

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
 Data File : BF082316.D
 Acq On : 15 Oct 2015 18:31
 Operator : UM/IZ
 Sample : G4015-06
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LF-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-BF101015.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	4.320	188	191	196	rBV	18673	28548	1.23%	0.129%
2	4.994	247	250	256	rBV	712671	1000982	43.02%	4.520%
3	5.428	285	288	290	rBV	1644279	1801318	77.41%	8.134%
4	5.634	304	306	311	rVB	27469	32078	1.38%	0.145%
5	6.469	376	379	382	rBV	1460057	1625886	69.87%	7.342%
6	6.594	387	390	392	rBV	2037445	1833156	78.78%	8.278%
7	6.823	408	410	413	rBV	562934	474530	20.39%	2.143%
8	6.983	421	424	426	rBV	1693409	1639818	70.47%	7.405%
9	7.394	457	460	463	rBV	1136164	1084492	46.61%	4.897%
10	8.114	520	523	525	rBV	734553	629874	27.07%	2.844%
11	8.675	569	572	573	rBV	73180	87069	3.74%	0.393%
12	9.200	615	618	620	rBV	1651640	2326902	100.00%	10.507%
13	9.589	650	652	655	rBV	391358	324924	13.96%	1.467%
14	9.875	674	677	679	rVB	767239	702083	30.17%	3.170%
15	10.149	697	701	703	rBV3	45601	62193	2.67%	0.281%
16	10.298	713	714	715	rBV	52590	39876	1.71%	0.180%
17	10.412	720	724	726	rBV3	12430	31618	1.36%	0.143%
18	10.526	731	734	735	rBV	51320	84033	3.61%	0.379%
19	10.663	744	746	751	rBV	1037680	1142266	49.09%	5.158%
20	10.789	754	757	759	rBV	203427	295764	12.71%	1.336%
21	10.869	760	764	766	rBV3	30932	89311	3.84%	0.403%
22	10.915	766	768	770	rBV3	25237	56071	2.41%	0.253%
23	10.983	772	774	777	rBV3	42468	103291	4.44%	0.466%
24	11.258	794	798	800	rBV	283104	475782	20.45%	2.148%
25	11.361	805	807	809	rBV	602908	675174	29.02%	3.049%
26	11.612	827	829	830	rBV	149001	153379	6.59%	0.693%
27	11.898	852	854	856	rBV3	244619	411549	17.69%	1.858%
28	12.961	945	947	950	rBV	1026693	1634237	70.23%	7.380%
29	14.058	1041	1043	1047	rVB	429147	574393	24.68%	2.594%
30	15.589	1175	1177	1186	rVB	474183	1147102	49.30%	5.180%
31	16.264	1234	1236	1239	rVB	347168	641362	27.56%	2.896%
32	16.687	1270	1273	1280	rVB	408103	936227	40.23%	4.228%

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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-BF101015.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

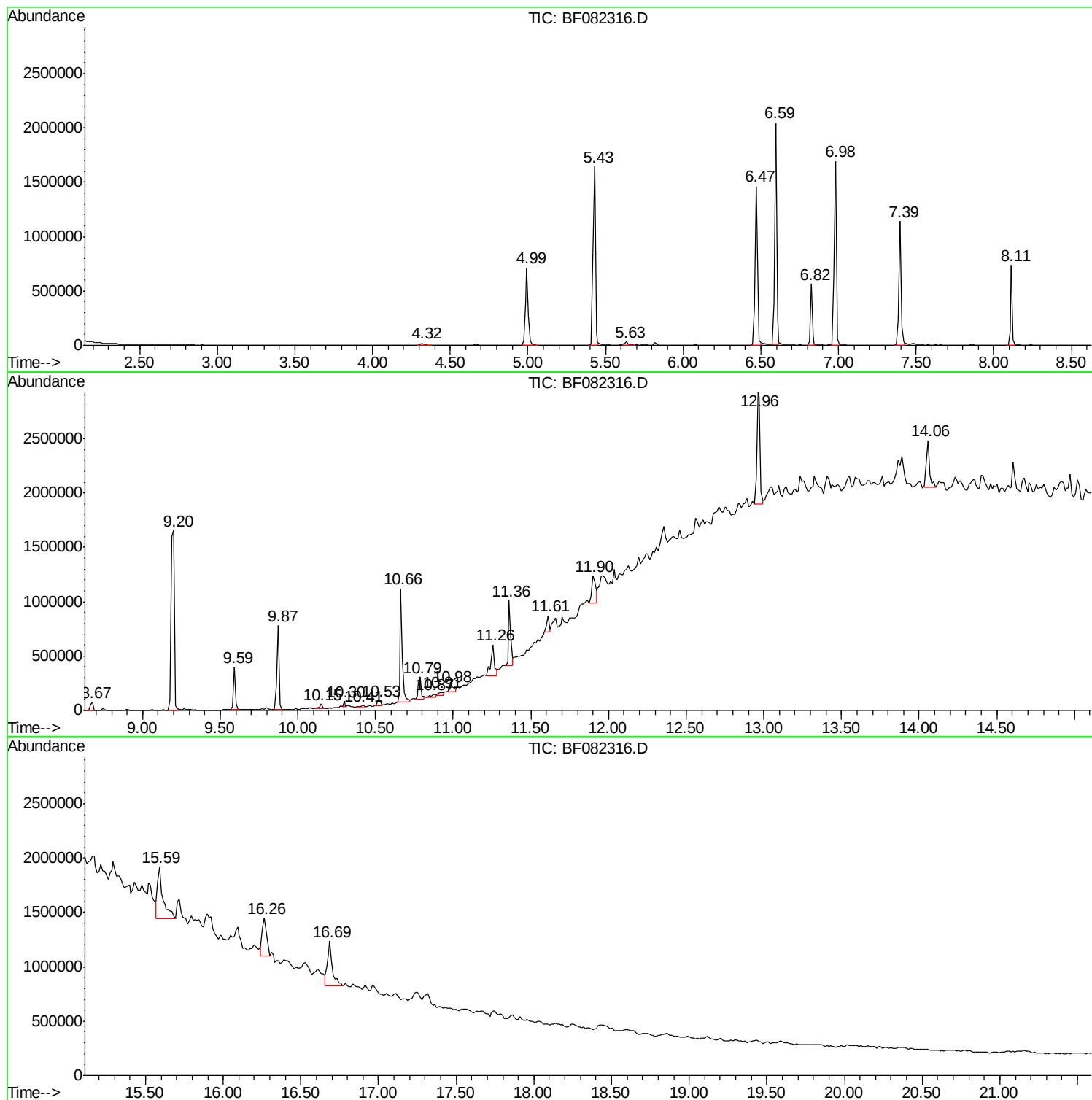
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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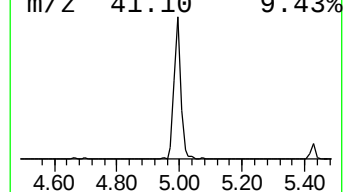
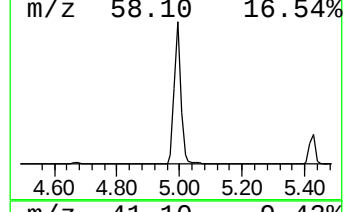
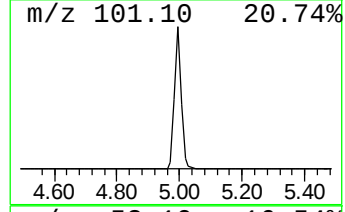
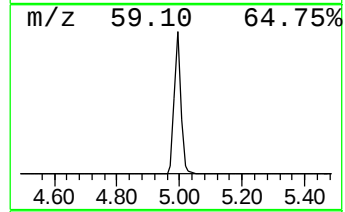
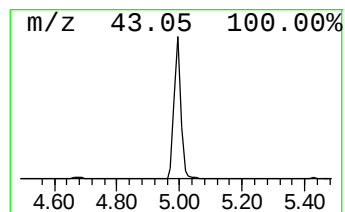
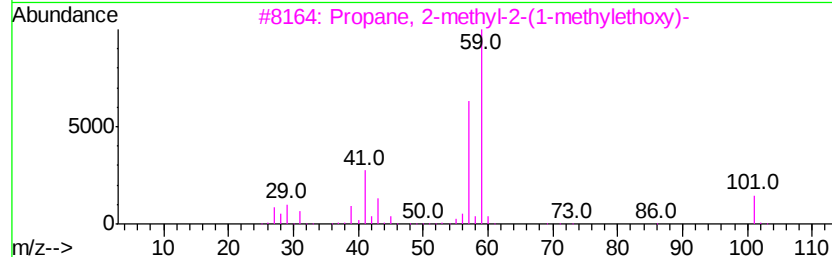
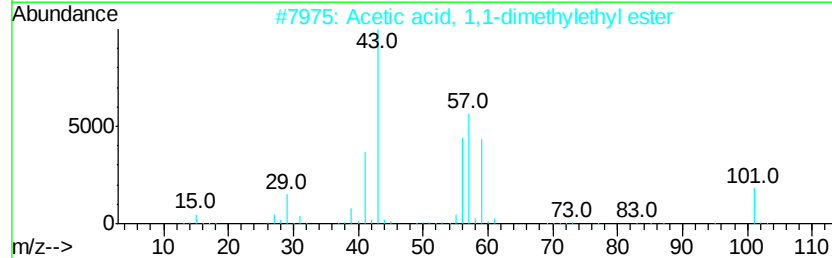
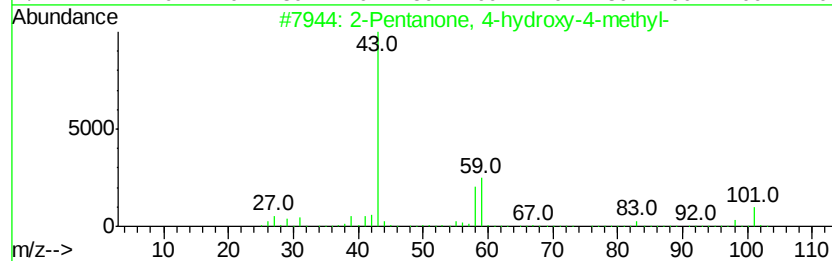
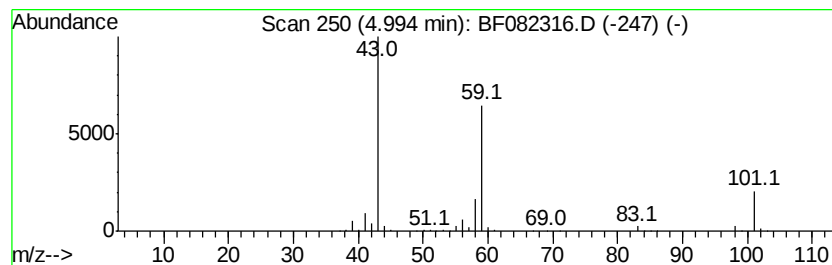
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	42.19 ng	1000980	1,4-Dichlorobenzene-d4	6.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25



