

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121516\  
 Data File : BF091651.D  
 Acq On : 16 Dec 2016 1:45  
 Operator : UM/SJ  
 Sample : H5940-09  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SS05-1FT

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-BF120916.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	4.956	248	251	262	rBV	1714141	2534838	37.65%	4.083%
2	5.162	268	269	275	rVB	53603	73429	1.09%	0.118%
3	5.390	286	289	291	rBV	6222389	6360373	94.47%	10.244%
4	6.430	377	380	383	rVB	5810150	6457499	95.91%	10.401%
5	6.556	388	391	393	rBV	6658584	6581958	97.76%	10.601%
6	6.785	409	411	413	rBV	1592142	1346637	20.00%	2.169%
7	6.945	422	425	427	rBV	5017720	4979757	73.96%	8.021%
8	7.356	457	461	463	rBV	3504844	4402598	65.39%	7.091%
9	7.459	466	470	474	rVB	128081	175442	2.61%	0.283%
10	8.065	521	523	526	rBV	1426409	1978609	29.39%	3.187%
11	9.151	615	618	620	rBV	7277765	6732700	100.00%	10.844%
12	9.551	650	653	655	rBV	1918688	2211641	32.85%	3.562%
13	9.825	674	677	679	rBV	2419646	2040533	30.31%	3.287%
14	10.259	714	715	717	rVB	121529	148295	2.20%	0.239%
15	10.476	731	734	738	rBV	51783	115850	1.72%	0.187%
16	10.614	743	746	748	rBV	3437733	3457769	51.36%	5.569%
17	10.739	755	757	761	rVB	162000	223728	3.32%	0.360%
18	11.185	794	796	797	rBV	109652	99927	1.48%	0.161%
19	11.299	804	806	809	rBV	1162778	1232884	18.31%	1.986%
20	11.539	824	827	830	rBV	194598	208454	3.10%	0.336%
21	11.848	852	854	858	rBV2	114453	166844	2.48%	0.269%
22	12.511	910	912	914	rVB	104935	95945	1.43%	0.155%
23	12.739	929	932	934	rBV	75521	86563	1.29%	0.139%
24	12.888	943	945	948	rBV	4111007	4977198	73.93%	8.017%
25	13.791	1022	1024	1030	rBV	421369	460428	6.84%	0.742%
26	13.940	1034	1037	1039	rBV	1425148	1649364	24.50%	2.657%
27	14.042	1044	1046	1048	rBV	79767	95882	1.42%	0.154%
28	14.420	1077	1079	1085	rVB3	71093	122526	1.82%	0.197%
29	14.854	1115	1117	1119	rVB	105133	106550	1.58%	0.172%
30	15.037	1130	1133	1138	rBV	223157	310247	4.61%	0.500%
31	15.345	1157	1160	1162	rBV	832109	950928	14.12%	1.532%
32	15.391	1162	1164	1166	rVB	57364	72965	1.08%	0.118%
33	15.574	1178	1180	1182	rBV	58651	79572	1.18%	0.128%
34	15.791	1197	1199	1202	rVB	291273	462621	6.87%	0.745%

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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	16.808	1285	1288	1291	rBV	67138	137536	2.04%	0.222%
36	17.243	1324	1326	1332	rVB6	35896	75835	1.13%	0.122%
37	17.448	1341	1344	1349	rVB3	61915	167159	2.48%	0.269%
38	17.677	1361	1364	1370	rVB4	63957	208209	3.09%	0.335%
39	19.163	1489	1494	1500	rVB4	145900	496949	7.38%	0.800%

