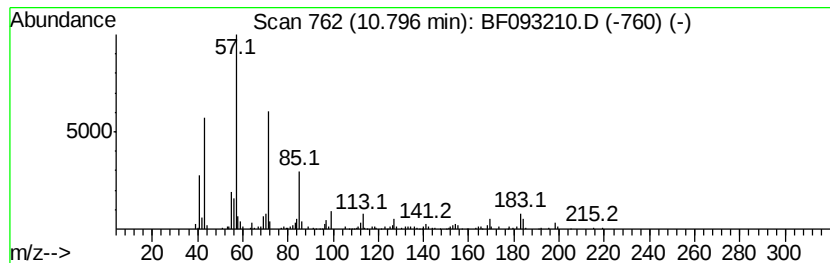
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Eicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.80	4.05 ng	277802	Phenanthrene-d10	11.38
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane	282 C20H42	000112-95-8	83
2	Tridecane, 1-iodo-	310 C13H27I	035599-77-0	80
3	Tetradecane	198 C14H30	000629-59-4	76
4	Heptadecane, 2,6,10,15-tetramethyl-	296 C21H44	054833-48-6	74
5	Pentadecane	212 C15H32	000629-62-9	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\
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Acq On : 27 Feb 2017 15:47
Operator : SJ/MA
Sample : I1972-02 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

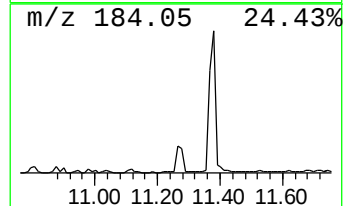
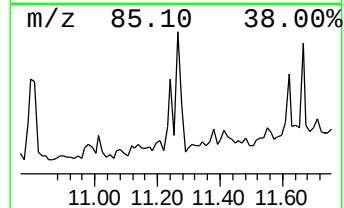
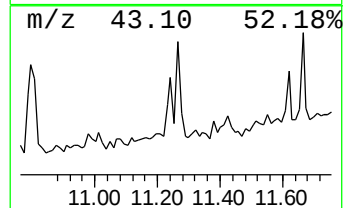
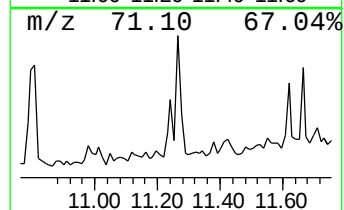
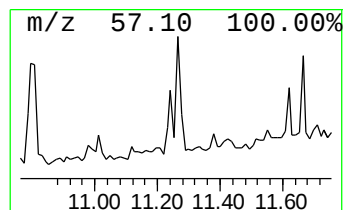
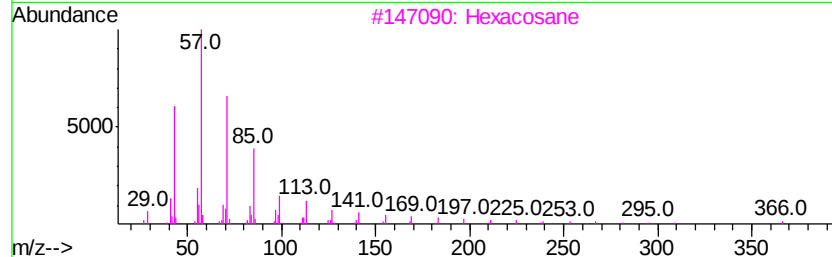
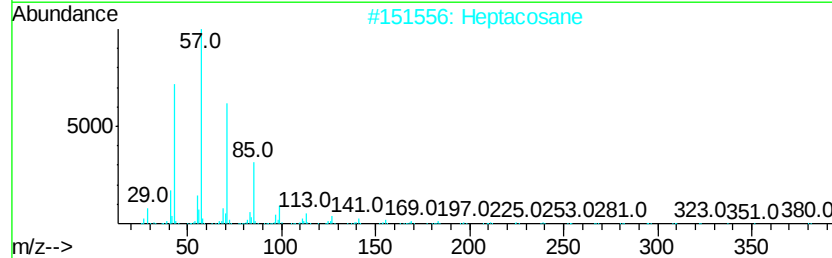
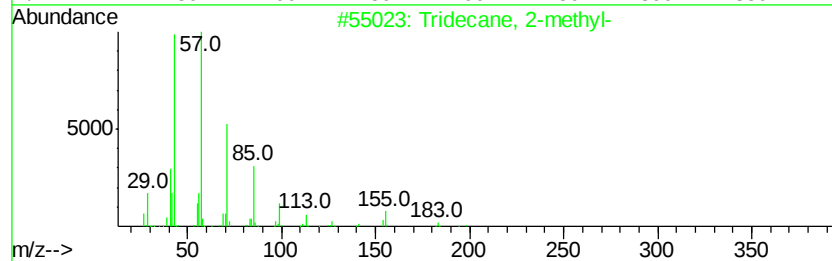
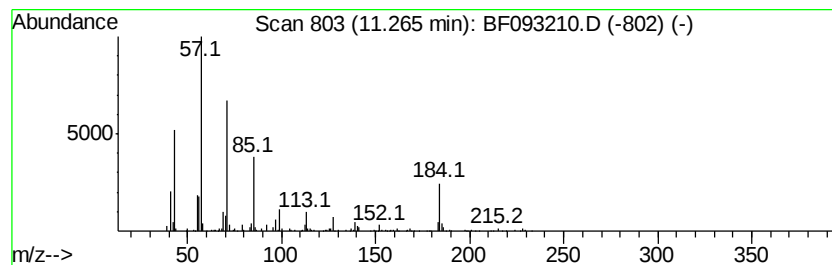
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tridecane, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.26	2.44 ng	167505	Phenanthrene-d10		11.38	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane, 2-methyl-	198	C14H30	001560-96-9	68
