

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033116\
 Data File : BF085949.D
 Acq On : 1 Apr 2016 1:37
 Operator : UM/SJ
 Sample : H2010-06
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AMBOY-SB25COMP0.5-15.0

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-BF032416.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	2.344	3	5	8	rBV	106444	90253	2.86%	0.334%
2	2.413	8	11	12	rVV2	137838	254017	8.04%	0.939%
3	2.436	12	13	17	rVV4	110994	276862	8.77%	1.024%
4	2.504	17	19	21	rVV3	95828	125438	3.97%	0.464%
5	2.596	23	27	30	rVB3	41025	113205	3.58%	0.419%
6	2.744	34	40	41	rBV4	244299	626054	19.82%	2.315%
7	2.778	41	43	45	rVV2	206728	358406	11.35%	1.325%
8	2.813	45	46	48	rVB2	139703	103026	3.26%	0.381%
9	3.681	119	122	128	rBV	18629	33365	1.06%	0.123%
10	4.973	234	235	252	rBV	52005	107274	3.40%	0.397%
11	5.384	269	271	274	rBV	1837279	2520661	79.81%	9.321%
12	6.436	360	363	365	rBV	2474437	2784621	88.17%	10.297%
13	6.562	371	374	376	rBV	2533432	2874612	91.02%	10.630%
14	6.790	392	394	396	rBV	643169	626708	19.84%	2.318%
15	6.939	405	407	410	rBV	1725131	2030181	64.28%	7.507%
16	7.362	441	444	446	rBV	1512185	1846226	58.46%	6.827%
17	8.070	504	506	509	rBV	822915	800258	25.34%	2.959%
18	9.156	598	601	603	rBV	2928000	3158160	100.00%	11.679%
19	9.545	633	635	638	rBV	397737	426823	13.51%	1.578%
20	9.762	652	654	656	rBV	120161	96318	3.05%	0.356%
21	9.831	658	660	662	rVB	856554	850907	26.94%	3.147%
22	9.991	671	674	678	rBV	21205	51132	1.62%	0.189%
23	10.265	695	698	699	rVB	100267	115704	3.66%	0.428%
24	10.482	714	717	720	rBV	45999	78191	2.48%	0.289%
25	10.619	726	729	731	rBV	1846146	1752253	55.48%	6.480%
26	10.733	737	739	743	rBV	143655	161972	5.13%	0.599%
27	10.928	754	756	760	rVB	15706	32914	1.04%	0.122%
28	11.179	776	778	779	rBV	66256	54623	1.73%	0.202%
29	11.305	787	789	792	rBV	585033	768191	24.32%	2.841%
30	12.722	910	913	915	rBV	73351	87698	2.78%	0.324%
31	12.894	925	928	931	rBV	2293579	2110688	66.83%	7.805%
32	13.374	968	970	972	rBV	27384	43487	1.38%	0.161%
33	13.419	972	974	976	rVB	82417	83380	2.64%	0.308%
34	13.797	1004	1007	1014	rBV	42784	79997	2.53%	0.296%

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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

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Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.945	1018	1020	1023	rVB	591551	534277	16.92%	1.976%
36	14.780	1090	1093	1094	rBV	37842	44385	1.41%	0.164%
37	15.363	1141	1144	1148	rBV	426396	611194	19.35%	2.260%
38	15.523	1155	1158	1164	rVB5	18780	45028	1.43%	0.167%
39	15.637	1166	1168	1169	rBV2	38149	39548	1.25%	0.146%
40	15.774	1178	1180	1184	rBV5	24247	69754	2.21%	0.258%
41	15.877	1187	1189	1194	rVB4	24497	41997	1.33%	0.155%
42	16.734	1261	1264	1266	rBV3	22325	38411	1.22%	0.142%
43	18.117	1381	1385	1387	rBV4	59704	62102	1.97%	0.230%
44	18.174	1387	1390	1394	rVV5	12335	31984	1.01%	0.118%

