

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020217\
 Data File : BF092742.D
 Acq On : 2 Feb 2017 18:32
 Operator : SJ/MA
 Sample : I1658-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-COMP-1A-0-2FT

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-BF013017.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	4.807	236	238	248	rBV	31967	56072	1.10%	0.116%
2	5.127	263	266	276	rBV	1722692	2693480	52.74%	5.581%
3	5.550	300	303	306	rVB	4278629	5106755	100.00%	10.581%
4	6.579	390	393	396	rBV	3664761	4735531	92.73%	9.812%
5	6.704	401	404	407	rVV	3785266	4522102	88.55%	9.370%
6	6.933	422	424	426	rBV	1136971	951027	18.62%	1.971%
7	7.093	435	438	440	rBV	3797725	3555748	69.63%	7.367%
8	7.505	470	474	476	rBV	2942618	3308532	64.79%	6.855%
9	8.225	534	537	539	rBV	1101362	1276438	25.00%	2.645%
10	8.808	586	588	590	rBV	77538	69475	1.36%	0.144%
11	9.299	628	631	633	rBV	4567767	4543702	88.97%	9.415%
12	9.379	635	638	642	rVV2	85544	163062	3.19%	0.338%
13	9.630	657	660	663	rVV	56463	93561	1.83%	0.194%
14	9.688	663	665	669	rVB	414303	415580	8.14%	0.861%
15	9.905	680	684	686	rBV	159089	168284	3.30%	0.349%
16	9.973	686	690	692	rBV	1504063	1367208	26.77%	2.833%
17	10.179	705	708	710	rVV3	48883	77718	1.52%	0.161%
18	10.225	710	712	714	rVB3	44791	64644	1.27%	0.134%
19	10.396	724	727	729	rBV	161835	226965	4.44%	0.470%
20	10.625	745	747	750	rVB2	82852	132290	2.59%	0.274%
21	10.762	756	759	762	rBV2	1918075	2676824	52.42%	5.546%
22	10.876	767	769	772	rBV	260380	384337	7.53%	0.796%
23	10.945	773	775	779	rBV	26680	59580	1.17%	0.123%
24	11.059	783	785	787	rBV2	27581	56041	1.10%	0.116%
25	11.219	795	799	801	rBV4	32652	76143	1.49%	0.158%
26	11.322	806	808	809	rVV	232008	211501	4.14%	0.438%
27	11.356	809	811	813	rVB	97494	134364	2.63%	0.278%
28	11.459	818	820	824	rBV2	1401566	1640015	32.11%	3.398%
29	11.539	824	827	828	rVB	60012	100739	1.97%	0.209%
30	11.699	837	841	843	rBV4	50001	106665	2.09%	0.221%
31	11.745	843	845	848	rVB	136374	127850	2.50%	0.265%
32	11.962	861	864	865	rBV2	51110	60883	1.19%	0.126%
33	11.985	865	866	868	rVV	66947	64028	1.25%	0.133%
34	12.076	872	874	878	rVB3	70896	145943	2.86%	0.302%

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35	12.156	878	881	883	rBV	99950	108026	2.12%	0.224%
36	12.259	887	890	893	rVV2	31463	62507	1.22%	0.130%
37	12.316	893	895	899	rVB2	34123	59739	1.17%	0.124%
38	12.408	900	903	904	rVB3	42960	60105	1.18%	0.125%
39	12.431	904	905	910	rBV3	33160	62222	1.22%	0.129%
40	12.511	910	912	913	rBV2	46452	57665	1.13%	0.119%
41	12.545	913	915	916	rVB	93068	90989	1.78%	0.189%
42	12.671	924	926	928	rVB	479340	396464	7.76%	0.821%
43	12.717	928	930	932	rVB2	40505	56651	1.11%	0.117%
44	12.899	943	946	949	rBV	380129	443451	8.68%	0.919%
45	13.048	956	959	961	rBV	3530564	3504305	68.62%	7.261%
46	13.231	973	975	977	rVB2	42335	59367	1.16%	0.123%
47	13.265	977	978	980	rBV	66357	84663	1.66%	0.175%
48	13.608	1006	1008	1011	rBV3	63796	109087	2.14%	0.226%
49	13.734	1016	1019	1023	rVB3	66298	123394	2.42%	0.256%
50	13.939	1035	1037	1043	rVB	196790	301401	5.90%	0.625%
51	14.100	1047	1051	1052	rBV2	923713	1209541	23.69%	2.506%
52	14.557	1089	1091	1092	rBV	65109	79130	1.55%	0.164%
53	15.128	1139	1141	1145	rBV	143545	279310	5.47%	0.579%
54	15.197	1145	1147	1149	rVB2	60463	71785	1.41%	0.149%
55	15.437	1165	1168	1170	rVB	72519	124805	2.44%	0.259%
56	15.494	1170	1173	1176	rVB	117343	149254	2.92%	0.309%
57	15.563	1176	1179	1185	rVB	572562	968750	18.97%	2.007%
58	15.700	1188	1191	1193	rBV2	62051	135745	2.66%	0.281%
59	16.008	1216	1218	1221	rBV	51805	77440	1.52%	0.160%
60	17.003	1302	1305	1309	rBV2	57884	116429	2.28%	0.241%
61	17.448	1341	1344	1347	rVB	45740	97417	1.91%	0.202%