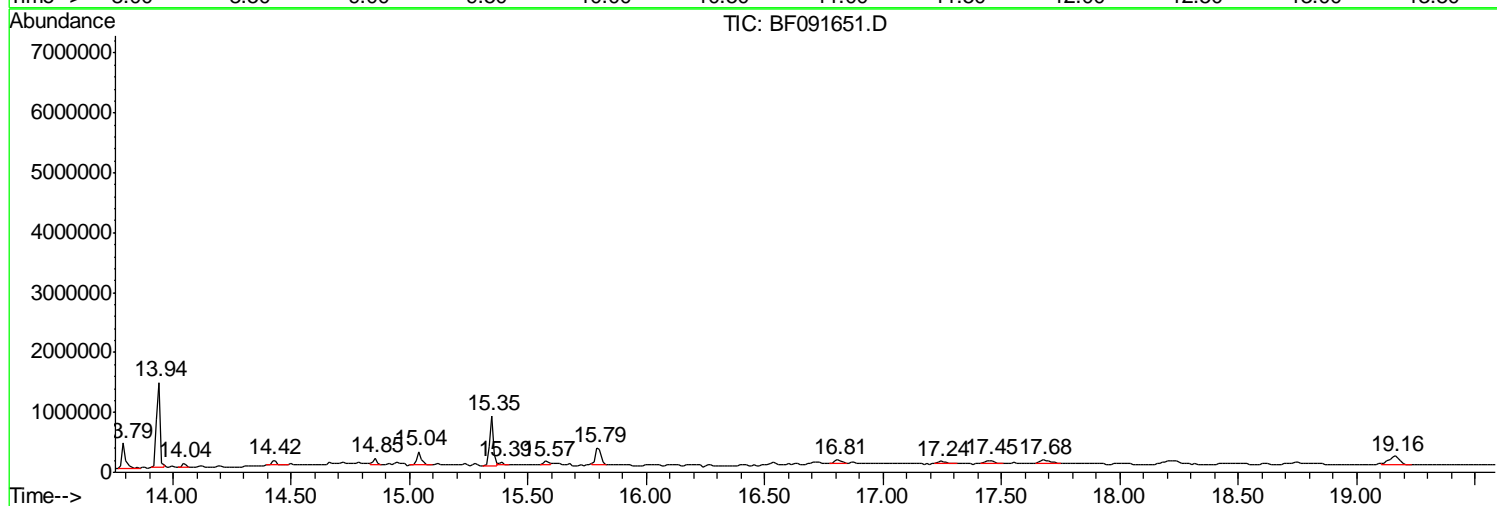
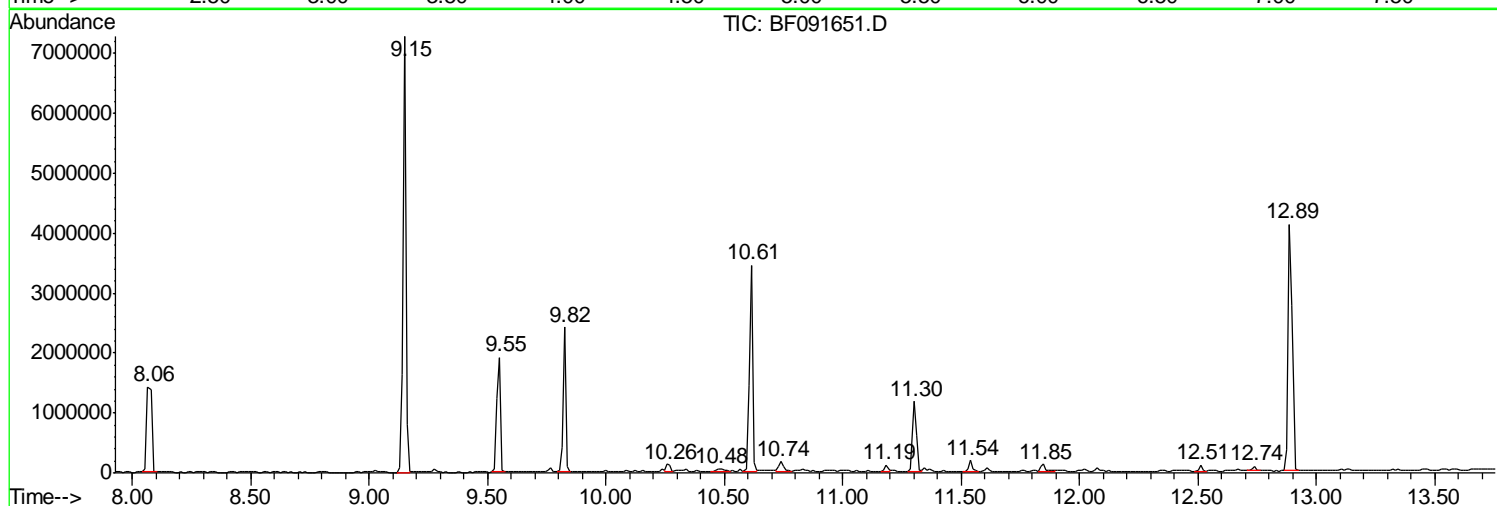
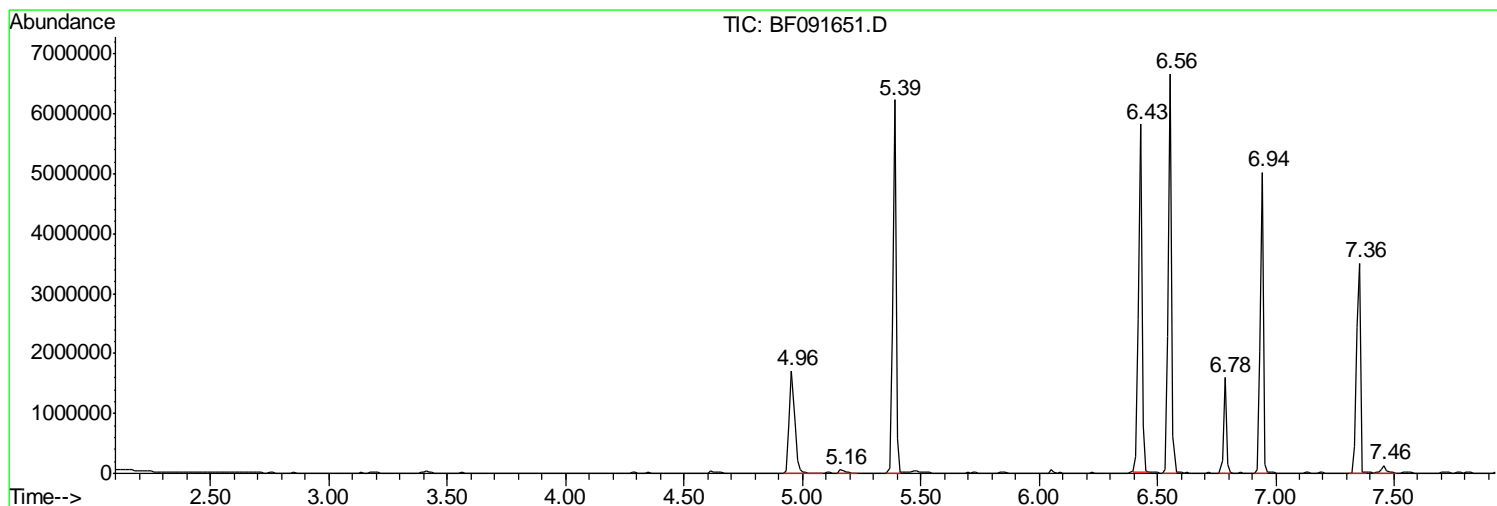
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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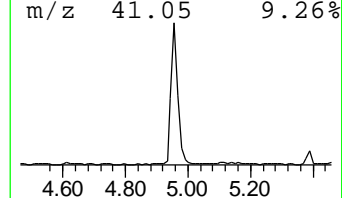
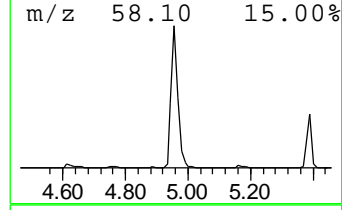
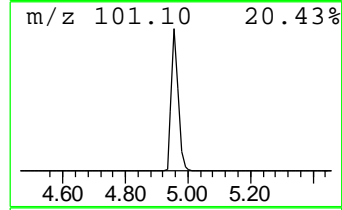
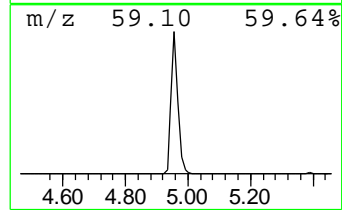
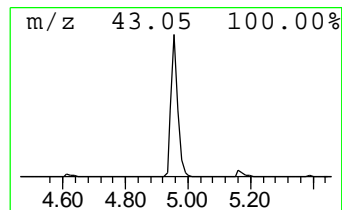
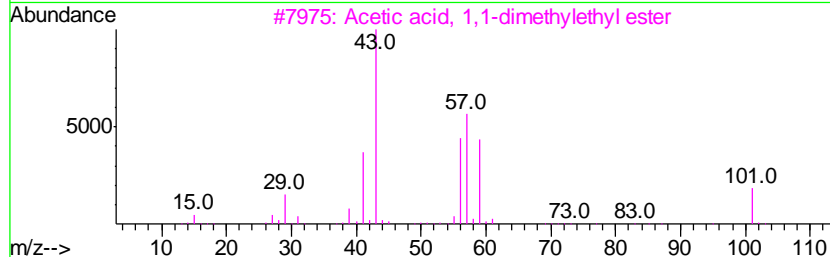
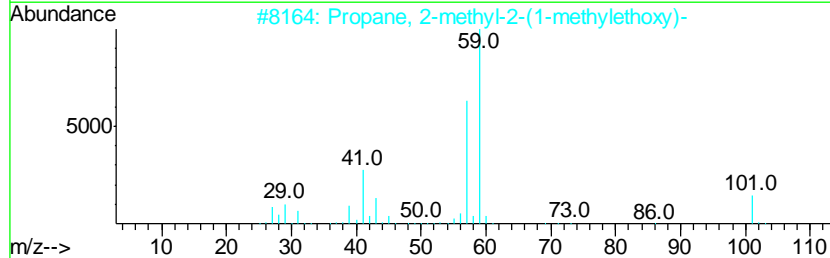
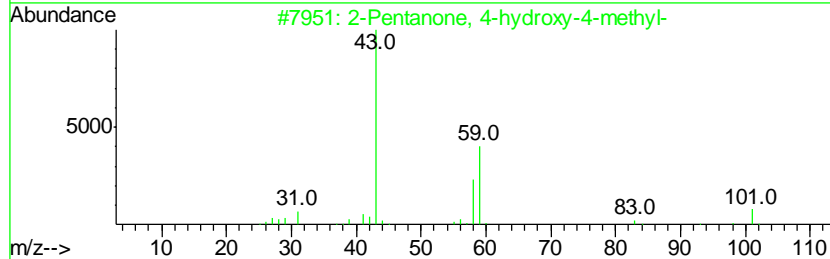
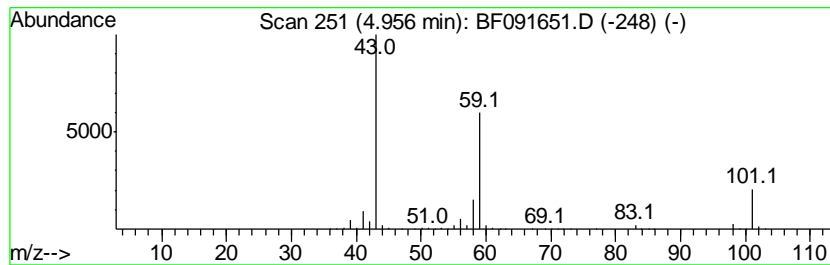
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	37.65 ng	2534840	1,4-Dichlorobenzene-d4	6.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	42
3		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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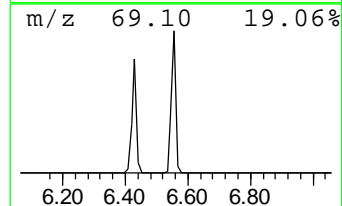
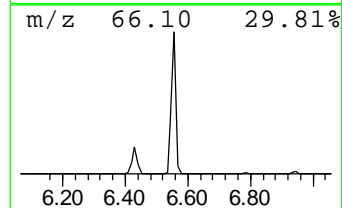
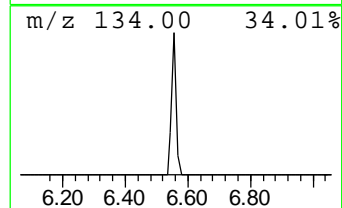
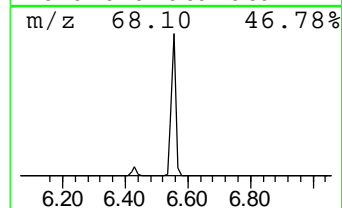
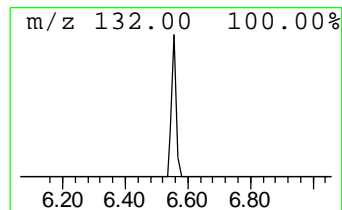
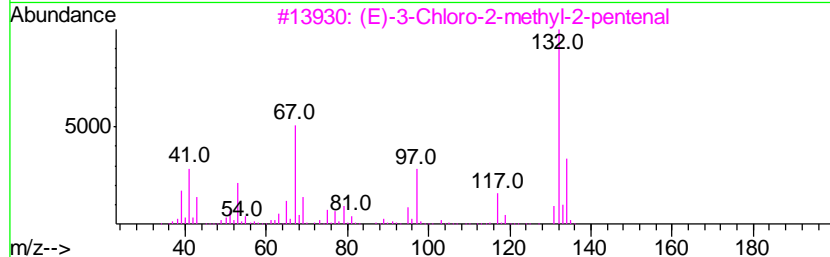
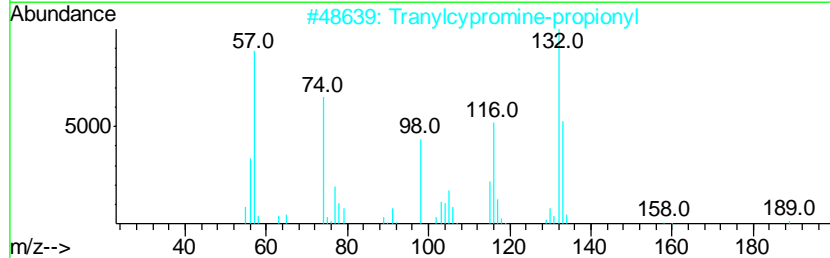
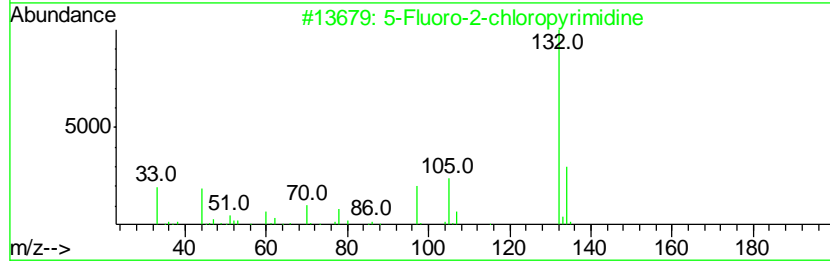
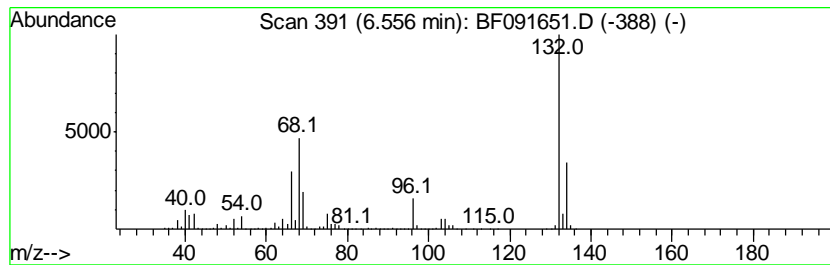
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 Peak Number 2 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	97.75 ng	6581960	1,4-Dichlorobenzene-d4	6.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		Tranlycypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