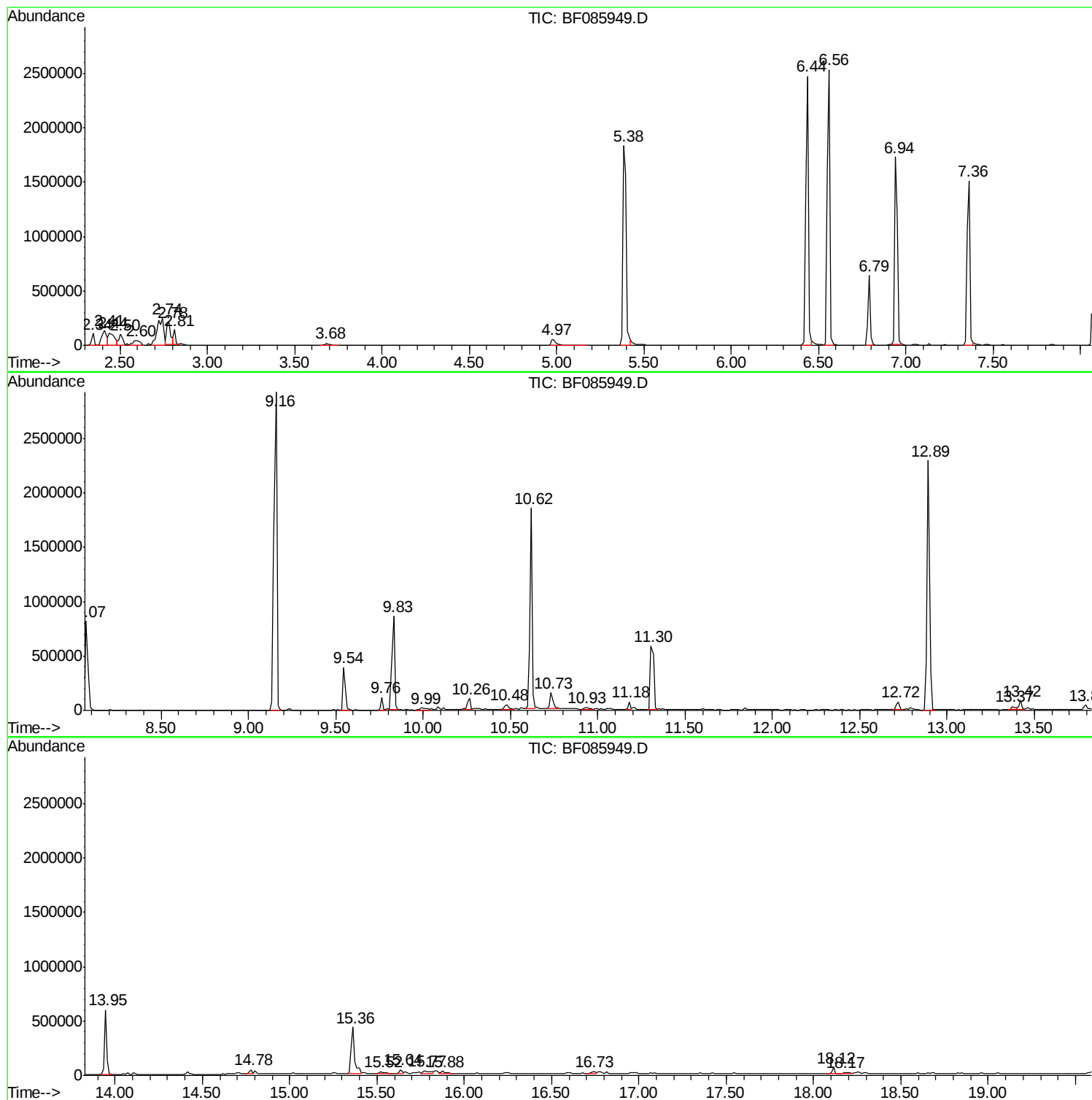
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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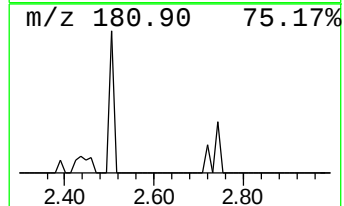
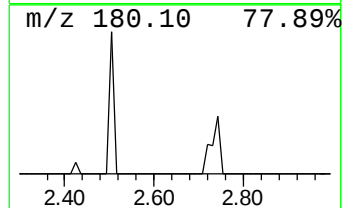
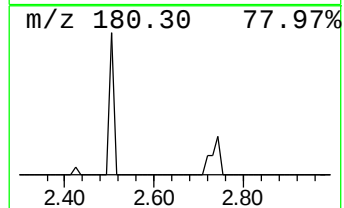
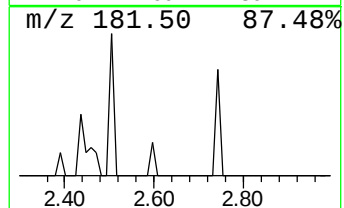
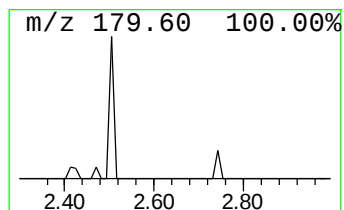
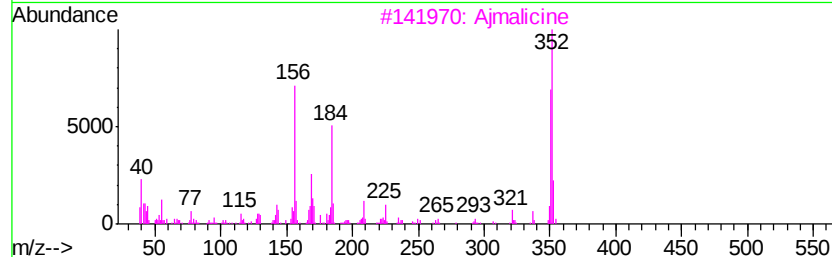
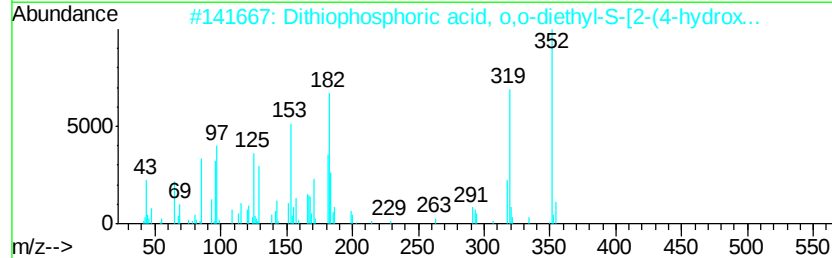
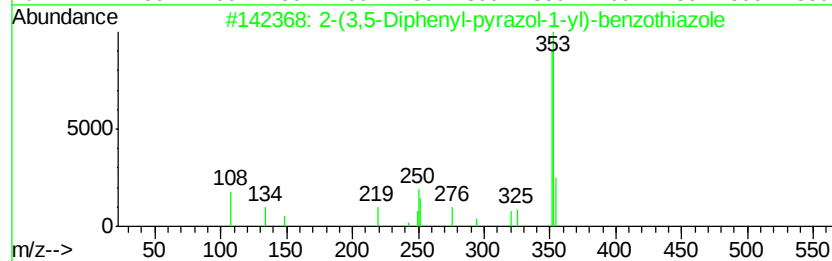
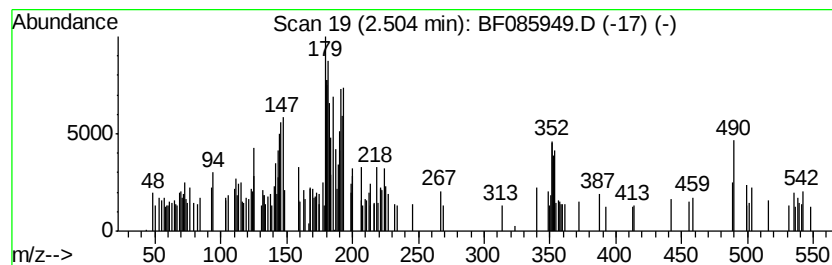
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 unknown2.50 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.50	4.00 ng	125438	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-(3,5-Diphenyl-pyrazol-1-yl)-be...	353	C22H15N3S	111199-59-8	38
2		Dithiophosphoric acid, o,o-dieth...	352	C12H17O6PS2	294874-89-8	25
3		Ajmalicine	352	C21H24N2O3	000483-04-5	18
4		Oxayohimban-16-carboxylic acid, ...	352	C21H24N2O3	006474-90-4	14
5		Yohimban-16-carboxylic acid, 17-...	352	C21H24N2O3	002671-57-0	14



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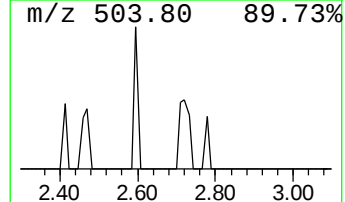
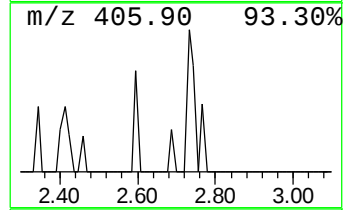
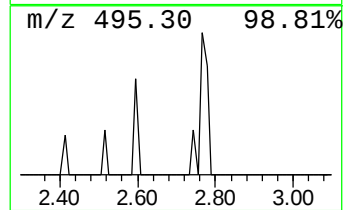
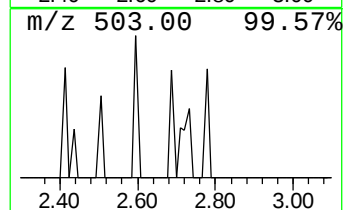
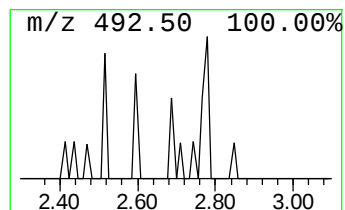
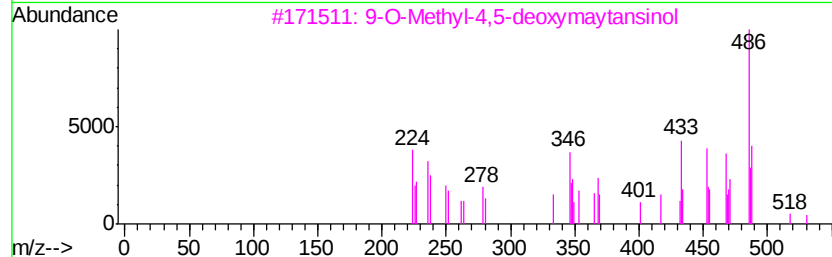
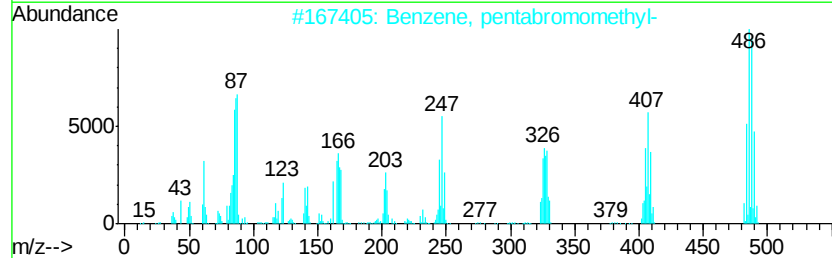
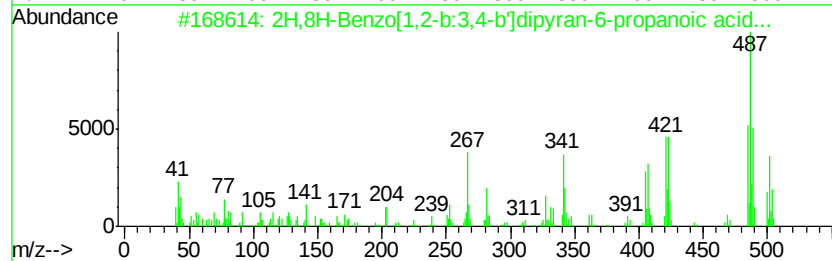
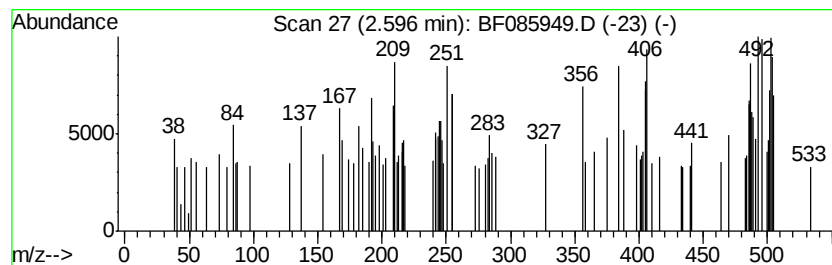
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TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown2.60 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.60	3.61 ng	113205	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2H,8H-Benzo[1,2-b:3,4-b']dipyran...	500	C20H22Br2O5	056336-18-6	23
2		Benzene, pentabromomethyl-	482	C7H3Br5	000087-83-2	9
3		9-O-Methyl-4,5-deoxymaytansinol	562	C29H39ClN2O7	075349-70-1	7
4		Dipyridamole	504	C24H40N8O4	000058-32-2	2
5		7-Chloro-3-[2,4-dichlorophenyl]-...	501	C26H26Cl3N3O	144155-58-8	1



