

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031616\
 Data File : BF085495.D
 Acq On : 17 Mar 2016 7:08
 Operator : UM/SJ
 Sample : H1768-02
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AG-1A

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-BF031516.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.584	25	26	27	rVB	61542	42187	1.04%	0.127%
2	2.653	30	32	33	rBV2	146013	106873	2.64%	0.322%
3	2.927	53	56	57	rBV	37560	43174	1.06%	0.130%
4	2.961	57	59	60	rVV	71783	62149	1.53%	0.187%
5	2.996	60	62	64	rVB	184522	141167	3.48%	0.425%
6	3.041	64	66	68	rVB3	156391	158013	3.90%	0.476%
7	3.384	93	96	99	rVB	25738	41678	1.03%	0.125%
8	3.567	110	112	114	rVV2	47823	43883	1.08%	0.132%
9	3.876	137	139	143	rBV	126366	212873	5.25%	0.641%
10	4.596	200	202	203	rBV	56340	40905	1.01%	0.123%
11	4.801	216	220	222	rVB2	90028	85186	2.10%	0.256%
12	5.110	245	247	250	rBV	60545	104618	2.58%	0.315%
13	5.499	278	281	284	rVV	2622944	3822502	94.28%	11.506%
14	6.379	354	358	360	rBV	33677	57786	1.43%	0.174%
15	6.424	360	362	364	rBV2	42989	45922	1.13%	0.138%
16	6.539	368	372	374	rVB	3041078	4054492	100.00%	12.204%
17	6.664	380	383	386	rBV	2750502	3968097	97.87%	11.944%
18	6.904	401	404	406	rBV	369060	458502	11.31%	1.380%
19	7.053	415	417	420	rBV	2321665	3018442	74.45%	9.086%
20	7.430	449	450	451	rBV	37782	45476	1.12%	0.137%
21	7.465	451	453	456	rVV	2223372	2338260	57.67%	7.038%
22	7.556	459	461	466	rVB	87330	131862	3.25%	0.397%
23	8.185	514	516	518	rBV	705887	581962	14.35%	1.752%
24	8.299	521	526	528	rVB2	33663	52307	1.29%	0.157%
25	8.447	537	539	541	rBV3	75641	57888	1.43%	0.174%
26	9.259	608	610	613	rBV	2946916	3927965	96.88%	11.823%
27	9.659	643	645	646	rBV	807661	840051	20.72%	2.529%
28	9.876	662	664	666	rVB	69170	83923	2.07%	0.253%
29	9.945	668	670	672	rVB	414675	557492	13.75%	1.678%
30	10.368	706	707	709	rVB	144147	125880	3.10%	0.379%
31	10.585	723	726	730	rBV	51462	89922	2.22%	0.271%
32	10.733	736	739	741	rBV	2345676	2549539	62.88%	7.674%
33	10.848	747	749	755	rBV	107111	191011	4.71%	0.575%
34	11.031	764	765	770	rBV3	18581	42880	1.06%	0.129%

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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	11.293	786	788	789 rBV	77606	80413	1.98%	0.242%
36	11.419	797	799	802 rBV	514763	525958	12.97%	1.583%
37	11.648	816	819	823 rBV	72920	94427	2.33%	0.284%
38	11.716	823	825	827 rVB	56081	46325	1.14%	0.139%
39	11.945	844	845	851 rBV	115844	143155	3.53%	0.431%
40	12.631	903	905	910 rBV	52663	53469	1.32%	0.161%
41	12.859	923	925	928 rVB	45790	50074	1.24%	0.151%
42	13.008	935	938	940 rBV	3020618	2742205	67.63%	8.254%
43	13.899	1014	1016	1025 rBV	194178	274280	6.76%	0.826%
44	14.059	1027	1030	1032 rBV	347758	377875	9.32%	1.137%
45	14.448	1057	1064	1065 rBV4	13266	41174	1.02%	0.124%
46	14.528	1069	1071	1078 rVB3	23252	54939	1.36%	0.165%
47	15.088	1117	1120	1123 rBV	27204	66773	1.65%	0.201%
48	15.500	1153	1156	1159 rBV	297319	402981	9.94%	1.213%
49	17.706	1345	1349	1357 rVB	20128	65998	1.63%	0.199%
50	18.517	1415	1420	1427 rBV2	20555	77342	1.91%	0.233%

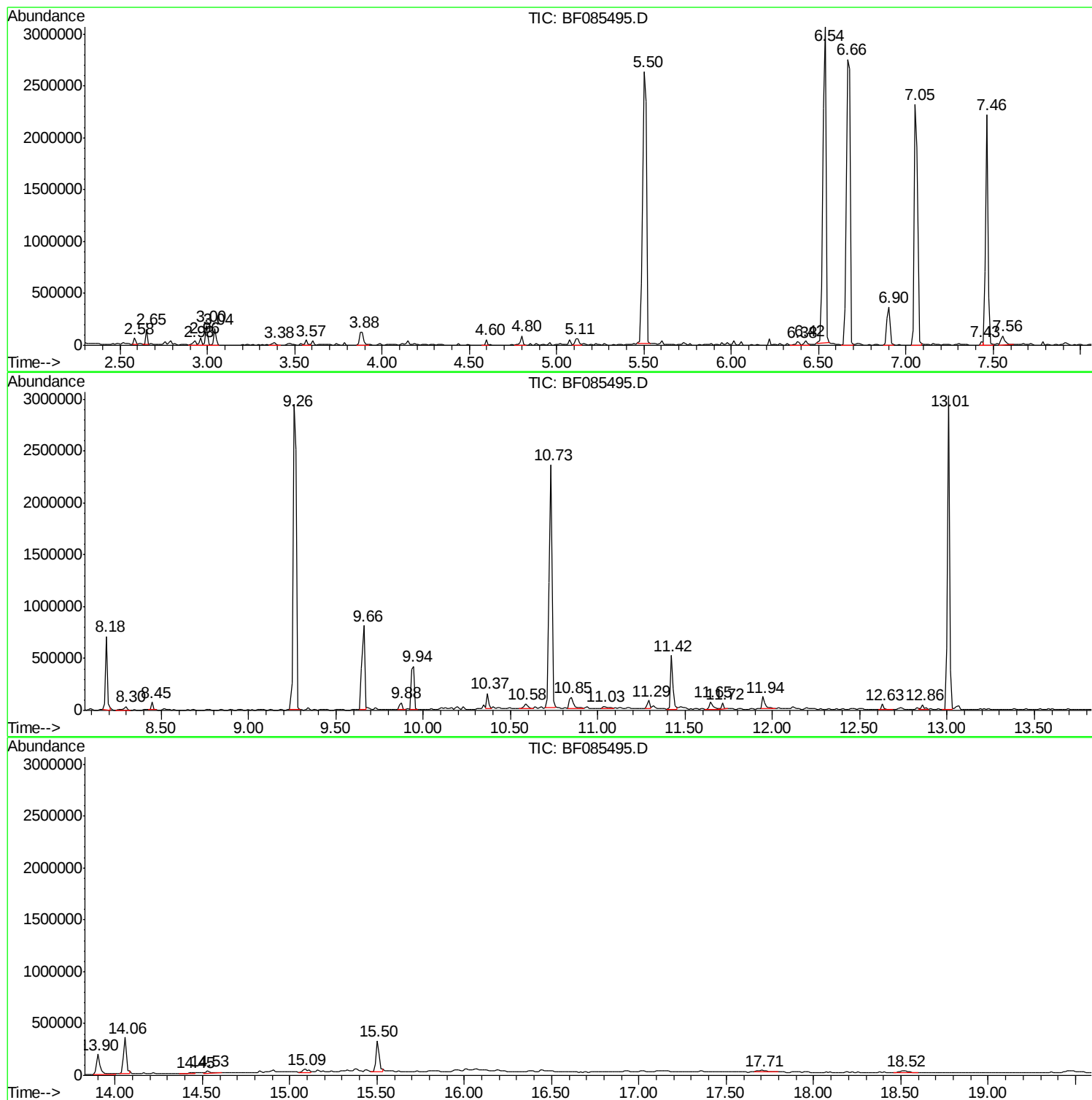
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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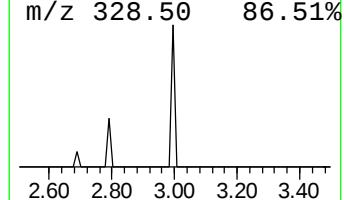
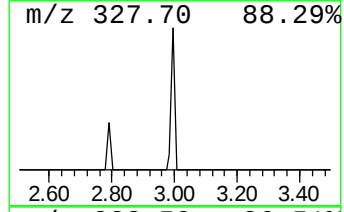
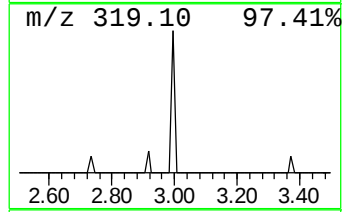
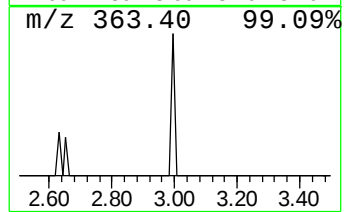
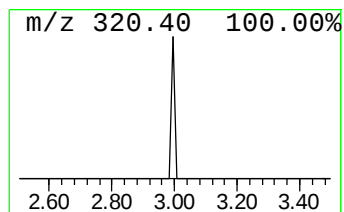
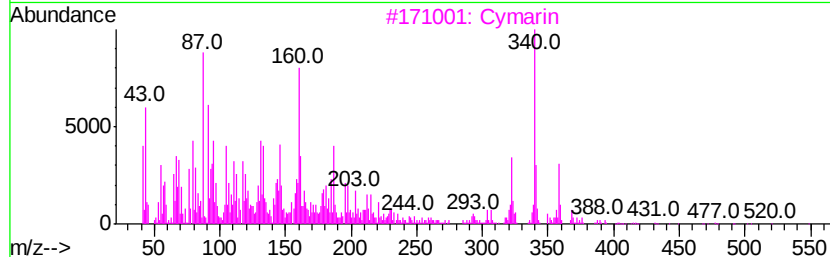
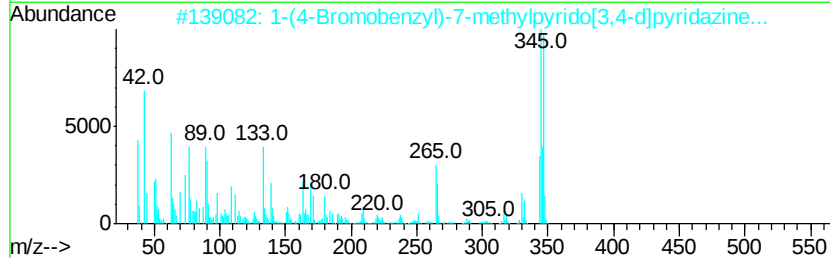
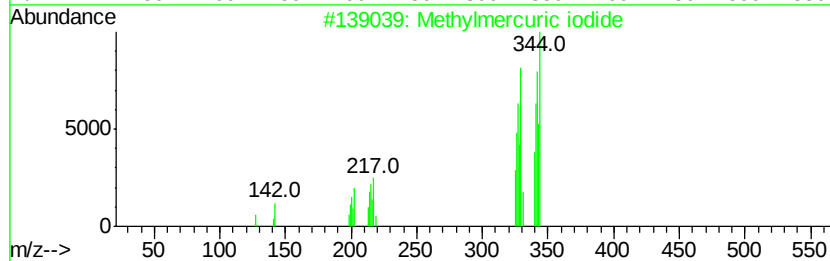
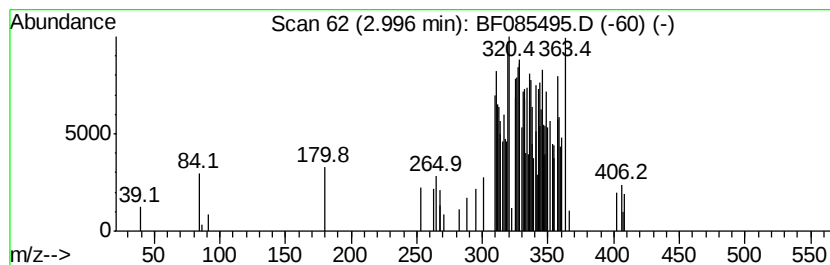
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown3.00 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.00	6.16 ng	141167	1,4-Dichlorobenzene-d4	6.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methylmercuric iodide	344	CH3HgI	000143-36-2	35
2		1-(4-Bromobenzyl)-7-methylpyrido...	345	C15H12BrN3O2	296798-69-1	18
3		Cymar	548	C30H44O9	000508-77-0	15
4		4H-1-Benzopyran-4-one, 2-(3,4-di...	342	C19H18O6	017093-86-6	11
5		A-Nor-5.alpha.-androstan-3-one, ...	344	C22H32O3	019594-91-3	11