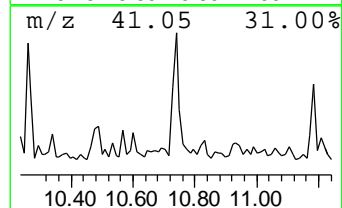
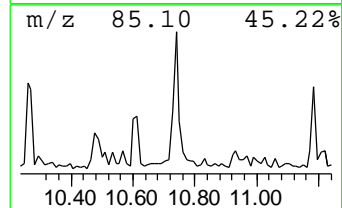
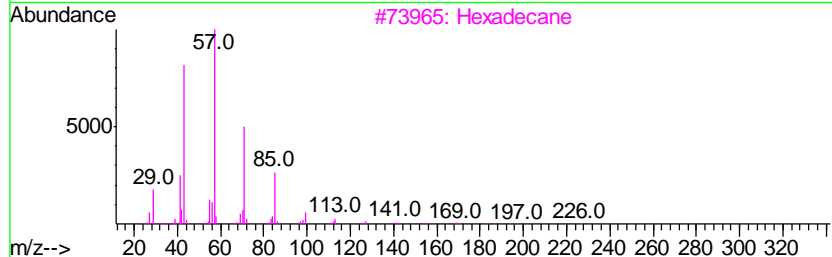
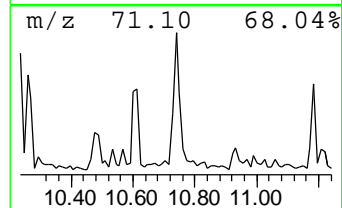
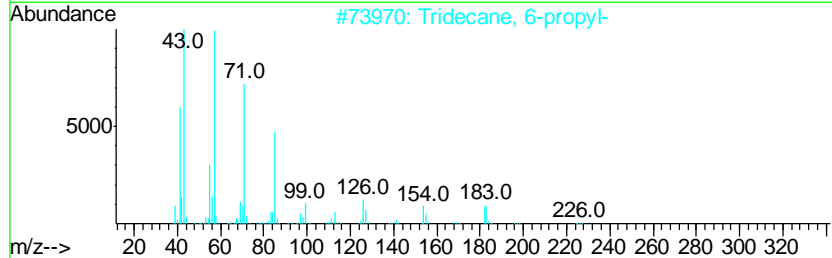
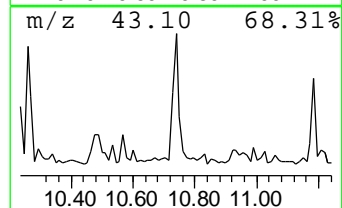
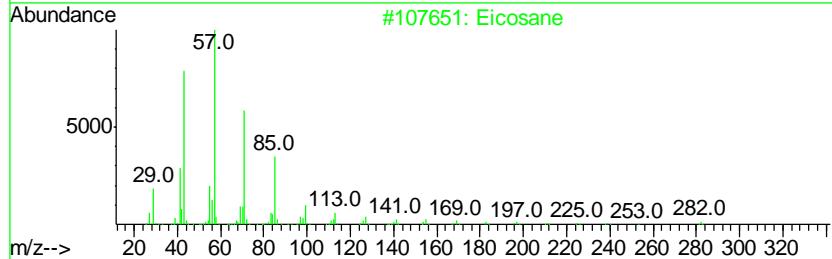
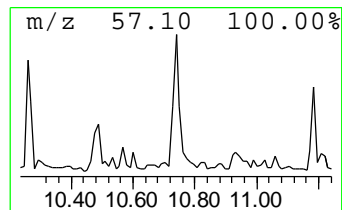
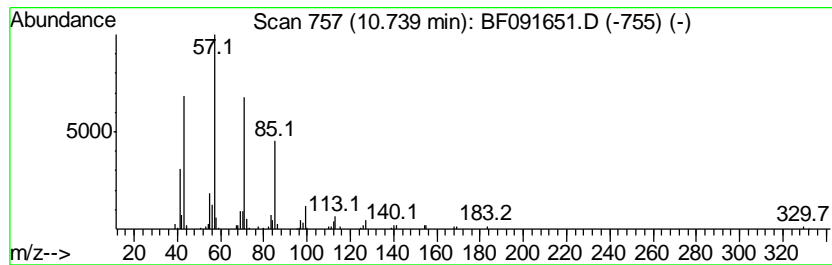
TIC Library : C:\DATABASE\NIST02.L  
 TIC Integration Parameters: LSCINT.P

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 Peak Number 4 Eicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.74	3.63 ng	223728	Phenanthrene-d10	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	90
2		Tridecane, 6-propyl-	226	C16H34	055045-10-8	90
3		Hexadecane	226	C16H34	000544-76-3	90
4		Tetracosane	338	C24H50	000646-31-1	86
5		Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	86



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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M  
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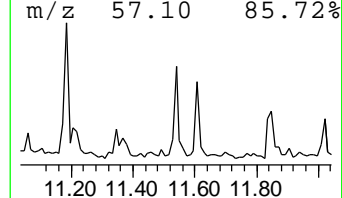
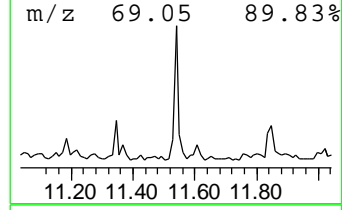
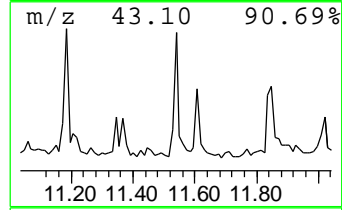
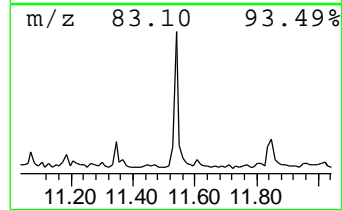
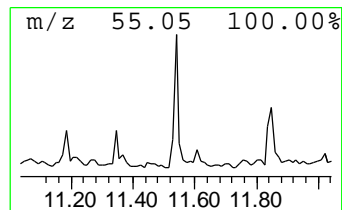
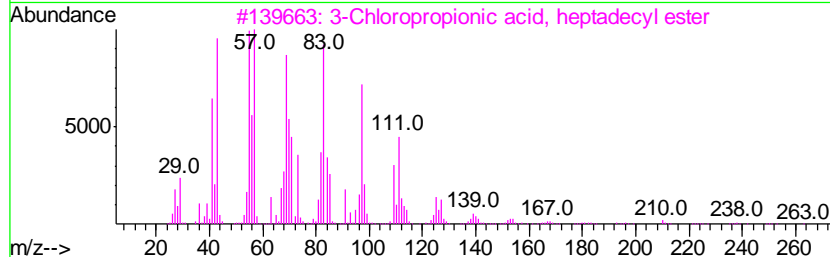
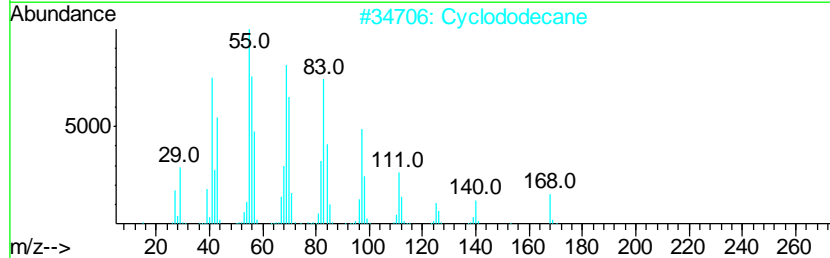
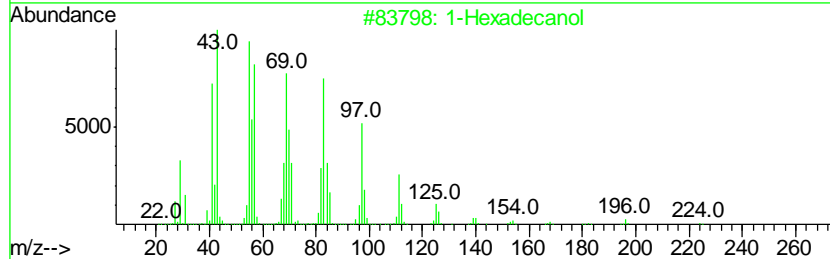
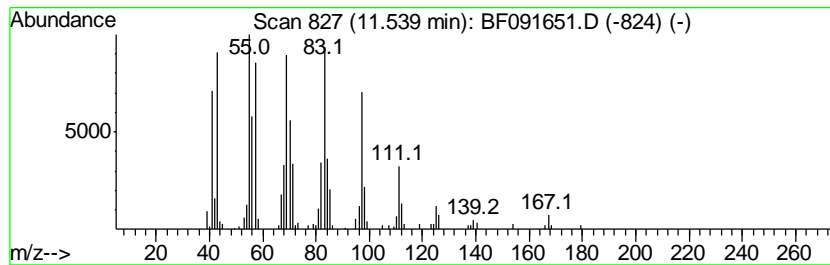
TIC Library : C:\DATABASE\NIST02.L  
 TIC Integration Parameters: LSCINT.P