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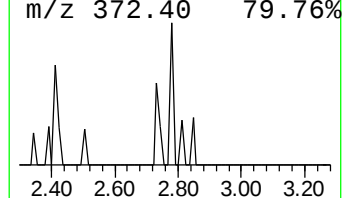
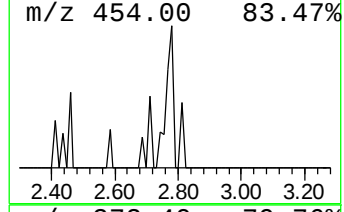
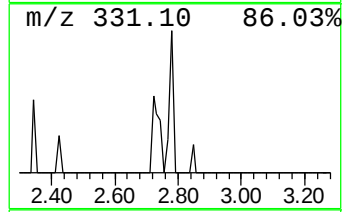
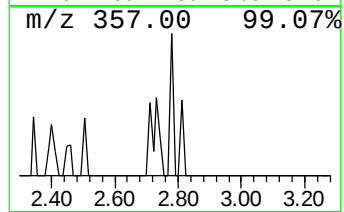
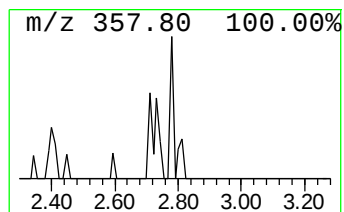
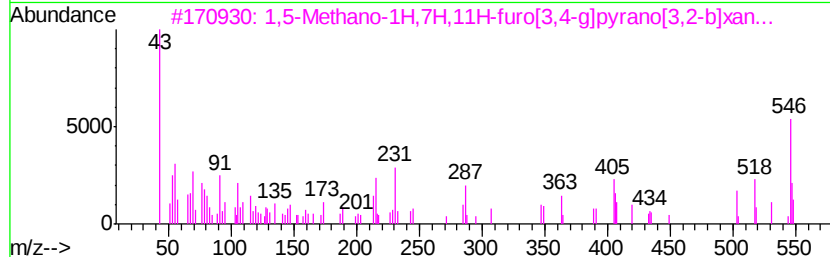
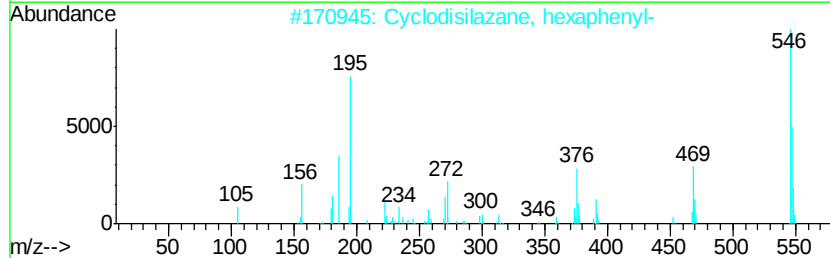
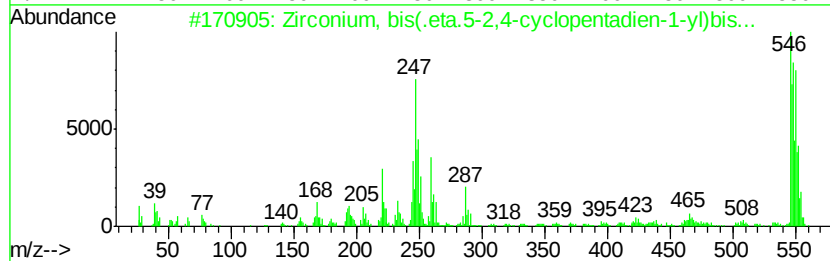
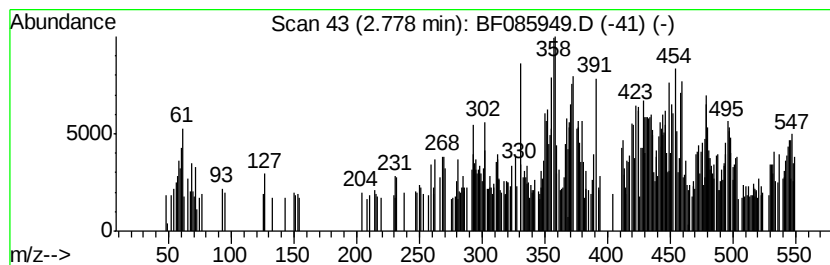
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown2.78 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.78	11.44 ng	358406	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Zirconium, bis(.eta.5-2,4-cyclop...	546	C28H30Zr2	119144-91-1	22
2		Cyclodisilazane, hexaphenyl-	546	C36H30N2Si2	003284-07-9	10
3		1,5-Methano-1H,7H,11H-furo[3,4-g...	546	C33H38O7	055452-65-8	10
4		Iron, (.eta.5-2,4-cyclopentadien...	546	C33H32FeP2	097959-85-8	9
5		Pregna-2,4-dien-20-one, 17-(acet...	546	C30H50O5Si2	069833-57-4	9