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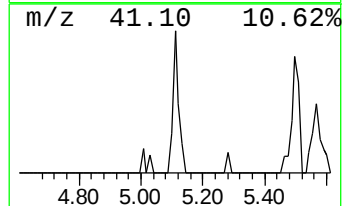
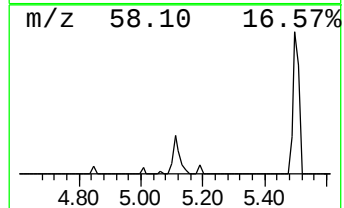
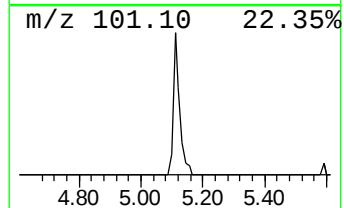
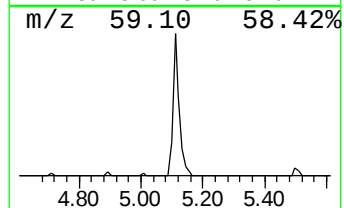
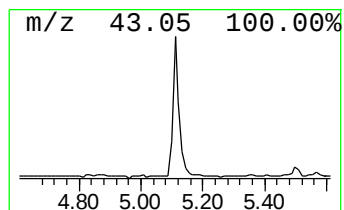
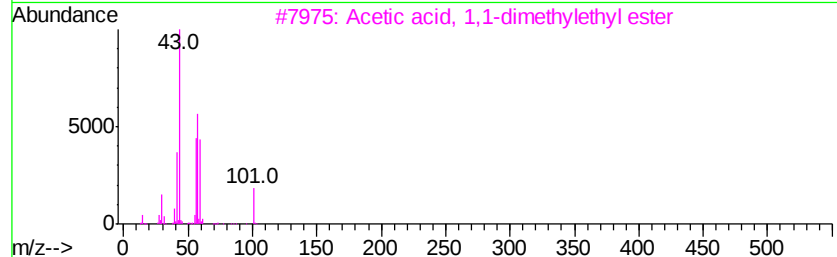
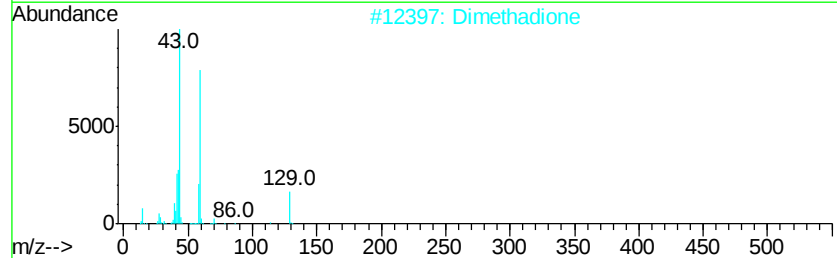
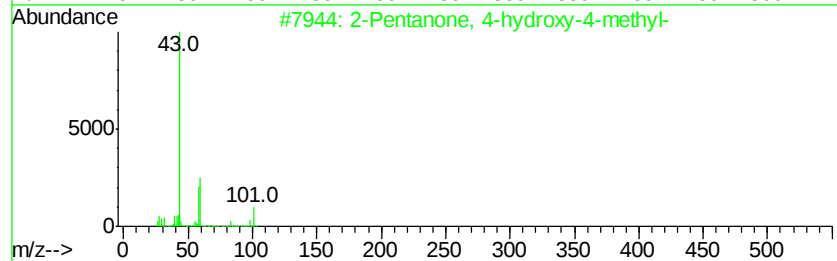
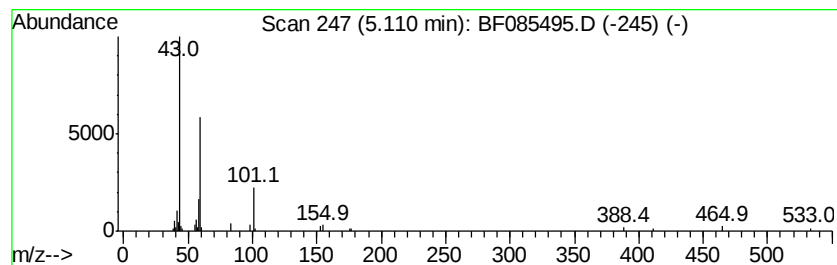
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 6 unknown5.11 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	4.56 ng	104618	1,4-Dichlorobenzene-d4	6.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45		
2	Dimethadione	129	C5H7NO3	000695-53-4	40		
3	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38		
4	Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	25		
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12		