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 Peak Number 5 1-Hexadecanol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.54	3.38 ng	208454	Phenanthrene-d10	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexadecanol	242	C16H34O	036653-82-4	95
2		Cyclododecane	168	C12H24	000294-62-2	94
3		3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	94
4		1-Tetradecanol	214	C14H30O	000112-72-1	91
5		5-Octadecene, (E)-	252	C18H36	007206-21-5	91



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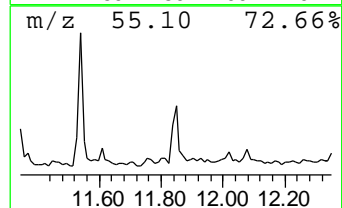
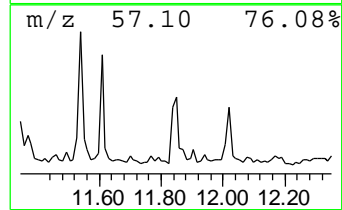
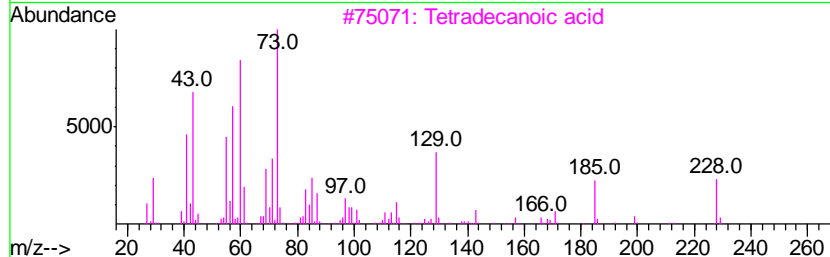
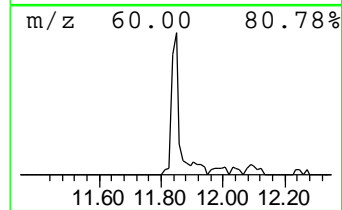
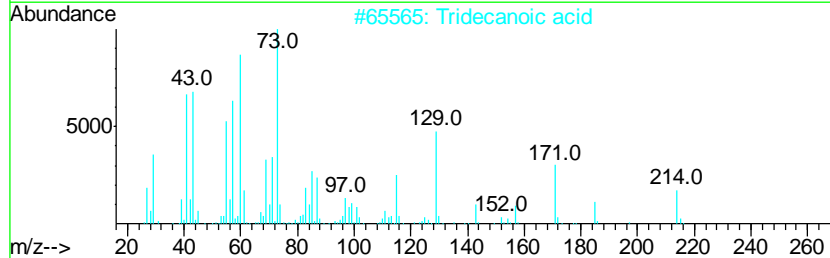
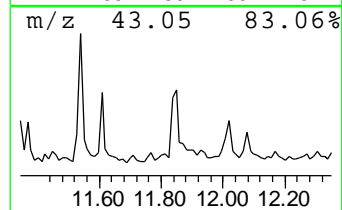
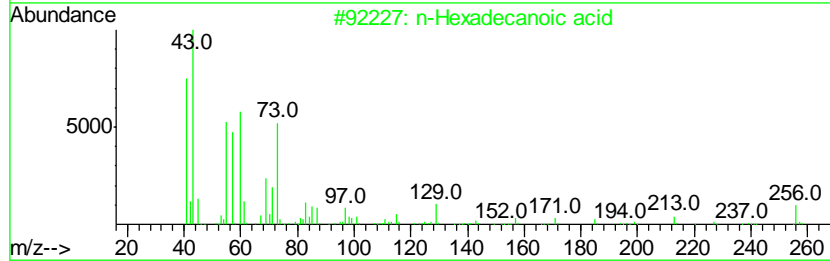
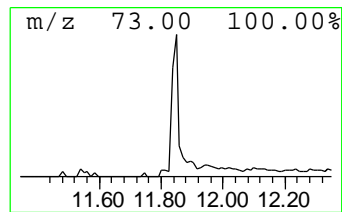
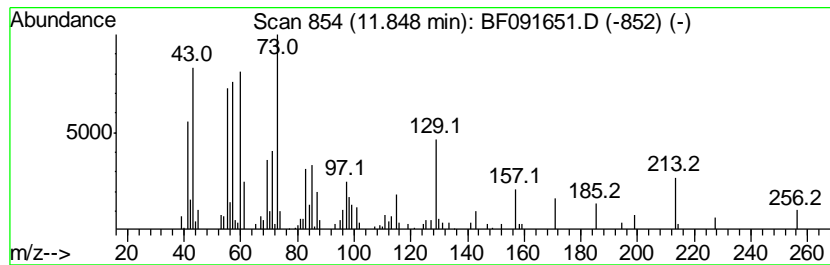
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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 Peak Number 6 n-Hexadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.85	2.71 ng	166844	Phenanthrene-d10	11.30		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
2		Tridecanoic acid	214	C13H26O2	000638-53-9	76
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	72
4		Dodecanoic acid	200	C12H24O2	000143-07-7	64
5		Amyl Nitrite	117	C5H11NO2	000110-46-3	27





Data Path : Z:\HPCHEM1\BNA F\DATA\BF121516\  
 Data File : BF091651.D  
 Acq On : 16 Dec 2016 1:45  
 Operator : UM/SJ  
 Sample : H5940-09  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleID :  
 SS05-1FT

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

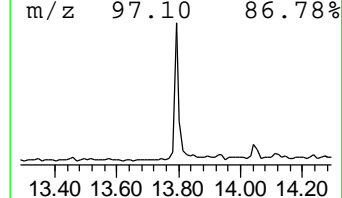
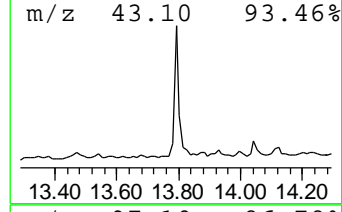
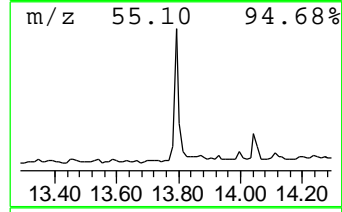
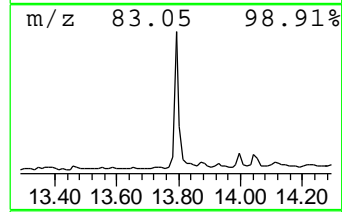
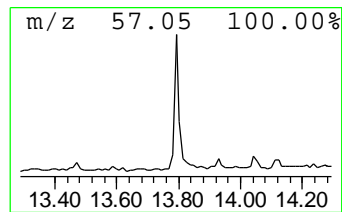
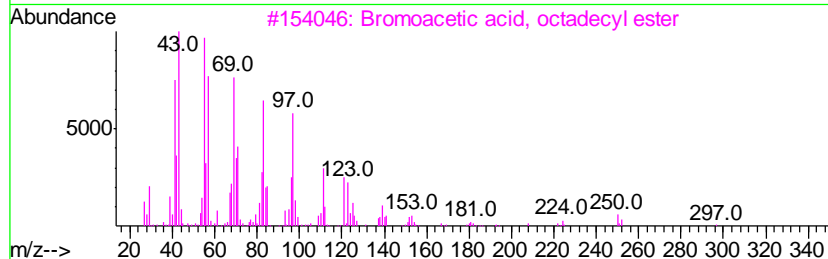
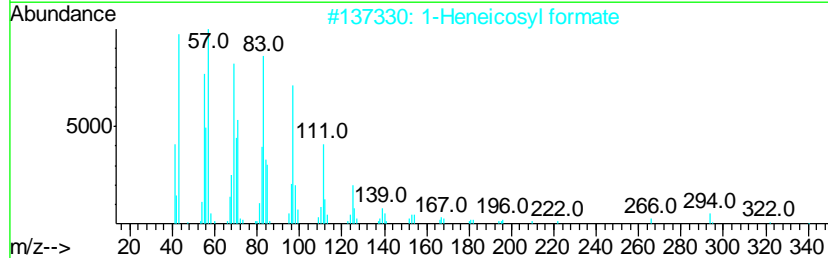
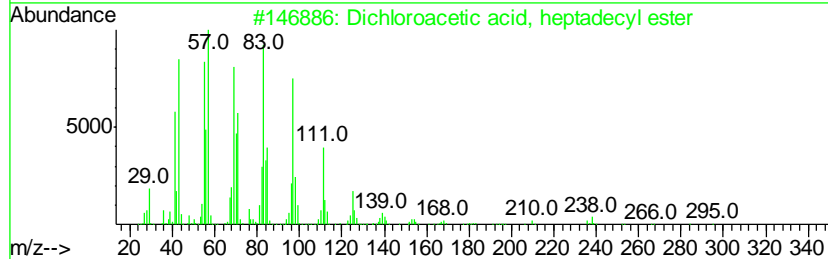
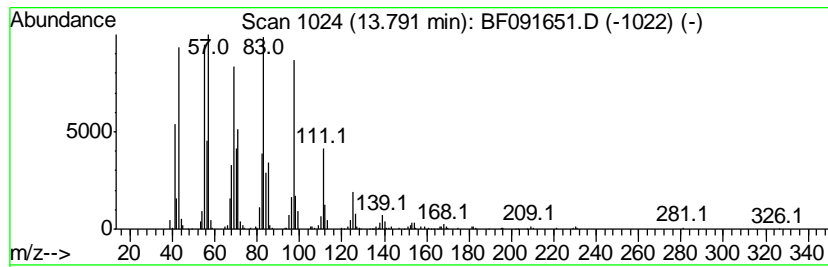
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 TIC Integration Parameters: LSCINT.P