2		Heptacosane	380	C27H56	000593-49-7	64
3		Hexacosane	366	C26H54	000630-01-3	64
4		Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	59
5		3,5-Dimethyldodecane	198	C14H30	107770-99-0	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022717\
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ClientSampleId :
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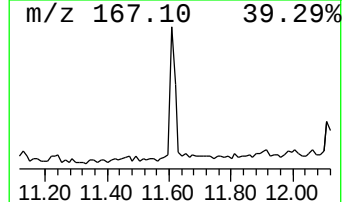
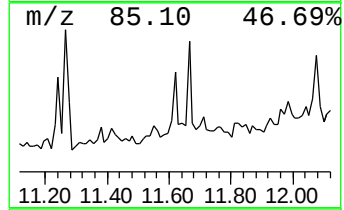
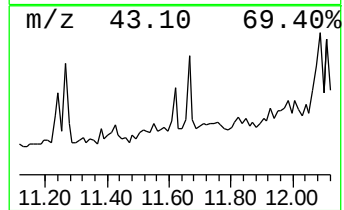
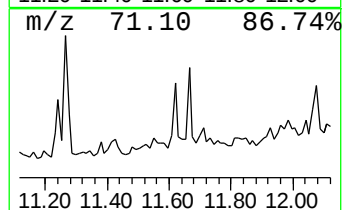
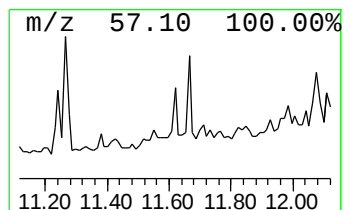
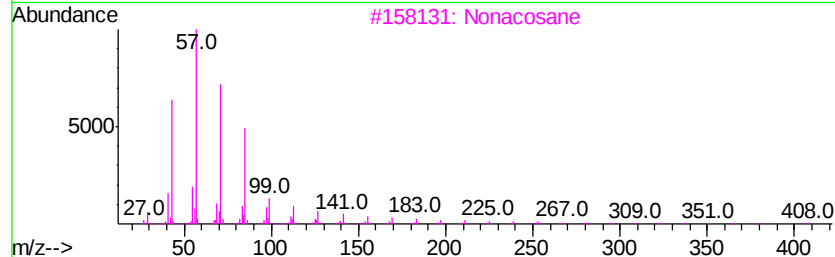
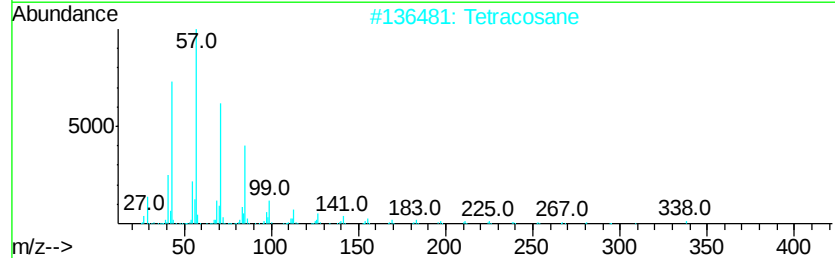
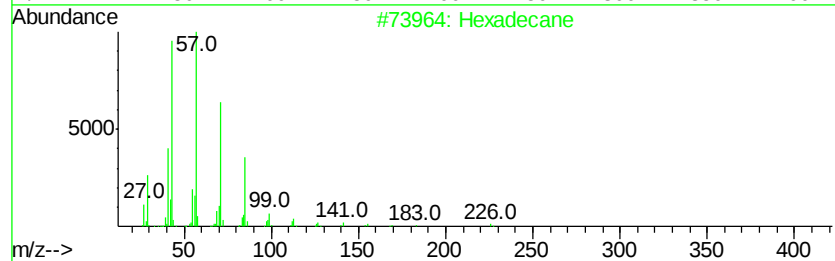
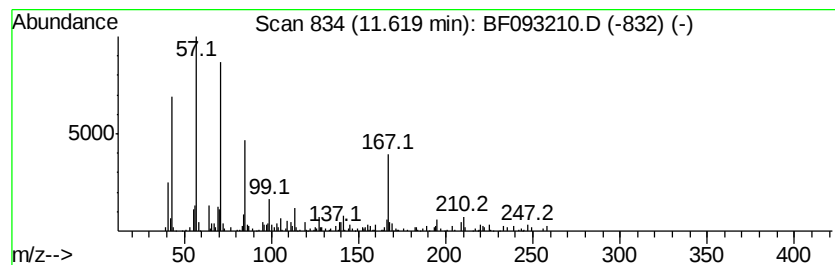
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.62	2.10 ng	144448	Phenanthrene-d10	11.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	70
2		Tetracosane	338	C24H50	000646-31-1	62
3		Nonacosane	408	C29H60	000630-03-5	58
4		Pentadecane	212	C15H32	000629-62-9	58
5		Heptacosane	380	C27H56	000593-49-7	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022717\
