

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040617\
 Data File : BF094222.D
 Acq On : 6 Apr 2017 12:25
 Operator : SJ/MA
 Sample : I2511-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-CONCRETE-SE

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-BF032217.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	2.810	152	157	165	rVB	179675	201233	2.47%	0.424%
2	3.469	261	269	272	rBV	4901503	8163502	100.00%	17.194%
3	4.016	352	362	365	rBV	3386784	6550854	80.25%	13.797%
4	4.381	421	424	434	rVB	697802	655385	8.03%	1.380%
5	5.451	601	606	617	rBV	2272891	2463638	30.18%	5.189%
6	5.592	626	630	634	rBV	1498086	1341030	16.43%	2.824%
7	5.851	670	674	678	rBV	1108074	988027	12.10%	2.081%
8	6.028	699	704	708	rBV	2961646	3087694	37.82%	6.503%
9	6.216	732	736	739	rBV	139579	119015	1.46%	0.251%
10	6.498	778	784	788	rBV	2161485	2622646	32.13%	5.524%
11	7.351	924	929	932	rBV	1374765	1369130	16.77%	2.884%
12	8.680	1147	1155	1158	rBV	4287218	4772221	58.46%	10.051%
13	8.816	1170	1178	1182	rBV2	71935	84116	1.03%	0.177%
14	9.174	1234	1239	1242	rBV	328856	325743	3.99%	0.686%
15	9.210	1242	1245	1253	rVB4	52426	102471	1.26%	0.216%
16	9.492	1285	1293	1297	rBV2	1541364	1713127	20.99%	3.608%
17	9.745	1328	1336	1341	rBV3	82462	145603	1.78%	0.307%
18	10.086	1390	1394	1398	rBV	246548	227106	2.78%	0.478%
19	10.163	1404	1407	1411	rVB	100445	99909	1.22%	0.210%
20	10.363	1436	1441	1444	rBV2	108402	129385	1.58%	0.273%
21	10.457	1454	1457	1460	rBV2	141497	150695	1.85%	0.317%
22	10.669	1490	1493	1496	rBV	258215	306506	3.75%	0.646%
23	11.227	1584	1588	1591	rBV	204966	203228	2.49%	0.428%
24	11.263	1591	1594	1598	rVV	80694	85738	1.05%	0.181%
25	11.316	1598	1603	1605	rVV	1460629	1501068	18.39%	3.162%
26	11.339	1605	1607	1611	rVV	496253	486709	5.96%	1.025%
27	11.404	1615	1618	1622	rVB	150246	141334	1.73%	0.298%
28	11.757	1675	1678	1682	rVB	126207	132312	1.62%	0.279%
29	11.810	1682	1687	1695	rVB2	73511	140228	1.72%	0.295%
30	11.939	1705	1709	1711	rBV	109751	112800	1.38%	0.238%
31	11.968	1711	1714	1719	rVB	134475	157538	1.93%	0.332%
32	12.068	1726	1731	1734	rBV2	119862	153362	1.88%	0.323%
33	12.257	1759	1763	1767	rVB2	115437	134827	1.65%	0.284%
34	12.504	1800	1805	1808	rBV2	99090	121790	1.49%	0.257%

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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-BF032217.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	12.804	1852	1856	1860	rVV	427538	407865	5.00%	0.859%
36	13.080	1898	1903	1907	rVB	358903	408854	5.01%	0.861%
37	13.298	1932	1940	1943	rBV	3596261	4295328	52.62%	9.047%
38	13.451	1961	1966	1970	rBV	329812	383985	4.70%	0.809%
39	14.462	2133	2138	2146	rBV2	133756	282901	3.47%	0.596%
40	14.562	2149	2155	2158	rVV	1248515	1407032	17.24%	2.963%
41	14.592	2158	2160	2163	rVB	111134	98249	1.20%	0.207%
42	14.856	2200	2205	2213	rVB4	49772	101277	1.24%	0.213%
43	15.674	2341	2344	2348	rVB	95644	94275	1.15%	0.199%
44	15.780	2359	2362	2366	rBV	75316	115938	1.42%	0.244%
45	16.186	2426	2431	2440	rBV	740422	893789	10.95%	1.882%

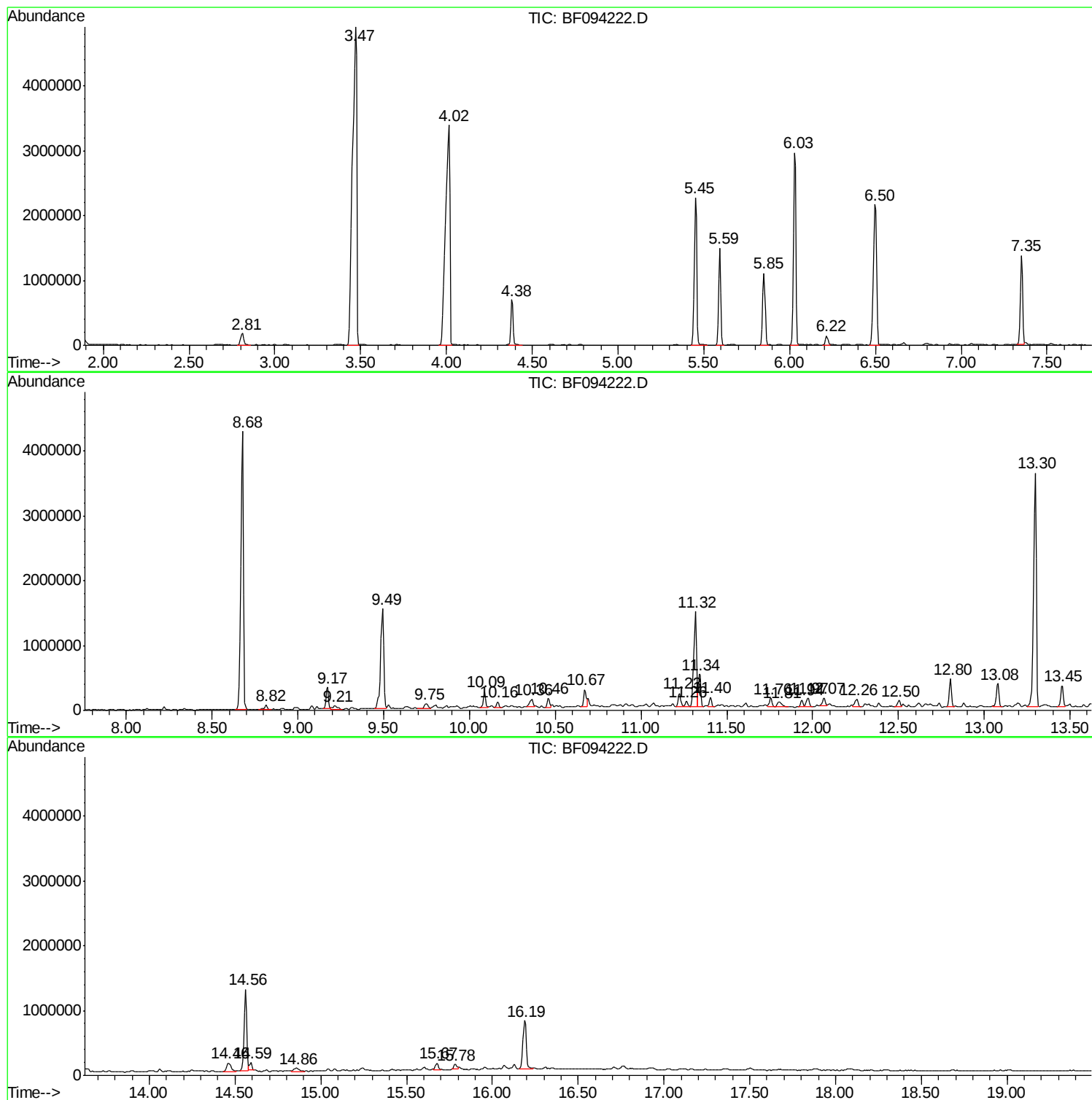
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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ALS Vial : 6 Sample Multiplier: 1