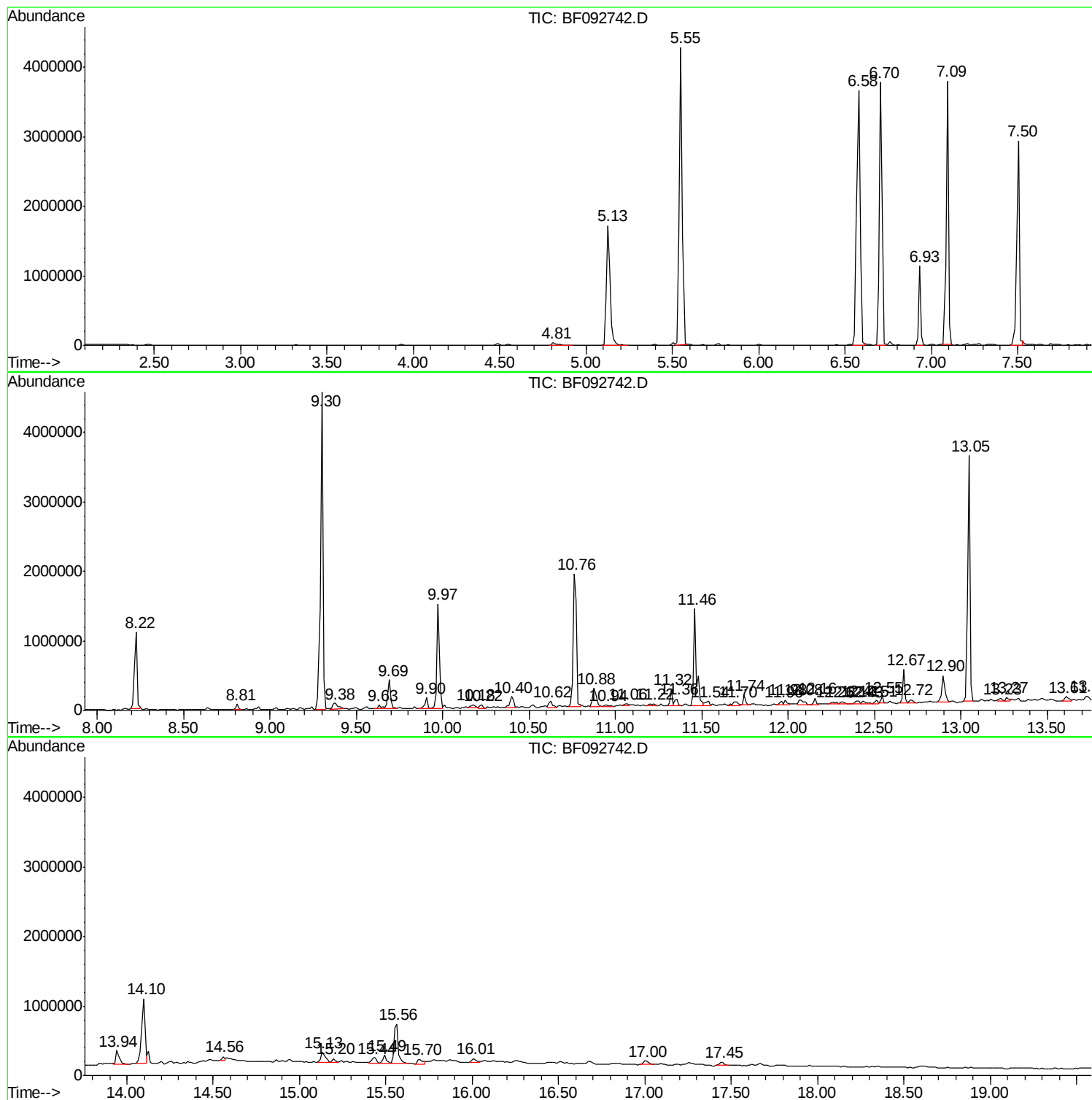
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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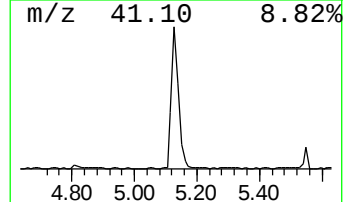
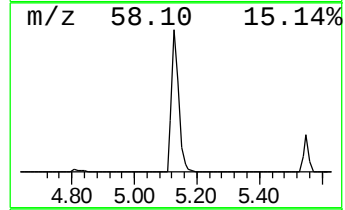
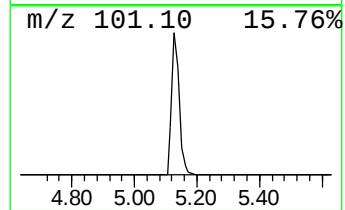
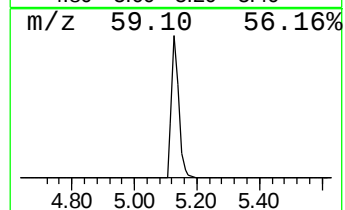
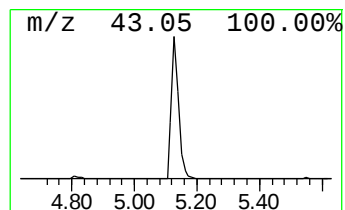
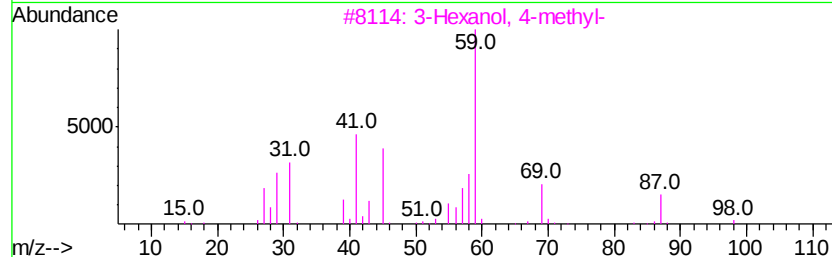
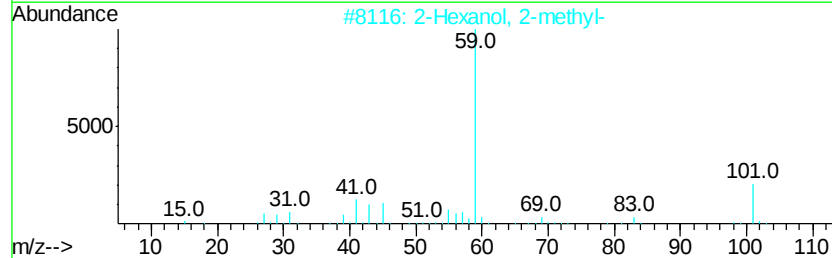
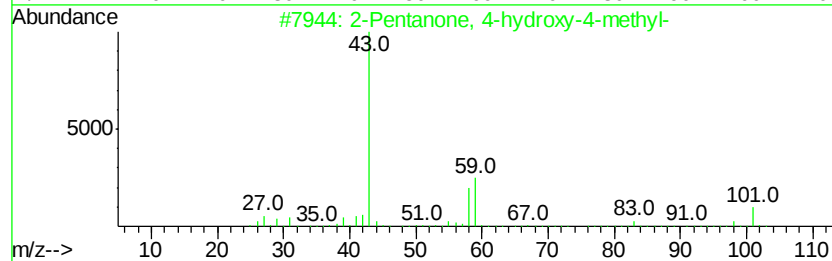
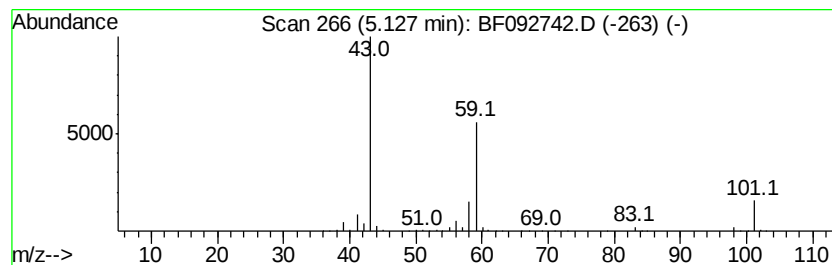
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	56.64 ng	2693480	1,4-Dichlorobenzene-d4	6.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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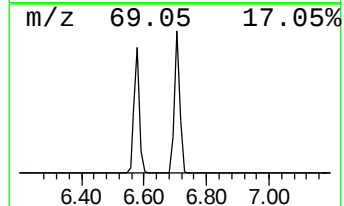
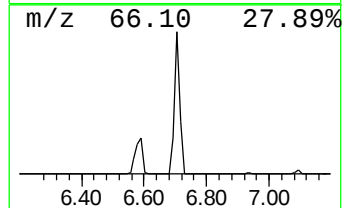
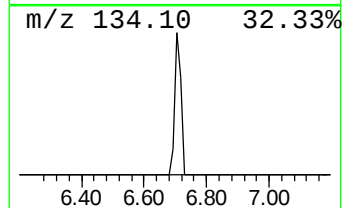
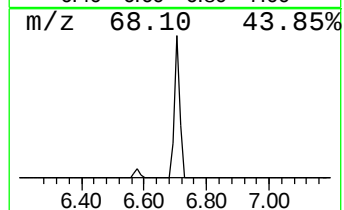
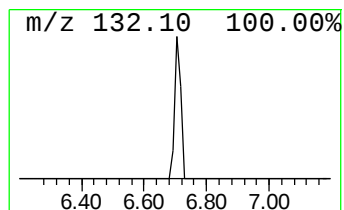
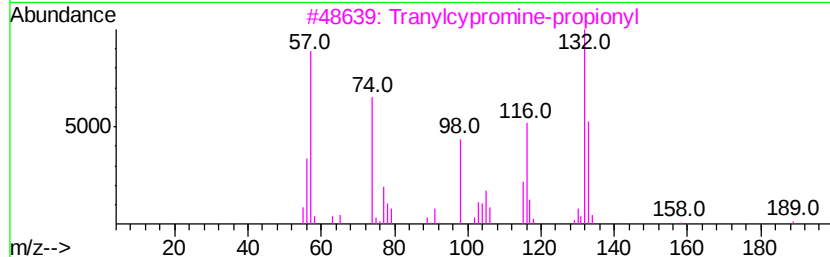
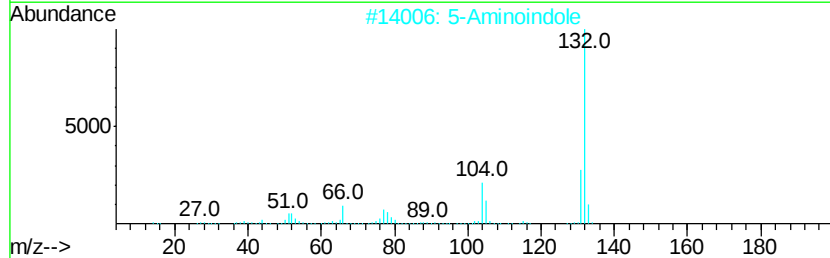
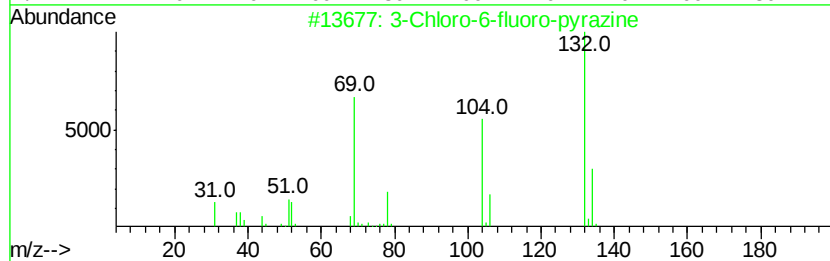
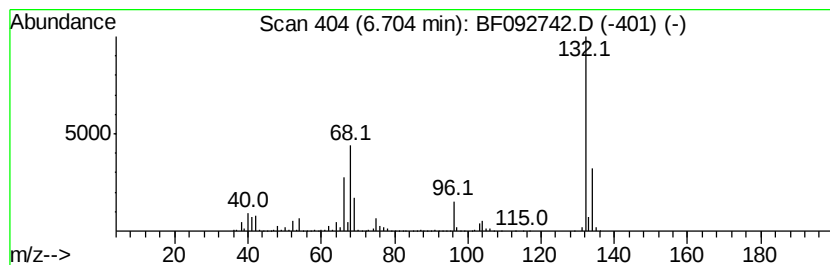
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 unknown6.70 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.70	95.10 ng	4522100	1,4-Dichlorobenzene-d4	6.93