Data File : BF093210.D
Acq On : 27 Feb 2017 15:47
Operator : SJ/MA
Sample : I1972-02 5X
Misc :
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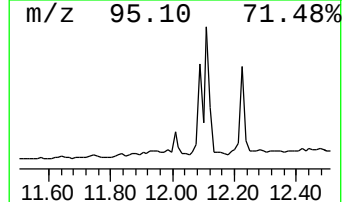
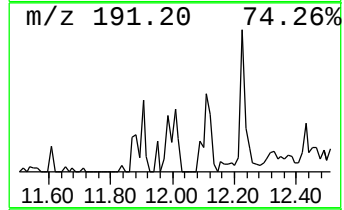
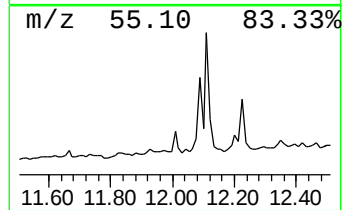
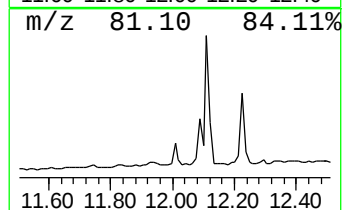
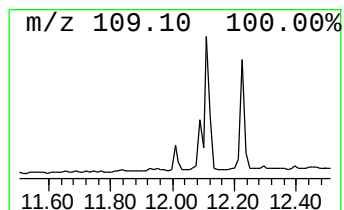
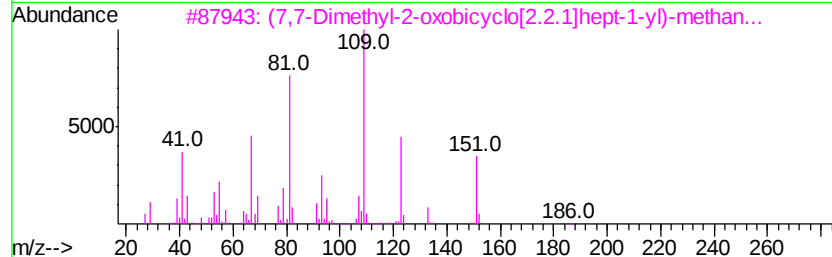
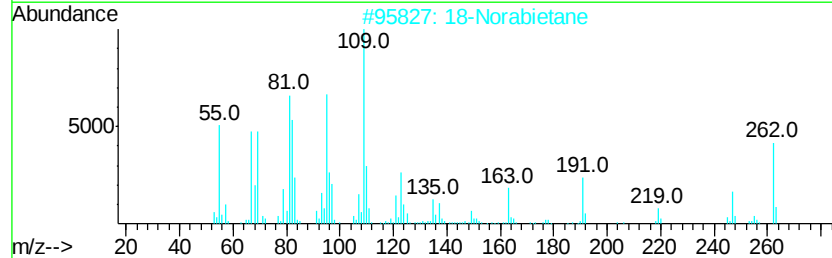
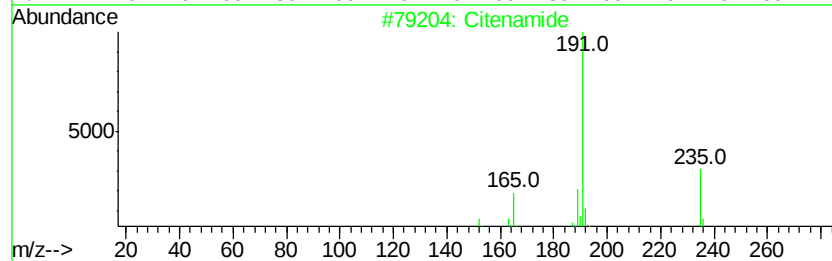
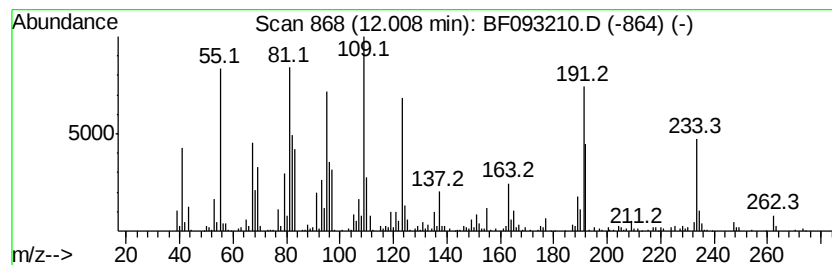
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown12.01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	6.43 ng	441660	Phenanthrene-d10	11.38

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Citenamide	235	C16H13NO	010423-37-7	35
2			18-Norabietane	262	C19H34	1000293-16-6	30
3			(7,7-Dimethyl-2-oxobicyclo[2.2.1...]	250	C10H15ClO3S	1000194-76-1	27
4			4-Triethylsilylbut-1-en-3-yne	166	C10H18Si	001693-70-5	27
5			Imidazole, 4-methyl-5-[3,3,3-tri...	234	C10H13F3N2O	1000129-15-8	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022717\
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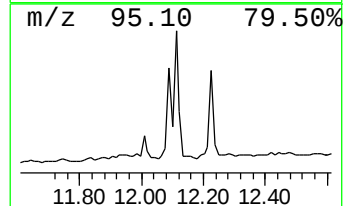