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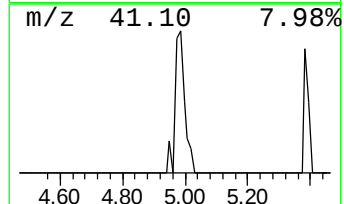
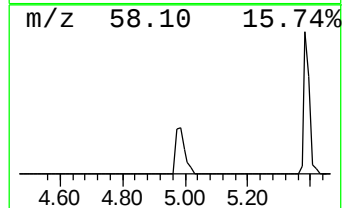
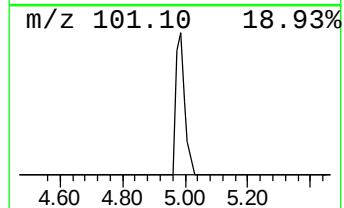
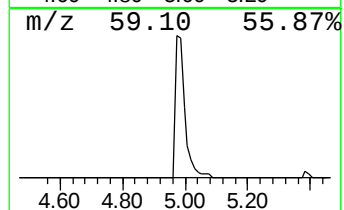
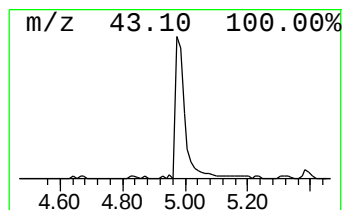
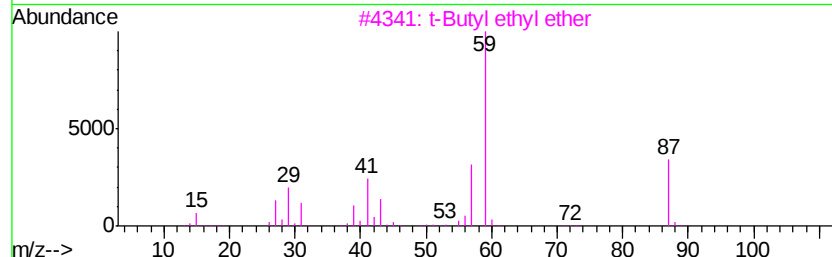
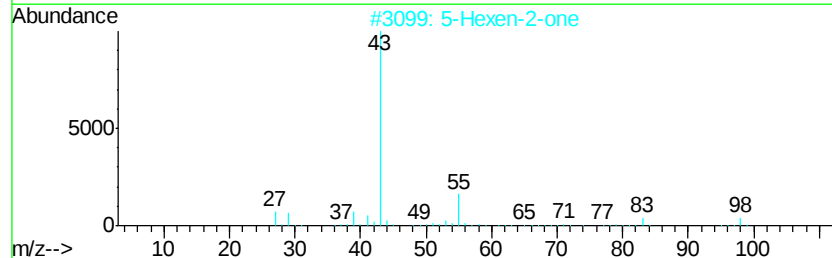
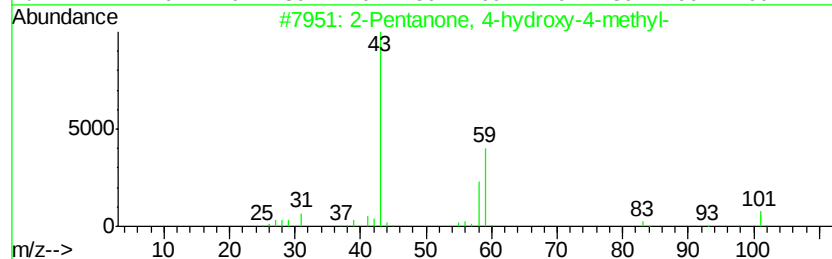
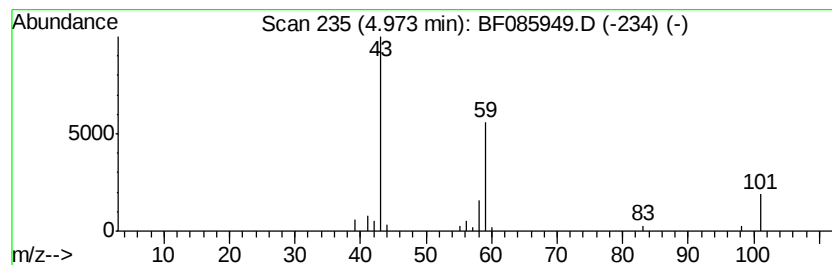
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.97	3.42 ng	107274	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		5-Hexen-2-one	98	C6H10O	000109-49-9	9
3		t-Butyl ethyl ether	102	C6H14O	1000118-18-4	9
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	9
5		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	9



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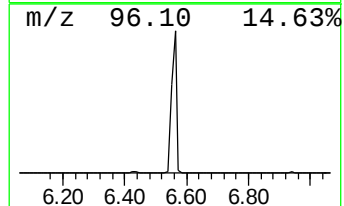
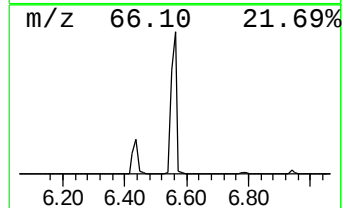
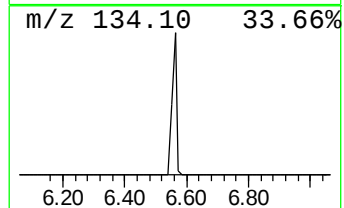
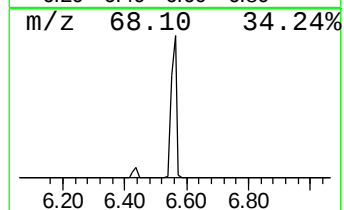
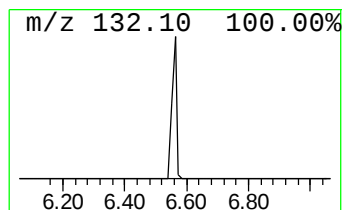
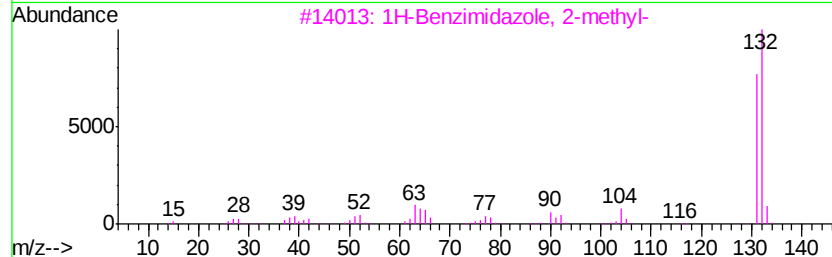
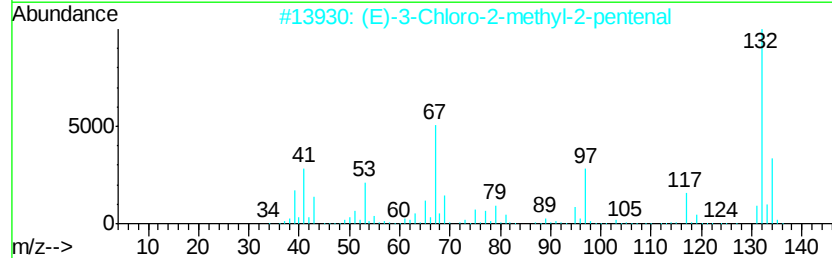
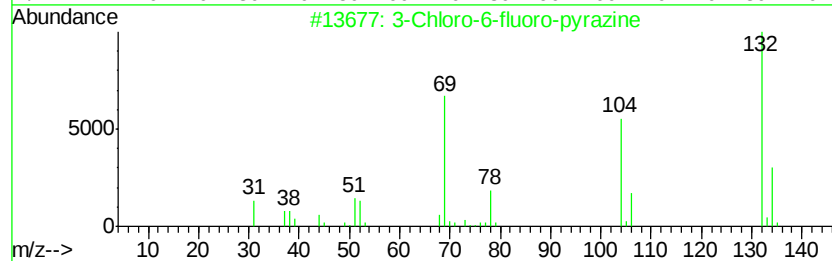
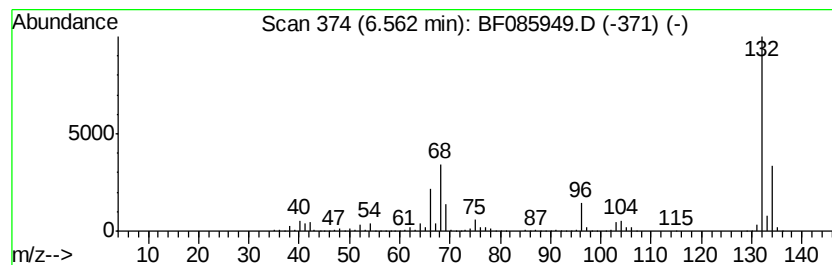
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	91.74 ng	2874610	1,4-Dichlorobenzene-d4	6.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3			1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
4			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5			4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033116\
Data File : BF085949.D
Acq On : 1 Apr 2016 1:37
Operator : UM/SJ
Sample : H2010-06
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AMBOY-SB25COMP0.5-15.0