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4			Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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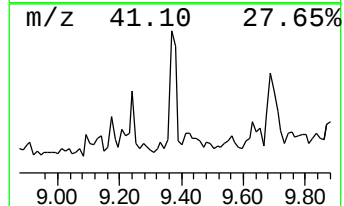
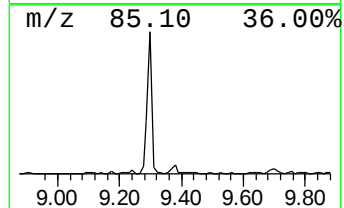
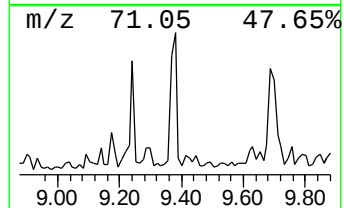
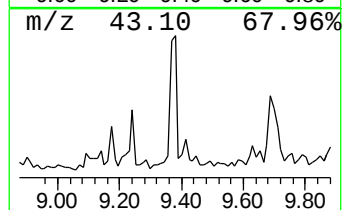
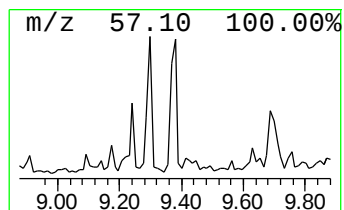
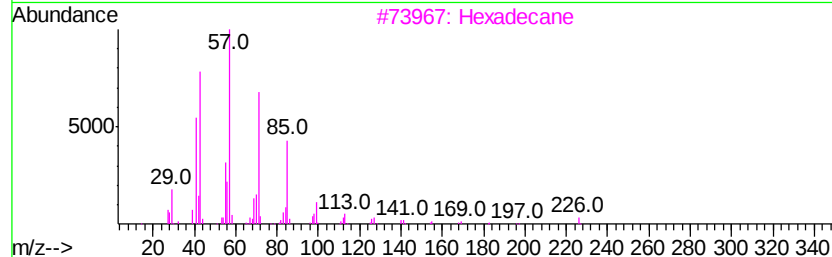
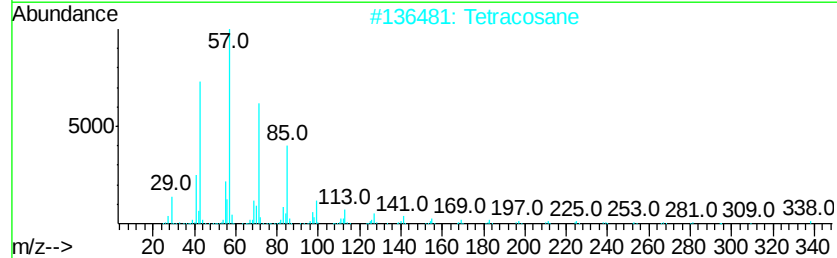
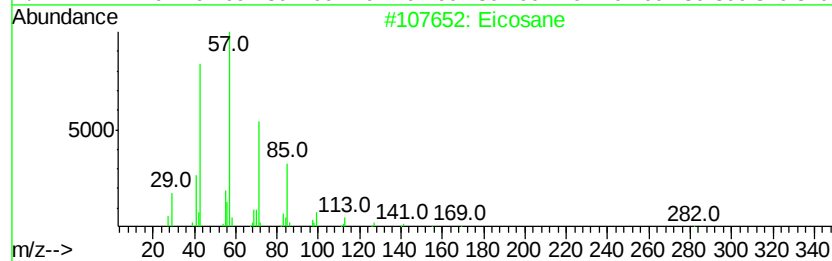
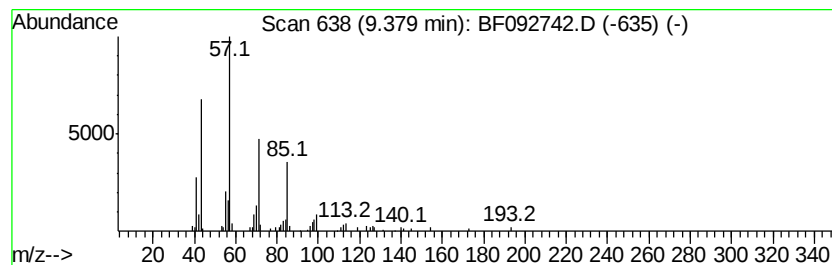
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 Eicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.38	2.39 ng	163062	Acenaphthene-d10	9.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	91
2		Tetracosane	338	C24H50	000646-31-1	90
3		Hexadecane	226	C16H34	000544-76-3	87
4		Dodecane	170	C12H26	000112-40-3	86
5		Nonadecane	268	C19H40	000629-92-5	86