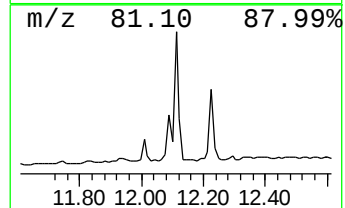
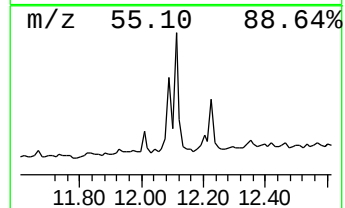
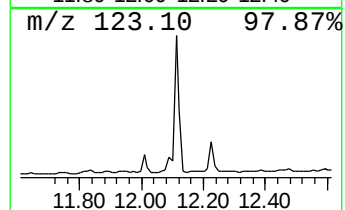
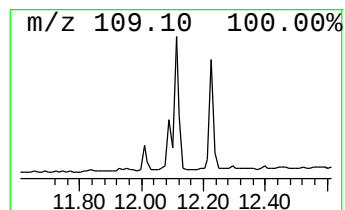
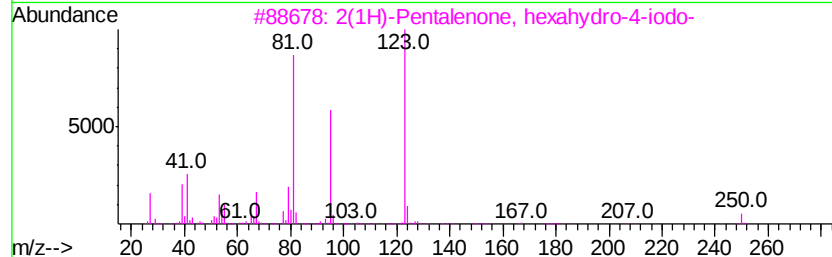
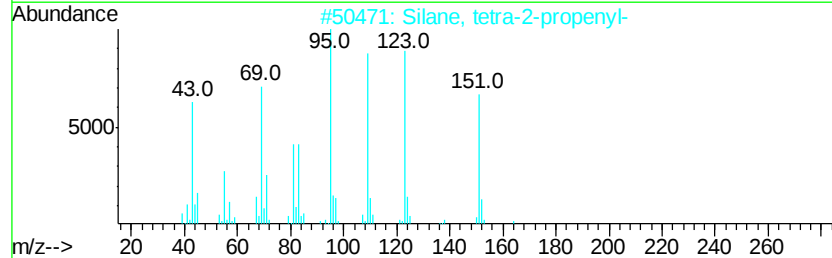
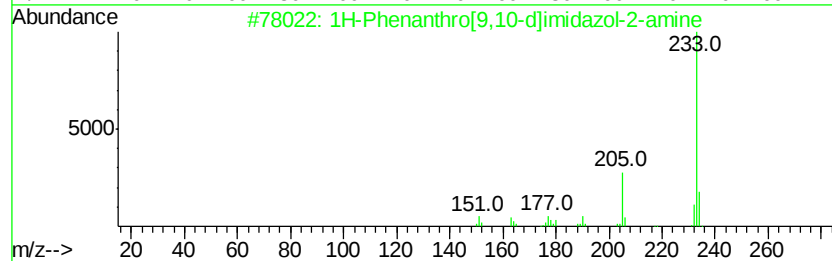
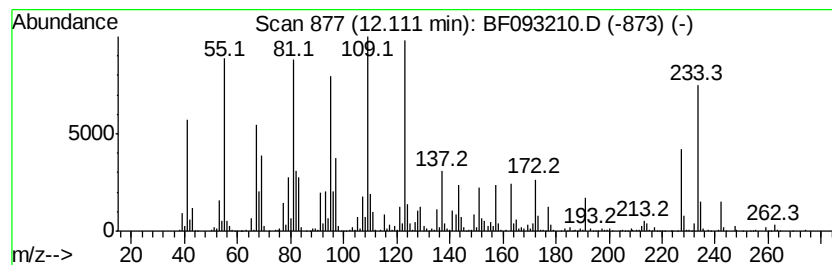
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown12.11 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.11	47.78 ng	3280690	Phenanthrene-d10	11.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Phenanthro[9,10-d]imidazol-2-...	233	C15H11N3	037052-13-4	38
2		Silane, tetra-2-propenyl-	192	C12H20Si	001112-66-9	35
3		2(1H)-Pentalenone, hexahydro-4-i...	250	C8H11IO	077070-56-5	35
4		Bicyclo[2.2.1]heptane, 1,3,3-tri...	138	C10H18	006248-88-0	30
5		1-(1-Ethyl-2,3-dimethyl-cyclopen...	166	C11H18O	1000185-18-6	27



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\
Data File : BF093210.D
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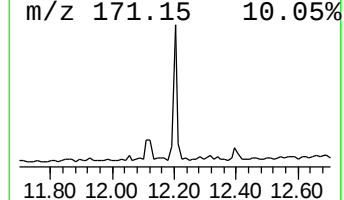
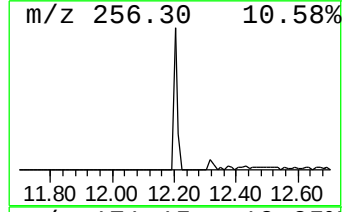
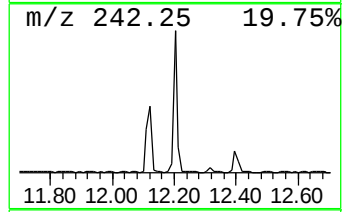
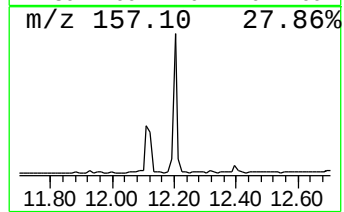
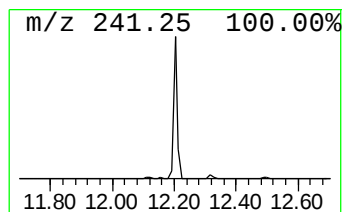