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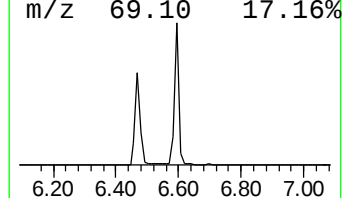
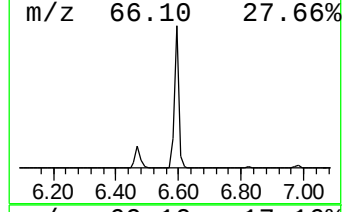
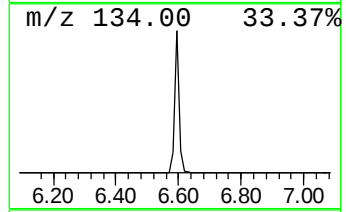
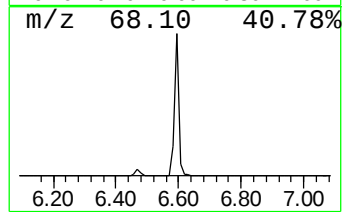
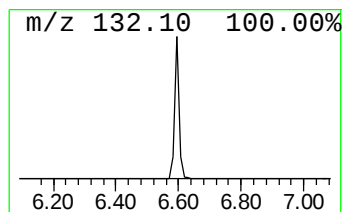
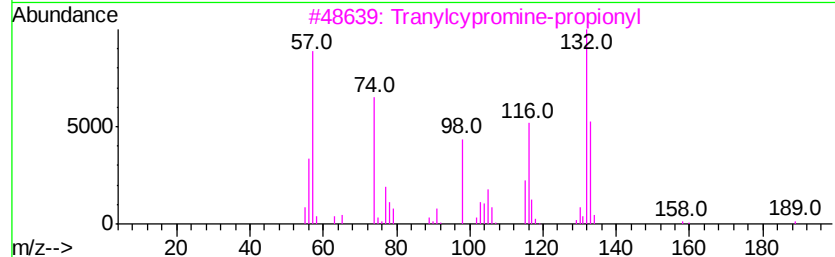
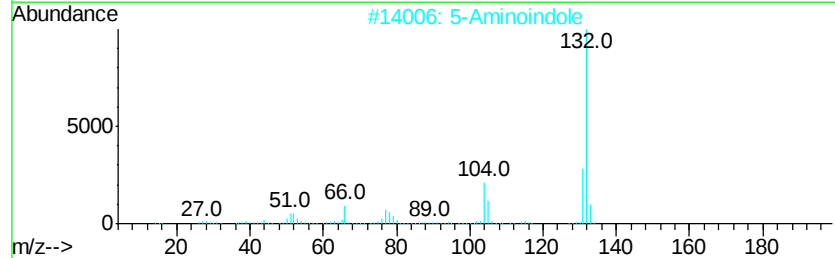
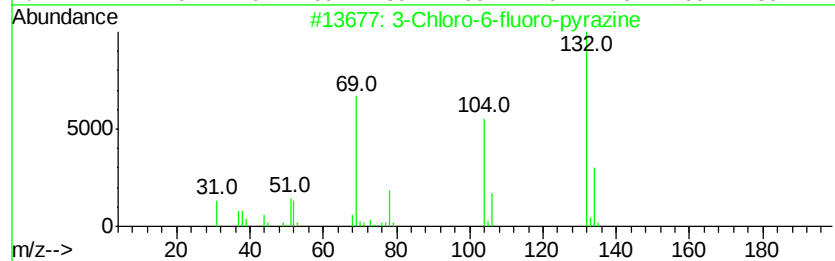
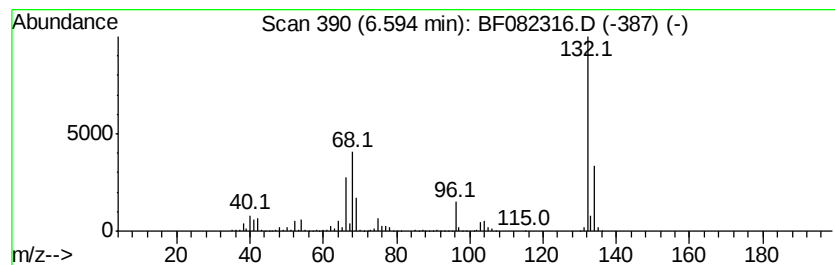
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.59	77.26 ng	1833160	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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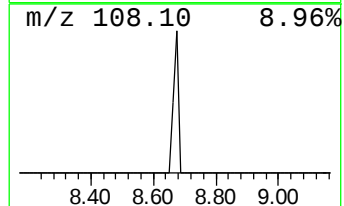
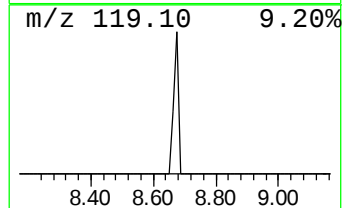
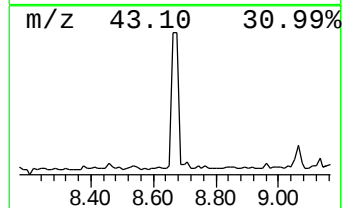
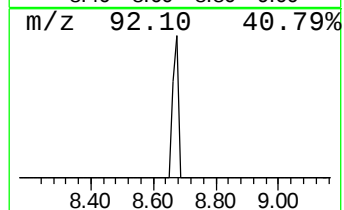
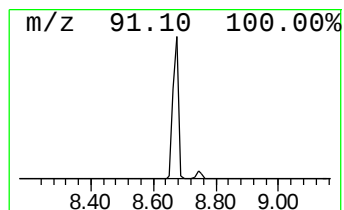
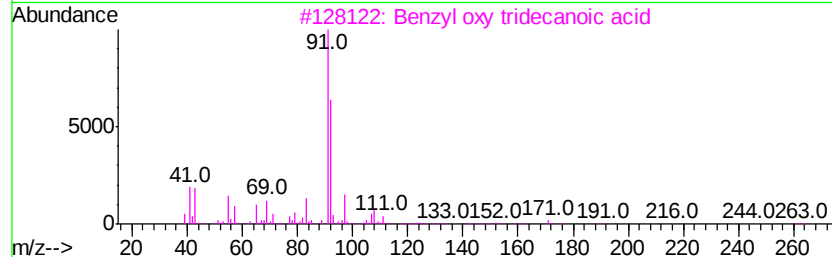
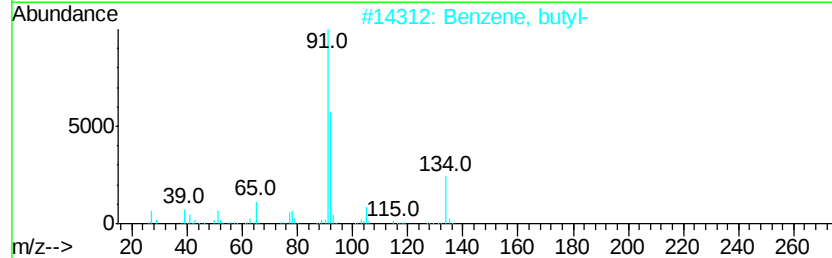
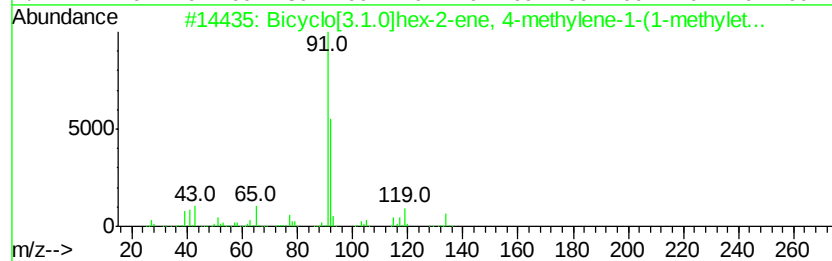
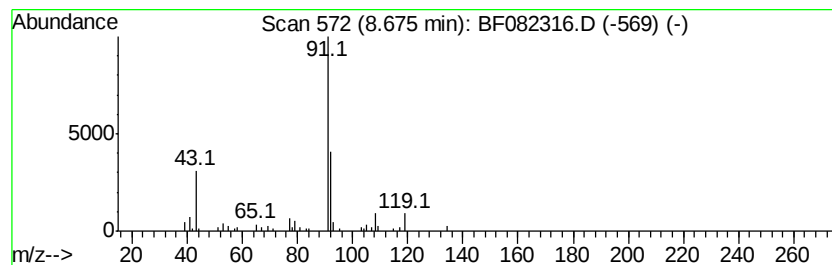
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Bicyclo[3.1.0]hex-2-ene, 4-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.67	2.76 ng	87069	Napthalene-d8	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[3.1.0]hex-2-ene, 4-methy...	134	C10H14	036262-09-6	50
2		Benzene, butyl-	134	C10H14	000104-51-8	42
3		Benzyl oxy tridecanoic acid	320	C20H32O3	1000289-36-6	38
4		Benzyl isopentyl ether	178	C12H18O	000122-73-6	28
5		trans-Verbenol	152	C10H16O	1000292-86-3	22