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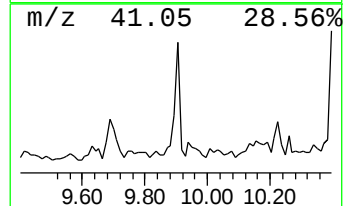
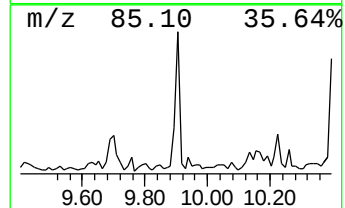
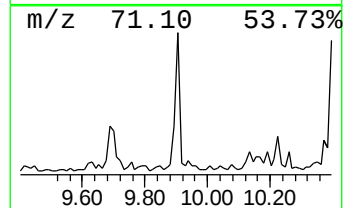
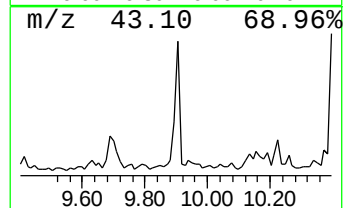
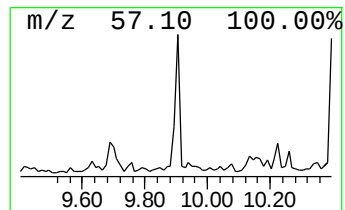
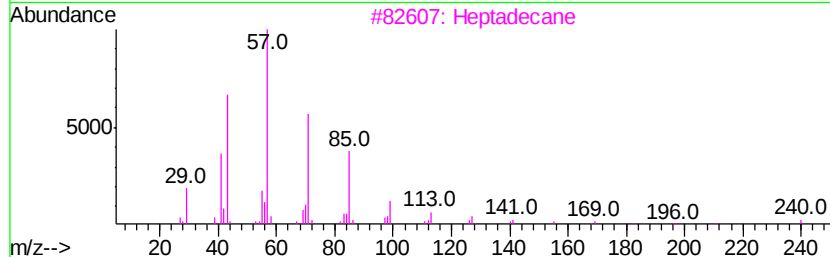
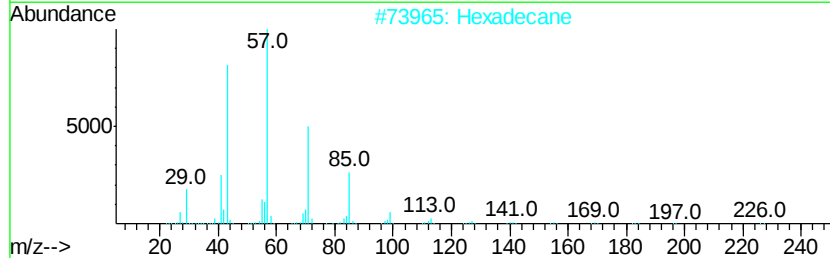
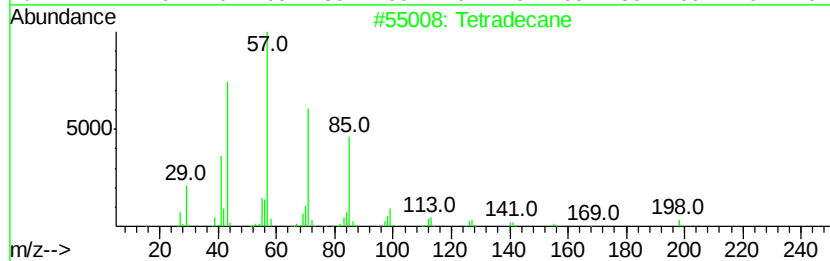
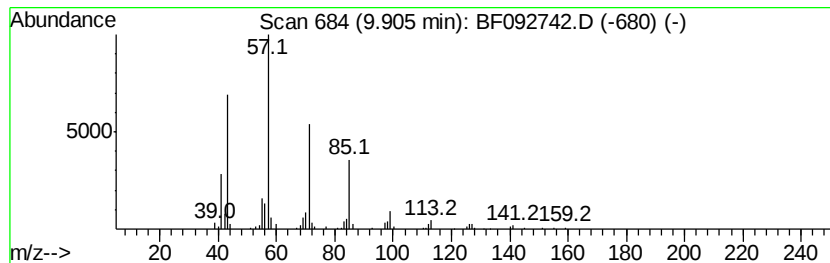
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Tetradecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.90	2.46 ng	168284	Acenaphthene-d10	9.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	91
2		Hexadecane	226	C16H34	000544-76-3	90
3		Heptadecane	240	C17H36	000629-78-7	90
4		Pentadecane	212	C15H32	000629-62-9	90
5		Tridecane	184	C13H28	000629-50-5	90