Instrument :
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E2-CONCRETE-SE

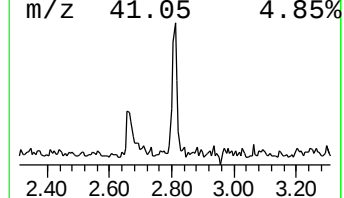
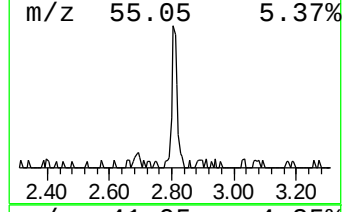
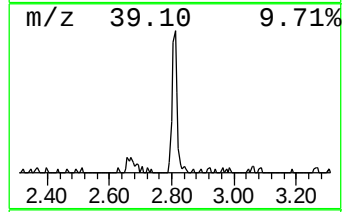
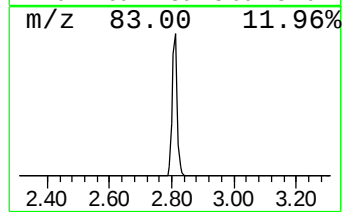
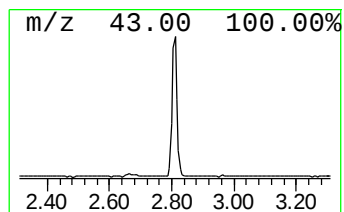
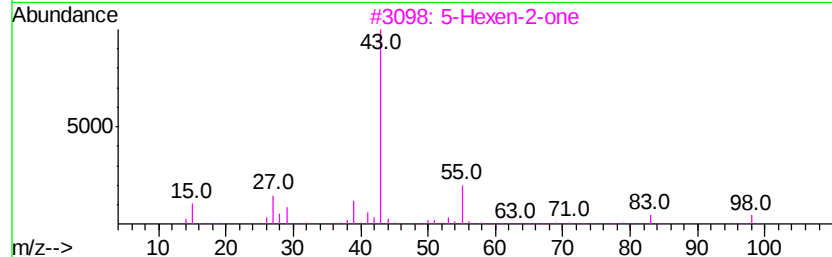
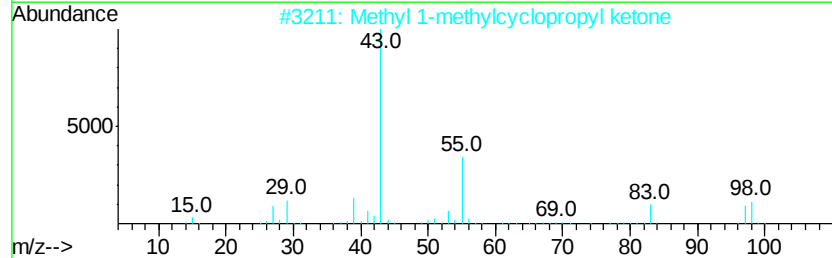
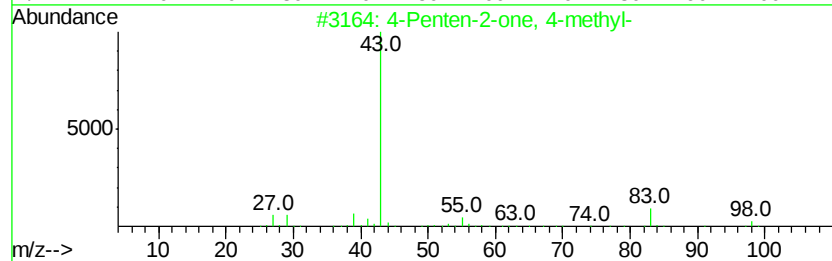
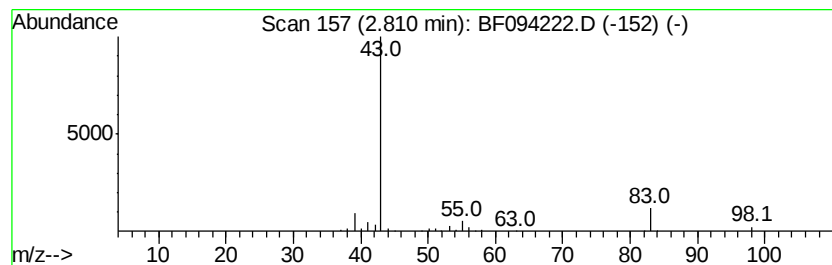
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 4-Penten-2-one, 4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.81	4.07 ng	201233	1,4-Dichlorobenzene-d4	5.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	80
2			Methyl 1-methylcyclopropyl ketone	98	C6H10O	001567-75-5	47
3			5-Hexen-2-one	98	C6H10O	000109-49-9	35
4			2-Pentanone, 3-methylene-	98	C6H10O	004359-77-7	23
5			2-Vinylethyl acetate	114	C6H10O2	001576-84-7	9



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Instrument :
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ClientSampleId :
E2-CONCRETE-SE

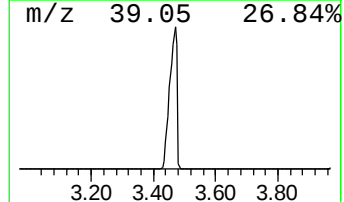
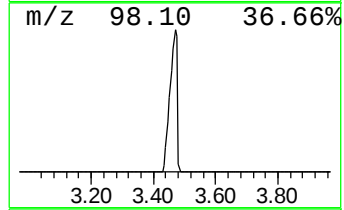
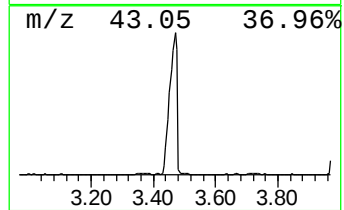
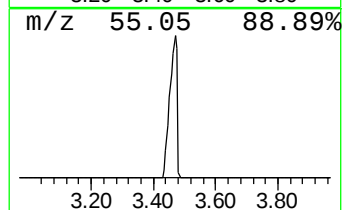
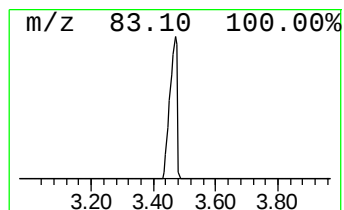
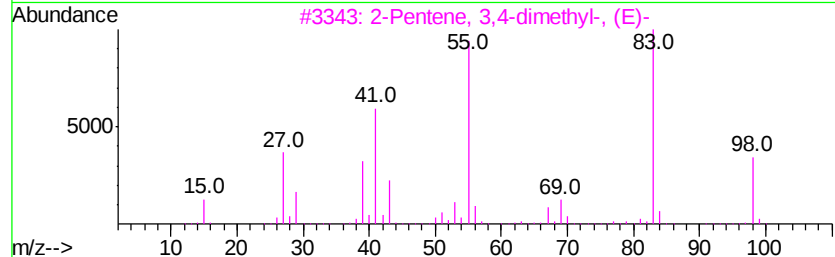
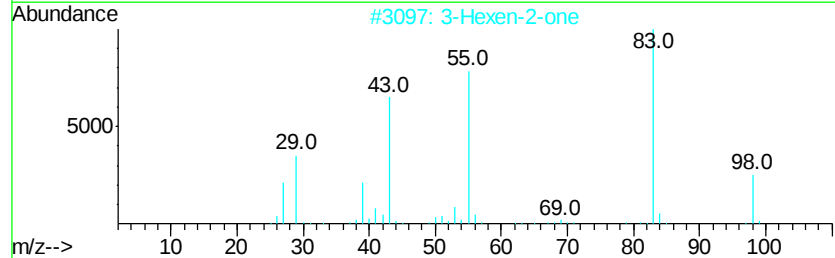
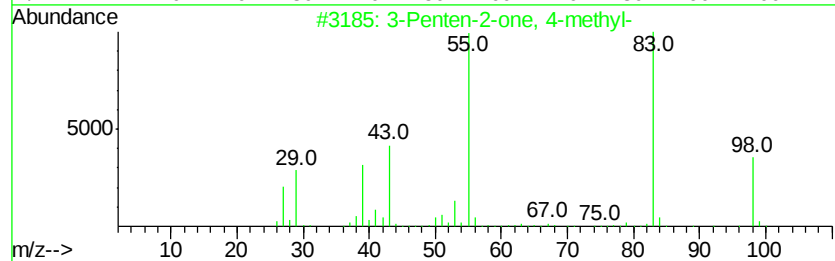
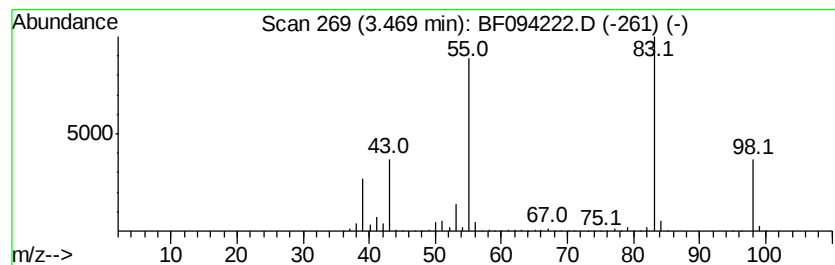
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 3-Penten-2-one, 4-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.47	165.25 ng	8163500	1,4-Dichlorobenzene-d4	5.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	95
2		3-Hexen-2-one	98	C6H10O	000763-93-9	91
3		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
4		2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	86
5		2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	80