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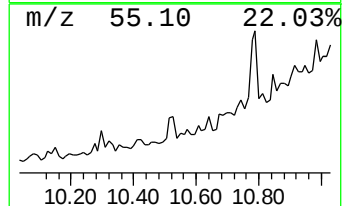
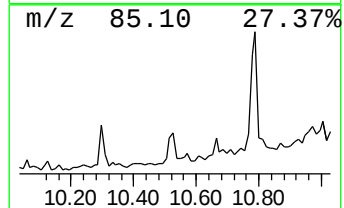
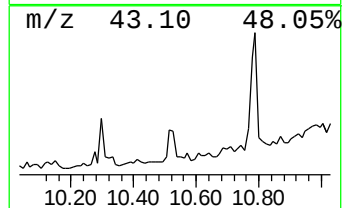
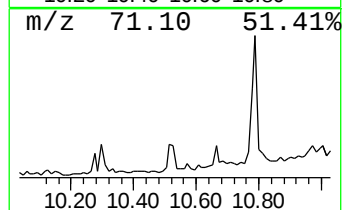
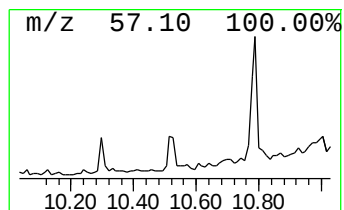
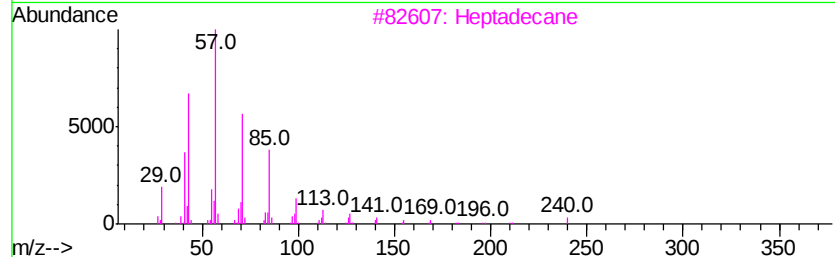
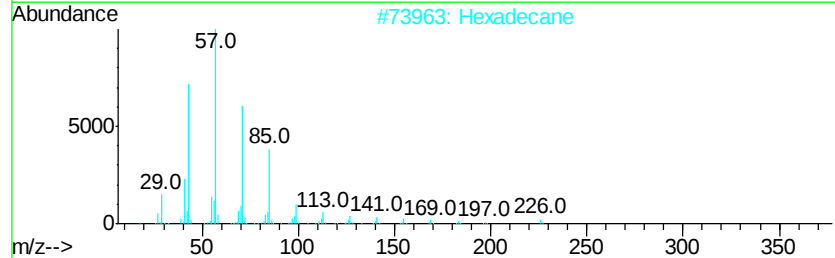
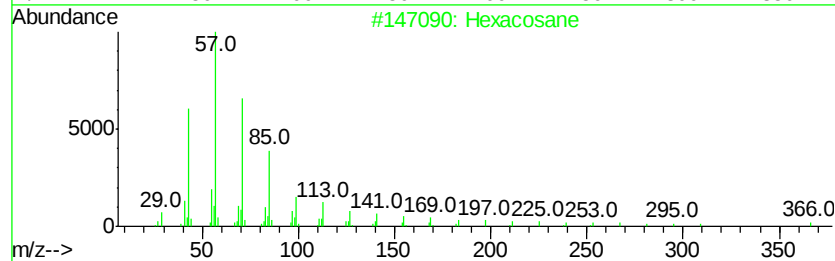
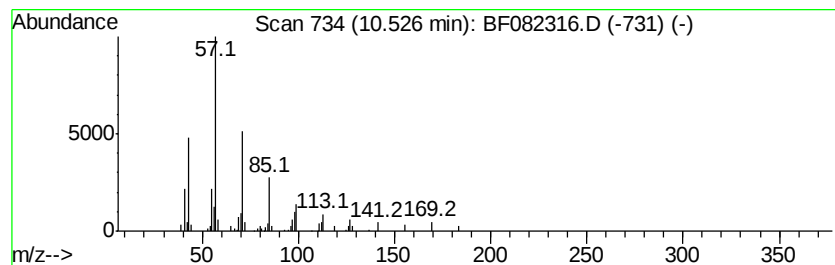
Instrument :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Hexacosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.53	2.39 ng	84033	Acenaphthene-d10	9.87
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexacosane	366 C26H54	000630-01-3	90
2	Hexadecane	226 C16H34	000544-76-3	86
3	Heptadecane	240 C17H36	000629-78-7	86
4	Tetratetracontane	619 C44H90	007098-22-8	86
5	Heptacosane	380 C27H56	000593-49-7	86