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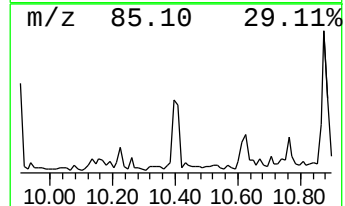
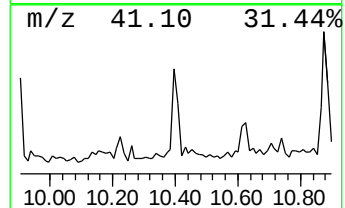
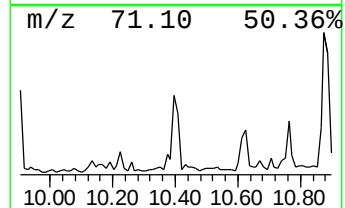
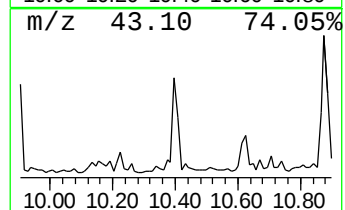
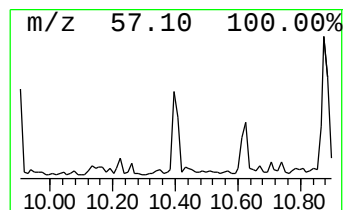
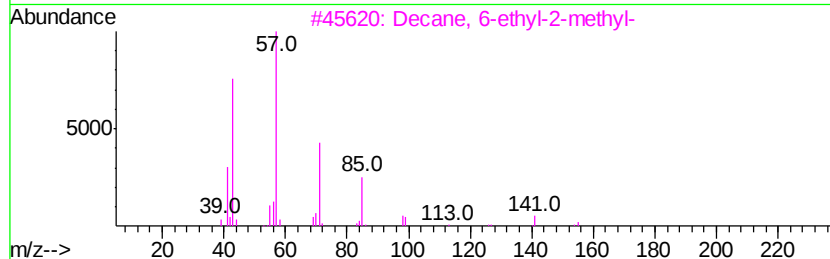
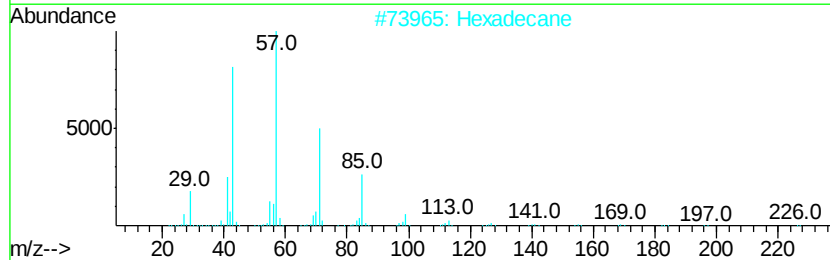
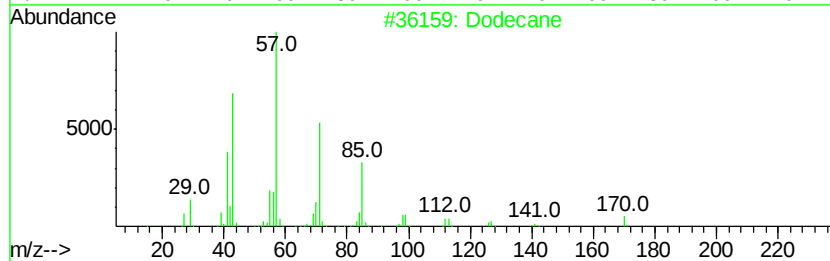
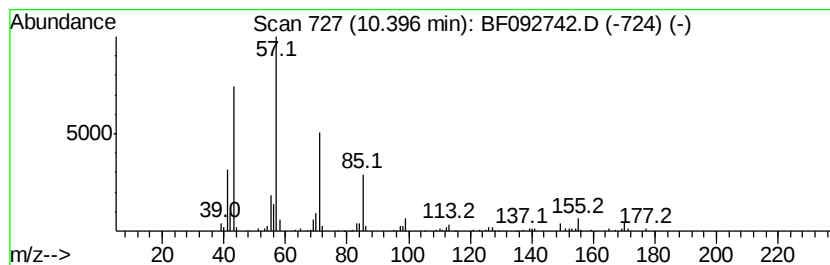
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ClientSampleId :
SP-COMP-1A-0-2FT

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Dodecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.40	3.32 ng	226965	Acenaphthene-d10	9.97
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Dodecane	170 C12H26	000112-40-3	83
2	Hexadecane	226 C16H34	000544-76-3	83
3	Decane, 6-ethyl-2-methyl-	184 C13H28	062108-21-8	78
4	Tridecane, 6-methyl-	198 C14H30	013287-21-3	72
5	Tetradecane	198 C14H30	000629-59-4	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020217\
Data File : BF092742.D
Acq On : 2 Feb 2017 18:32
Operator : SJ/MA
Sample : I1658-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-COMP-1A-0-2FT

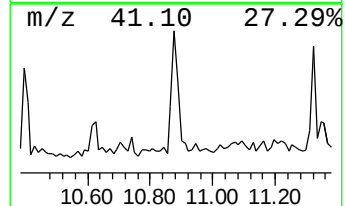
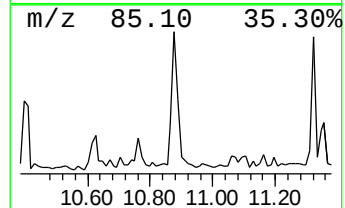
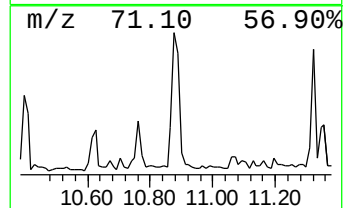
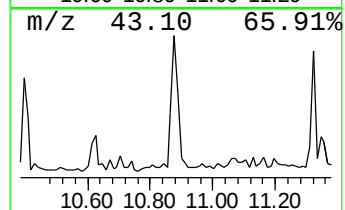
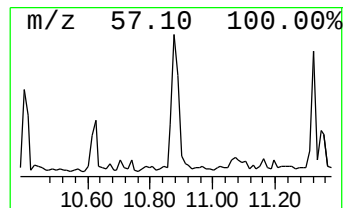
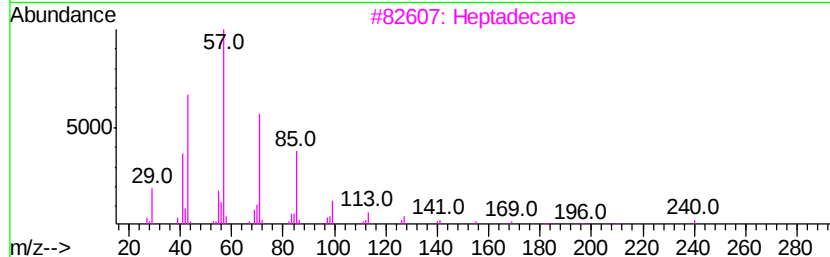
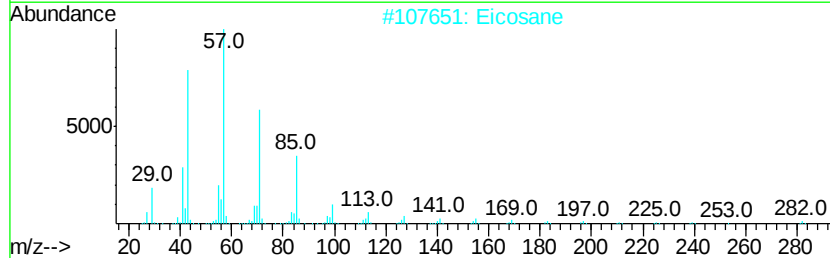
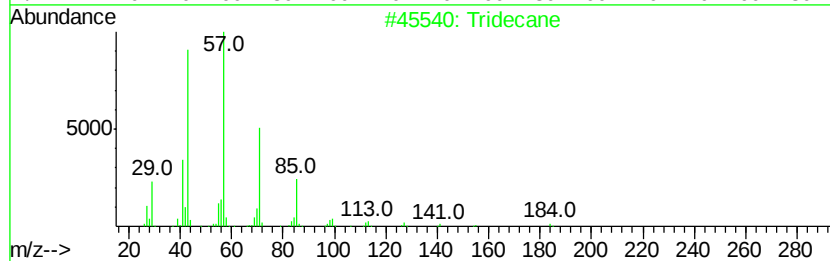
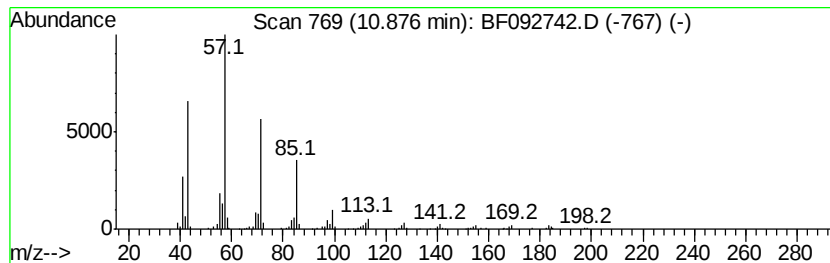
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tridecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.88	4.69 ng	384337	Phenanthrene-d10	11.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane	184	C13H28	000629-50-5	91
2		Eicosane	282	C20H42	000112-95-8	91
3		Heptadecane	240	C17H36	000629-78-7	90
4		Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	90
5		Tetradecane	198	C14H30	000629-59-4	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020217\
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Acq On : 2 Feb 2017 18:32
Operator : SJ/MA
Sample : I1658-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