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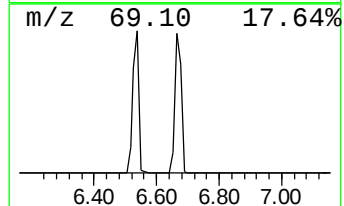
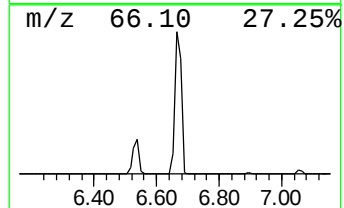
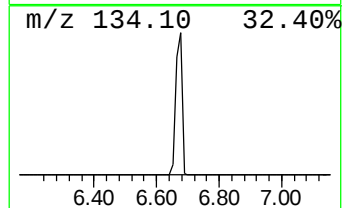
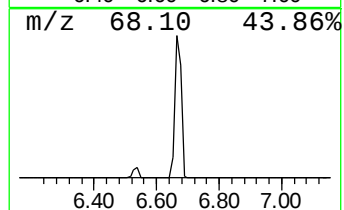
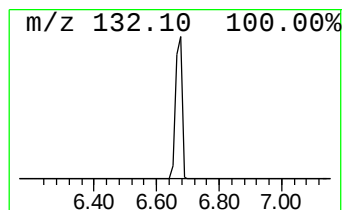
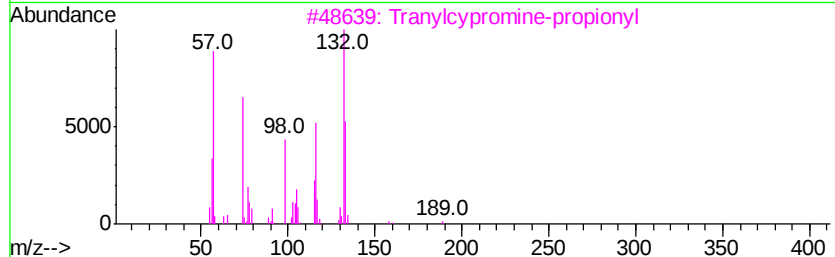
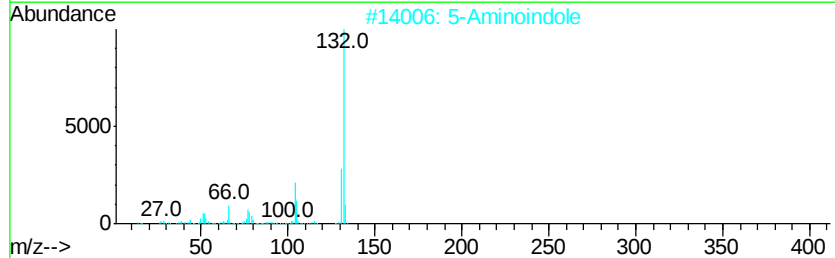
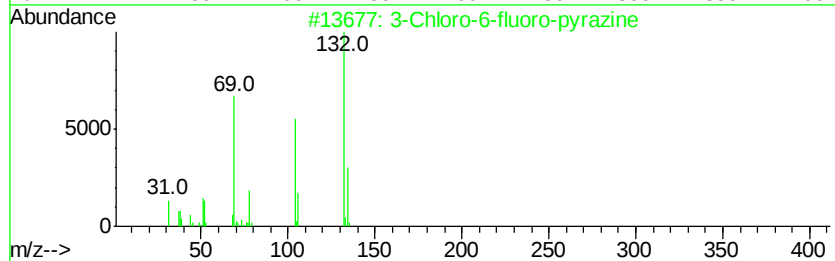
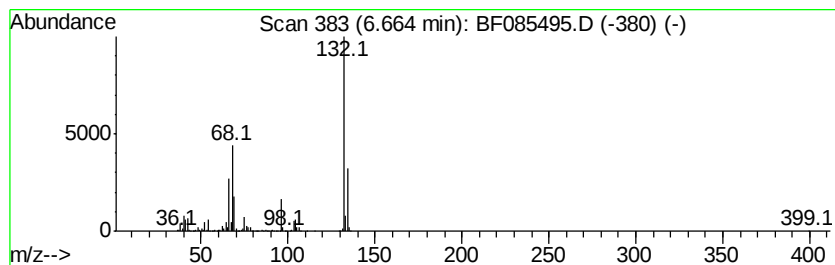
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown6.66 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.66	173.09 ng	3968100	1,4-Dichlorobenzene-d4	6.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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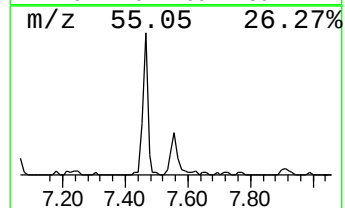
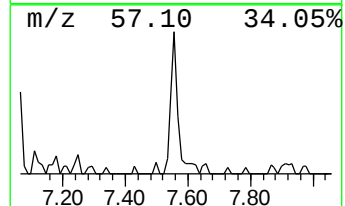
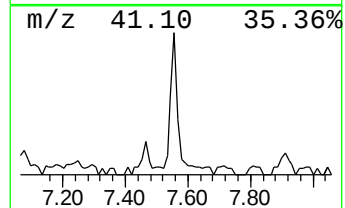
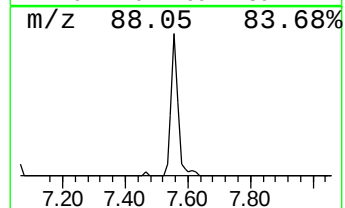
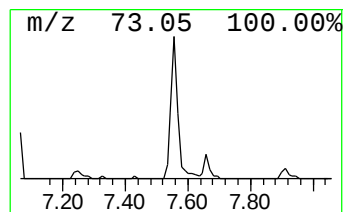
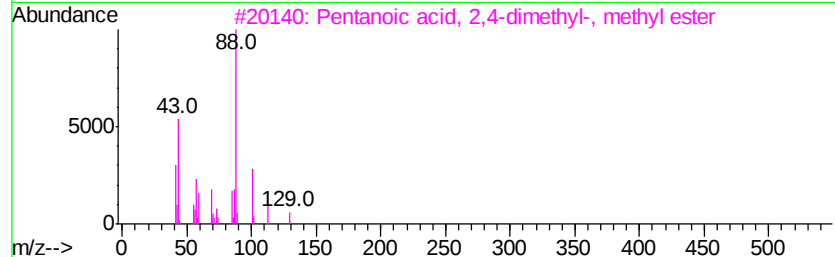
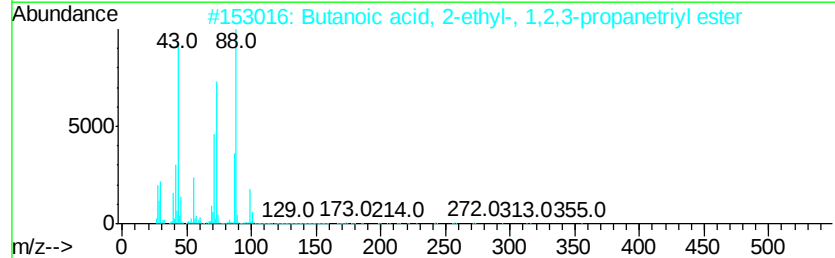
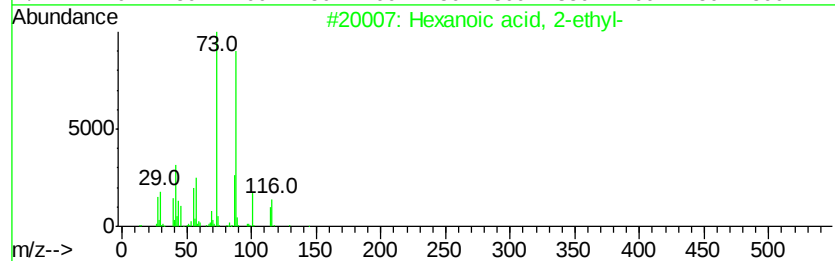
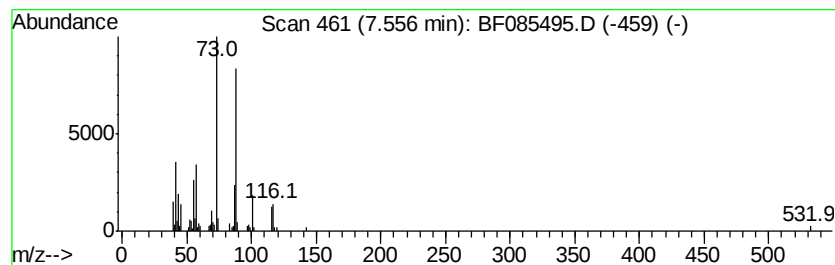
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Peak Number 9 Hexanoic acid, 2-ethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	4.53 ng	131862	Naphthalene-d8	8.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
2		Butanoic acid, 2-ethyl-, 1,2,3-p...	386	C21H38O6	056554-54-2	59
3		Pentanoic acid, 2,4-dimethyl-, m...	144	C8H16O2	071672-33-8	53
4		Octanoic acid, ethyl ester	172	C10H20O2	000106-32-1	47
5		Dithiocarbamic acid, N,N-dimethy...	477	C13H12F9N3S3	1000268-72-7	45



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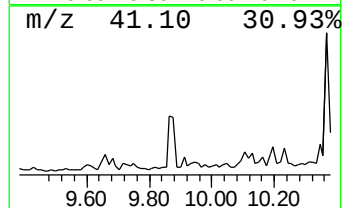
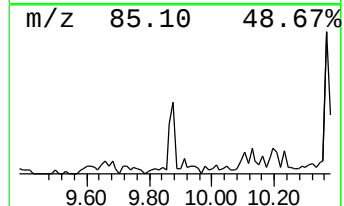
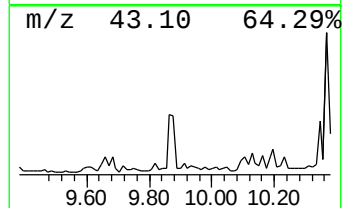
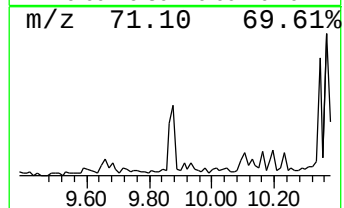
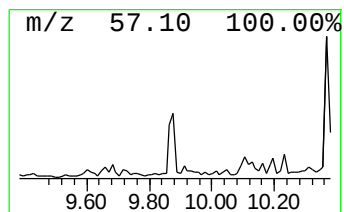
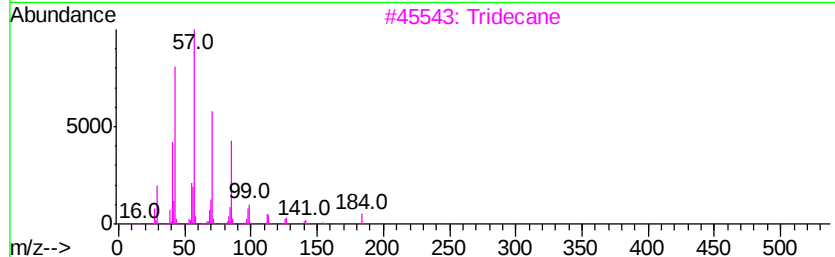
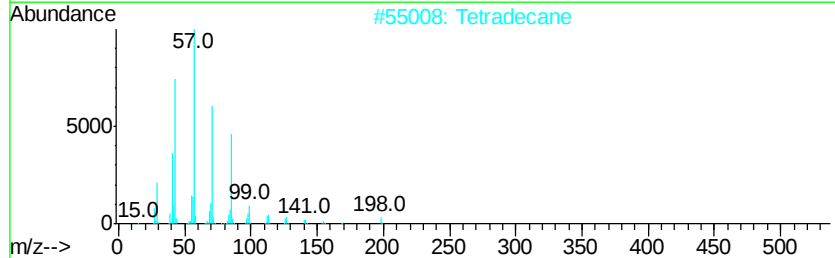
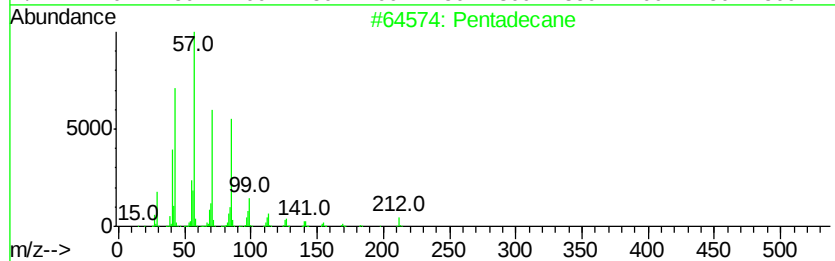
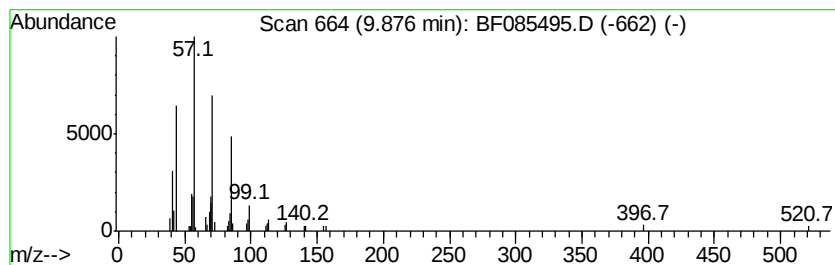
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Pentadecane Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.88	3.01 ng	83923	Acenaphthene-d10	9.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane	212	C15H32	000629-62-9	87
2			Tetradecane	198	C14H30	000629-59-4	87
3			Tridecane	184	C13H28	000629-50-5	86
4			Eicosane	282	C20H42	000112-95-8	83
5			Nonadecane	268	C19H40	000629-92-5	80