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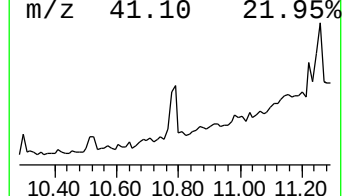
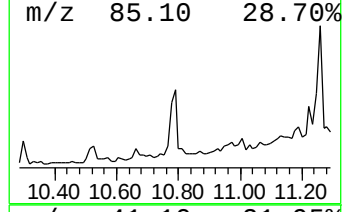
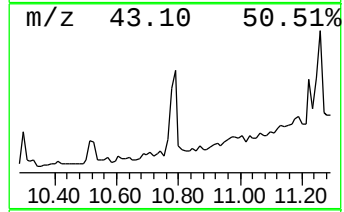
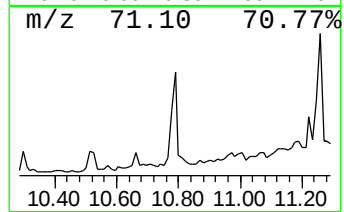
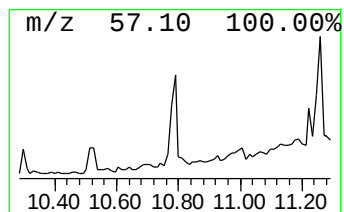
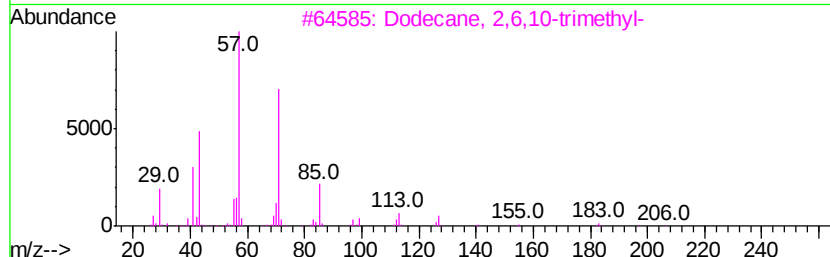
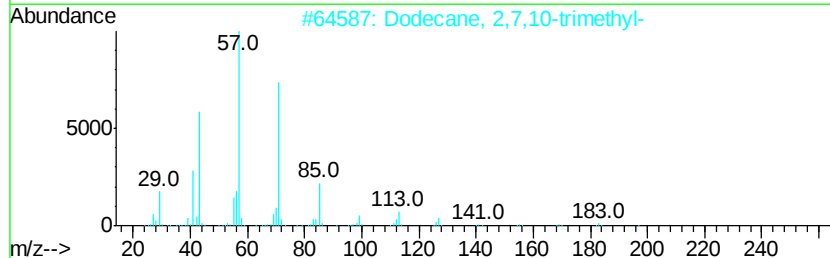
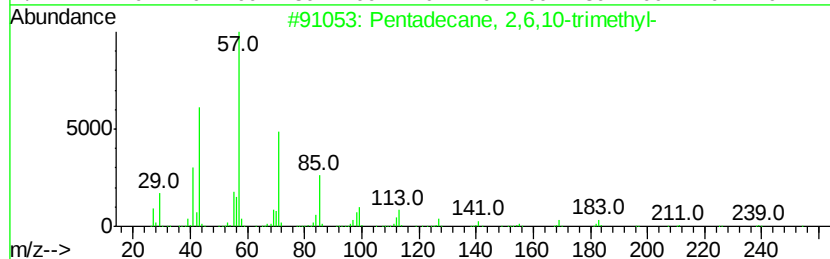
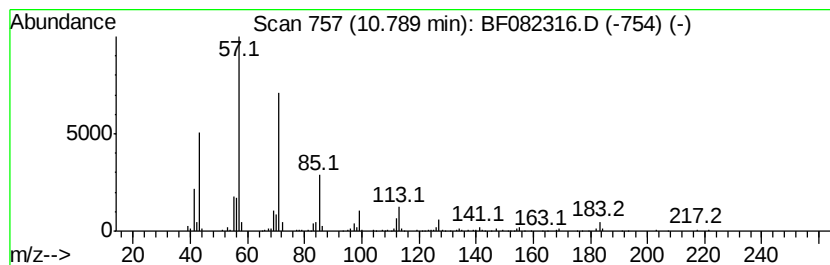
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Pentadecane, 2,6,10-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.79	8.76 ng	295764	Phenanthrene-d10	11.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	90
2		Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	83
3		Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	78
4		Undecane, 6-ethyl-	184	C13H28	017312-60-6	76
5		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	72



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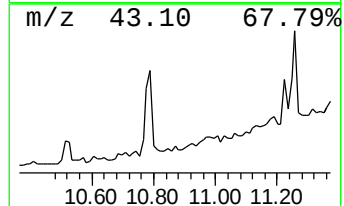
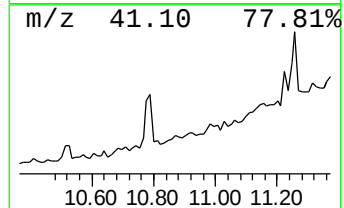
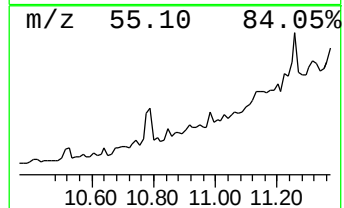
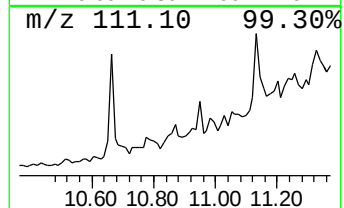
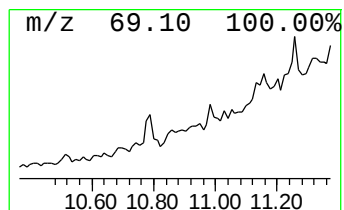
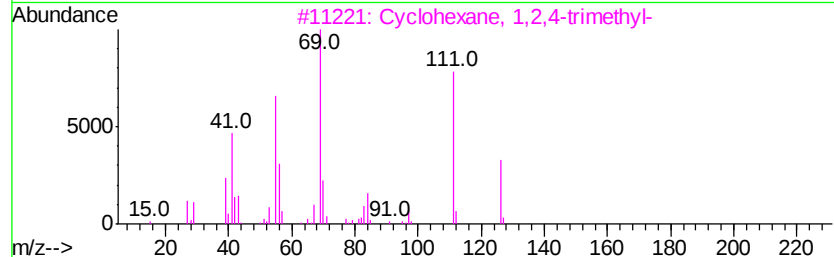
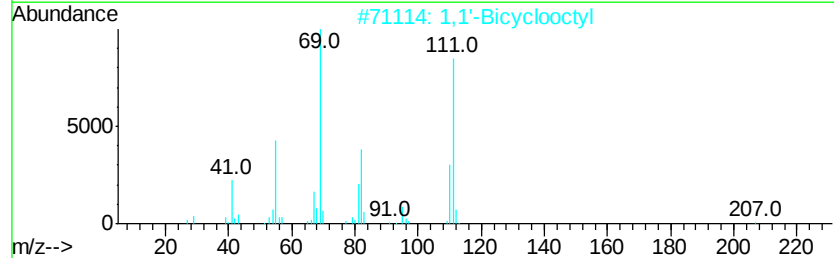
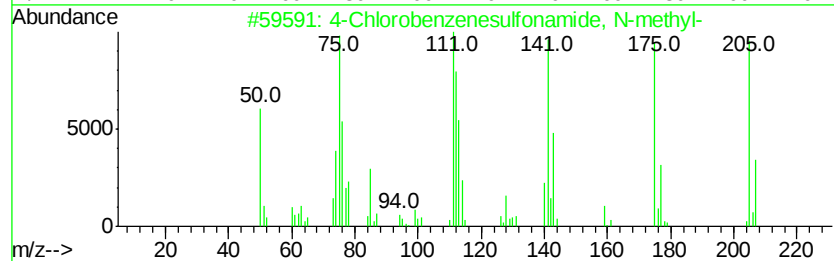
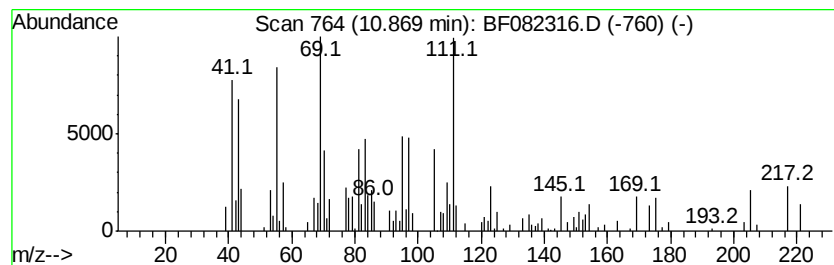
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 4-Chlorobenzenesulfonamide,... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.87	2.65 ng	89311	Phenanthrene-d10	11.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Chlorobenzenesulfonamide, N-me...	205	C7H8ClN02S	006333-79-5	59
2	1,1'-Bicyclooctyl	222	C16H30	006708-17-4	35
3	Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	35
4	6-Octenal, 3,7-dimethyl-	154	C10H18O	000106-23-0	30
5	1,1,3-Trimethyl-1-silacyclo-3-pe...	126	C7H14Si	003528-14-1	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
Data File : BF082316.D
Acq On : 15 Oct 2015 18:31
Operator : UM/IZ
Sample : G4015-06
Misc :
ALS Vial : 42 Sample Multiplier: 1