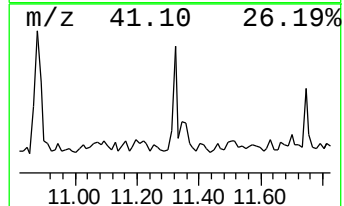
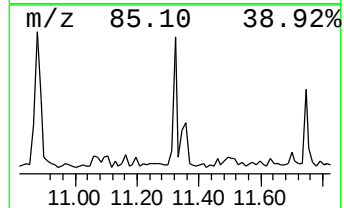
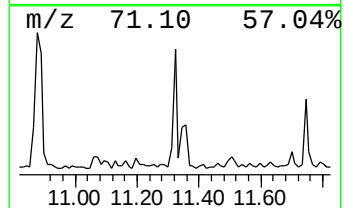
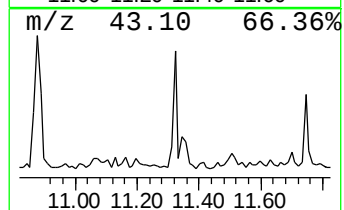
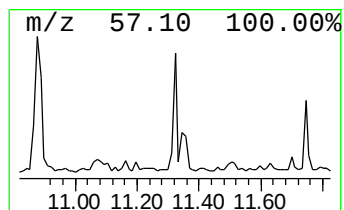
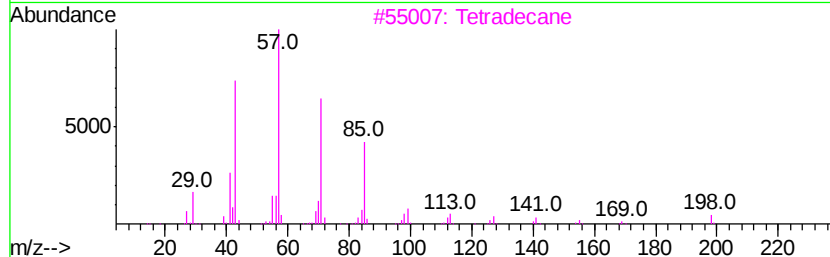
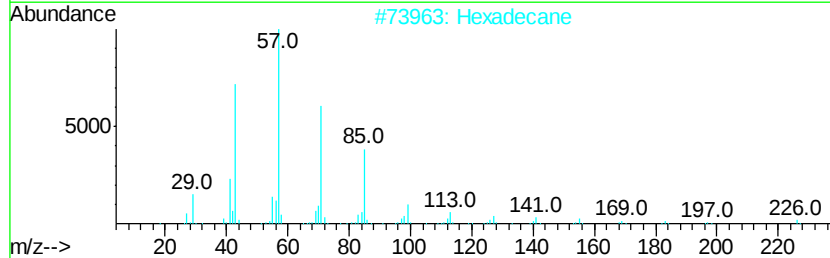
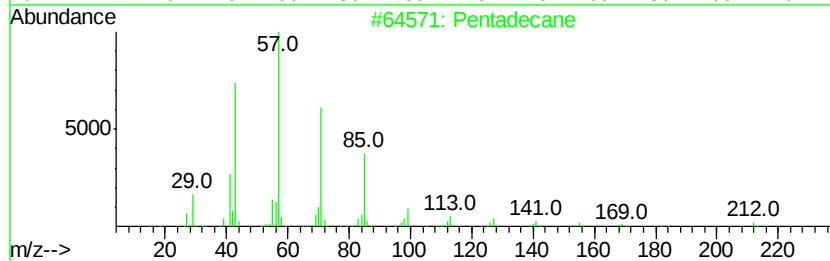
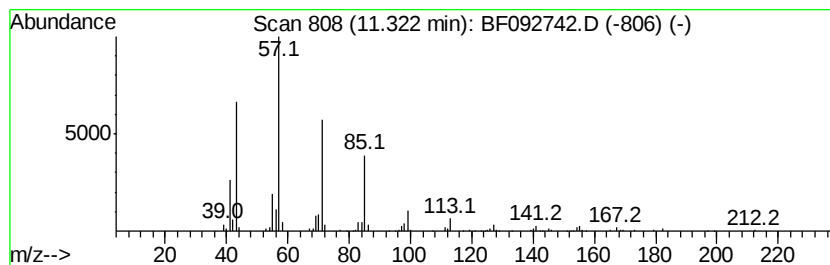
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ClientSampleId :
SP-COMP-1A-0-2FT

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Pentadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.		
11.32	2.58 ng	211501	Phenanthrene-d10		11.46		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane	212	C15H32	000629-62-9	94
2			Hexadecane	226	C16H34	000544-76-3	90
3			Tetradecane	198	C14H30	000629-59-4	80
4			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	80
5			Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	78



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020217\
Data File : BF092742.D
Acq On : 2 Feb 2017 18:32
Operator : SJ/MA
Sample : I1658-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-COMP-1A-0-2FT