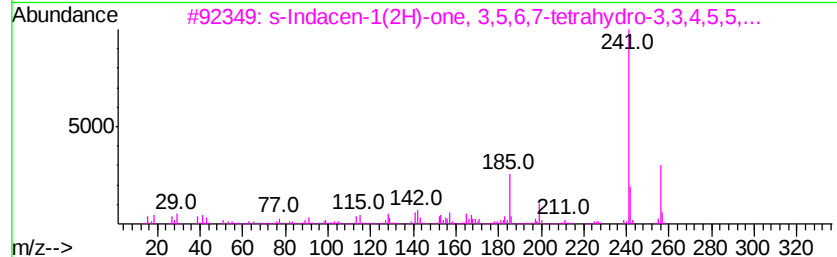
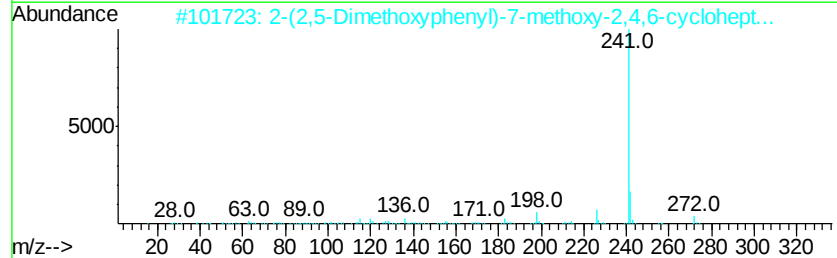
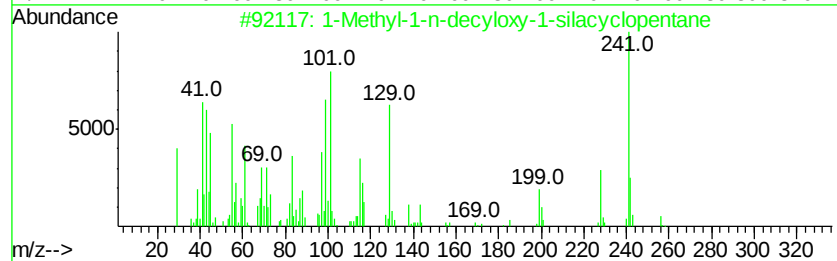
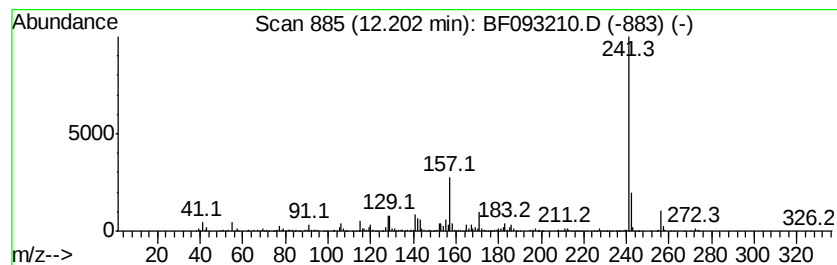
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1-Methyl-1-n-decyloxy-1-sil... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.20	12.12 ng	832487	Phenanthrene-d10	11.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Methyl-1-n-decyloxy-1-silacycl...	256	C15H32OSi	1000216-91-2	53
2		2-(2,5-Dimethoxyphenyl)-7-methox...	272	C16H16O4	003525-17-5	42
3		s-Indacen-1(2H)-one, 3,5,6,7-tet...	256	C18H24O	038754-94-8	42
4		1,3,5,7-Tetrasilatricyclo[3.3.1....	256	C10H24Si4	017995-33-4	40
5		Dibenzo(b,def)carbazole	241	C18H11N	104313-09-9	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\
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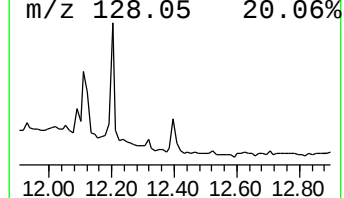
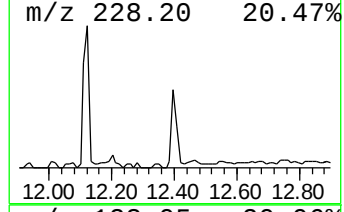
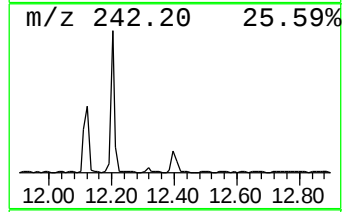
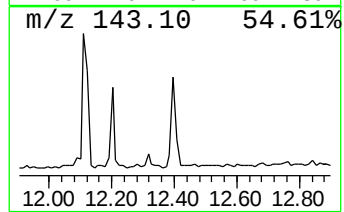
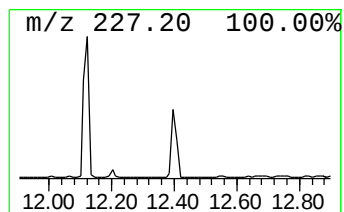
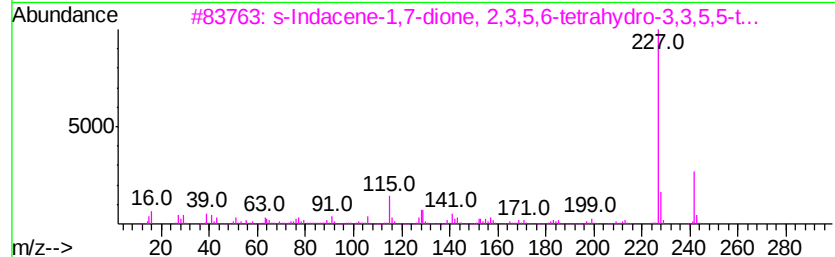
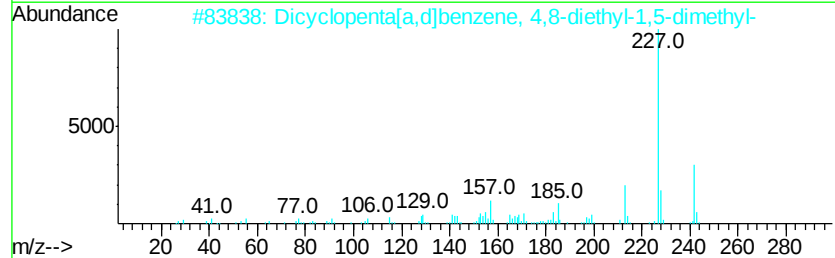
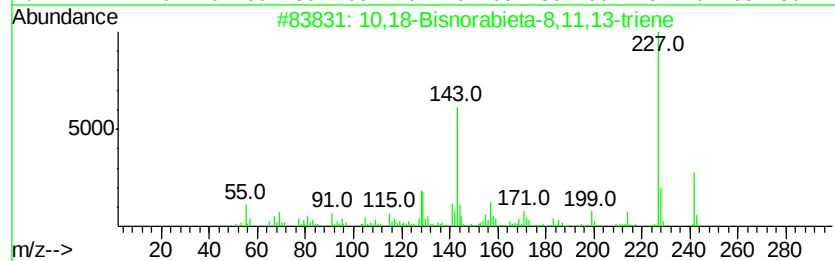
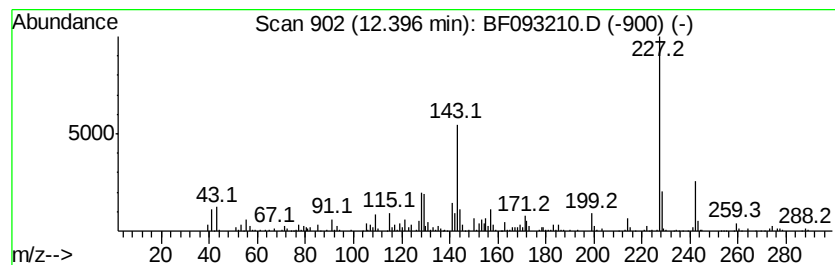