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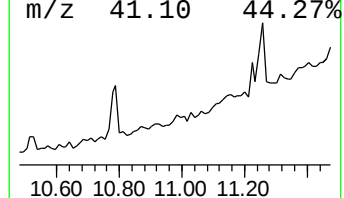
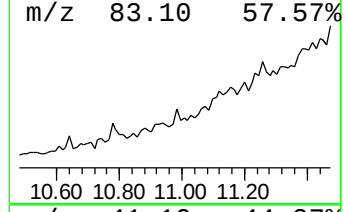
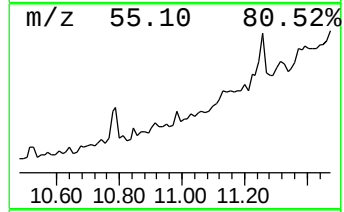
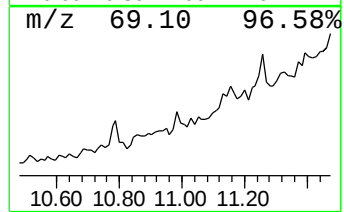
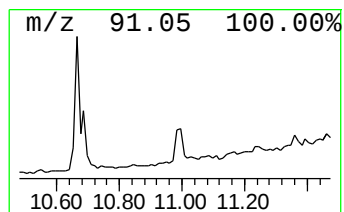
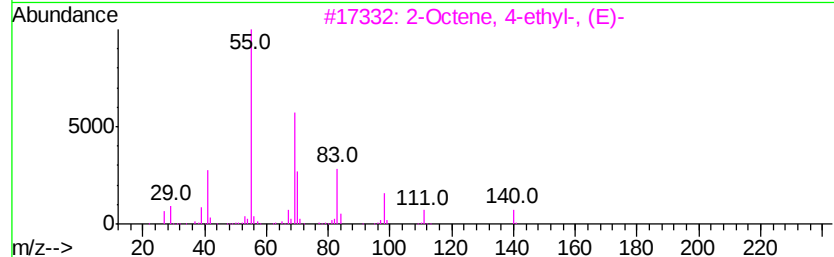
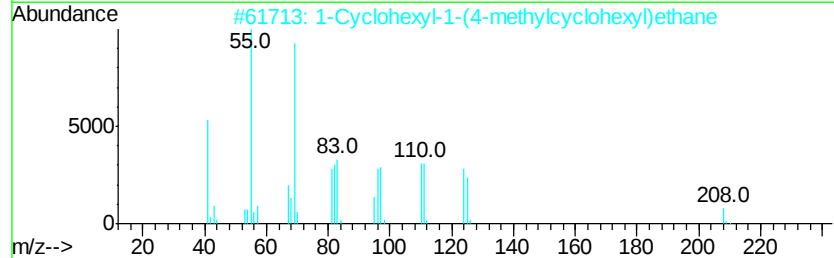
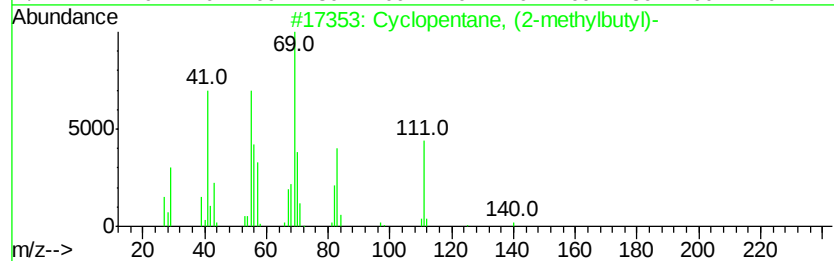
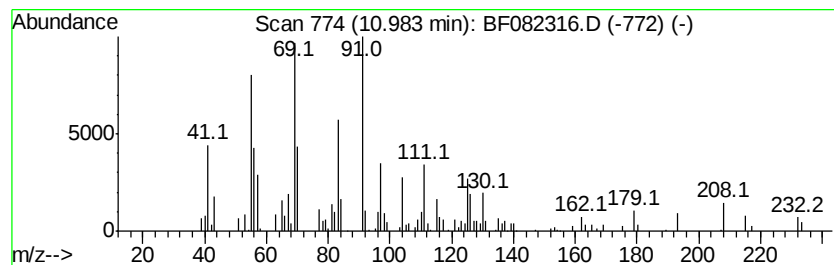
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Cyclopentane, (2-methylbutyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.98	3.06 ng	103291	Phenanthrene-d10	11.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, (2-methylbutyl)-	140	C10H20	053366-38-4	58
2		1-Cyclohexyl-1-(4-methylcyclohexyl)ethane	208	C15H28	002027-12-5	46
3		2-Octene, 4-ethyl-, (E)-	140	C10H20	074630-09-4	46
4		Cyclohexanone, 3,5-dimethyl-, cis-	126	C8H14O	007214-52-0	43
5		Triallylsilane	152	C9H16Si	001116-62-7	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
Data File : BF082316.D
Acq On : 15 Oct 2015 18:31
Operator : UM/IZ
Sample : G4015-06
Misc :
ALS Vial : 42 Sample Multiplier: 1