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 Peak Number 7 Dichloroacetic acid, heptad... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.79	5.58 ng	460428	Chrysene-d12	13.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	94
2		1-Heneicosyl formate	340	C22H44O2	077899-03-7	91
3		Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	91
4		1-Nonadecene	266	C19H38	018435-45-5	90
5		Trichloroacetic acid, tetradecyl...	358	C16H29Cl3O2	074339-52-9	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121516\  
 Data File : BF091651.D  
 Acq On : 16 Dec 2016 1:45  
 Operator : UM/SJ  
 Sample : H5940-09  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleID :  
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

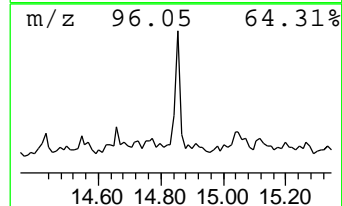
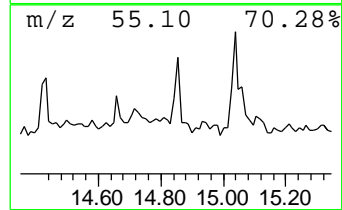
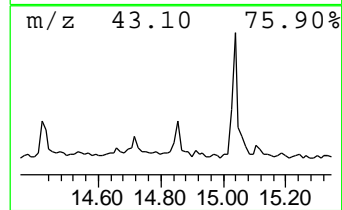
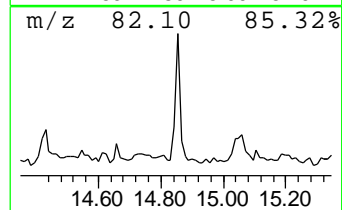
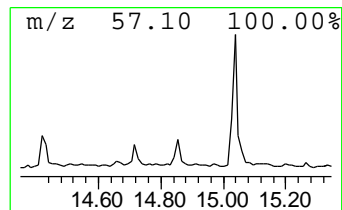
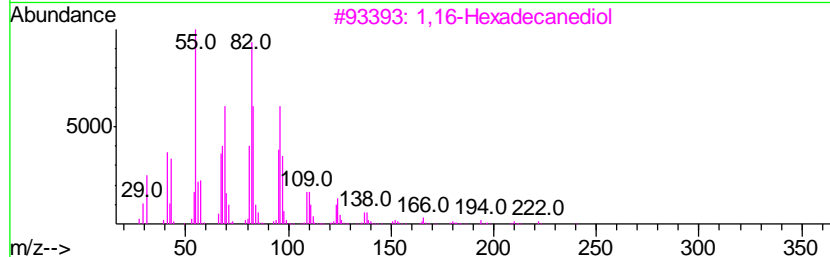
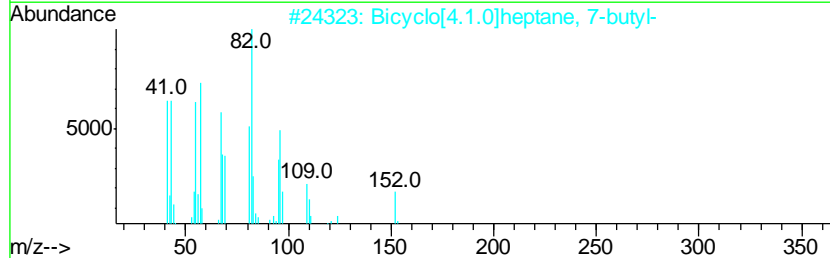
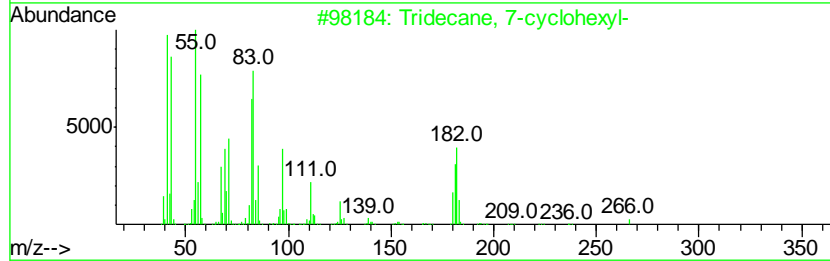
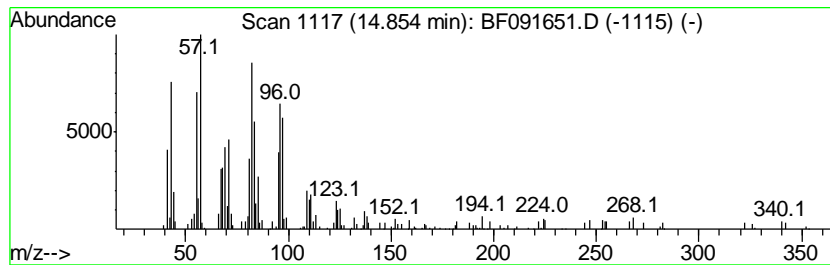
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 Peak Number 8 Tridecane, 7-cyclohexyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	2.24 ng	106550	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 7-cyclohexyl-	266	C19H38	013151-92-3	83
2		Bicyclo[4.1.0]heptane, 7-butyl-	152	C11H20	018645-10-8	64
3		1,16-Hexadecanediol	258	C16H34O2	007735-42-4	62
4		Oxirane, heptadecyl-	282	C19H38O	067860-04-2	58
5		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	58



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121516\  
 Data File : BF091651.D  
 Acq On : 16 Dec 2016 1:45  
 Operator : UM/SJ  
 Sample : H5940-09  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleID :  
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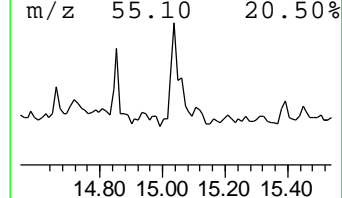
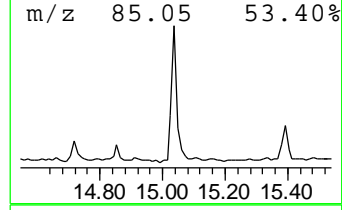
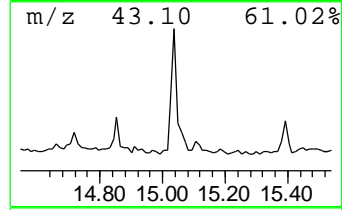
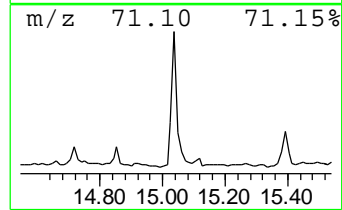
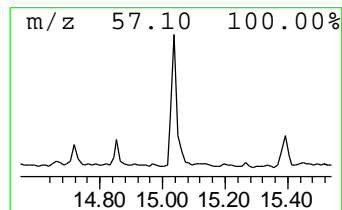
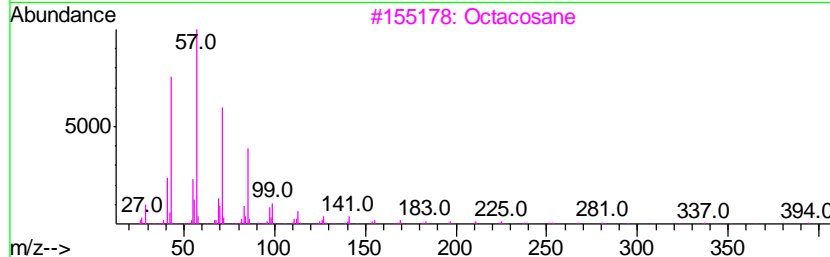
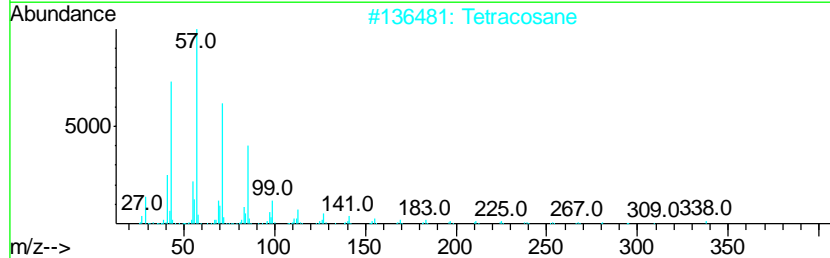
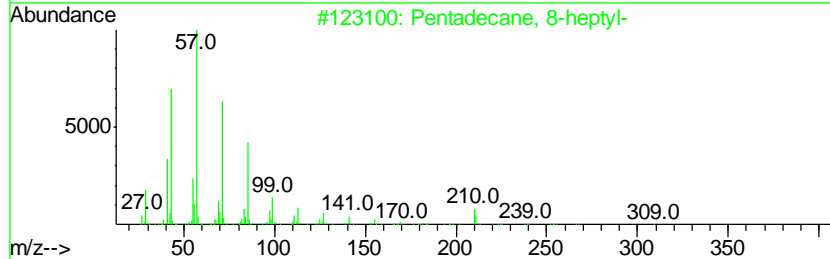
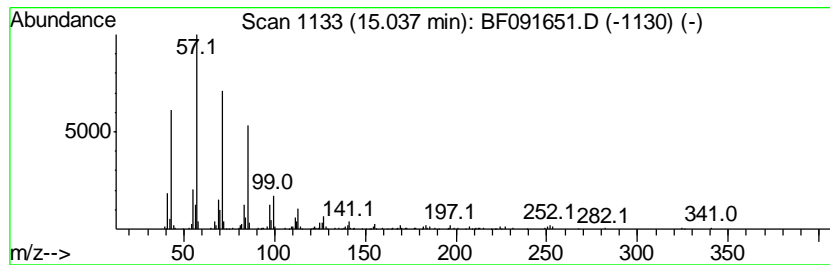
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 TIC Integration Parameters: LSCINT.P