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 10,18-Bisnorabieta-8,11,13-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.40	2.92 ng	200702	Phenanthrene-d10	11.38	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,18-Bisnorabieta-8,11,13-triene	242	C18H26	032624-67-2	95
2	Dicyclopenta[a,d]benzene, 4,8-di...	242	C18H26	1000156-41-6	47
3	s-Indacene-1,7-dione, 2,3,5,6-te...	242	C16H18O2	055591-17-8	46
4	5-Methoxy-2-methyl-4-oxo-1,2,3,4...	242	C14H14N2O2	1000227-06-0	43
5	2-(4'-Hydroxyphenyl)-2-(4'-metho...	242	C16H18O2	016530-58-8	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022717\
Data File : BF093210.D
Acq On : 27 Feb 2017 15:47
Operator : SJ/MA
Sample : I1972-02 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

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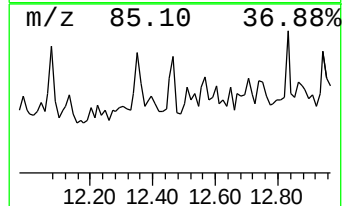
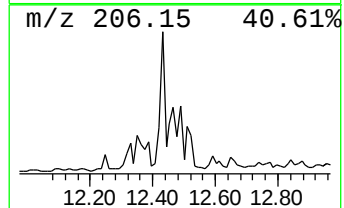
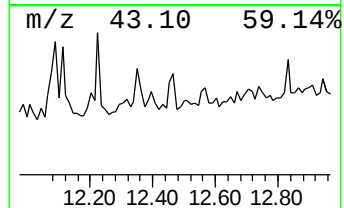
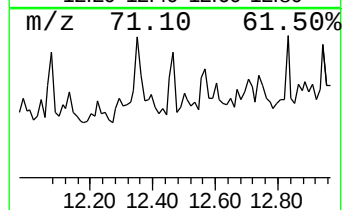
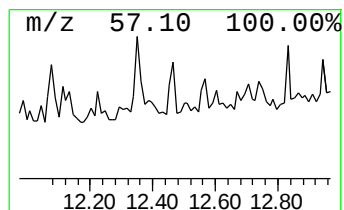
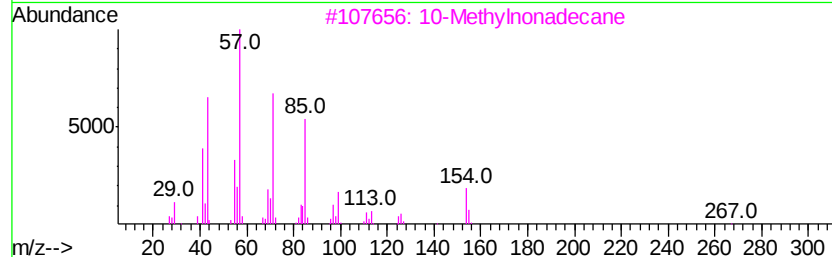
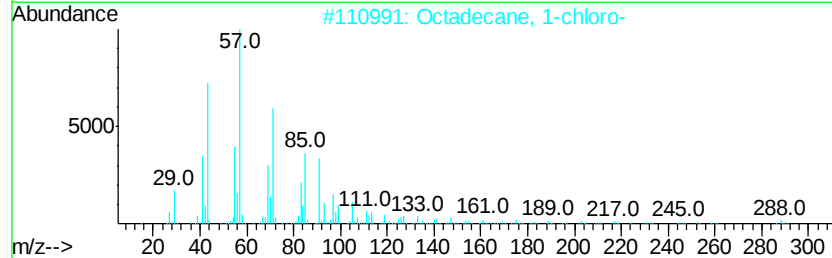
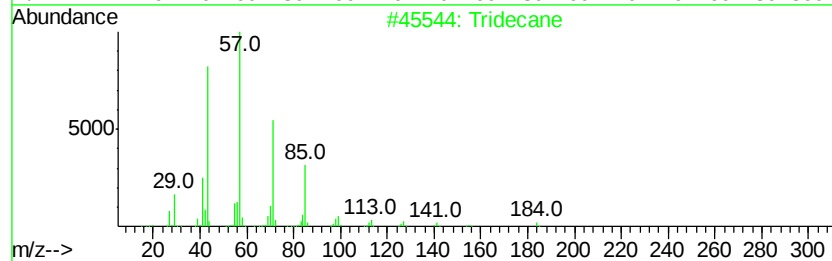
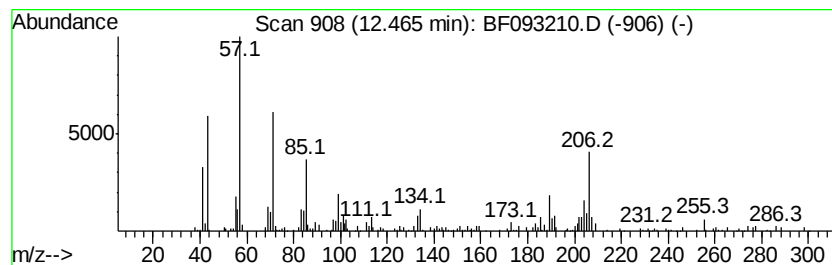