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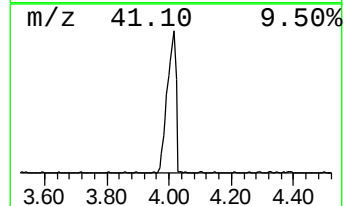
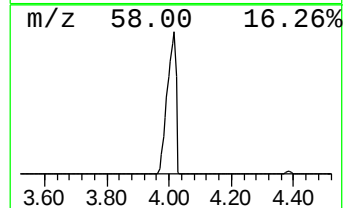
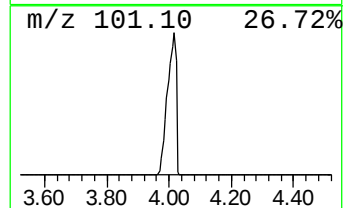
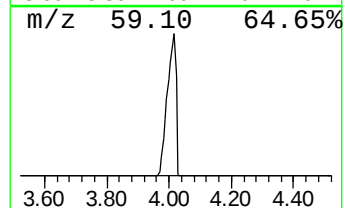
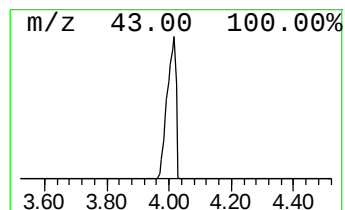
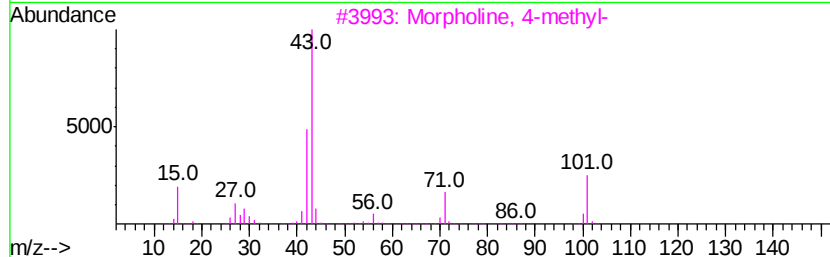
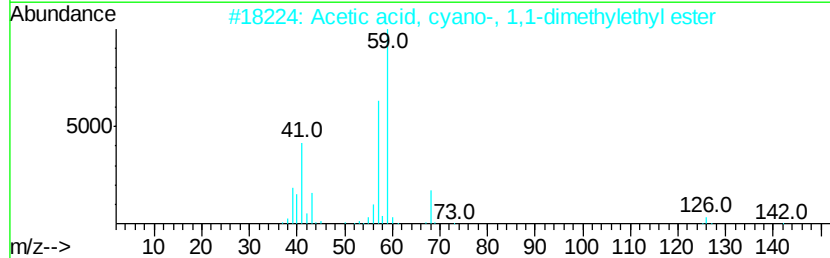
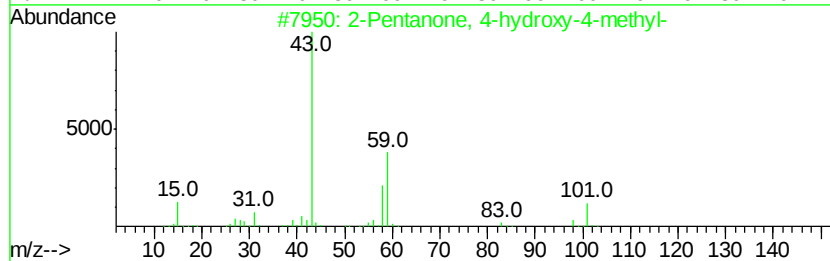
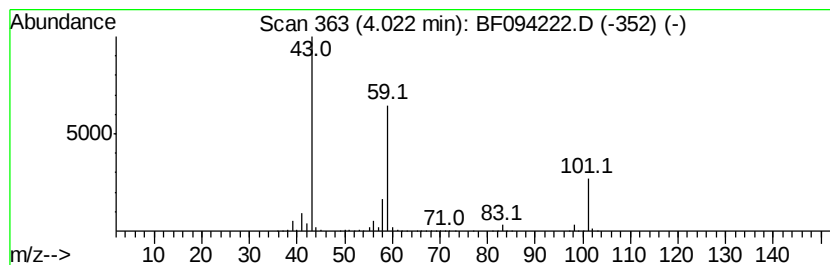
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.02	132.61 ng	6550850	1,4-Dichlorobenzene-d4	5.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