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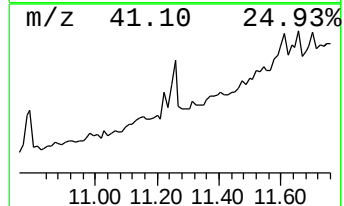
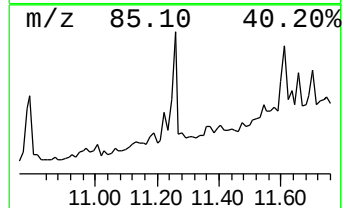
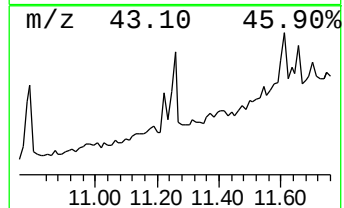
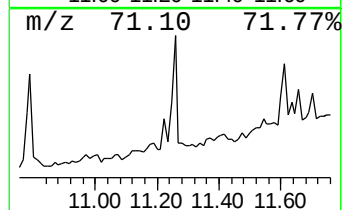
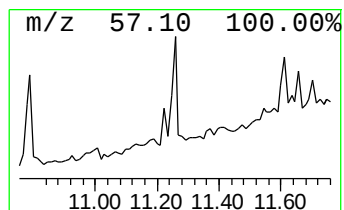
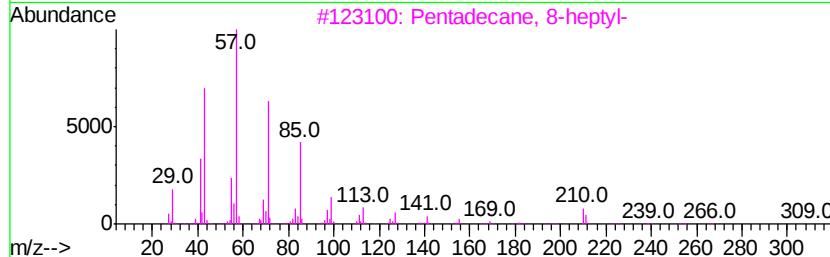
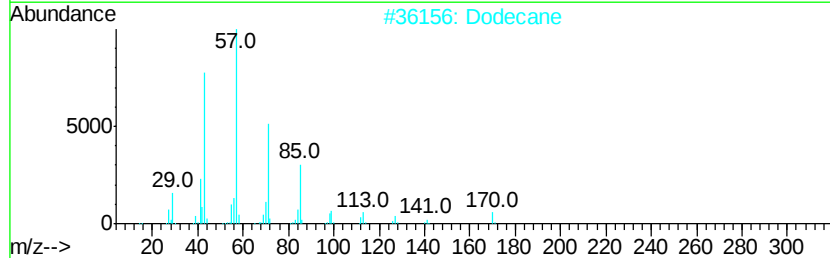
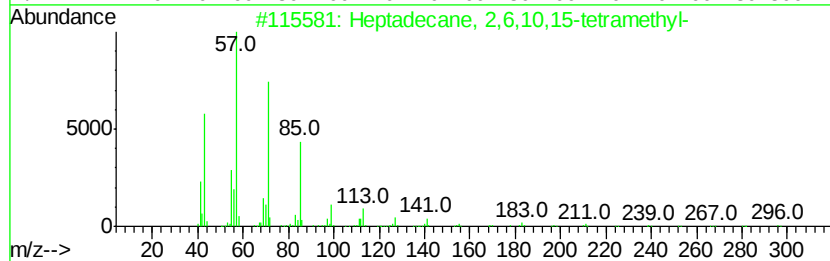
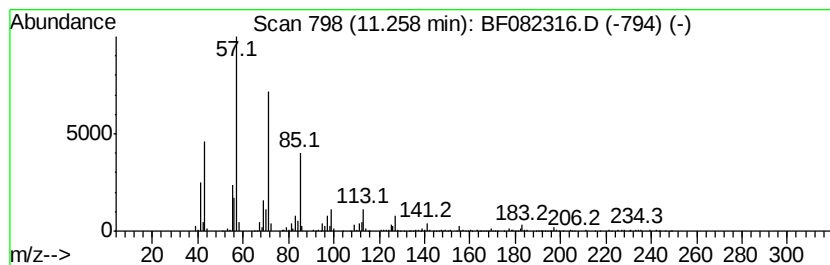
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptadecane, 2,6,10,15-tetr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.26	14.09 ng	475782	Phenanthrene-d10	11.36

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
2		Dodecane	170	C12H26	000112-40-3	90
3		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	90
4		Octadecane	254	C18H38	000593-45-3	90
5		Tridecane, 2-methyl-	198	C14H30	001560-96-9	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
Data File : BF082316.D
Acq On : 15 Oct 2015 18:31
Operator : UM/IZ
Sample : G4015-06
Misc :
ALS Vial : 42 Sample Multiplier: 1

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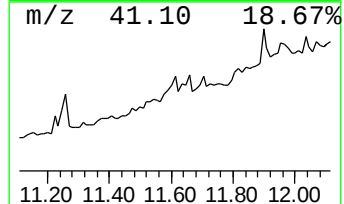
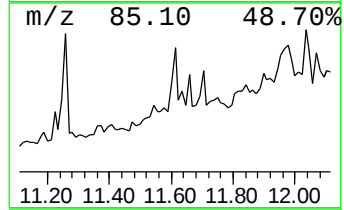
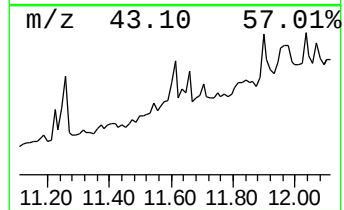
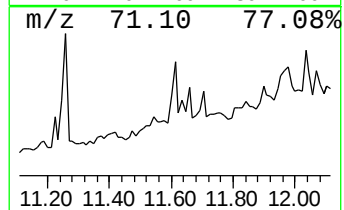
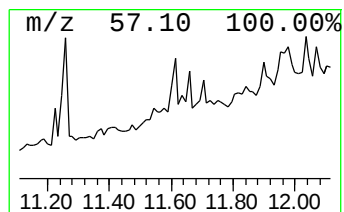
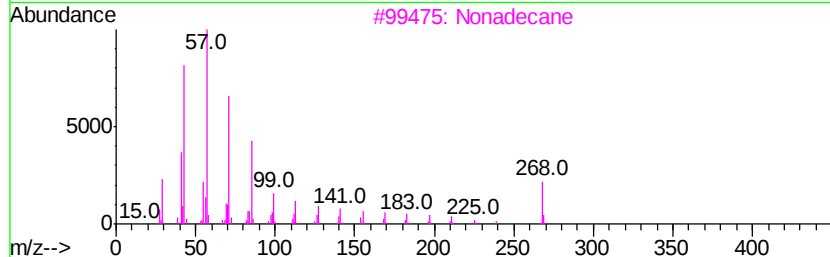
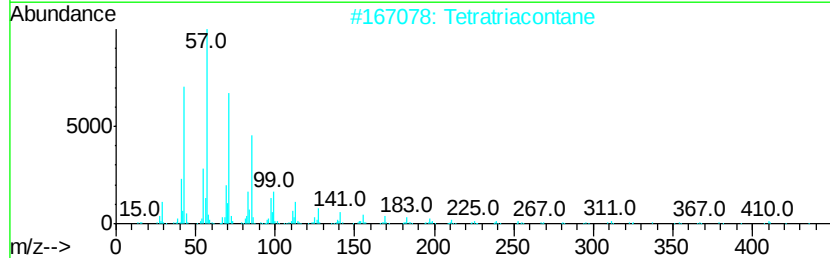
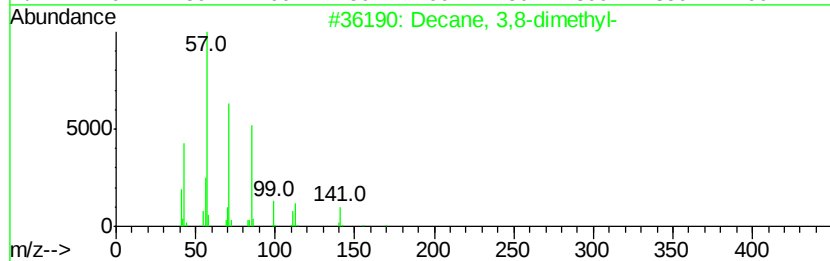
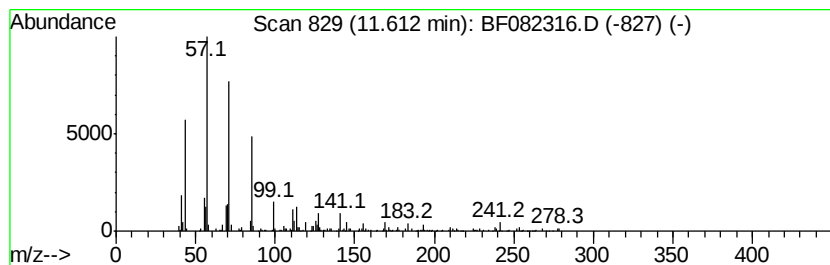
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Decane, 3,8-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.61	4.54 ng	153379	Phenanthrene-d10		11.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	94
2			Tetratriacontane	479	C34H70	014167-59-0	91
3			Nonadecane	268	C19H40	000629-92-5	90
4			Octadecane	254	C18H38	000593-45-3	87
5			Hexacosane	366	C26H54	000630-01-3	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
Data File : BF082316.D
Acq On : 15 Oct 2015 18:31
Operator : UM/IZ
Sample : G4015-06
Misc :
ALS Vial : 42 Sample Multiplier: 1