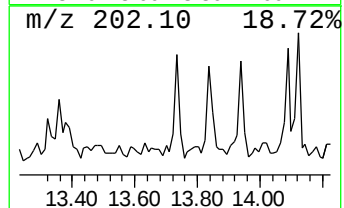
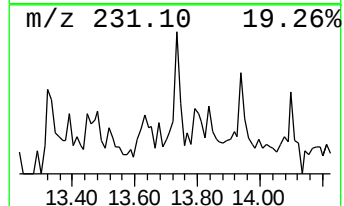
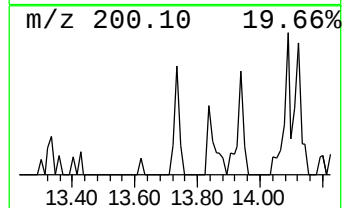
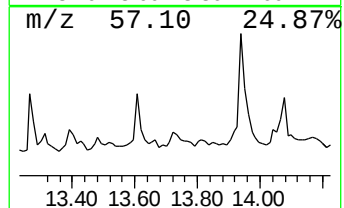
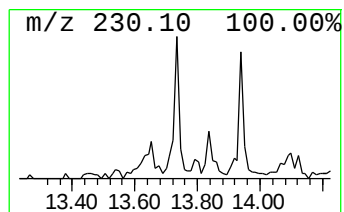
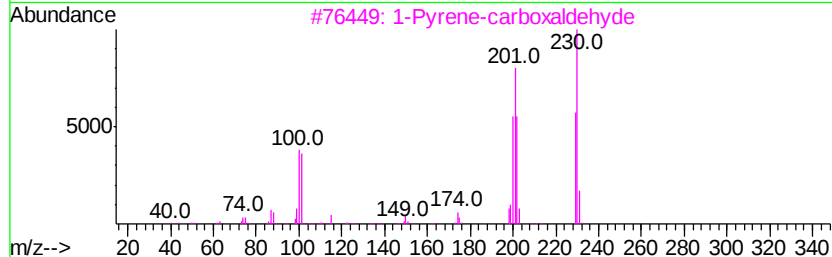
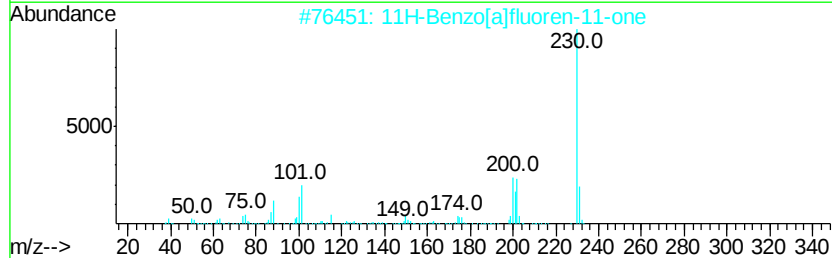
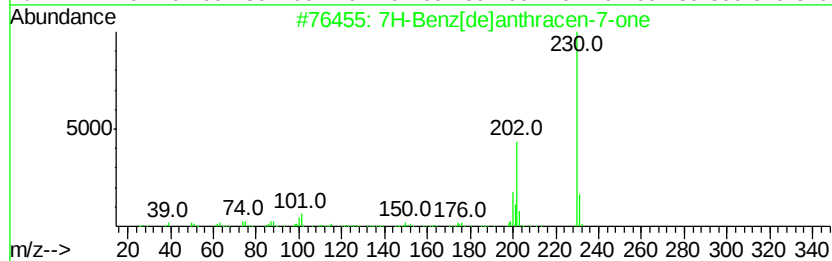
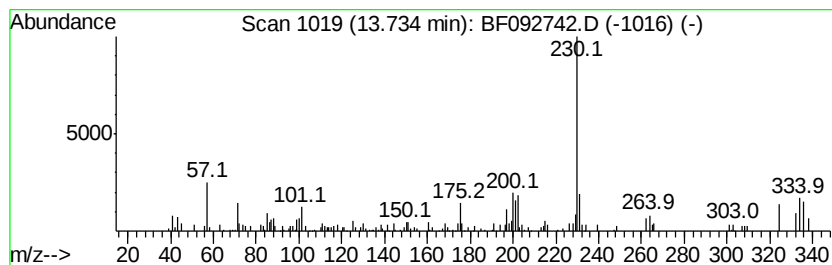
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 7H-Benz[de]anthracen-7-one Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	2.04 ng	123394	Chrysene-d12	14.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	64
2		11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	64
3		1-Pyrene-carboxaldehyde	230	C17H10O	003029-19-4	50
4		4-Chloro-indeno[1,2-b]pyridin-5-...	230	C12H7ClN2O	1000294-18-6	50
5		Pyrene, 1,3-dimethyl-	230	C18H14	064401-21-4	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020217\
Data File : BF092742.D
Acq On : 2 Feb 2017 18:32
Operator : SJ/MA
Sample : I1658-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SP-COMP-1A-0-2FT

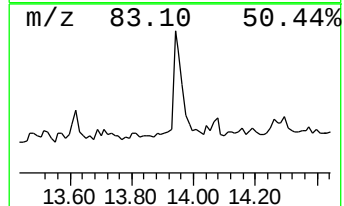
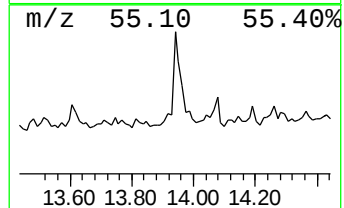
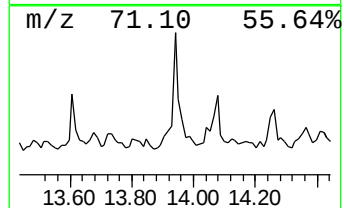
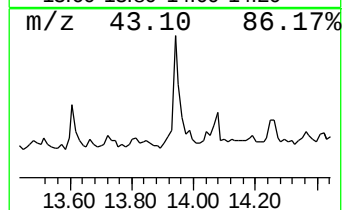
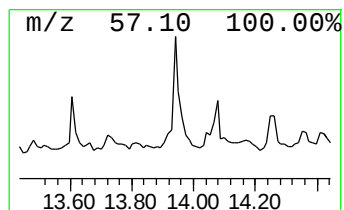
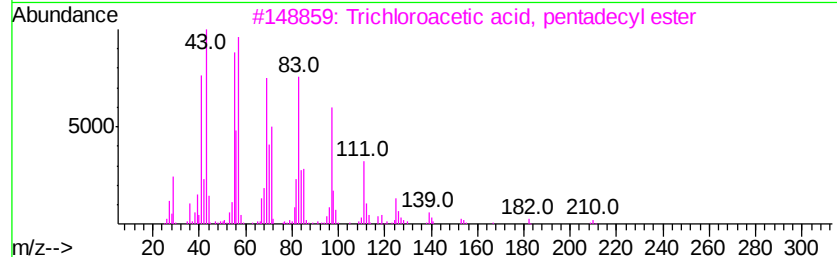
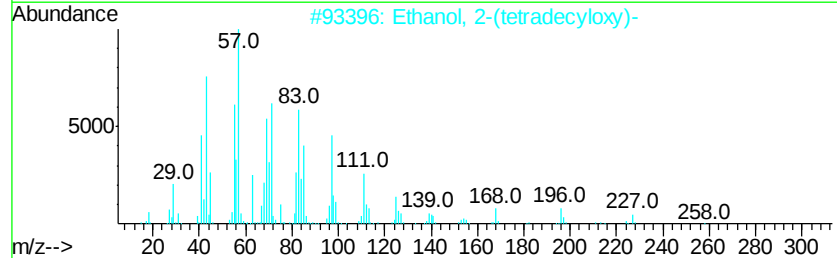
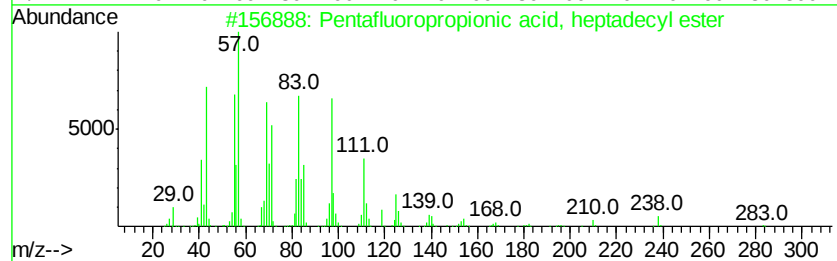
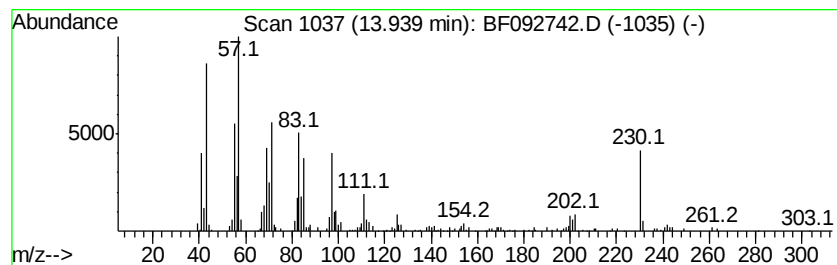
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Pentafluoropropionic acid, ... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.94	4.98 ng	301401	Chrysene-d12	14.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	74
2		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	68
3		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	64
4		Cyclododecane	168	C12H24	000294-62-2	60
5		3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	60