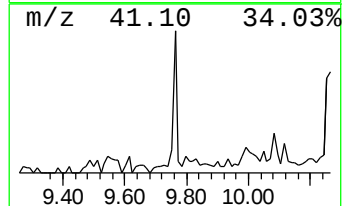
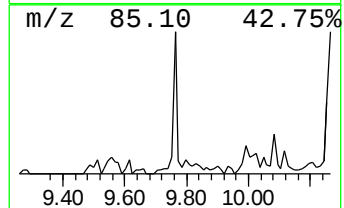
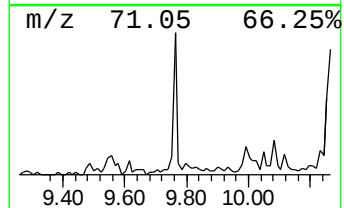
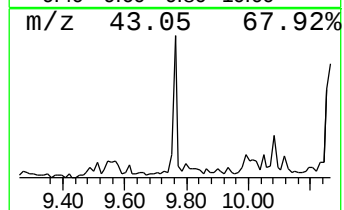
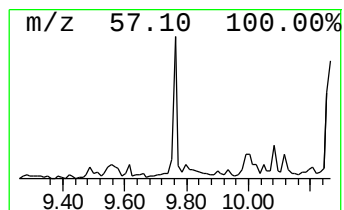
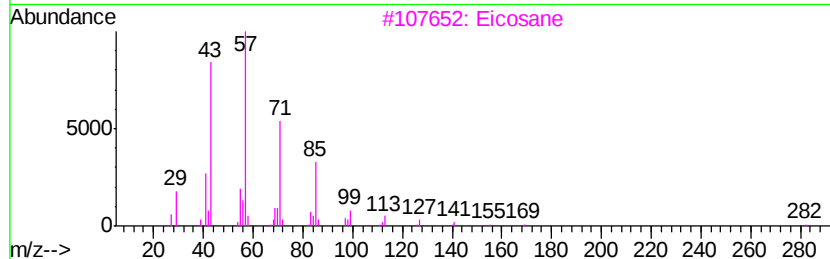
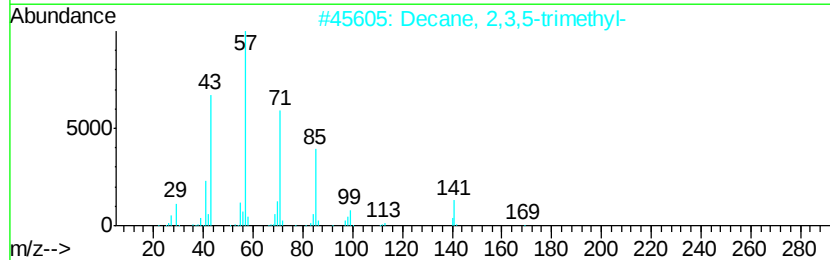
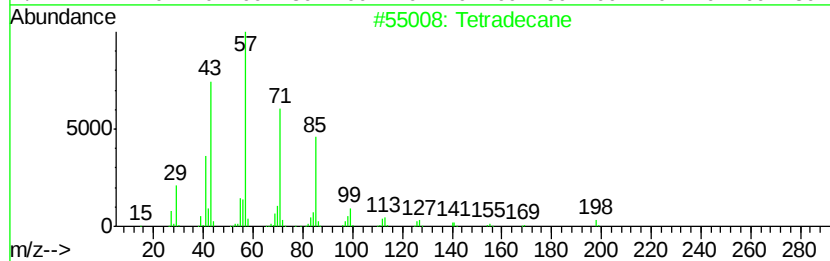
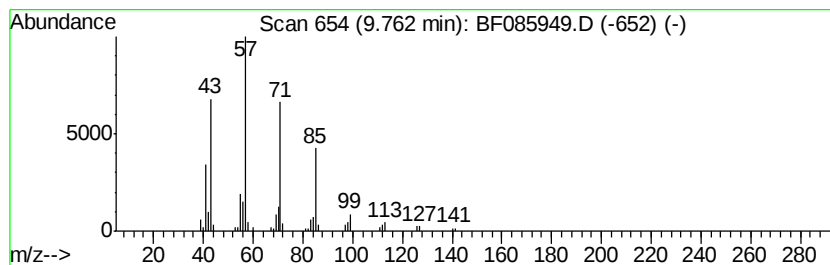
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tetradecane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.76	2.26 ng	96318	Acenaphthene-d10	9.83

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	86
2			Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	83
3			Eicosane	282	C20H42	000112-95-8	83
4			Dodecane, 2-methyl-	184	C13H28	001560-97-0	80
5			Hexadecane	226	C16H34	000544-76-3	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033116\
Data File : BF085949.D
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Operator : UM/SJ
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Instrument :
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ClientSampleId :
AMBOY-SB25COMP0.5-15.0