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031616\
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Acq On : 17 Mar 2016 7:08
Operator : UM/SJ
Sample : H1768-02
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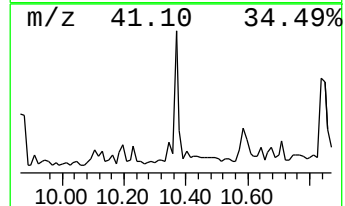
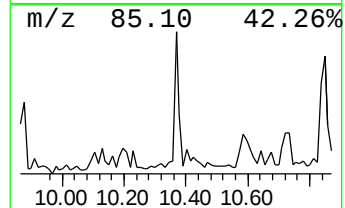
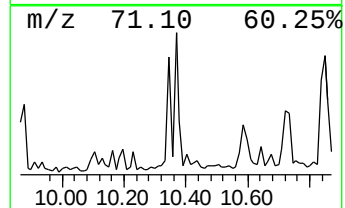
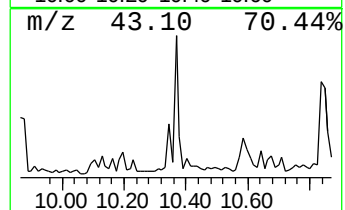
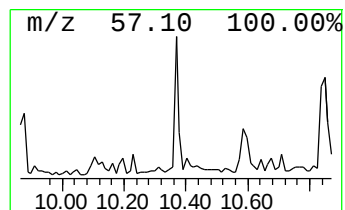
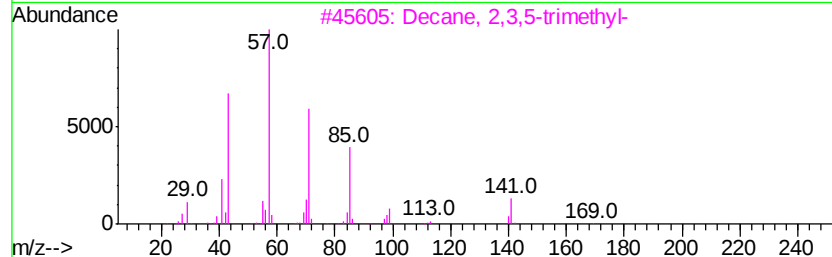
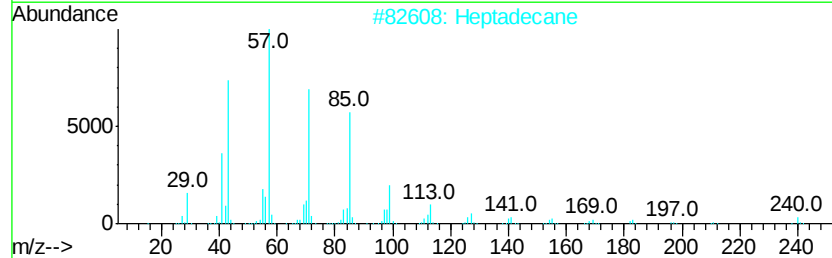
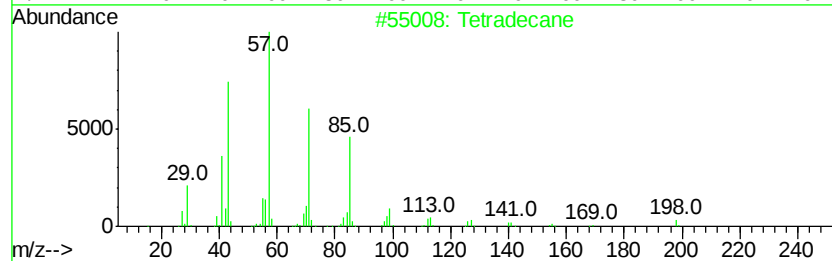
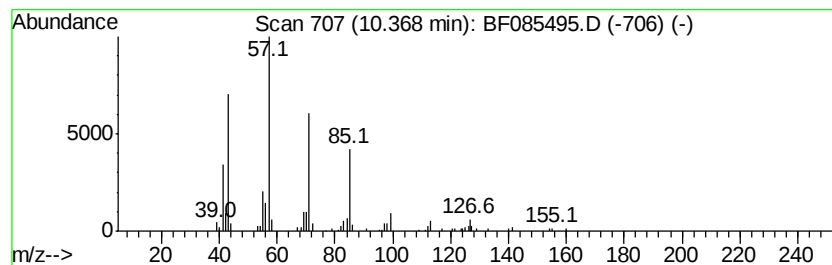
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tetradecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.37	4.52 ng	125880	Acenaphthene-d10		9.94
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	87
2	Heptadecane	240	C17H36	000629-78-7	83
3	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	83
4	Heptacosane	380	C27H56	000593-49-7	83
5	Eicosane	282	C20H42	000112-95-8	74



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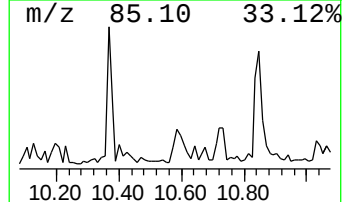
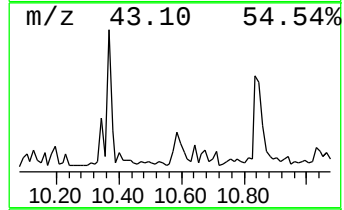
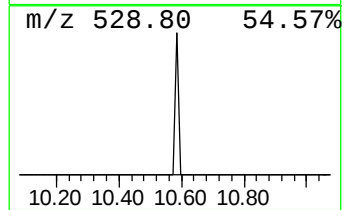
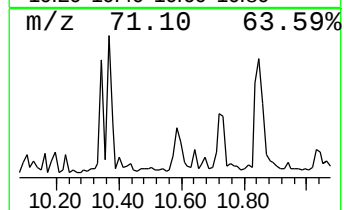
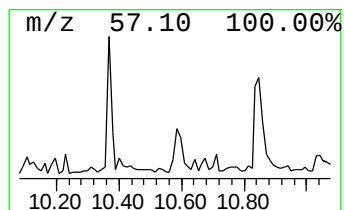
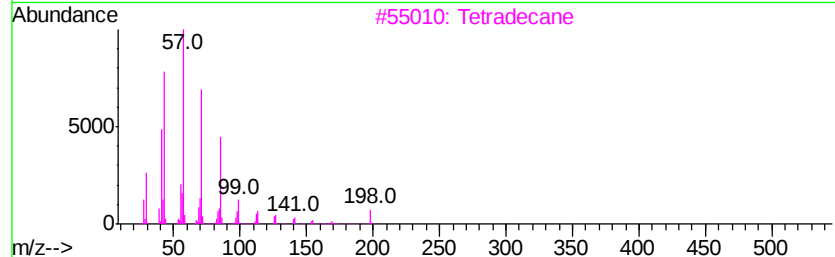
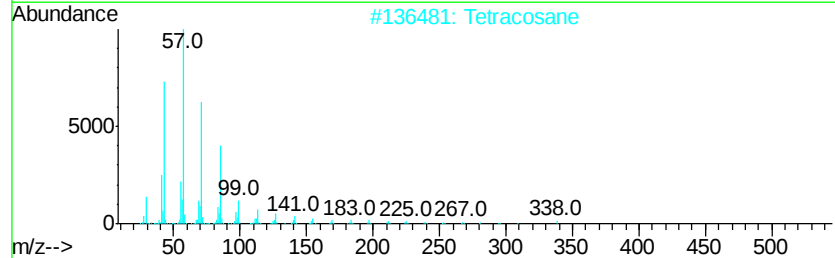
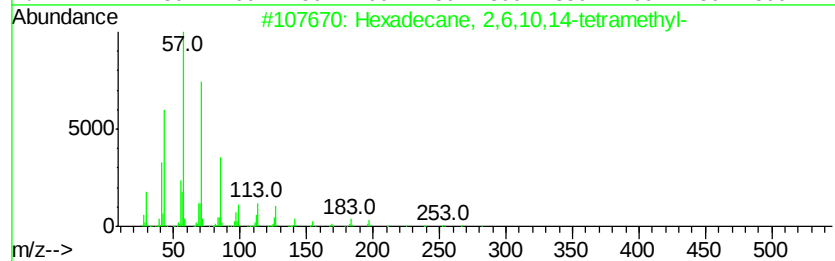
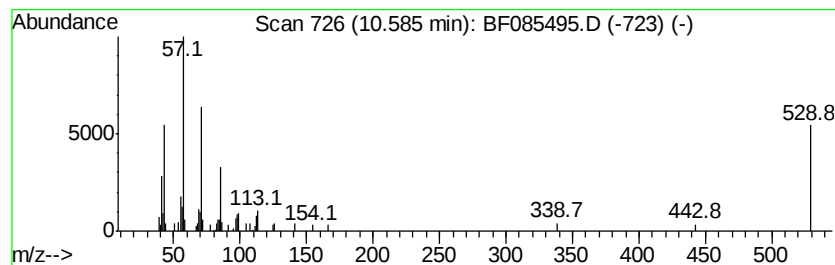
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hexadecane, 2,6,10,14-tetra... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.58	3.23 ng	89922	Acenaphthene-d10	9.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	58
2		Tetracosane	338	C24H50	000646-31-1	52
3		Tetradecane	198	C14H30	000629-59-4	52
4		Nonadecane	268	C19H40	000629-92-5	52
5		Eicosane	282	C20H42	000112-95-8	52