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown12.47 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.47	2.03 ng	139185	Phenanthrene-d10	11.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane	184	C13H28	000629-50-5	42
2		Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	42
3		10-Methylnonadecane	282	C20H42	056862-62-5	38
4		Heptadecane	240	C17H36	000629-78-7	35
5		Tetradecane	198	C14H30	000629-59-4	30



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\
Data File : BF093210.D
Acq On : 27 Feb 2017 15:47
Operator : SJ/MA
Sample : I1972-02 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
E2B15-BASE-2

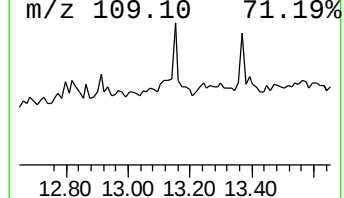
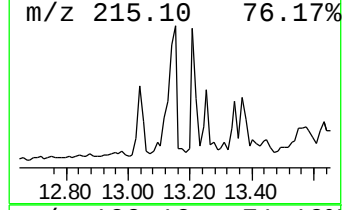
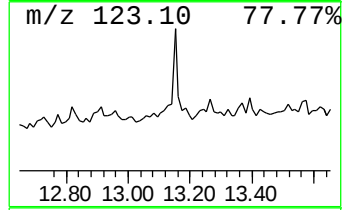
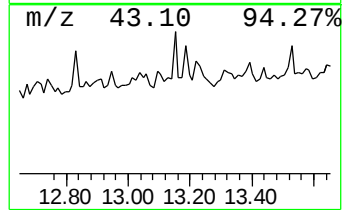
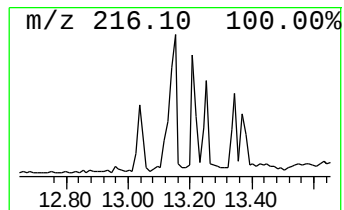
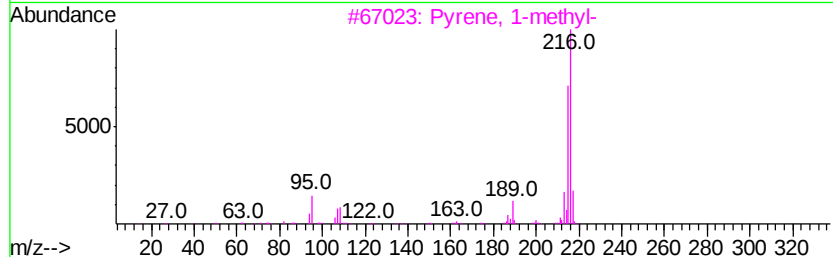
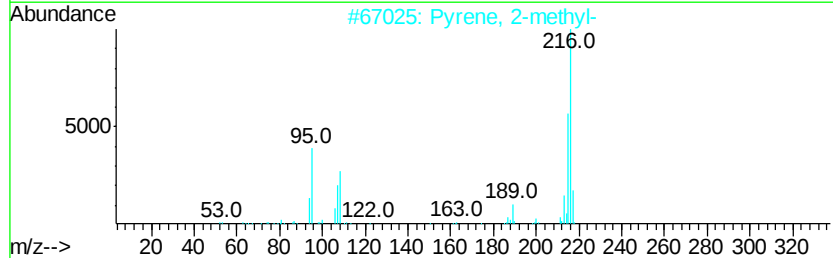
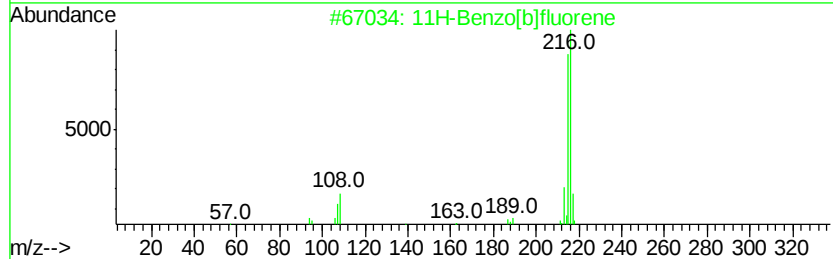
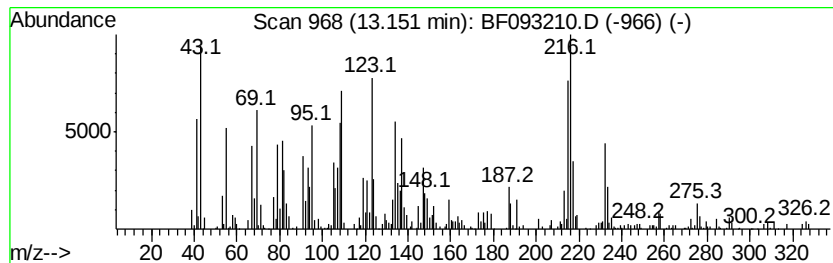
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown13.15 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.15	4.56 ng	277896	Chrysene-d12	14.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	49