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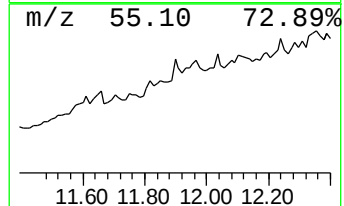
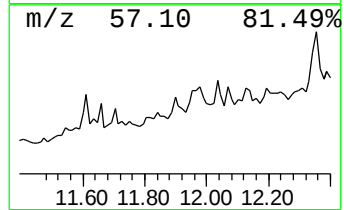
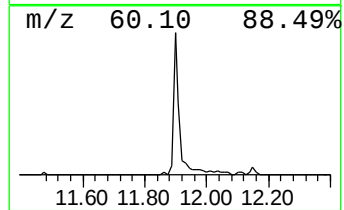
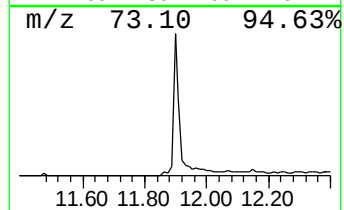
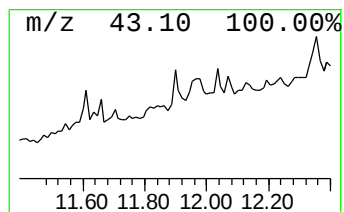
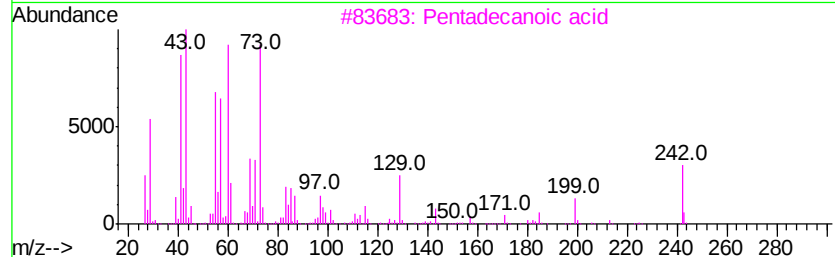
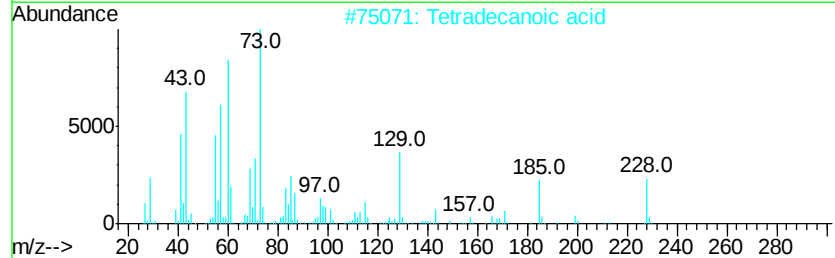
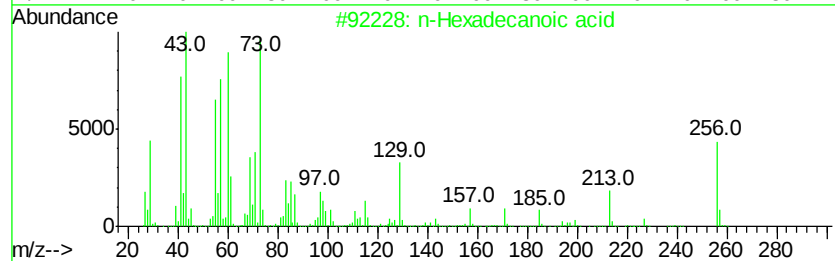
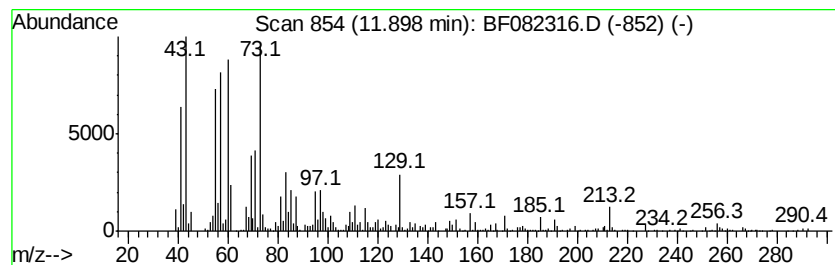
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.90	12.19 ng	411549	Phenanthrene-d10	11.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	90
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	83
4			Tridecanoic acid	214	C13H26O2	000638-53-9	70
5			Dodecanoic acid	200	C12H24O2	000143-07-7	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
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Acq On : 15 Oct 2015 18:31
Operator : UM/IZ
Sample : G4015-06
Misc :
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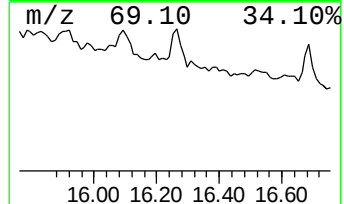
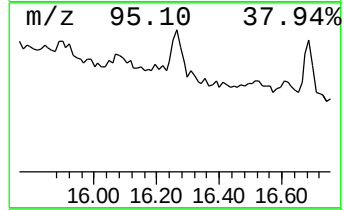
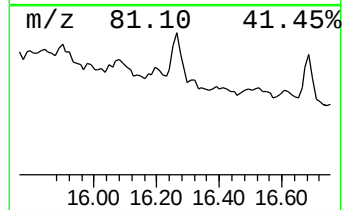
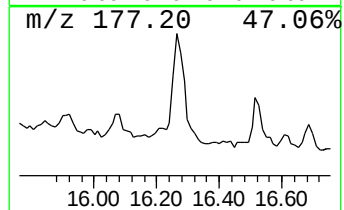
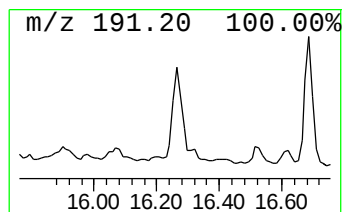
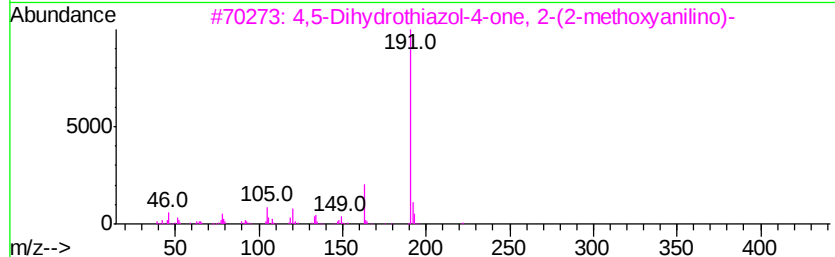
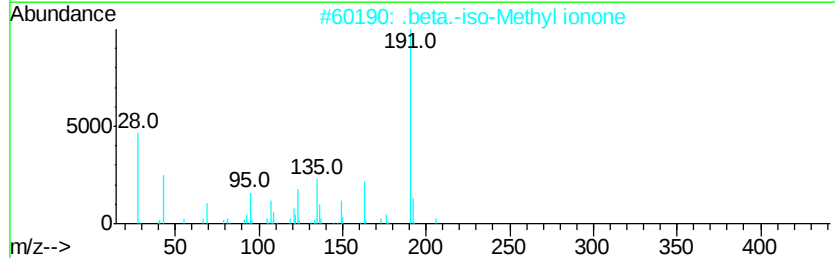
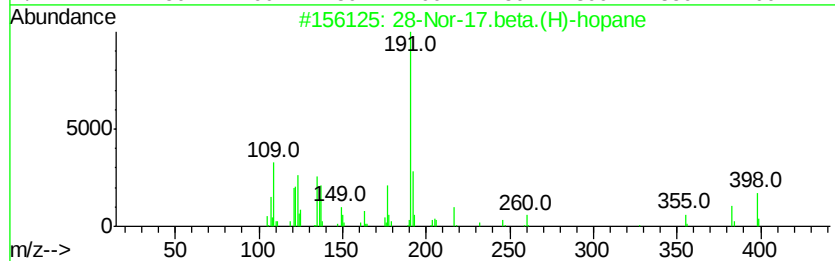
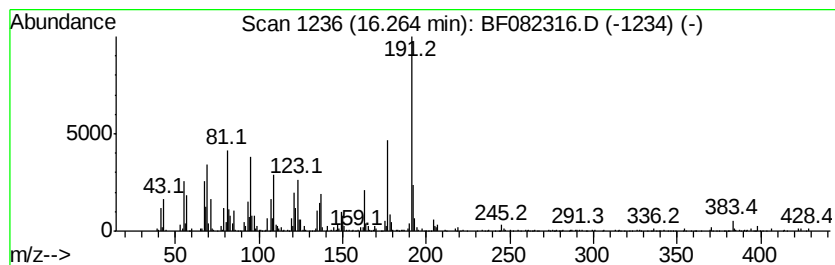
Instrument :
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ClientSampled :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 28-Nor-17.beta.(H)-hopane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.26	11.18 ng	641362	Perylene-d12	15.59
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		28-Nor-17.beta.(H)-hopane	398 C29H50	036728-72-0 53
2		.beta.-iso-Methyl ionone	206 C14H22O	1000285-40-2 43
3		4,5-Dihydrothiazol-4-one, 2-(2-m...	222 C10H10N2O2S	022476-61-5 38
4		Anthracene, 9-butyltetradecahydro-	248 C18H32	055133-89-6 38
5		1-Penten-3-one, 1-(2,6,6-trimeth...	206 C14H22O	000127-43-5 38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
Data File : BF082316.D
Acq On : 15 Oct 2015 18:31
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Sample : G4015-06
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Instrument :
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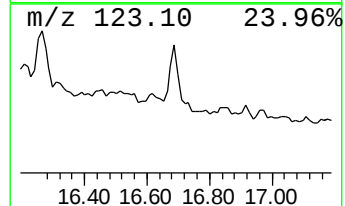
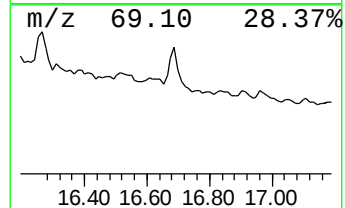
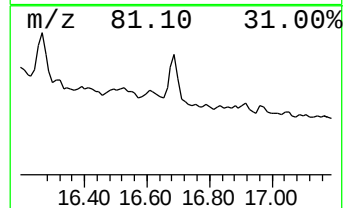
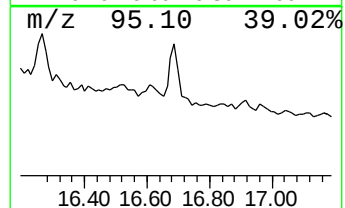
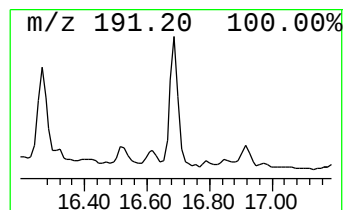
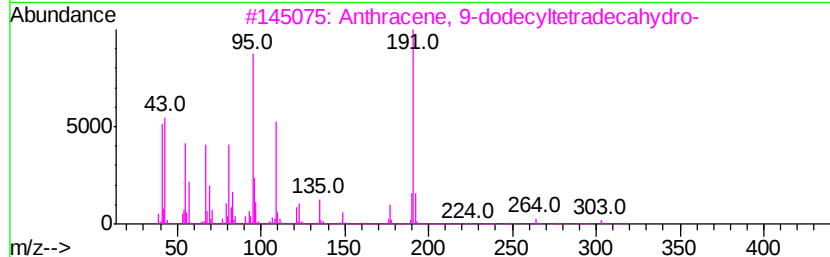
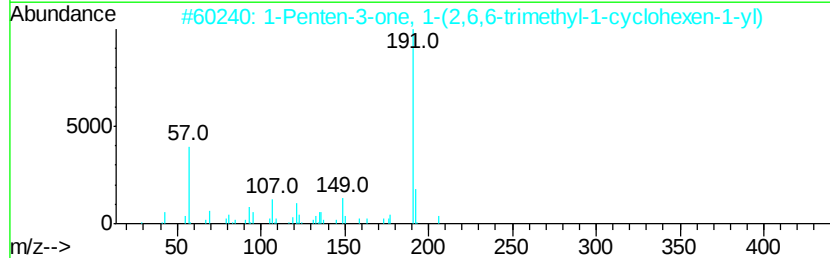
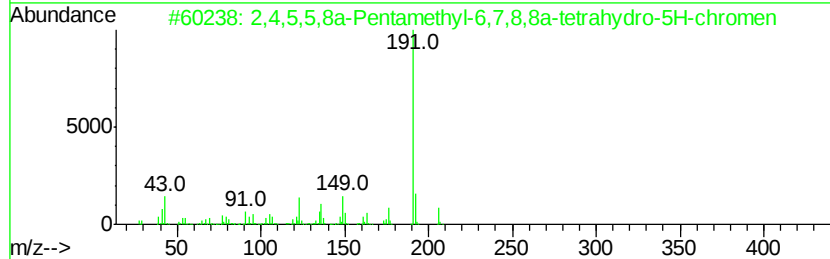
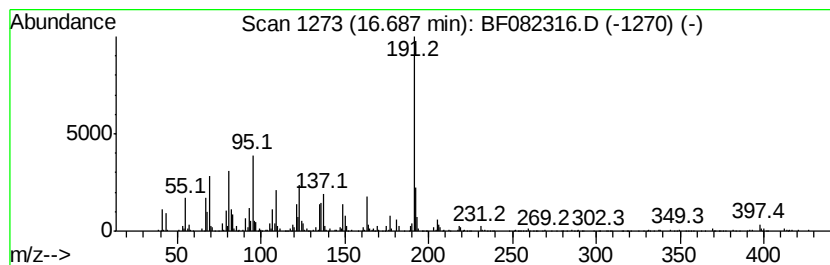
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown16.69 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.69	16.32 ng	936227	Perylene-d12	15.59