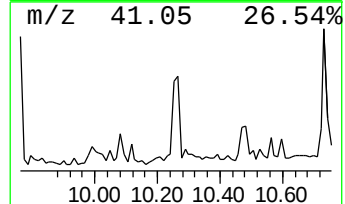
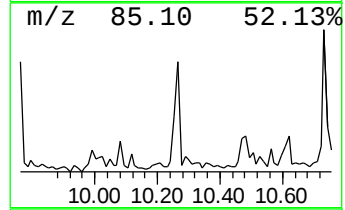
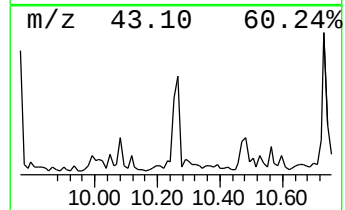
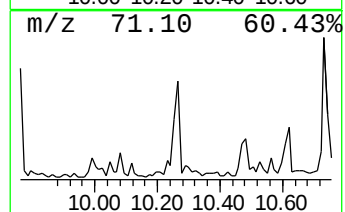
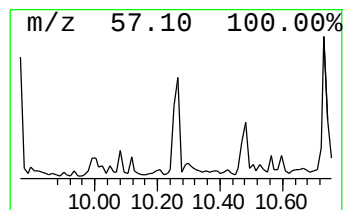
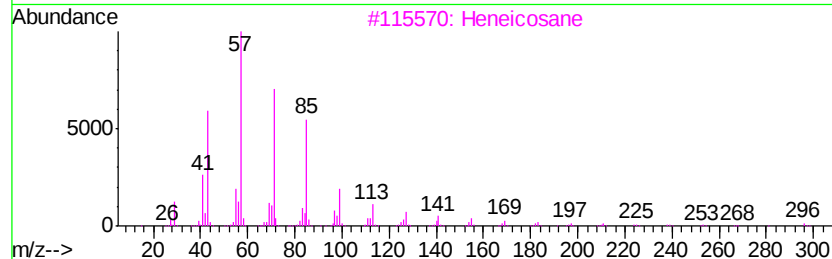
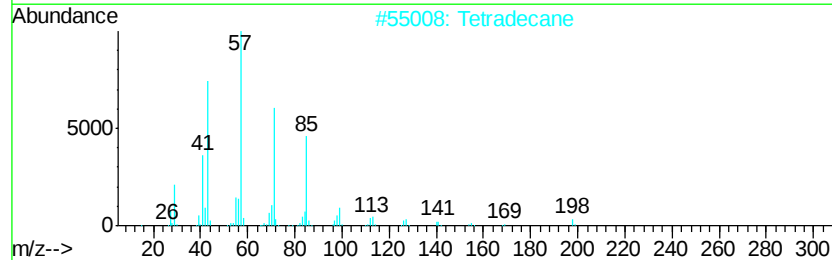
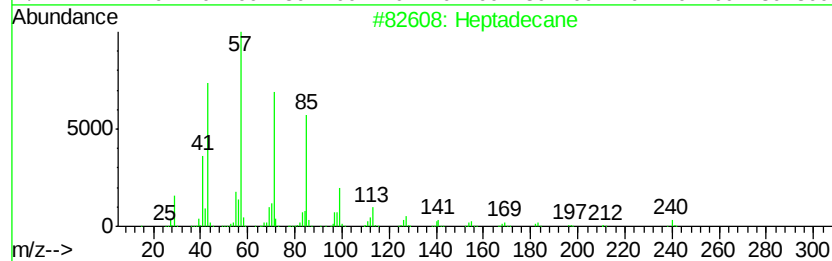
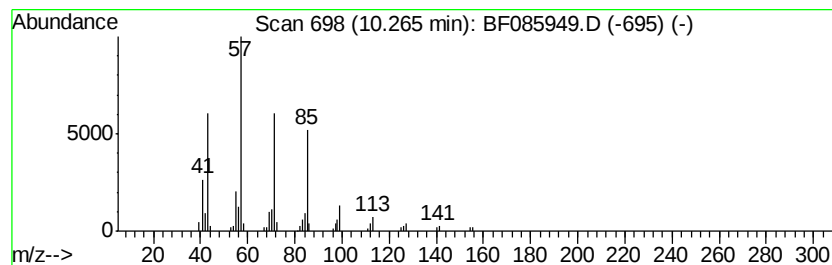
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heptadecane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.26	2.72 ng	115704	Acenaphthene-d10	9.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	91
2		Tetradecane	198	C14H30	000629-59-4	87
3		Heneicosane	296	C21H44	000629-94-7	86
4		Hexadecane	226	C16H34	000544-76-3	86
5		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033116\
Data File : BF085949.D
Acq On : 1 Apr 2016 1:37
Operator : UM/SJ
Sample : H2010-06
Misc :
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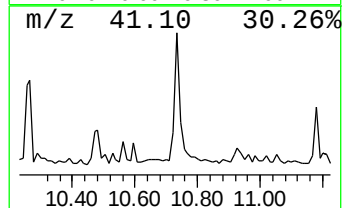
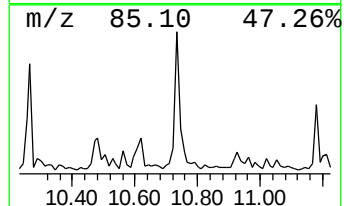
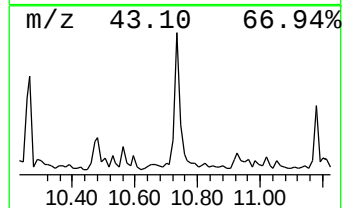
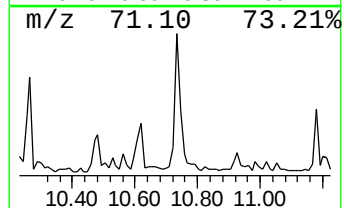
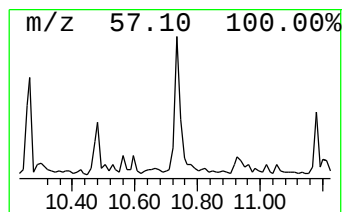
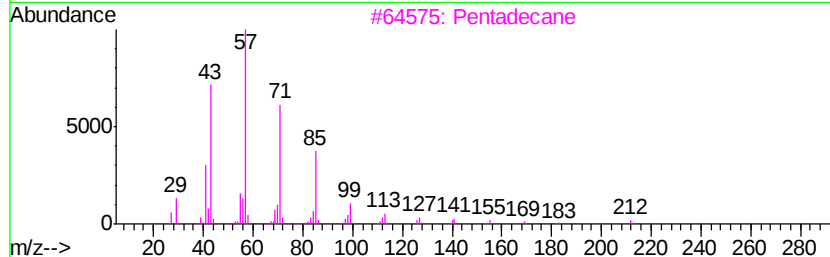
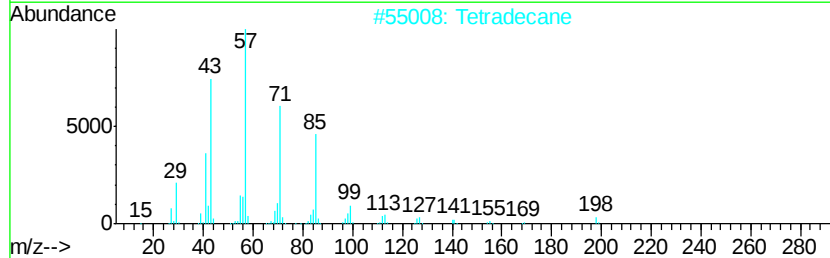
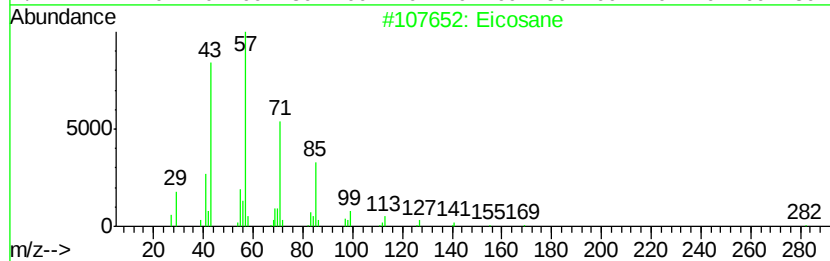
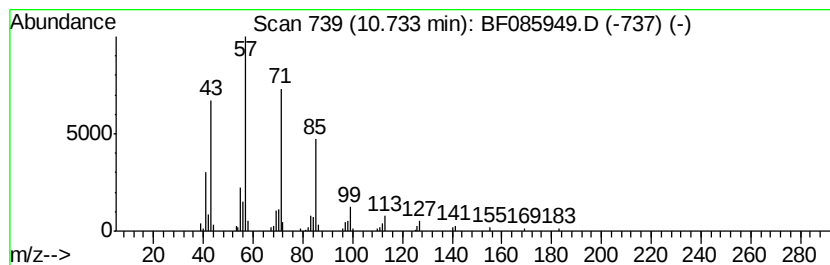
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Eicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.73	4.22 ng	161972	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	91
2		Tetradecane	198	C14H30	000629-59-4	91
3		Pentadecane	212	C15H32	000629-62-9	91
4		Heptadecane	240	C17H36	000629-78-7	91
5		Hexadecane	226	C16H34	000544-76-3	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033116\
Data File : BF085949.D
Acq On : 1 Apr 2016 1:37
Operator : UM/SJ
Sample : H2010-06
Misc :
ALS Vial : 24 Sample Multiplier: 1