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020217\
Data File : BF092742.D
Acq On : 2 Feb 2017 18:32
Operator : SJ/MA
Sample : I1658-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SP-COMP-1A-0-2FT

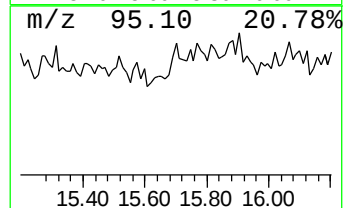
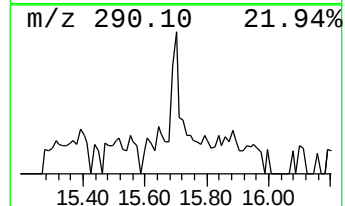
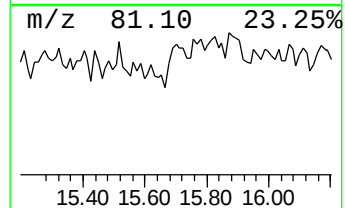
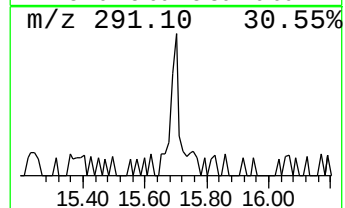
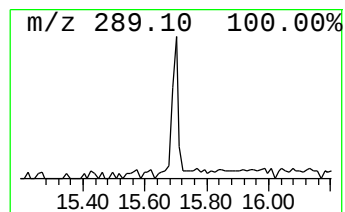
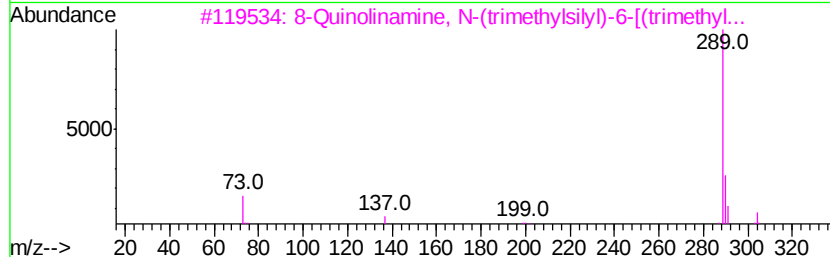
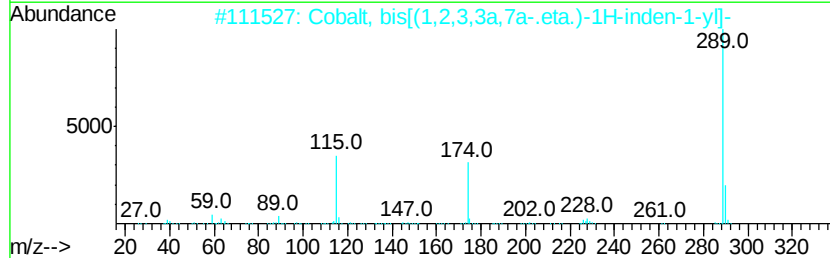
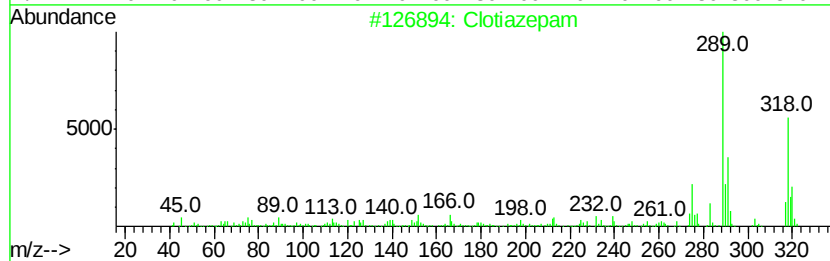
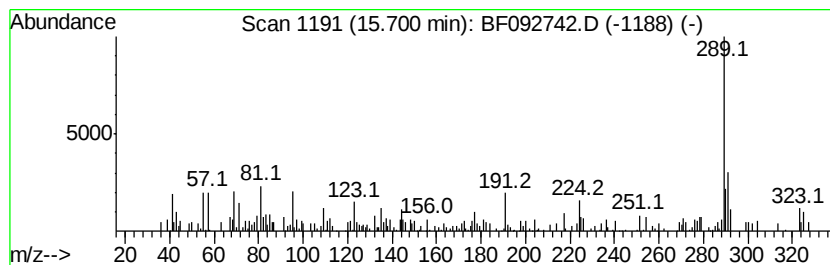
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown15.70 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.70	2.80 ng	135745	Perylene-d12	15.56

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Clotiazepam	318	C16H15ClN2OS	033671-46-4	38
2			Cobalt, bis[(1,2,3,3a,7a-.eta.)-...	289	C18H14Co	033056-04-1	38
3			8-Quinolinamine, N-(trimethylsil...	304	C15H24N2OSi2	036972-87-9	38
4			1H-Indole-2(3H)-one, 3-(4-triflu...	289	C16H10F3NO	035315-56-1	38
5			3-(4-Chlorophenyl)-7-methyl-6-pr...	318	C16H15ClN2OS	283170-54-7	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020217\
Data File : BF092742.D
Acq On : 2 Feb 2017 18:32
Operator : SJ/MA
Sample : I1658-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.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
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unknown6.70	6.70	95.1	ng	4522100	1	6.93	951027	20.0
Eicosane	9.38	2.4	ng	163062	3	9.97	1367210	20.0
Tetradecane	9.90	2.5	ng	168284	3	9.97	1367210	20.0
Dodecane	10.40	3.3	ng	226965	3	9.97	1367210	20.0
Tridecane	10.88	4.7	ng	384337	4	11.46	1640020	20.0
Pentadecane	11.32	2.6	ng	211501	4	11.46	1640020	20.0
7H-Benz[de]anthra...	13.73	2.0	ng	123394	5	14.10	1209540	20.0
Pentafluoropropio...	13.94	5.0	ng	301401	5	14.10	1209540	20.0
unknown15.70	15.70	2.8	ng	135745	6	15.56	968750	20.0