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031616\
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Sample : H1768-02
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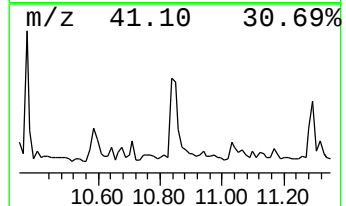
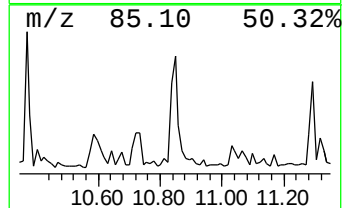
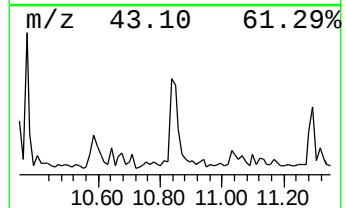
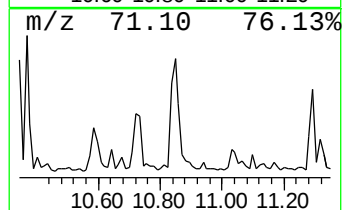
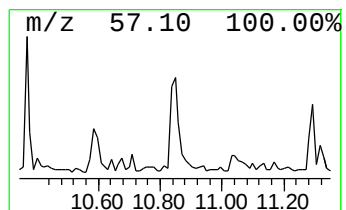
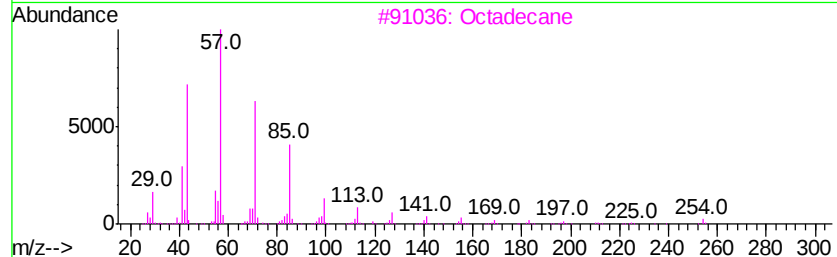
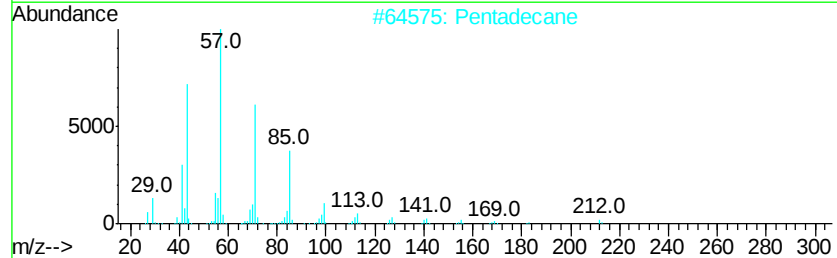
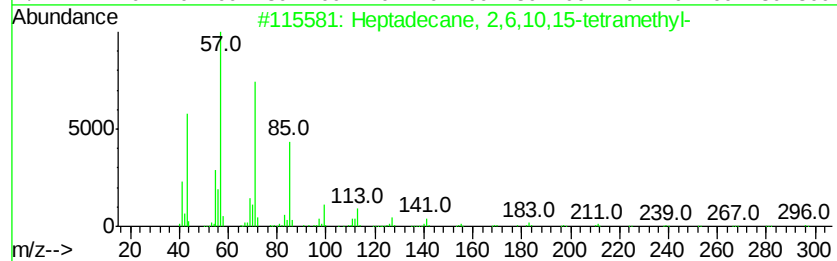
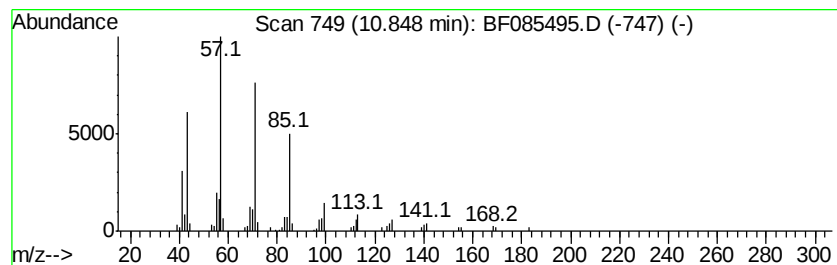
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heptadecane, 2,6,10,15-tetr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.85	7.26 ng	191011	Phenanthrene-d10	11.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
2		Pentadecane	212	C15H32	000629-62-9	91
3		Octadecane	254	C18H38	000593-45-3	91
4		Heptadecane	240	C17H36	000629-78-7	91
5		Hexadecane	226	C16H34	000544-76-3	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031616\
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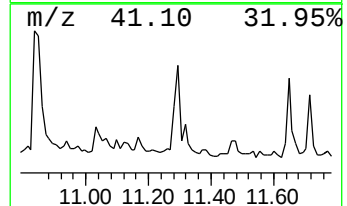
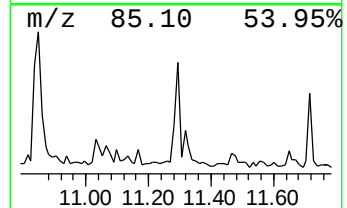
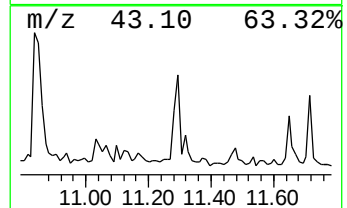
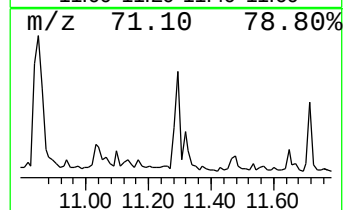
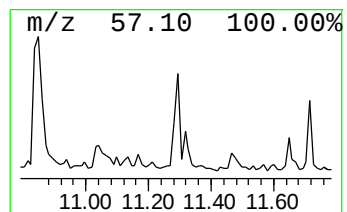
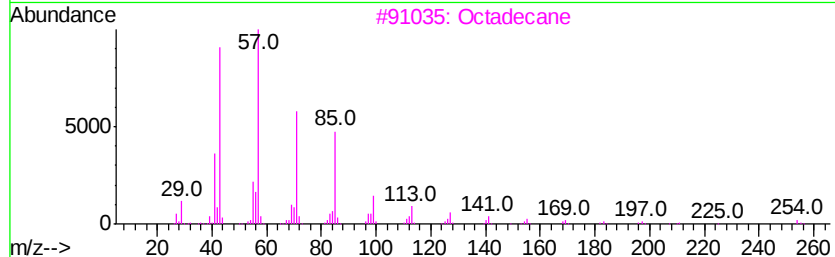
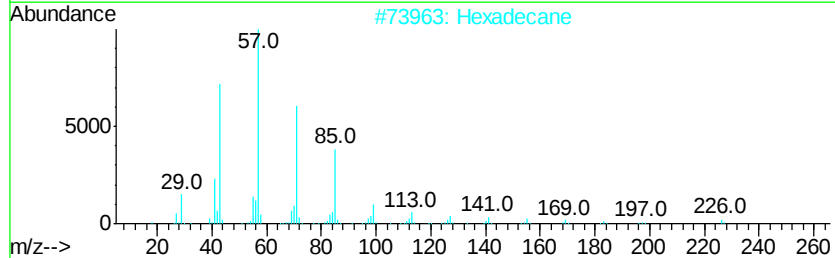
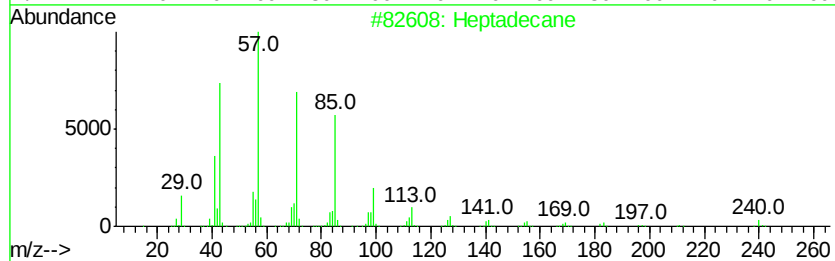
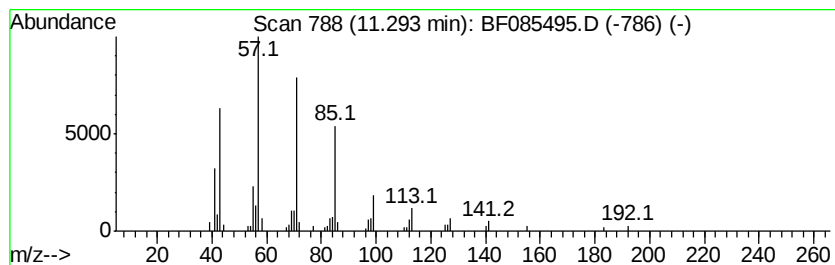
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Heptadecane Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.29	3.06 ng	80413	Phenanthrene-d10	11.42