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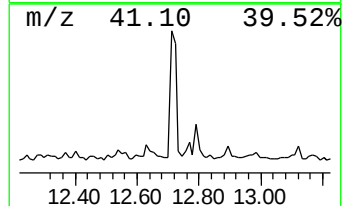
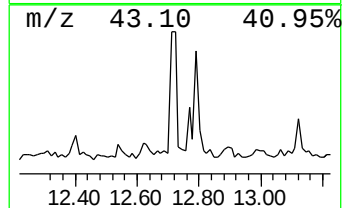
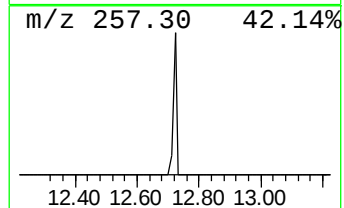
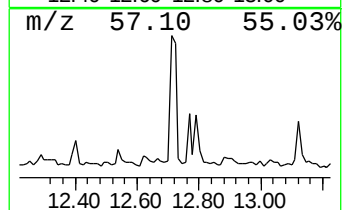
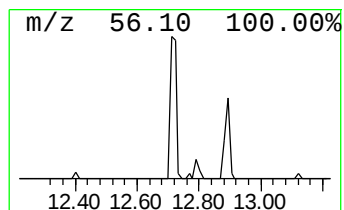
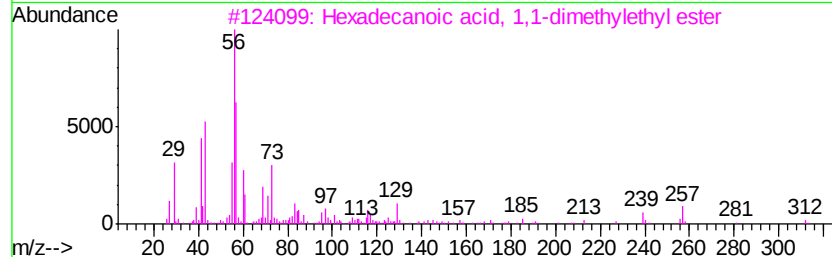
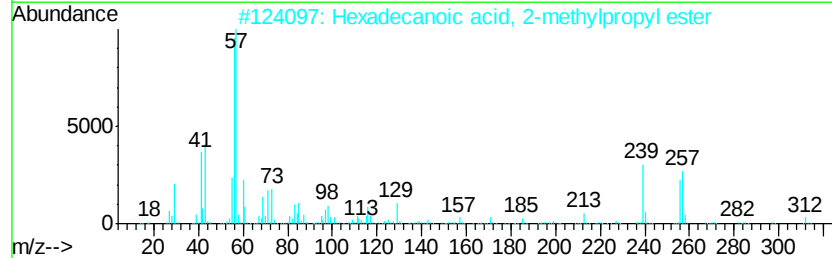
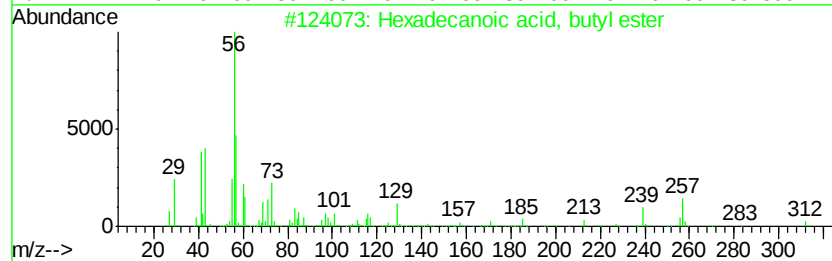
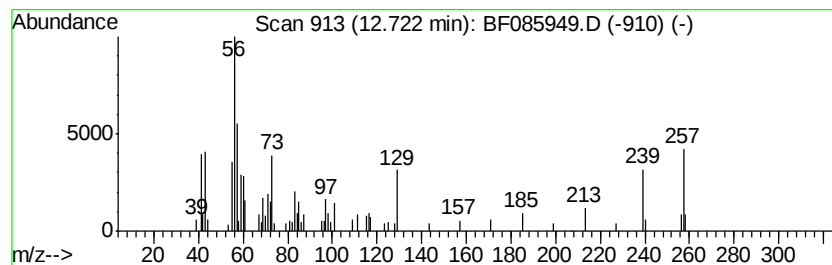
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Hexadecanoic acid, butyl ester Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.72	3.28 ng	87698	Chrysene-d12	13.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	64
2		Hexadecanoic acid, 2-methylpropy...	312	C20H40O2	000110-34-9	53
3		Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	43
4		1-Piperidinyloxy, 4-(hydroxyimin...	185	C9H17N2O2	003229-75-2	25
5		Cyclohexanol,2-amino-,trans-	115	C6H13NO	006982-39-4	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033116\
Data File : BF085949.D
Acq On : 1 Apr 2016 1:37
Operator : UM/SJ
Sample : H2010-06
Misc :
ALS Vial : 24 Sample Multiplier: 1