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	47		
2	1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	43		
3	Anthracene, 9-dodecyltetradecahv...	360	C26H48	055401-75-7	42		
4	28-Nor-17.beta.(H)-hopane	398	C29H50	036728-72-0	37		
5	5-(p-Aminophenyl)-2-thiazolamine	191	C9H9N3S	090349-87-4	35		



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
 Data File : BF082316.D
 Acq On : 15 Oct 2015 18:31
 Operator : UM/IZ
 Sample : G4015-06
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LF-2

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.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...	4.99	42.2	ng	1000980	1	6.82	474530	20.0
unknown6.59	6.59	77.3	ng	1833160	1	6.82	474530	20.0
Bicyclo[3.1.0]hex...	8.67	2.8	ng	87069	2	8.11	629874	20.0
Hexacosane	10.53	2.4	ng	84033	3	9.87	702083	20.0
Pentadecane, 2,6,...	10.79	8.8	ng	295764	4	11.36	675174	20.0
4-Chlorobenzenesu...	10.87	2.6	ng	89311	4	11.36	675174	20.0
Cyclopentane, (2-...	10.98	3.1	ng	103291	4	11.36	675174	20.0
Heptadecane, 2,6,...	11.26	14.1	ng	475782	4	11.36	675174	20.0
Decane, 3,8-dimet...	11.61	4.5	ng	153379	4	11.36	675174	20.0
n-Hexadecanoic acid	11.90	12.2	ng	411549	4	11.36	675174	20.0
28-Nor-17.beta.(H...	16.26	11.2	ng	641362	6	15.59	1147100	20.0
unknown16.69	16.69	16.3	ng	936227	6	15.59	1147100	20.0