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	91
2			Hexadecane	226	C16H34	000544-76-3	90
3			Octadecane	254	C18H38	000593-45-3	90
4			Heneicosane	296	C21H44	000629-94-7	90
5			Pentadecane	212	C15H32	000629-62-9	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031616\
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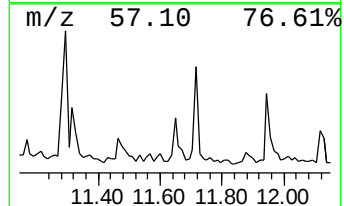
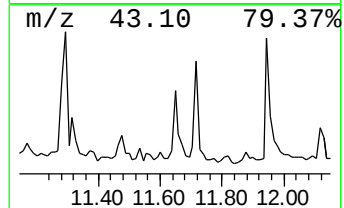
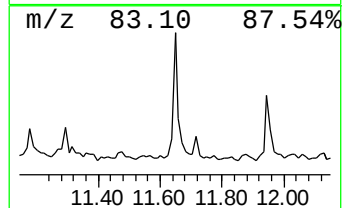
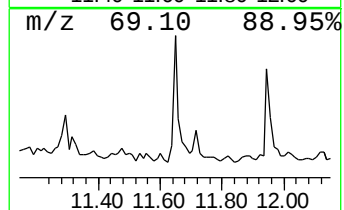
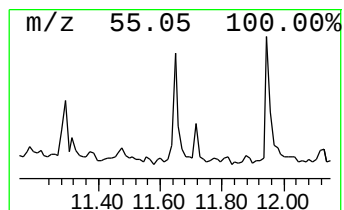
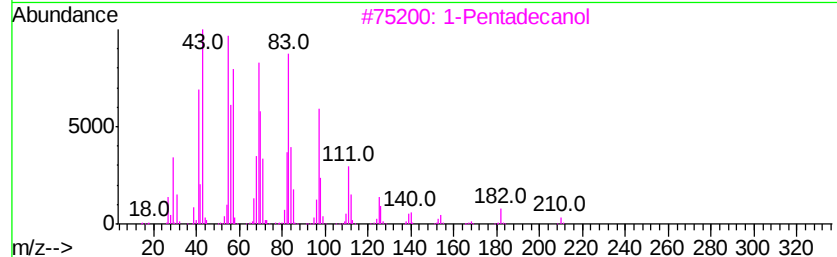
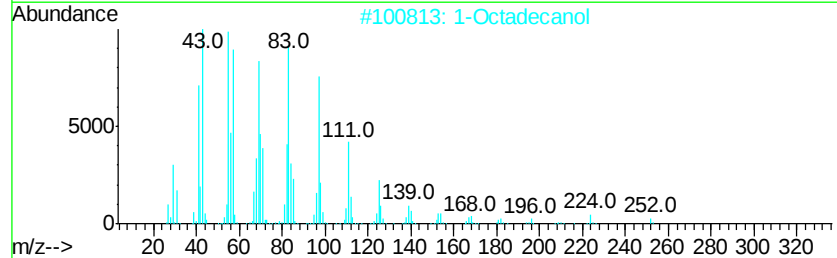
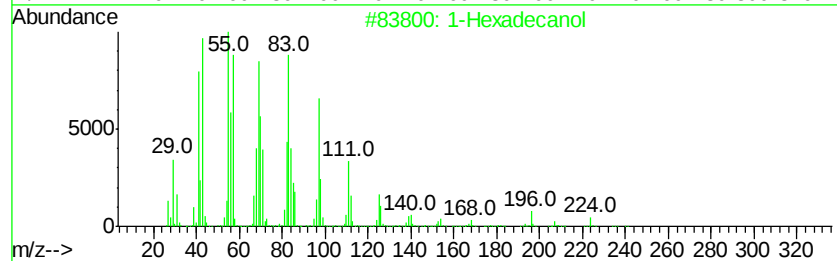
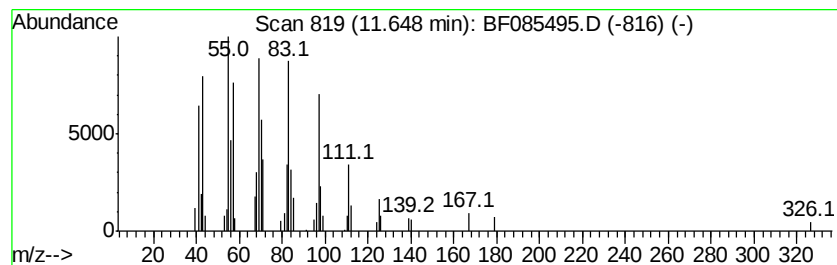
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 1-Hexadecanol Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD		R.T.		
11.65	3.59 ng	94427	Phenanthrene-d10		11.42		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexadecanol	242	C16H34O	036653-82-4	95
2			1-Octadecanol	270	C18H38O	000112-92-5	93
3			1-Pentadecanol	228	C15H32O	000629-76-5	91
4			4-Heptafluorobutylxyloxyhexadecane	438	C20H33F7O2	1000282-97-2	91
5			5-Octadecene, (E)-	252	C18H36	007206-21-5	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031616\
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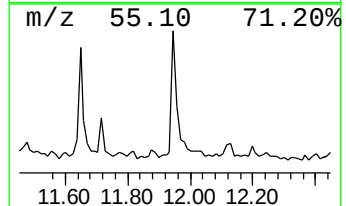
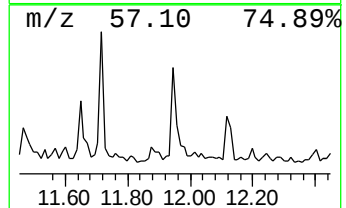
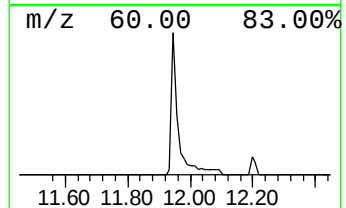
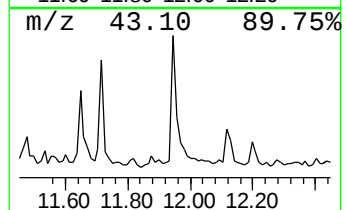
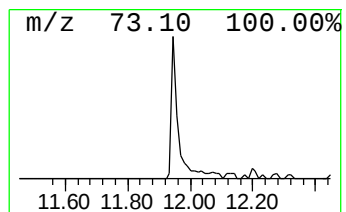
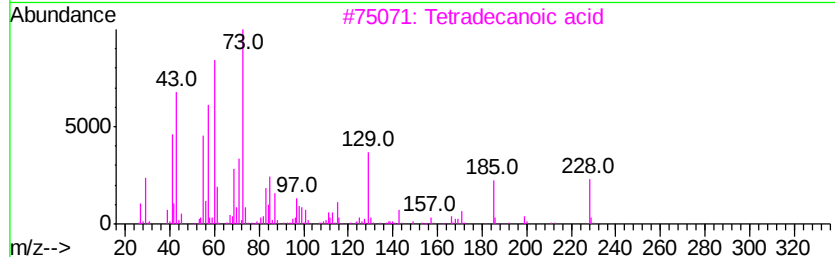
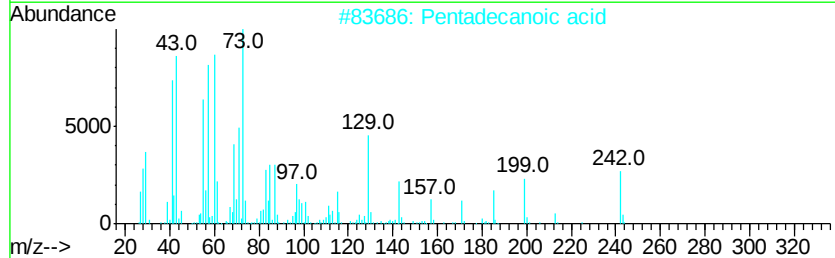
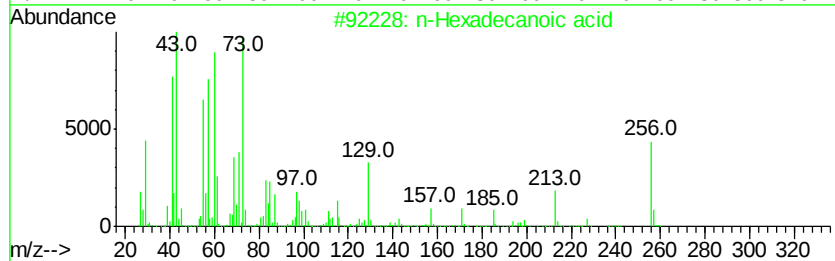
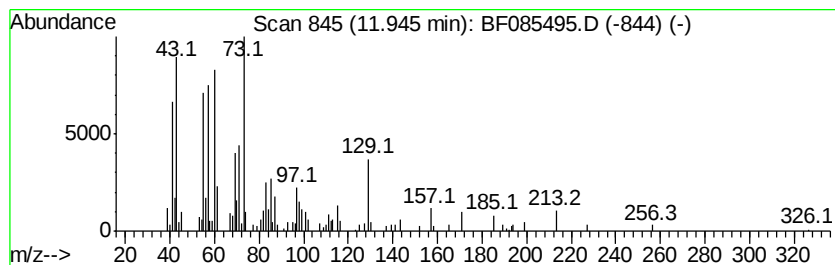
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.95	5.44 ng	143155	Phenanthrene-d10	11.42	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	80
4	Tridecanoic acid	214	C13H26O2	000638-53-9	64
5	Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	11