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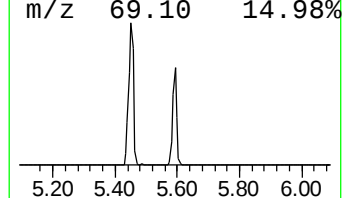
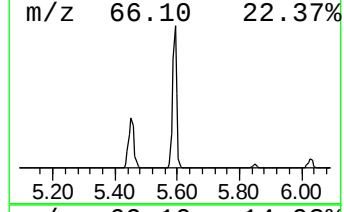
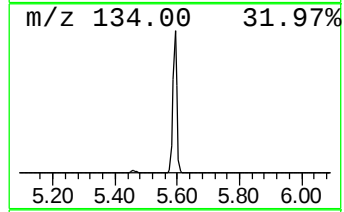
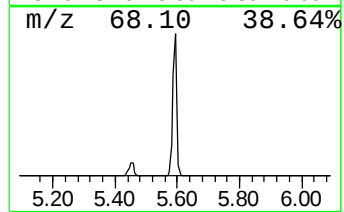
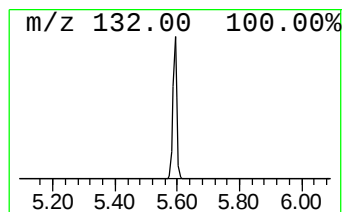
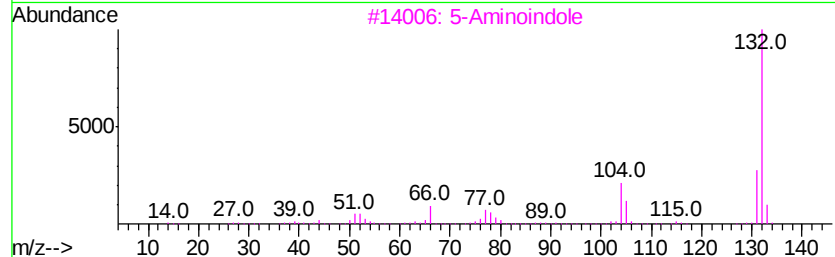
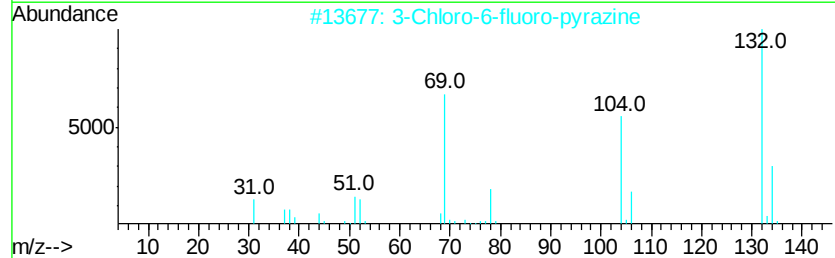
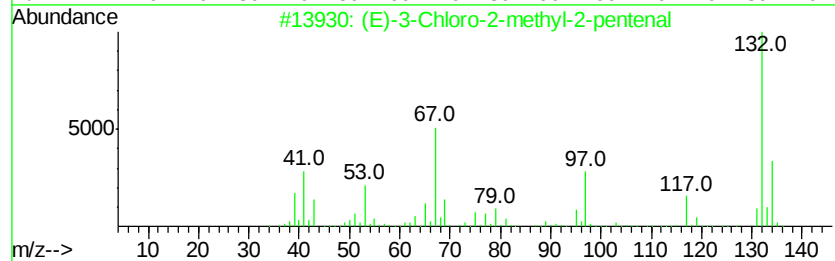
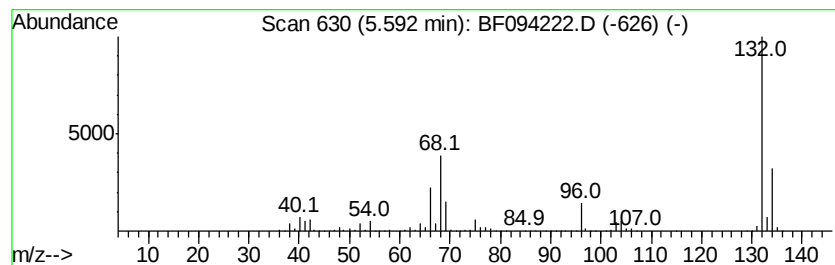
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown5.59 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.59	27.15 ng	1341030	1,4-Dichlorobenzene-d4	5.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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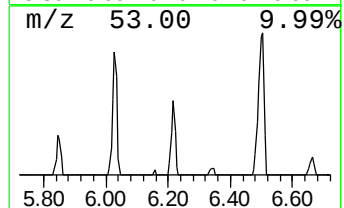
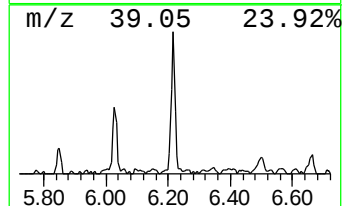
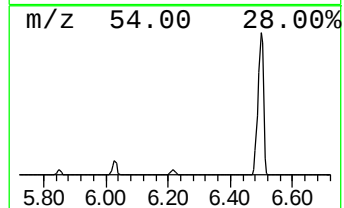
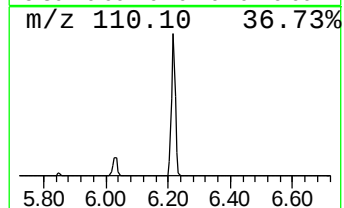
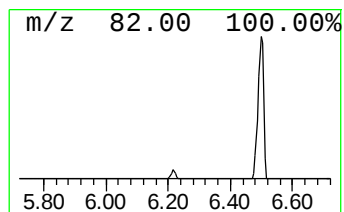
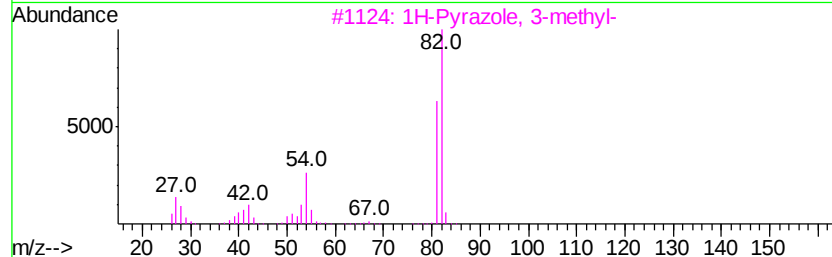
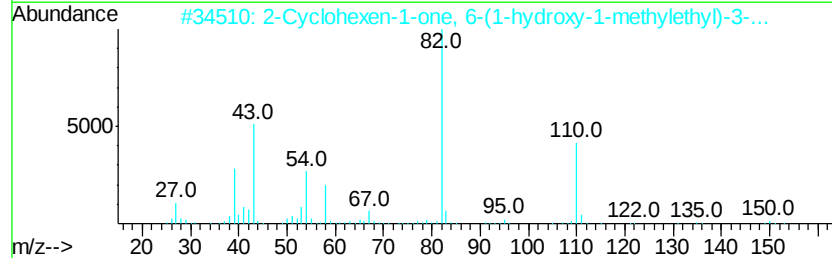
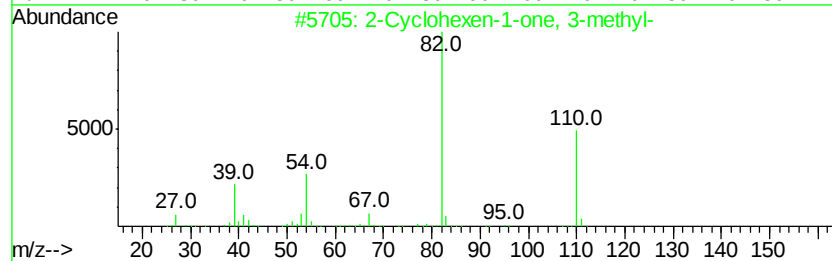
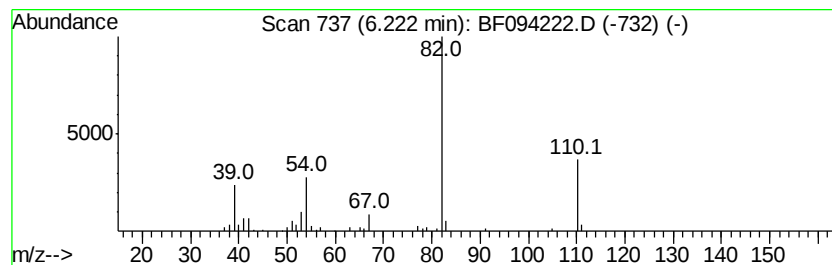
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2-Cyclohexen-1-one, 3-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.22	2.41 ng	119015	1,4-Dichlorobenzene-d4	5.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Cyclohexen-1-one, 3-methyl-	110	C7H10O	001193-18-6	91
2			2-Cyclohexen-1-one, 6-(1-hydroxy...	168	C10H16O2	087791-00-2	43
3			1H-Pyrazole, 3-methyl-	82	C4H6N2	001453-58-3	42
4			1'-Hydroxy-4,3'-dimethyl-bicyclo...	220	C14H20O2	1000189-12-5	40
5			1H-Imidazole, 4-methyl-	82	C4H6N2	000822-36-6	40



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040617\
Data File : BF094222.D
Acq On : 6 Apr 2017 12:25
Operator : SJ/MA
Sample : I2511-02
Misc :
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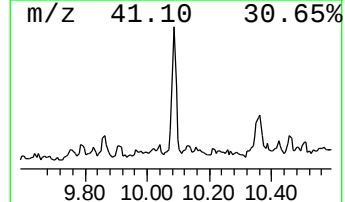
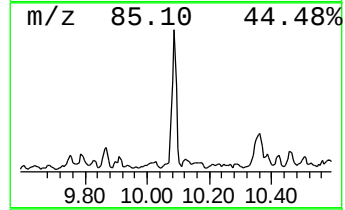
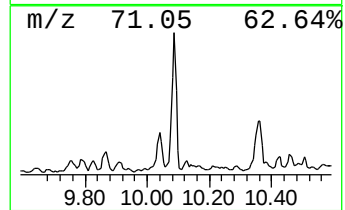
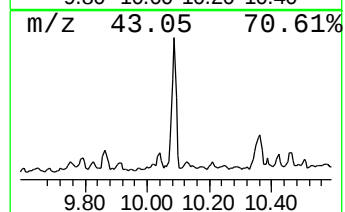
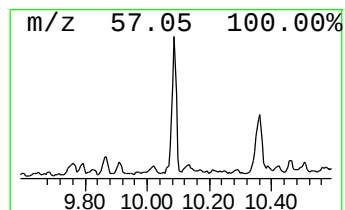
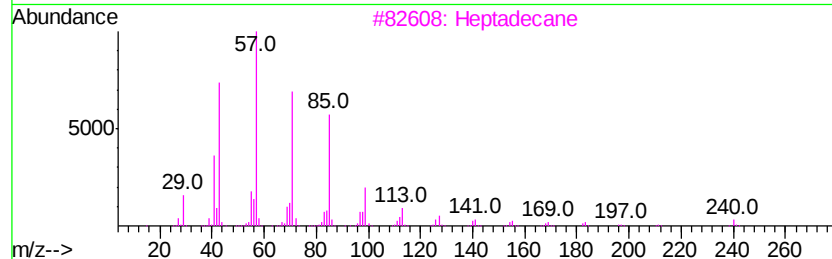
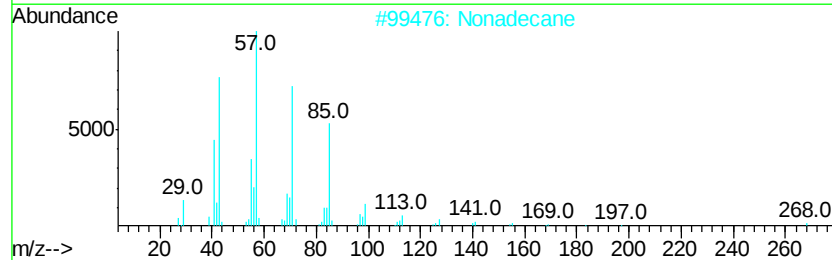
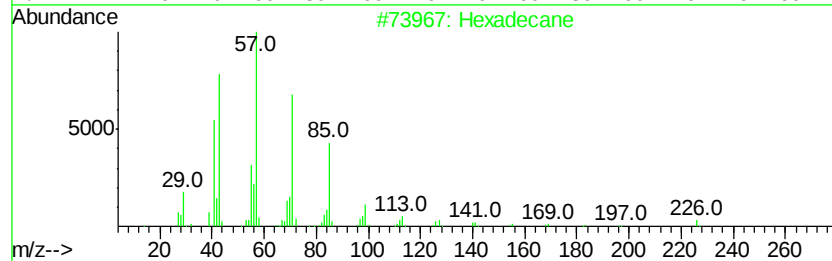
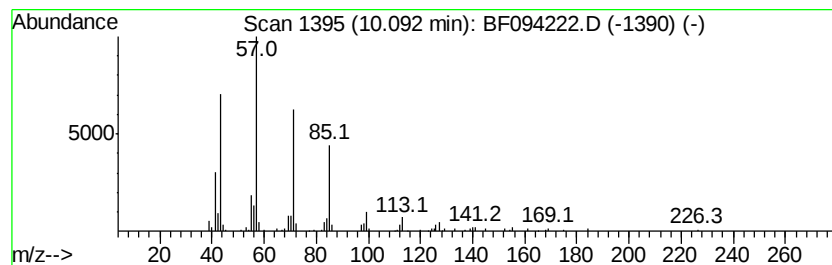
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.09	2.65 ng	227106	Acenaphthene-d10		9.49
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	94
2	Nonadecane	268	C19H40	000629-92-5	90
3	Heptadecane	240	C17H36	000629-78-7	90
4	Tetradecane	198	C14H30	000629-59-4	87
5	Eicosane	282	C20H42	000112-95-8	86