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 Peak Number 9 Pentadecane, 8-heptyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.04	6.53 ng	310247	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
2		Tetracosane	338	C24H50	000646-31-1	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Heneicosane	296	C21H44	000629-94-7	90
5		Eicosane	282	C20H42	000112-95-8	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121516\  
 Data File : BF091651.D  
 Acq On : 16 Dec 2016 1:45  
 Operator : UM/SJ  
 Sample : H5940-09  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleID :  
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M  
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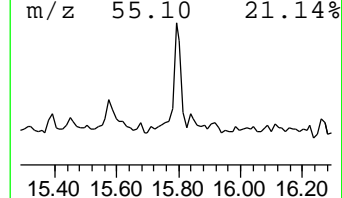
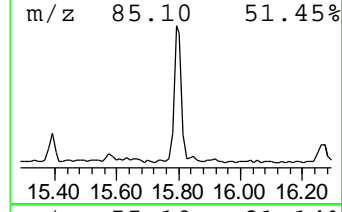
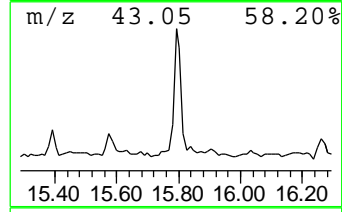
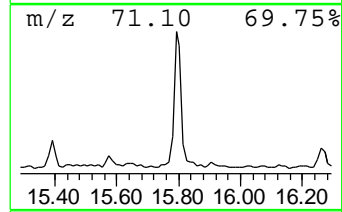
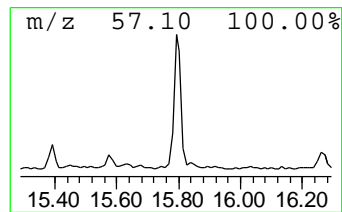
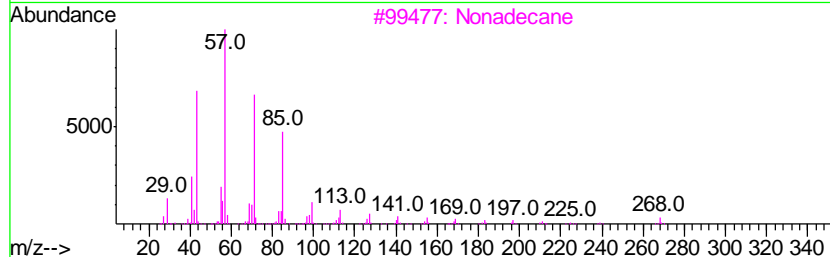
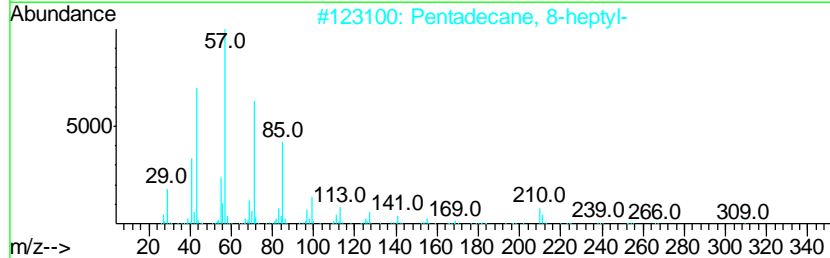
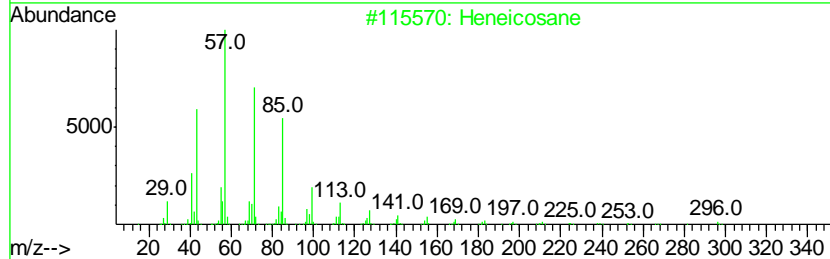
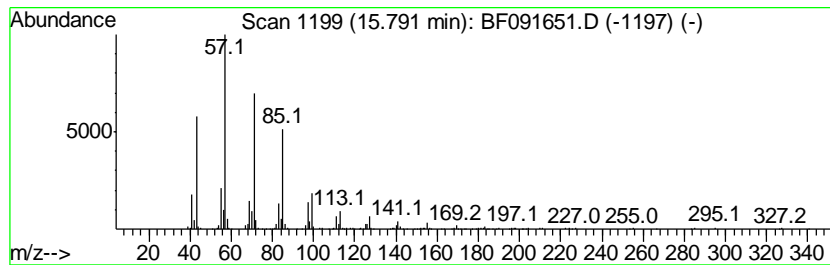
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 TIC Integration Parameters: LSCINT.P

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 Peak Number 10 Heneicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.79	9.73 ng	462621	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
3		Nonadecane	268	C19H40	000629-92-5	86
4		Tetratetracontane	619	C44H90	007098-22-8	80
5		Octacosane	394	C28H58	000630-02-4	74



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121516\  
 Data File : BF091651.D  
 Acq On : 16 Dec 2016 1:45  
 Operator : UM/SJ  
 Sample : H5940-09  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
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 ClientSampleID :  
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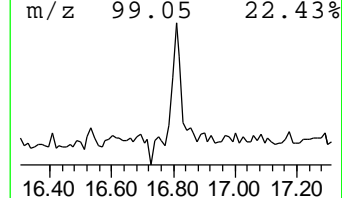
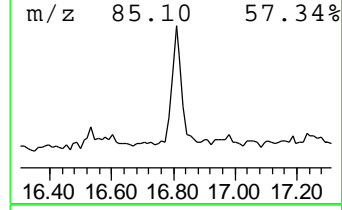
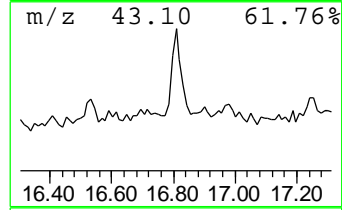
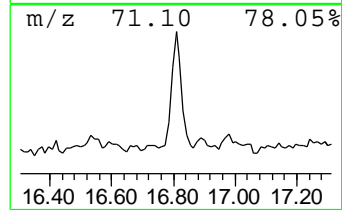
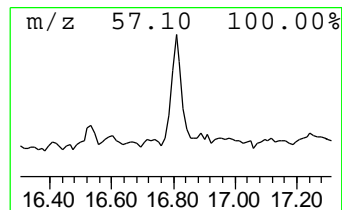
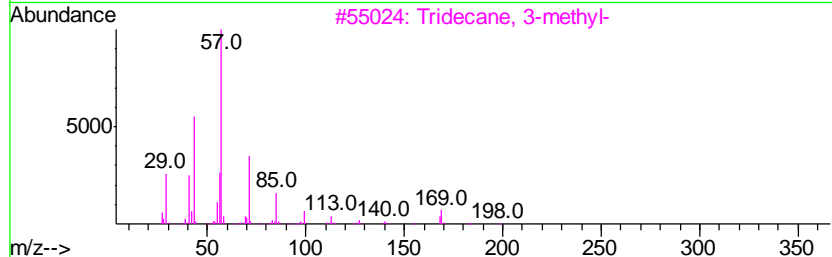
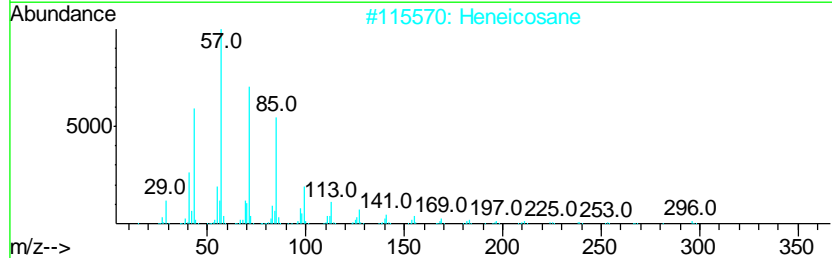
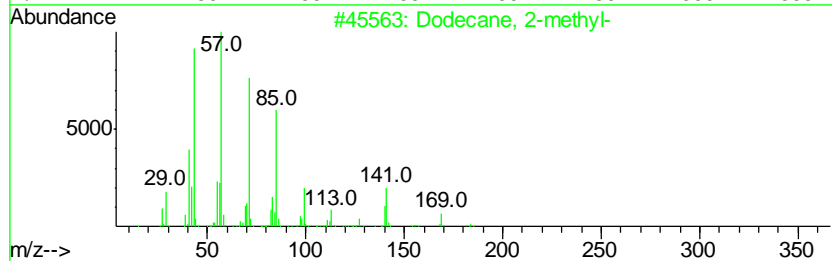
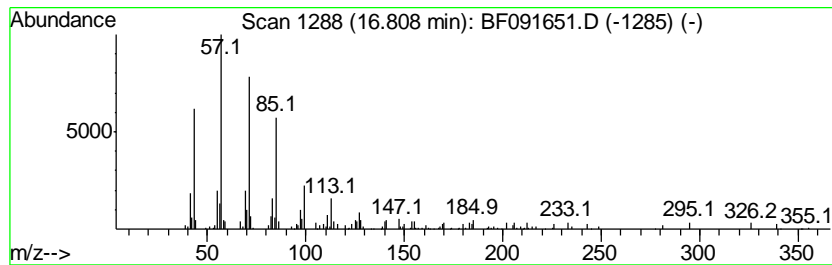
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 TIC Integration Parameters: LSCINT.P