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031616\
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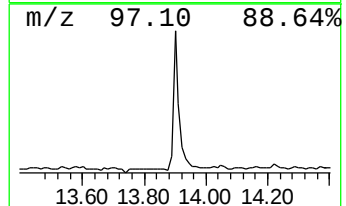
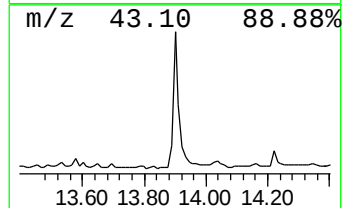
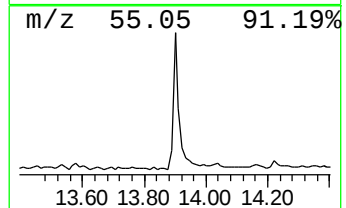
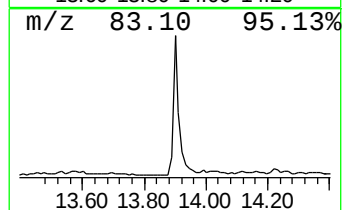
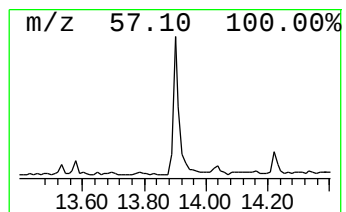
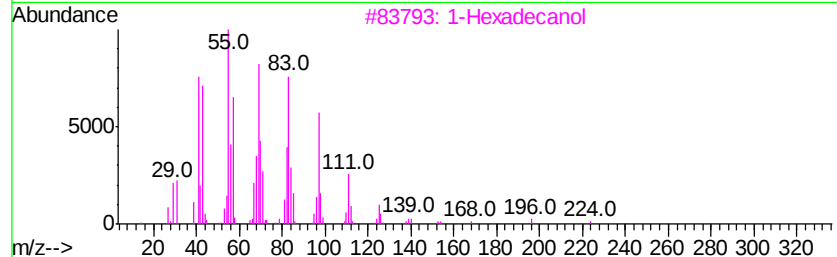
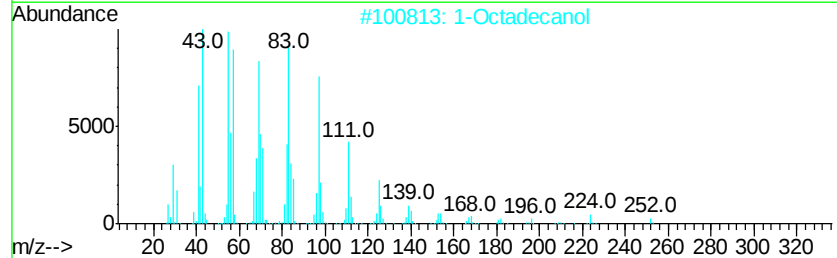
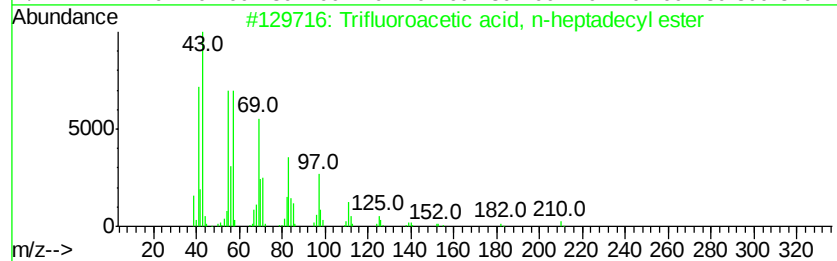
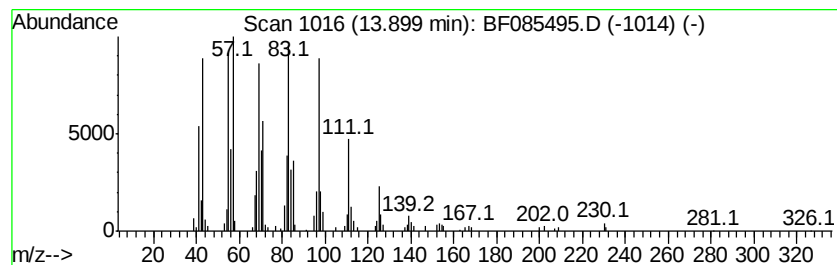
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Trifluoroacetic acid, n-hep... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	14.52 ng	274280	Chrysene-d12	14.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trifluoroacetic acid, n-heptadec...	324	C17H31F3O2	1000216-79-2	91
2		1-Octadecanol	270	C18H38O	000112-92-5	91
3		1-Hexadecanol	242	C16H34O	036653-82-4	91
4		1-Heptadecanol	256	C17H36O	001454-85-9	91
5		Cyclopentadecane	210	C15H30	000295-48-7	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031616\
Data File : BF085495.D
Acq On : 17 Mar 2016 7:08
Operator : UM/SJ
Sample : H1768-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AG-1A

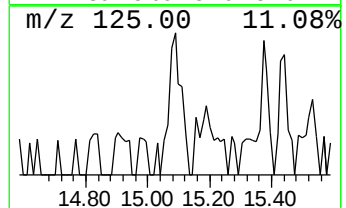
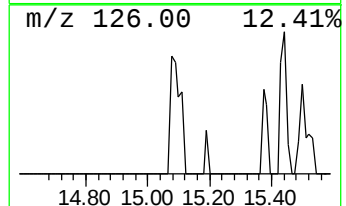
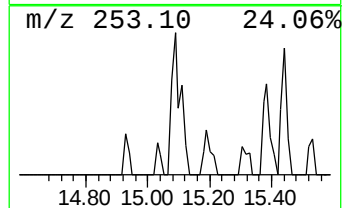
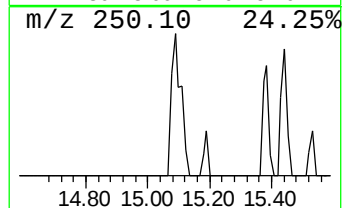
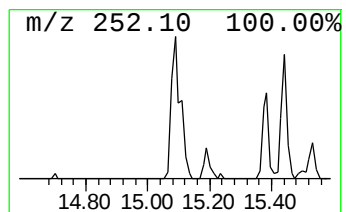
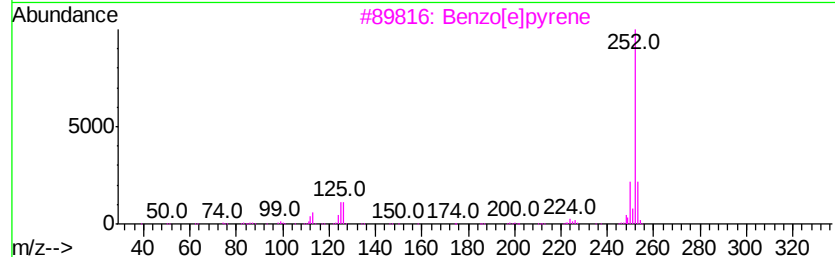
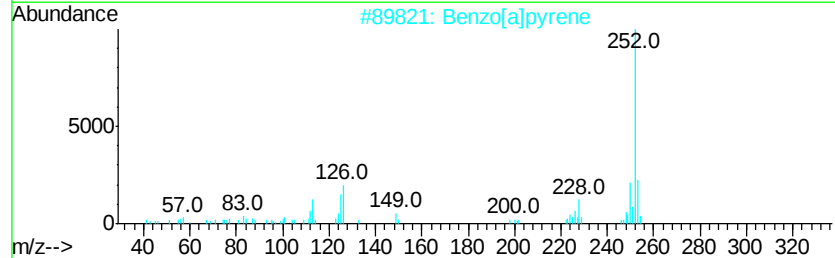
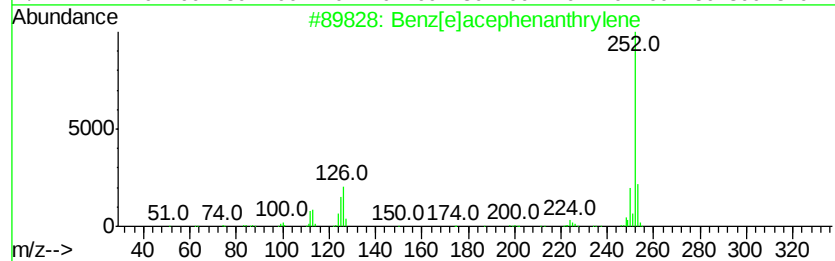
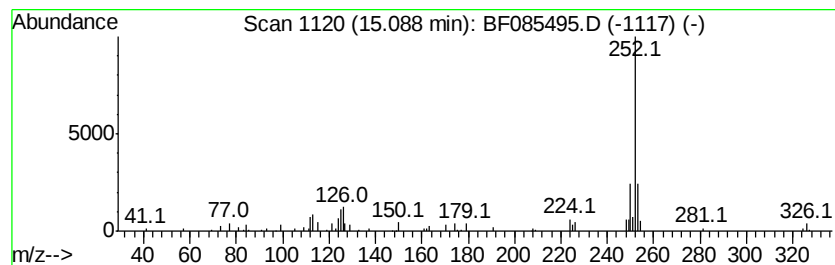
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Benz[e]acephenanthrylene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.09	3.31 ng	66773	Perylene-d12	15.50

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benz[e]acephenanthrylene	252	C20H12	000205-99-2	90
2			Benzo[a]pyrene	252	C20H12	000050-32-8	89
3			Benzo[e]pyrene	252	C20H12	000192-97-2	76
4			Benzo[k]fluoranthene	252	C20H12	000207-08-9	76
5			Benzo[j]fluoranthene	252	C20H12	000205-82-3	68



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031616\
Data File : BF085495.D
Acq On : 17 Mar 2016 7:08
Operator : UM/SJ
Sample : H1768-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AG-1A