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040617\
Data File : BF094222.D
Acq On : 6 Apr 2017 12:25
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Misc :
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Instrument :
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ClientSampleId :
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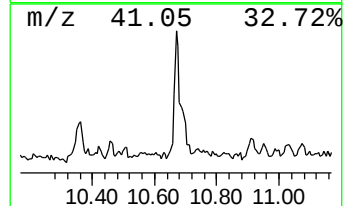
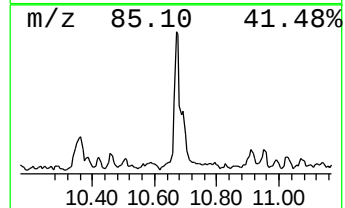
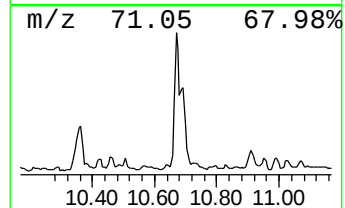
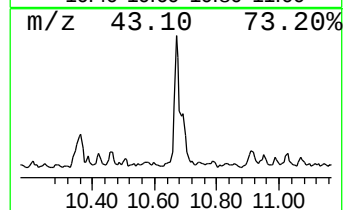
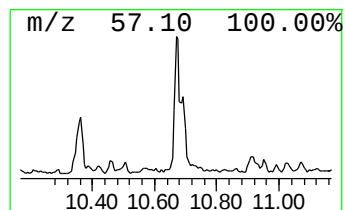
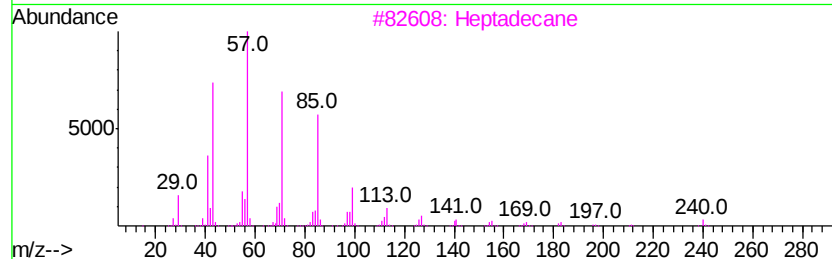
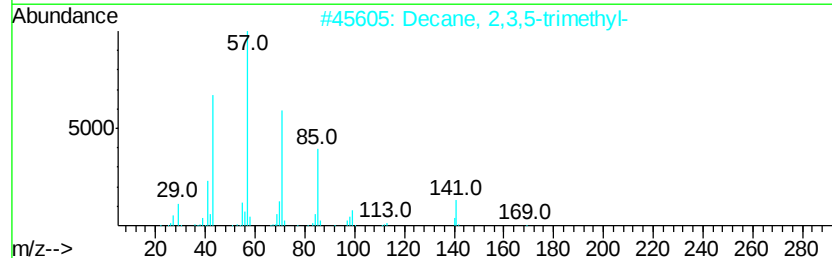
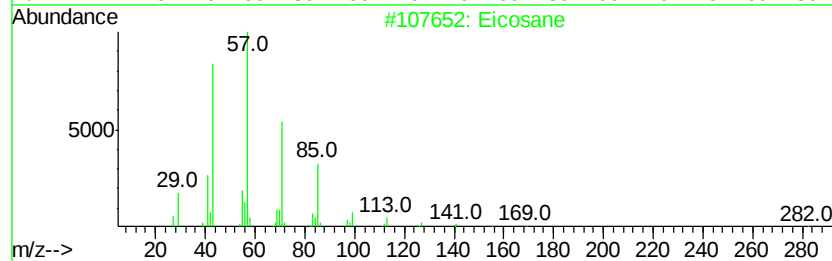
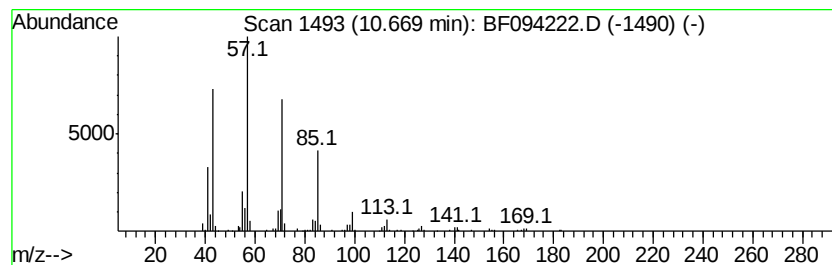
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Eicosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.67	4.08 ng	306506	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	91
2		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	90
3		Heptadecane	240	C17H36	000629-78-7	90
4		Tridecane	184	C13H28	000629-50-5	86
5		10-Methylnonadecane	282	C20H42	056862-62-5	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040617\
Data File : BF094222.D
Acq On : 6 Apr 2017 12:25
Operator : SJ/MA
Sample : I2511-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

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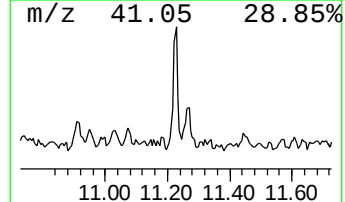
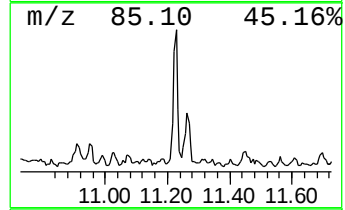
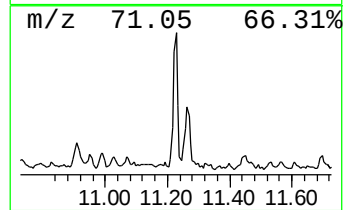
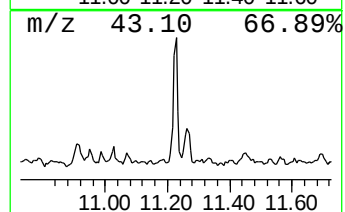
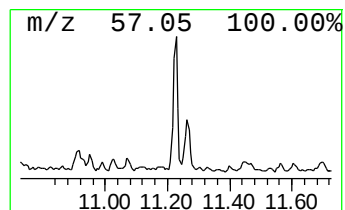
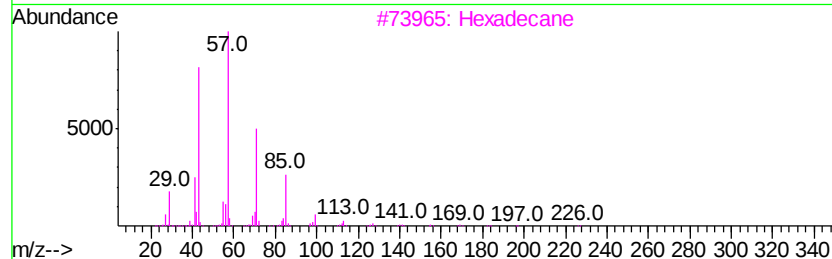
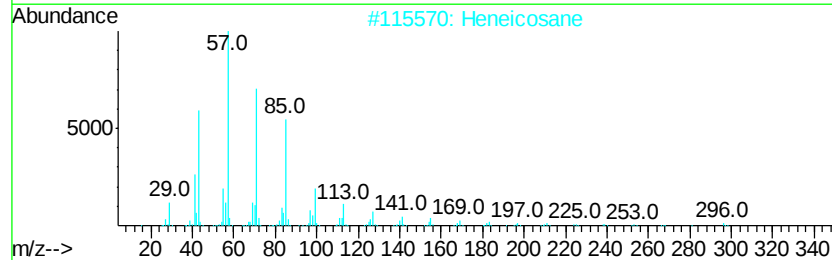
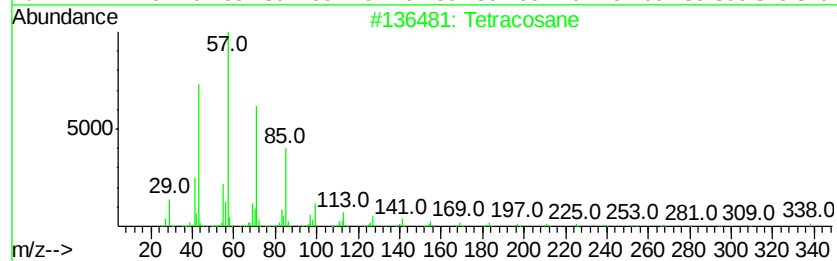
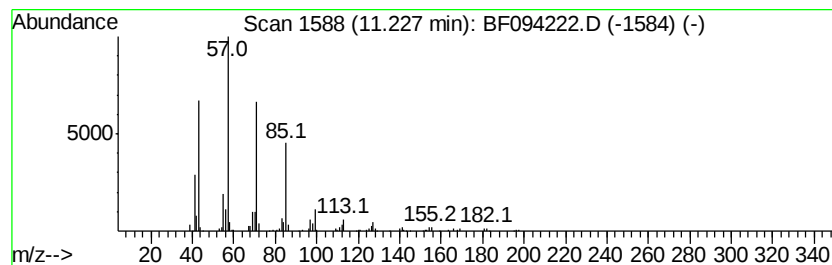
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetracosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.23	2.71 ng	203228	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	91
2		Heneicosane	296	C21H44	000629-94-7	90
3		Hexadecane	226	C16H34	000544-76-3	90
4		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	87
5		Octadecane	254	C18H38	000593-45-3	86