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 Peak Number 11 Dodecane, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.81	2.89 ng	137536	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane, 2-methyl-	184	C13H28	001560-97-0	90
2		Heneicosane	296	C21H44	000629-94-7	87
3		Tridecane, 3-methyl-	198	C14H30	006418-41-3	86
4		Heptadecane	240	C17H36	000629-78-7	86
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	83



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121516\  
 Data File : BF091651.D  
 Acq On : 16 Dec 2016 1:45  
 Operator : UM/SJ  
 Sample : H5940-09  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

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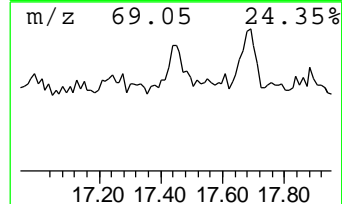
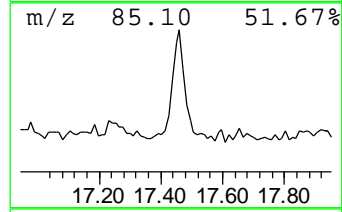
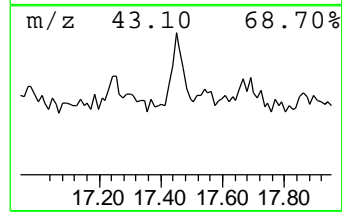
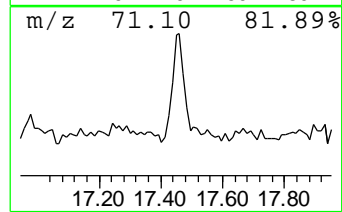
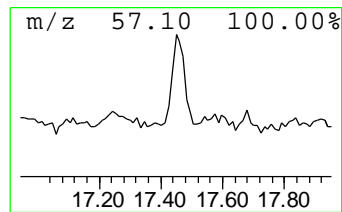
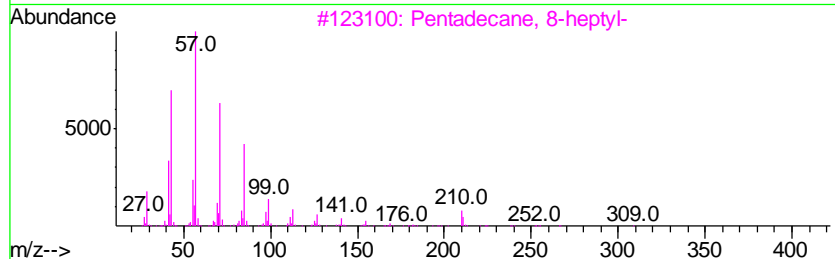
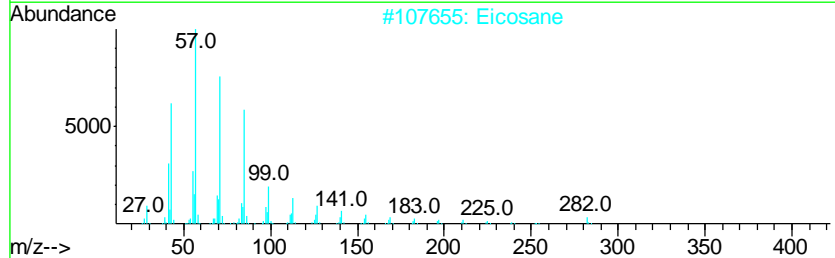
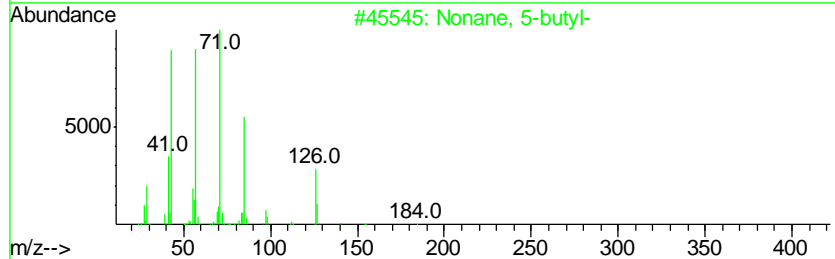
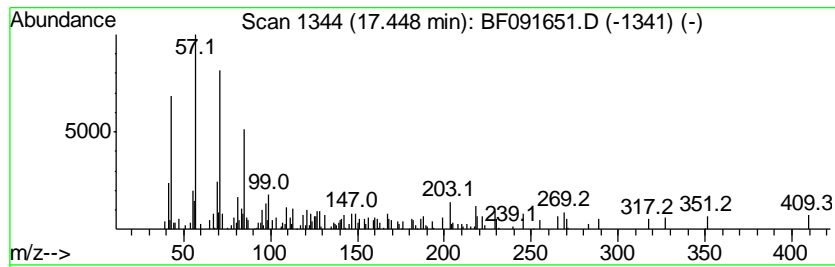
TIC Library : C:\DATABASE\NIST02.L  
 TIC Integration Parameters: LSCINT.P

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 Peak Number 12 Nonane, 5-butyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.45	3.52 ng	167159	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonane, 5-butyl-	184	C13H28	017312-63-9	76
2		Eicosane	282	C20H42	000112-95-8	68
3		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	64
4		Hexadecane	226	C16H34	000544-76-3	64
5		Nonacosane	408	C29H60	000630-03-5	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121516\  
 Data File : BF091651.D  
 Acq On : 16 Dec 2016 1:45  
 Operator : UM/SJ  
 Sample : H5940-09  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleID :  
 SS05-1FT

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

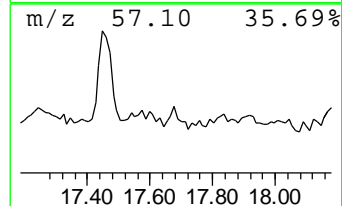
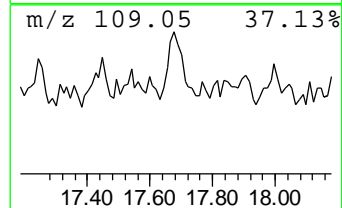
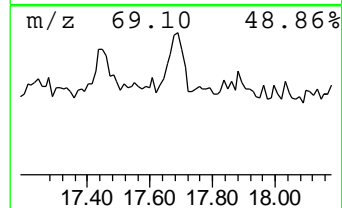
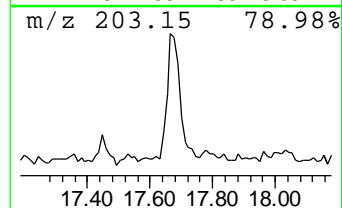
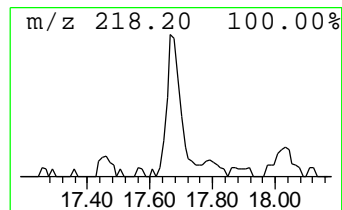
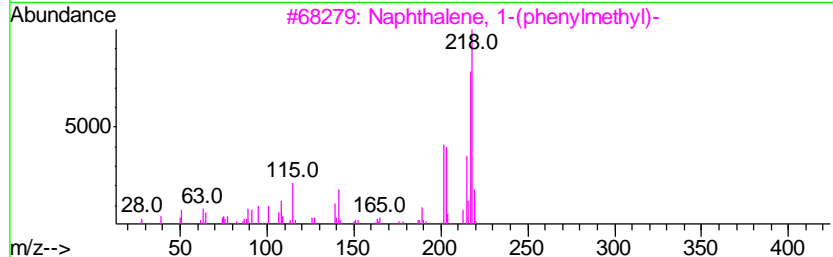
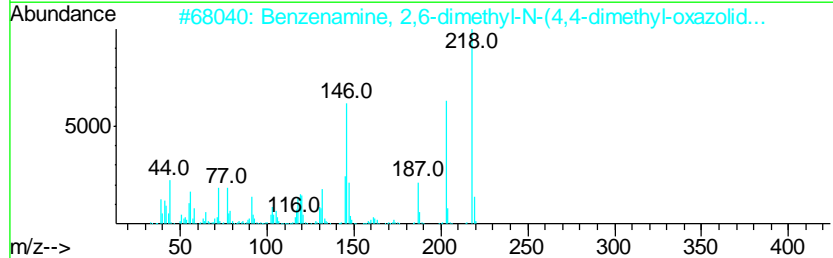
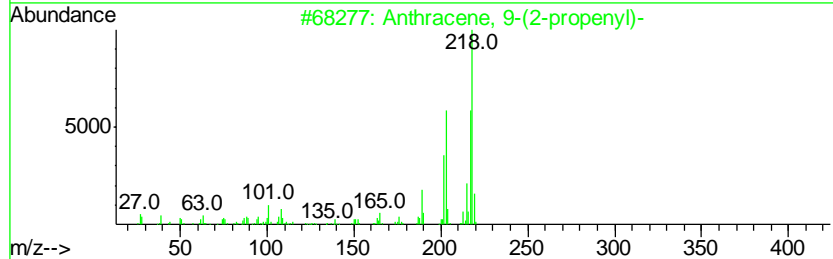
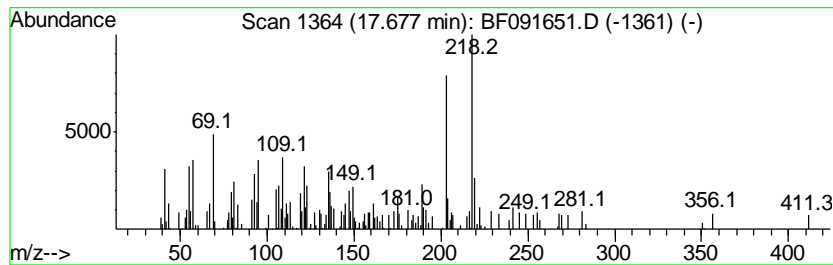
TIC Library : C:\DATABASE\NIST02.L  
 TIC Integration Parameters: LSCINT.P