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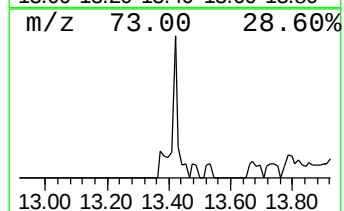
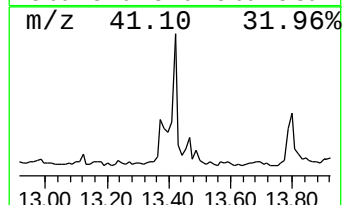
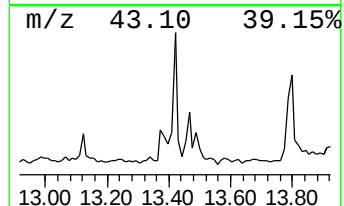
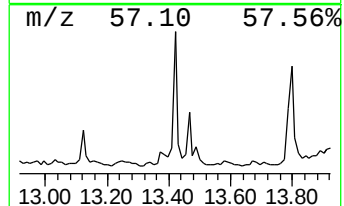
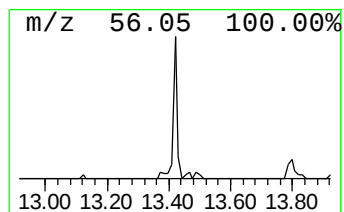
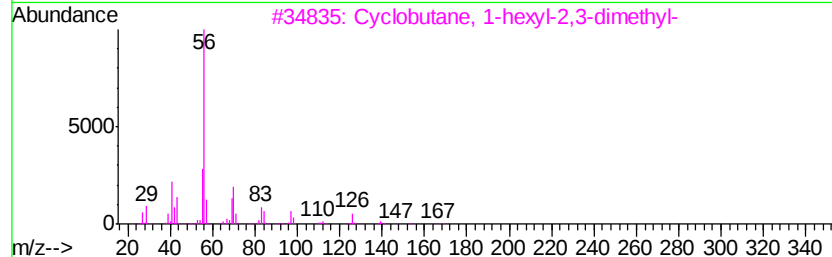
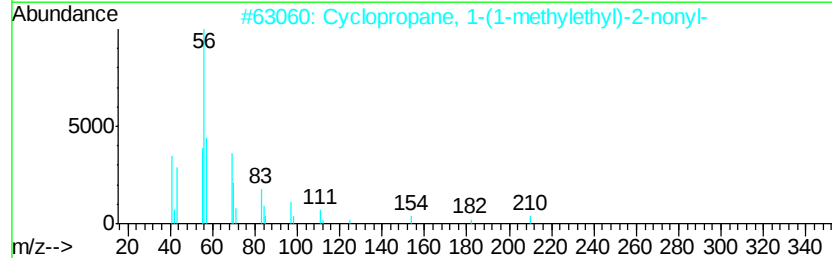
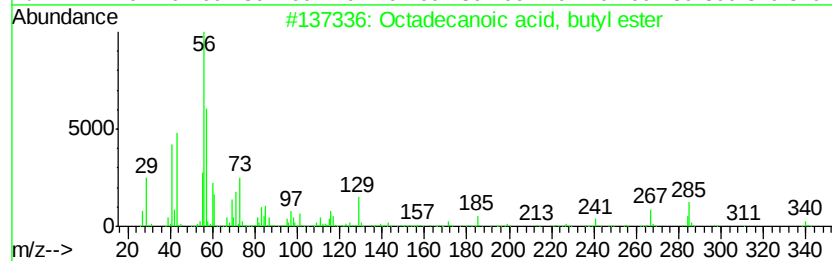
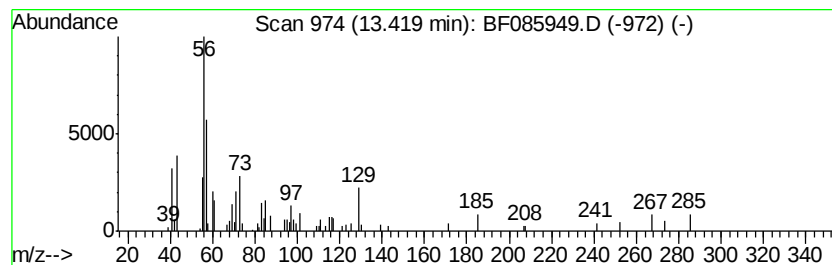
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Octadecanoic acid, butyl ester Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.42	3.12 ng	83380	Chrysene-d12	13.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	74
2			Cyclopropane, 1-(1-methylethyl)-...	210	C15H30	041977-39-3	35
3			Cyclobutane, 1-hexyl-2,3-dimethyl-	168	C12H24	055170-84-8	35
4			1-Hexene, 4-methyl-	98	C7H14	003769-23-1	30
5			Cyclopentanone, 2,2-dimethyl-	112	C7H12O	004541-32-6	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033116\
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Sample : H2010-06
Misc :
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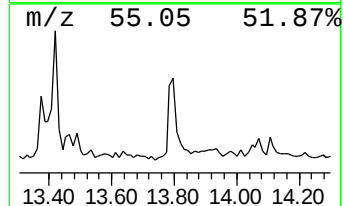
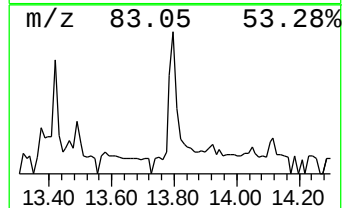
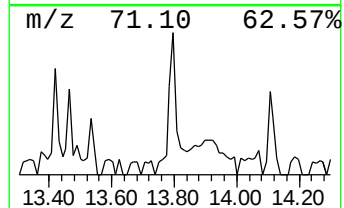
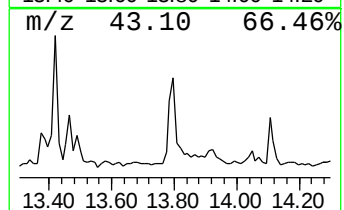
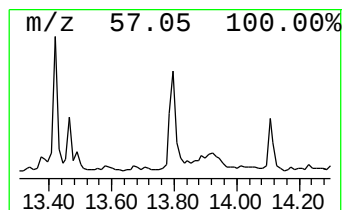
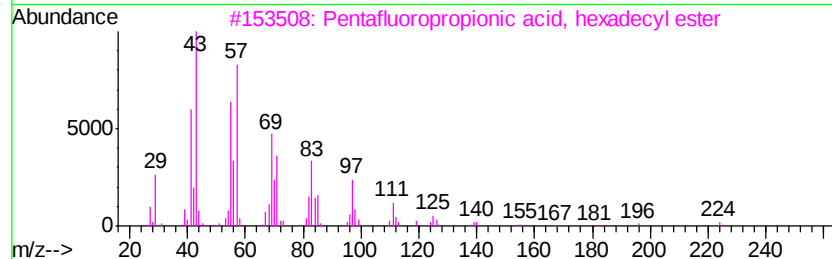
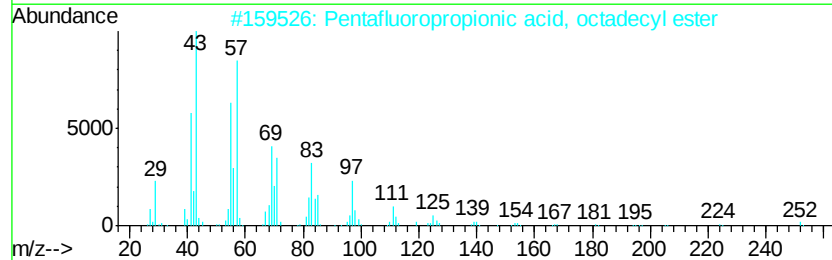
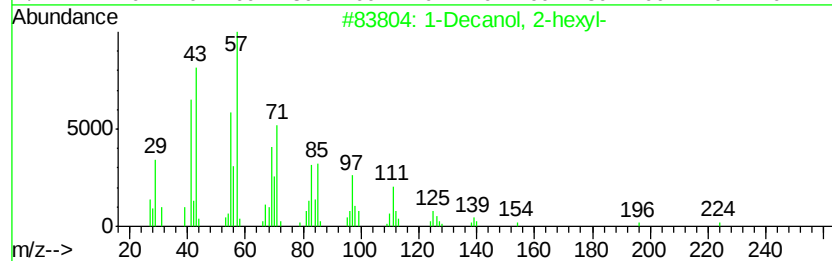
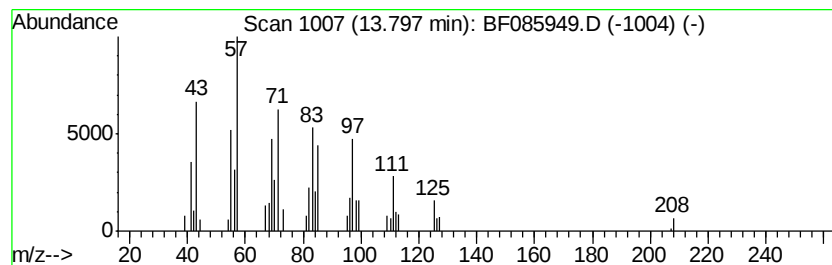
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 1-Decanol, 2-hexyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.80	2.99 ng	79997	Chrysene-d12	13.95
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Decanol, 2-hexyl-	242 C16H34O	002425-77-6	74
2	Pentafluoropropionic acid, octadecyl ester	416 C21H37F5O2	1000280-07-7	72
3	Pentafluoropropionic acid, hexadecyl ester	388 C19H33F5O2	006222-07-7	68
4	17-Pentatriacontene	491 C35H70	006971-40-0	64
5	1-Octadecene	252 C18H36	000112-88-9	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033116\
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 Acq On : 1 Apr 2016 1:37
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 Sample : H2010-06
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 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AMBOY-SB25COMP0.5-15.0

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.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
unknown2.50	2.50	4.0	ng	125438	1	6.79	626708	20.0
unknown2.60	2.60	3.6	ng	113205	1	6.79	626708	20.0
unknown2.78	2.78	11.4	ng	358406	1	6.79	626708	20.0
2-Pentanone, 4-hy...	4.97	3.4	ng	107274	1	6.79	626708	20.0
unknown6.56	6.56	91.7	ng	2874610	1	6.79	626708	20.0
Tetradecane	9.76	2.3	ng	96318	3	9.83	850907	20.0
Heptadecane	10.26	2.7	ng	115704	3	9.83	850907	20.0
Eicosane	10.73	4.2	ng	161972	4	11.32	768191	20.0
Hexadecanoic acid...	12.72	3.3	ng	87698	5	13.95	534277	20.0
Octadecanoic acid...	13.42	3.1	ng	83380	5	13.95	534277	20.0
1-Decanol, 2-hexyl-	13.80	3.0	ng	79997	5	13.95	534277	20.0