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040617\
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Acq On : 6 Apr 2017 12:25
Operator : SJ/MA
Sample : I2511-02
Misc :
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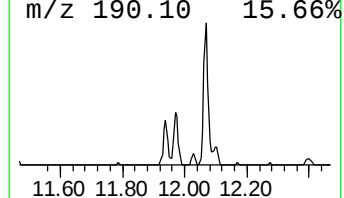
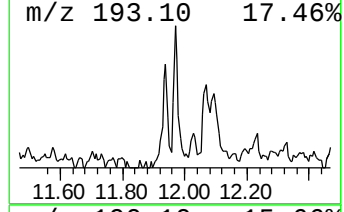
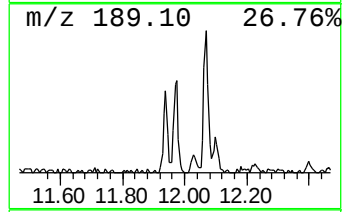
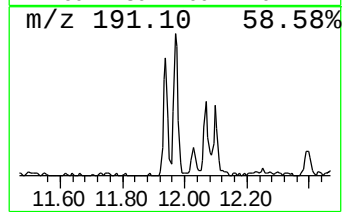
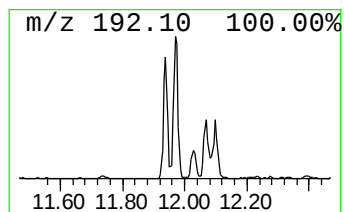
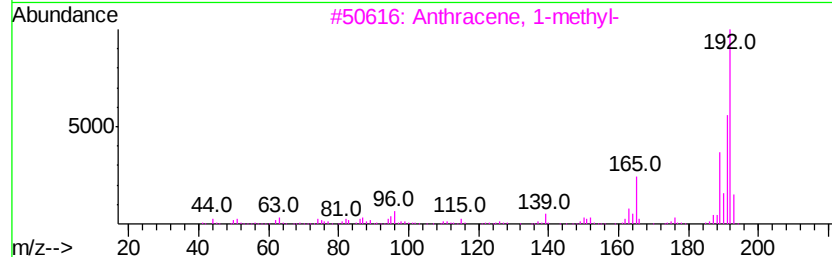
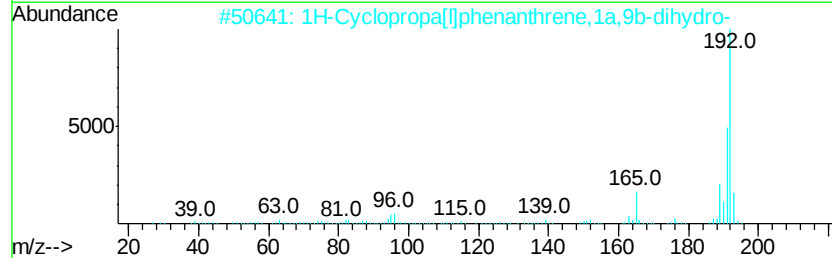
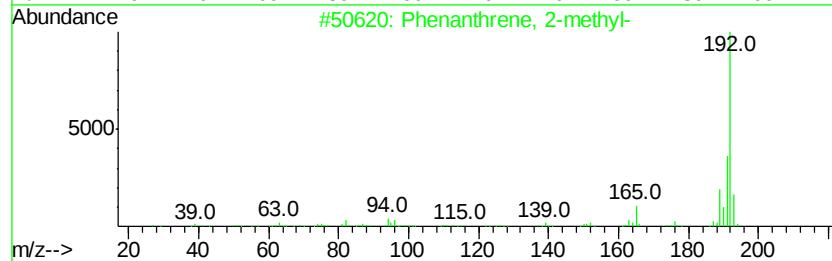
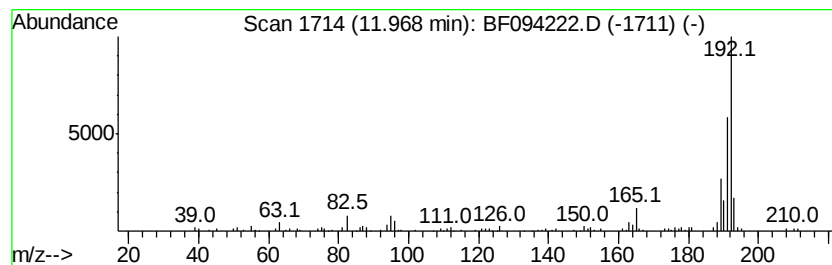
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Phenanthrene, 2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.97	2.10 ng	157538	Phenanthrene-d10	11.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
2	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	94
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	92



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040617\
Data File : BF094222.D
Acq On : 6 Apr 2017 12:25
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ALS Vial : 6 Sample Multiplier: 1

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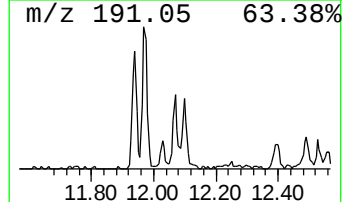
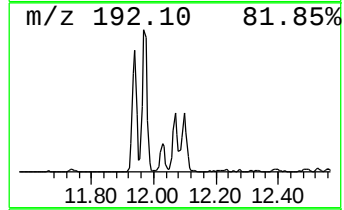
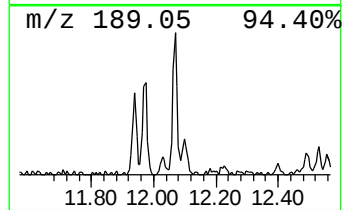
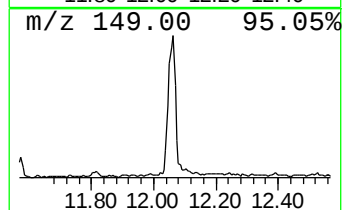
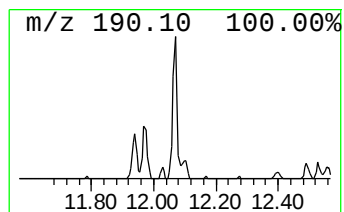
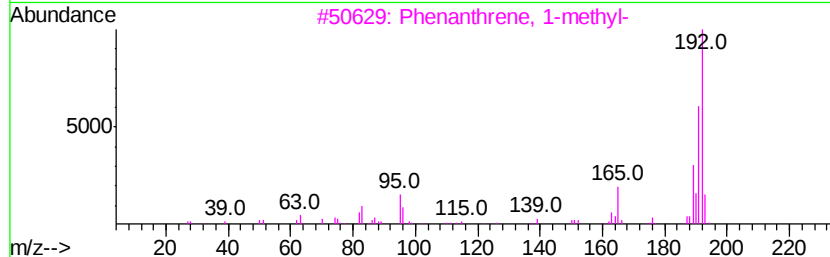
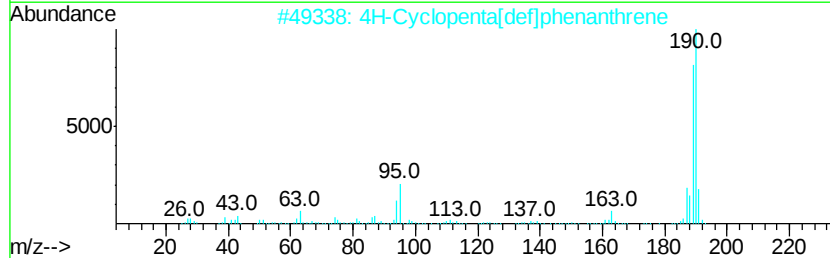
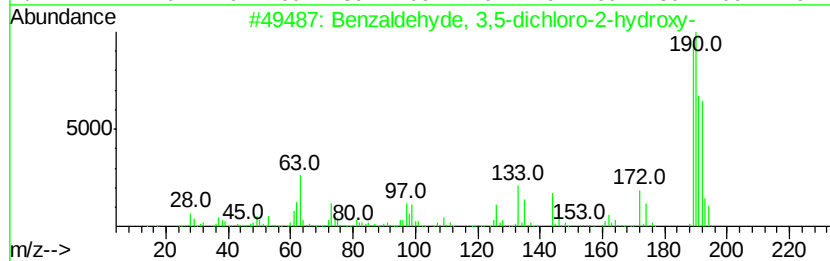
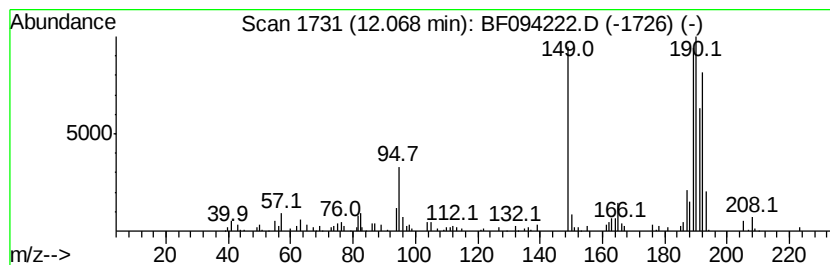
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzaldehyde, 3,5-dichloro-... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.07	2.04 ng	153362	Phenanthrene-d10	11.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehydve, 3,5-dichloro-2-hvd...	190	C7H4Cl2O2	000090-60-8	50
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	46
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	35
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	30
5			Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	20