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 Peak Number 13 Anthracene, 9-(2-propenyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.68	4.38 ng	208209	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 9-(2-propenyl)-	218	C17H14	023707-65-5	62
2		Benzenamine, 2,6-dimethyl-N-(4,4...	218	C13H18N2O	281211-68-5	53
3		Naphthalene, 1-(phenylmethyl)-	218	C17H14	000611-45-0	38
4		Naphthalene, 2-(phenylmethyl)-	218	C17H14	000613-59-2	38
5		2H-Cyclopropa[g]benzofuran, 4,5,...	218	C15H22O	102681-49-2	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121516\  
 Data File : BF091651.D  
 Acq On : 16 Dec 2016 1:45  
 Operator : UM/SJ  
 Sample : H5940-09  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampled :  
 SS05-1FT

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

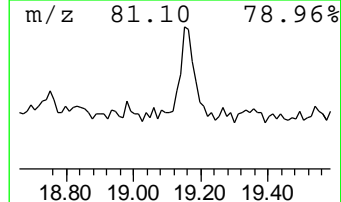
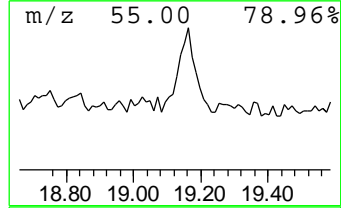
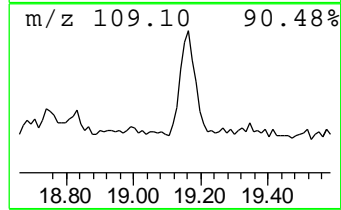
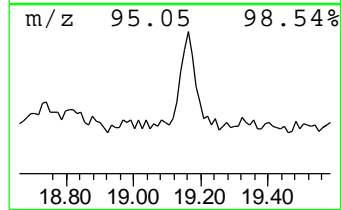
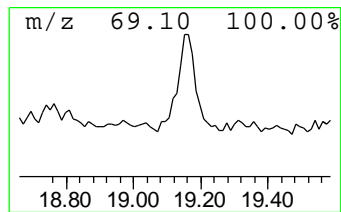
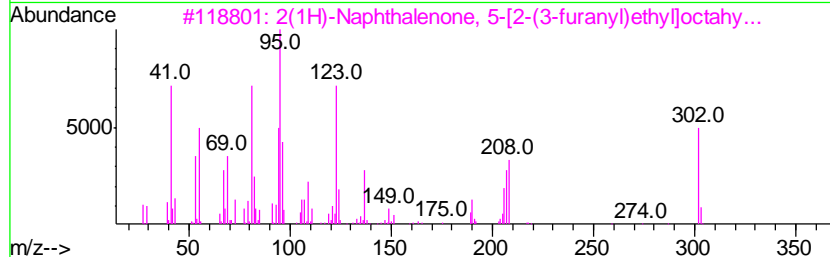
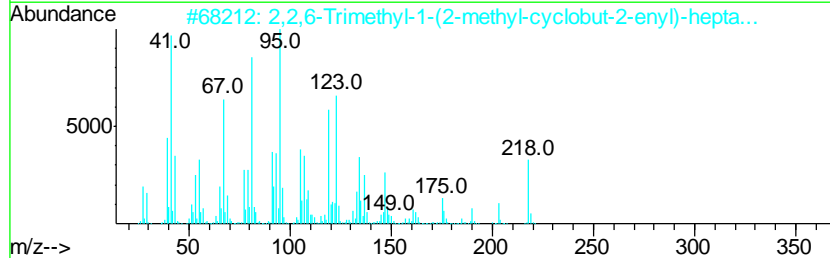
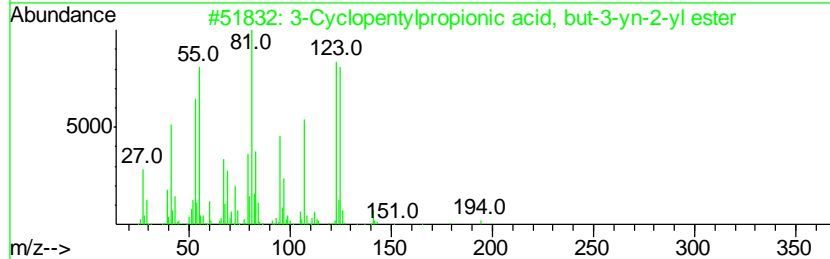
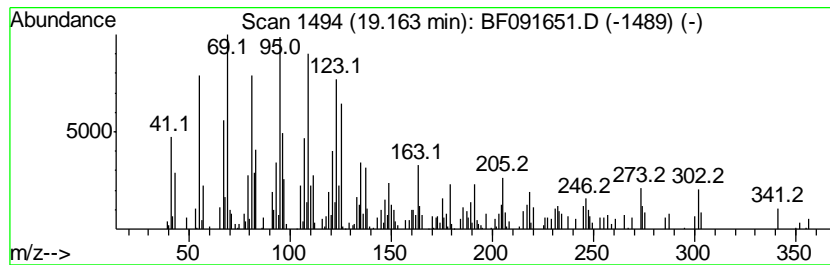
TIC Library : C:\DATABASE\NIST02.L  
 TIC Integration Parameters: LSCINT.P

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 Peak Number 14 3-Cyclopentylpropionic acid... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.16	10.45 ng	496949	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Cyclopentylpropionic acid, but...	194	C12H18O2	1000292-46-4	60
2		2,2,6-Trimethyl-1-(2-methyl-cycl...	218	C15H22O	1000188-72-8	50
3		2(1H)-Naphthalenone, 5-[2-(3-fur...	302	C20H30O2	059742-40-4	38
4		Cyclopentane, 1-methyl-1-(2-meth...	138	C10H18	074764-47-9	27
5		1H-Imidazole, 2-ethyl-	96	C5H8N2	001072-62-4	25





Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF121516\  
 Data File : BF091651.D  
 Acq On : 16 Dec 2016 1:45  
 Operator : UM/SJ  
 Sample : H5940-09  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SS05-1FT

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF120916.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...	4.96	37.6	ng	2534840	1	6.78	1346640	20.0
unknown6.56	6.56	97.8	ng	6581960	1	6.78	1346640	20.0
Eicosane	10.74	3.6	ng	223728	4	11.30	1232880	20.0
1-Hexadecanol	11.54	3.4	ng	208454	4	11.30	1232880	20.0
n-Hexadecanoic acid	11.85	2.7	ng	166844	4	11.30	1232880	20.0
Dichloroacetic ac...	13.79	5.6	ng	460428	5	13.94	1649360	20.0
Tridecane, 7-cycl...	14.85	2.2	ng	106550	6	15.35	950928	20.0
Pentadecane, 8-he...	15.04	6.5	ng	310247	6	15.35	950928	20.0
Heneicosane	15.79	9.7	ng	462621	6	15.35	950928	20.0
Dodecane, 2-methyl-	16.81	2.9	ng	137536	6	15.35	950928	20.0
Nonane, 5-butyl-	17.45	3.5	ng	167159	6	15.35	950928	20.0
Anthracene, 9-(2-...	17.68	4.4	ng	208209	6	15.35	950928	20.0
3-Cyclopentylprop...	19.16	10.4	ng	496949	6	15.35	950928	20.0