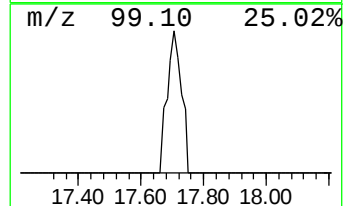
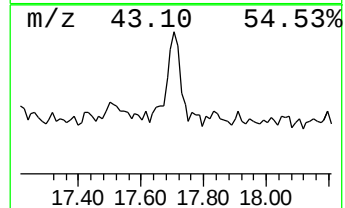
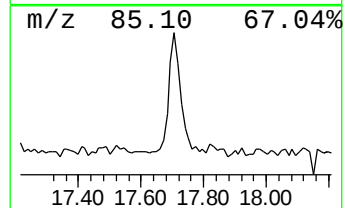
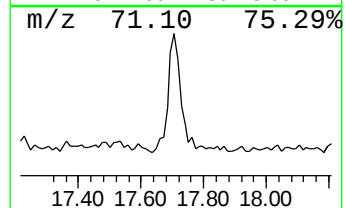
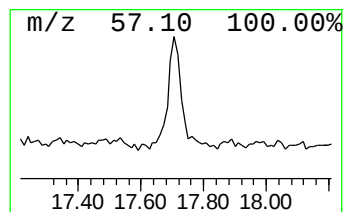
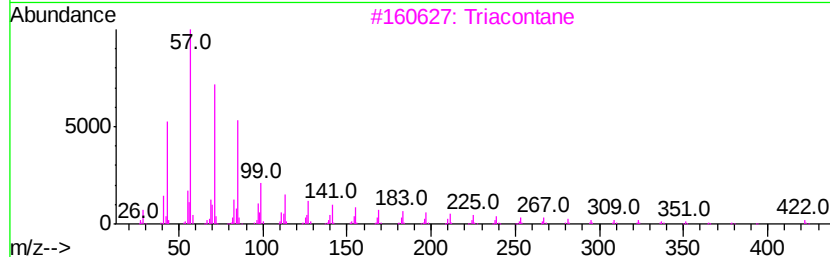
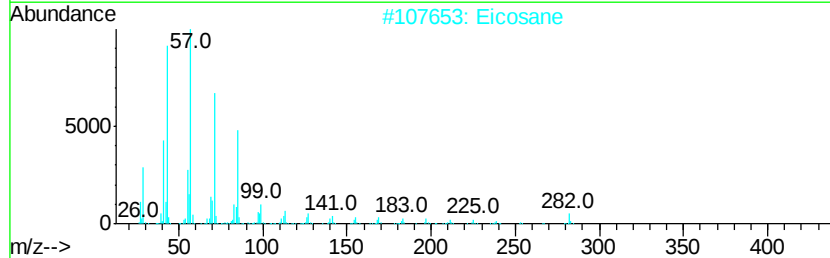
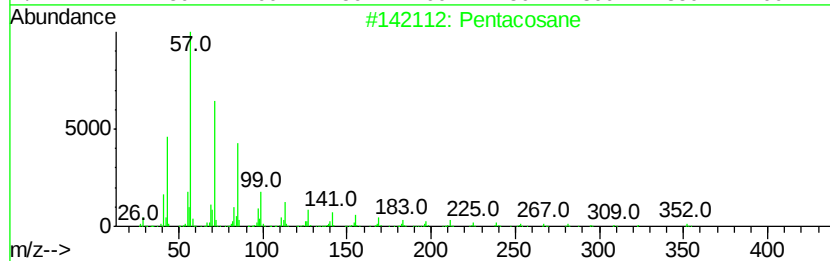
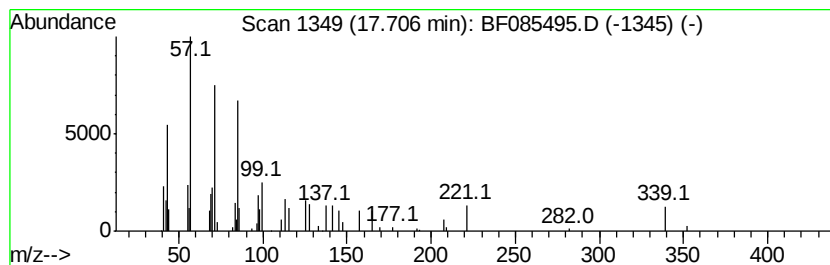
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Pentacosane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.71	3.28 ng	65998	Perylene-d12	15.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentacosane	352	C25H52	000629-99-2	86
2		Eicosane	282	C20H42	000112-95-8	64
3		Triacontane	422	C30H62	000638-68-6	53
4		Tetracosane	338	C24H50	000646-31-1	53
5		Docosane, 9-butyl-	366	C26H54	055282-14-9	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031616\
Data File : BF085495.D
Acq On : 17 Mar 2016 7:08
Operator : UM/SJ
Sample : H1768-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

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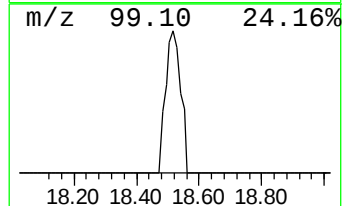
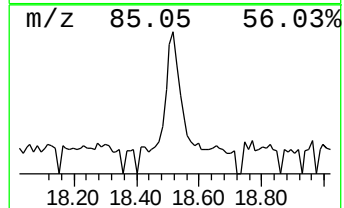
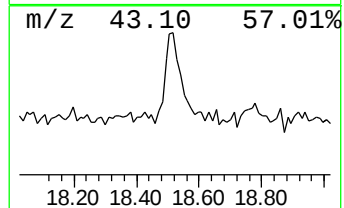
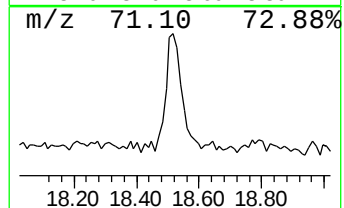
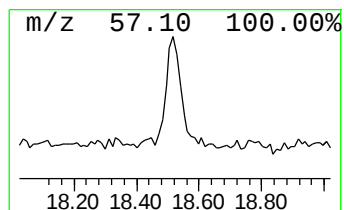
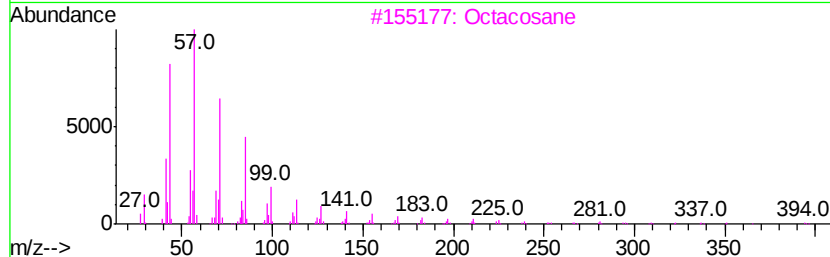
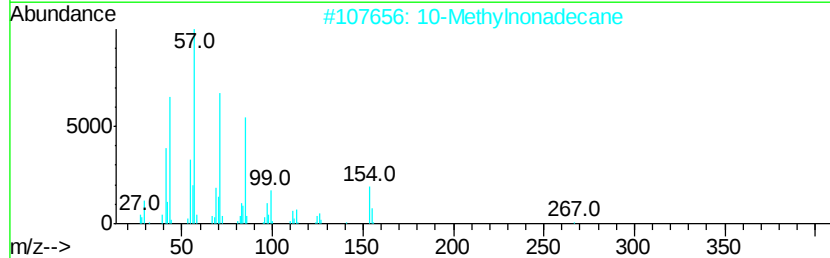
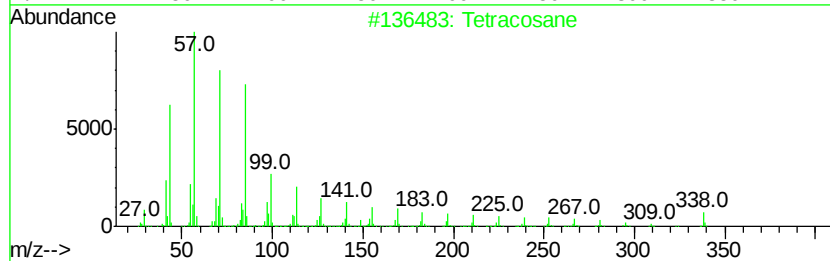
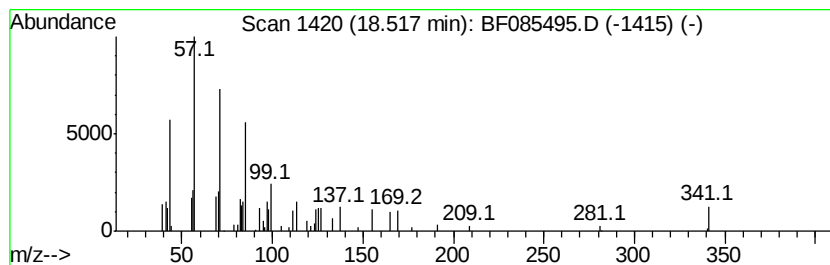
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Tetracosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.52	3.84 ng	77342	Perylene-d12	15.50

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	59
2			10-Methylnonadecane	282	C20H42	056862-62-5	58
3			Octacosane	394	C28H58	000630-02-4	58
4			Disulfide, di-tert-dodecyl	402	C24H50S2	027458-90-8	58
5			Pentacosane	352	C25H52	000629-99-2	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031616\
 Data File : BF085495.D
 Acq On : 17 Mar 2016 7:08
 Operator : UM/SJ
 Sample : H1768-02
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AG-1A

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031516.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
unknown3.00	3.00	6.2	ng	141167	1	6.90	458502	20.0
unknown5.11	5.11	4.6	ng	104618	1	6.90	458502	20.0
unknown6.66	6.66	173.1	ng	3968100	1	6.90	458502	20.0
Hexanoic acid, 2-...	7.56	4.5	ng	131862	2	8.18	581962	20.0
Pentadecane	9.88	3.0	ng	83923	3	9.94	557492	20.0
Tetradecane	10.37	4.5	ng	125880	3	9.94	557492	20.0
Hexadecane, 2,6,1...	10.58	3.2	ng	89922	3	9.94	557492	20.0
Heptadecane, 2,6,...	10.85	7.3	ng	191011	4	11.42	525958	20.0
Heptadecane	11.29	3.1	ng	80413	4	11.42	525958	20.0
1-Hexadecanol	11.65	3.6	ng	94427	4	11.42	525958	20.0
n-Hexadecanoic acid	11.95	5.4	ng	143155	4	11.42	525958	20.0
Trifluoroacetic a...	13.90	14.5	ng	274280	5	14.06	377875	20.0
Benz[e]acephenant...	15.09	3.3	ng	66773	6	15.50	402981	20.0
Pentacosane	17.71	3.3	ng	65998	6	15.50	402981	20.0
Tetracosane	18.52	3.8	ng	77342	6	15.50	402981	20.0