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040617\
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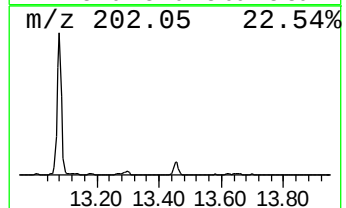
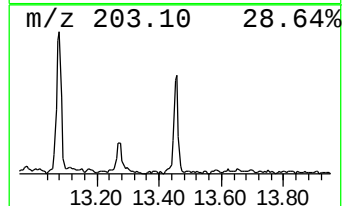
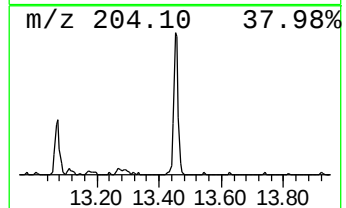
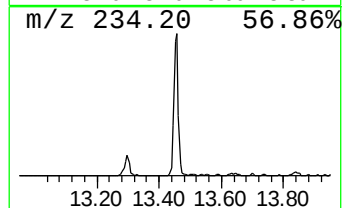
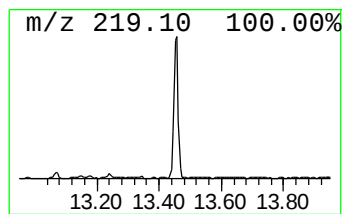
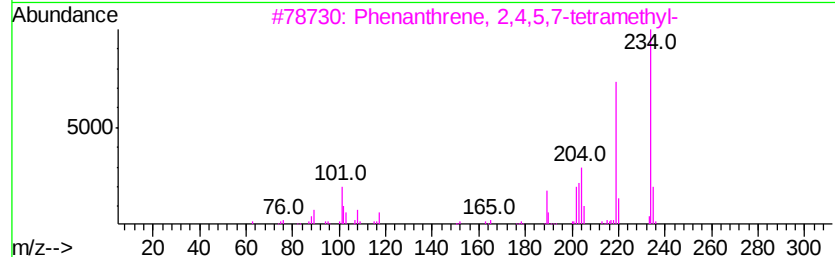
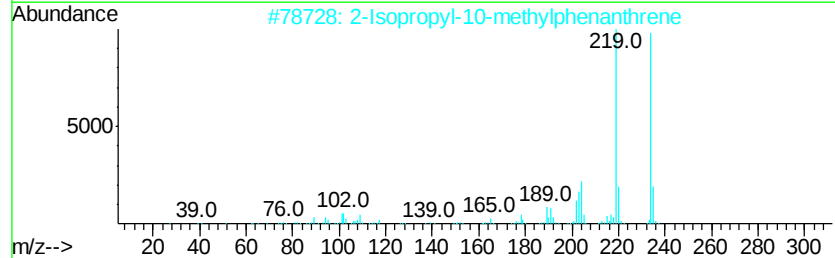
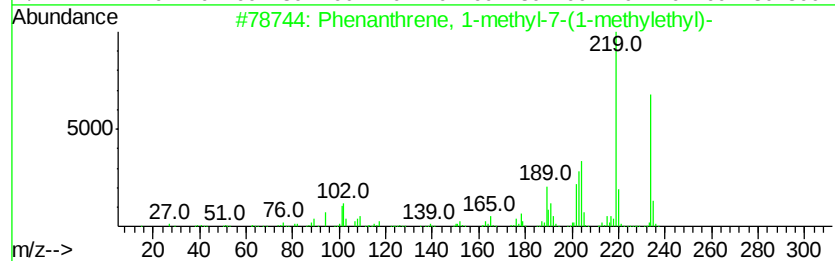
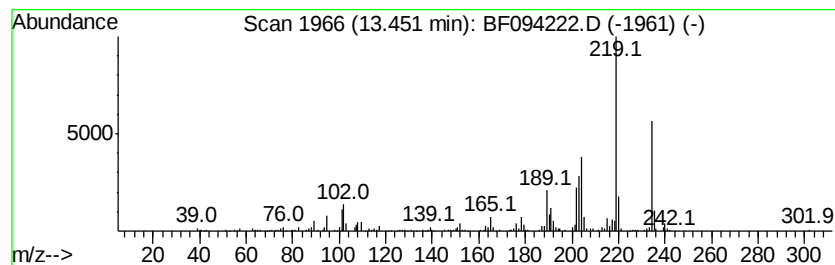
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Phenanthrene, 1-methyl-7-(1... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.45	5.46 ng	383985	Chrysene-d12	14.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 1-methyl-7-(1-meth...	234	C18H18	000483-65-8	99
2		2-Isopropyl-10-methylphenanthrene	234	C18H18	066552-97-4	94
3		Phenanthrene, 2,4,5,7-tetramethyl-	234	C18H18	007396-38-5	89
4		2,5-Diphenyl-2,4-hexadiene	234	C18H18	016819-47-9	68
5		1-(10-Methylantracen-9-yl)ethanone	234	C17H14O	036778-18-4	62



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040617\
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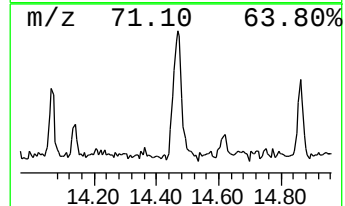
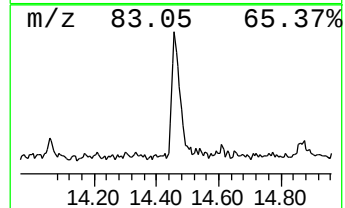
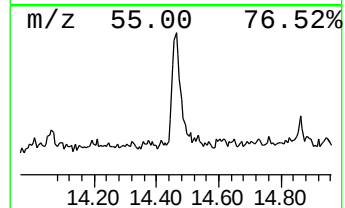
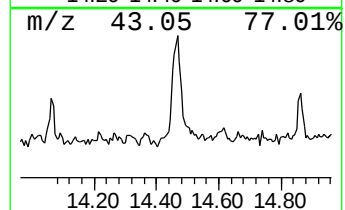
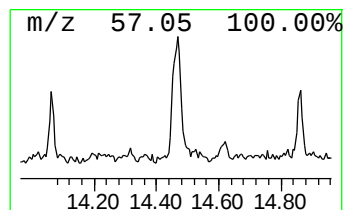
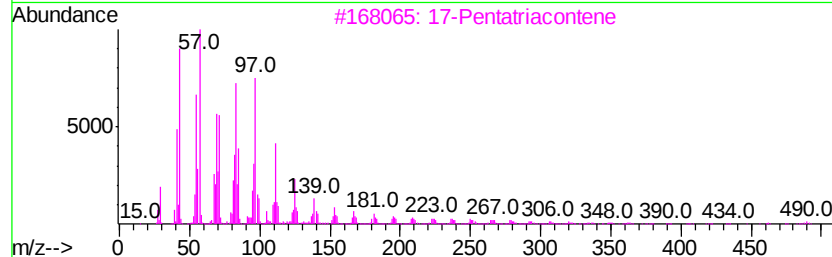
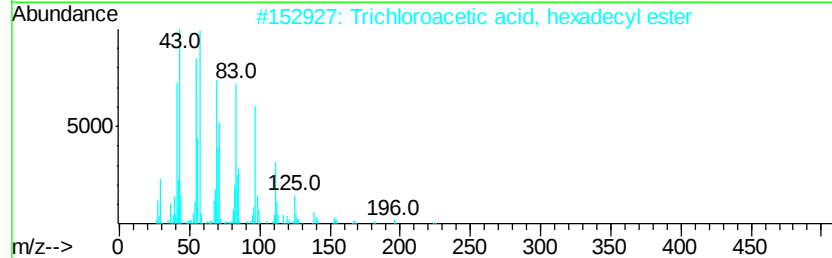
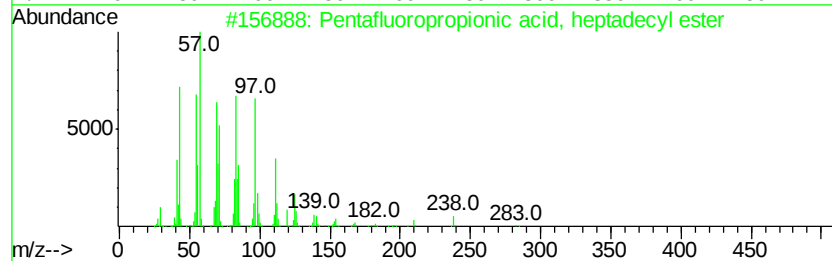
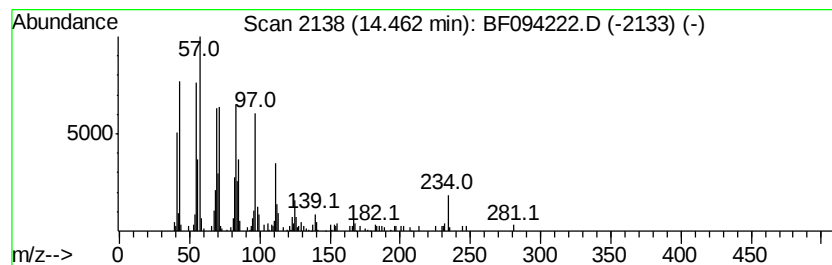
Instrument :
BNA_F
ClientSampleId :
E2-CONCRETE-SE