2		Pyrene, 2-methyl-	216	C17H12	003442-78-2	46
3		Pyrene, 1-methyl-	216	C17H12	002381-21-7	38
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	30
5		3-Phenanthrenol, tetradecahydro-...	250	C17H30O	054155-13-4	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022717\
 Data File : BF093210.D
 Acq On : 27 Feb 2017 15:47
 Operator : SJ/MA
 Sample : I1972-02 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2B15-BASE-2

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.98	7.2 ng		355287	1	6.83	985179	20.0
Hexyl chloroformate	5.36	3.0 ng		145940	1	6.83	985179	20.0
unknown6.60	6.60	20.9 ng		1027000	1	6.83	985179	20.0
Tridecane	10.53	2.0 ng		117535	3	9.89	1160790	20.0
Eicosane	10.80	4.0 ng		277802	4	11.38	1373350	20.0
Tridecane, 2-methyl-	11.26	2.4 ng		167505	4	11.38	1373350	20.0
Hexadecane	11.62	2.1 ng		144448	4	11.38	1373350	20.0
unknown12.01	12.01	6.4 ng		441660	4	11.38	1373350	20.0
unknown12.11	12.11	47.8 ng		3280690	4	11.38	1373350	20.0
1-Methyl-1-n-decy...	12.20	12.1 ng		832487	4	11.38	1373350	20.0
10,18-Bisnorabiet...	12.40	2.9 ng		200702	4	11.38	1373350	20.0
unknown12.47	12.47	2.0 ng		139185	4	11.38	1373350	20.0
unknown13.15	13.15	4.6 ng		277896	5	14.02	1219190	20.0