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Pentafluoropropionic acid, ... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
14.46	4.02 ng	282901	Chrysene-d12	14.56		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	87
2		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	87
3		17-Pentatriacontene	491	C35H70	006971-40-0	83
4		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	81
5		3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	81



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040617\
Data File : BF094222.D
Acq On : 6 Apr 2017 12:25
Operator : SJ/MA
Sample : I2511-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
E2-CONCRETE-SE

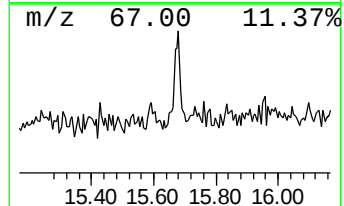
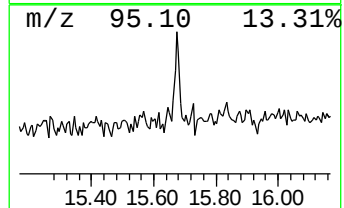
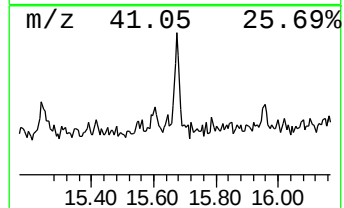
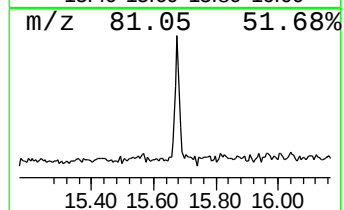
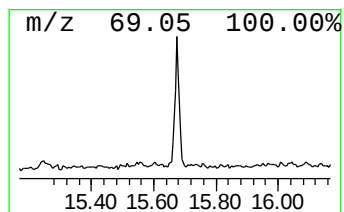
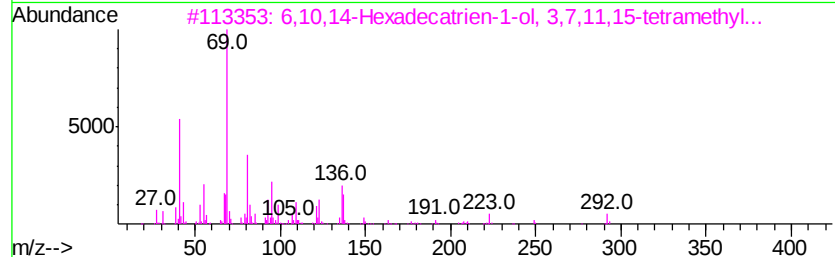
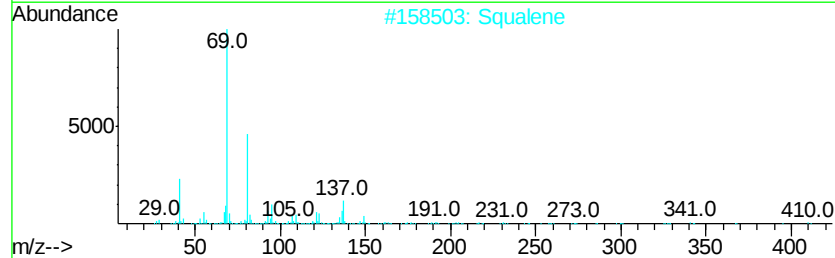
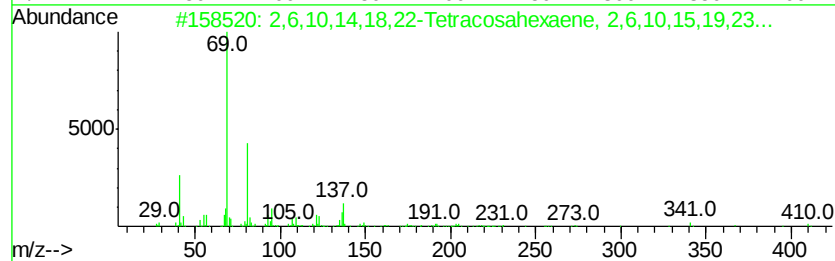
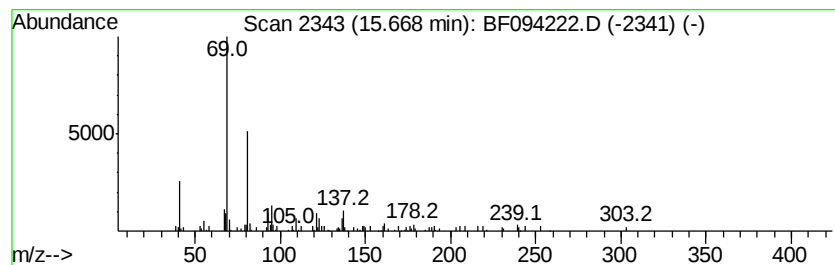
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 2,6,10,14,18,22-Tetracosahexaene... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.67	2.11 ng	94275	Perylene-d12	16.19

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,10,14,18,22-Tetracosahexaene...	410	C30H50	000111-02-4	83		
2	Squalene	410	C30H50	007683-64-9	83		
3	6,10,14-Hexadecatrien-1-ol, 3,7,...	292	C20H36O	036237-66-8	72		
4	7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24	125529-10-4	72		
5	1,5,9-Undecatriene, 2,6,10-trime...	192	C14H24	062951-96-6	64		



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040617\
 Data File : BF094222.D
 Acq On : 6 Apr 2017 12:25
 Operator : SJ/MA
 Sample : I2511-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-CONCRETE-SE

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.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
4-Penten-2-one, 4...	2.81	4.1 ng		201233	1	5.85	988027	20.0
3-Penten-2-one, 4...	3.47	165.3 ng		8163500	1	5.85	988027	20.0
2-Pentanone, 4-hy...	4.02	132.6 ng		6550850	1	5.85	988027	20.0
unknown5.59	5.59	27.1 ng		1341030	1	5.85	988027	20.0
2-Cyclohexen-1-on...	6.22	2.4 ng		119015	1	5.85	988027	20.0
Hexadecane	10.09	2.6 ng		227106	3	9.49	1713130	20.0
Eicosane	10.67	4.1 ng		306506	4	11.32	1501070	20.0
Tetracosane	11.23	2.7 ng		203228	4	11.32	1501070	20.0
Phenanthrene, 2-m...	11.97	2.1 ng		157538	4	11.32	1501070	20.0
Benzaldehyde, 3,5...	12.07	2.0 ng		153362	4	11.32	1501070	20.0
Phenanthrene, 1-m...	13.45	5.5 ng		383985	5	14.56	1407030	20.0
Pentafluoropropio...	14.46	4.0 ng		282901	5	14.56	1407030	20.0
2,6,10,14,18,22-T...	15.67	2.1 ng		94275	6	16.19	893789	20.0