

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121516\
 Data File : BF091638.D
 Acq On : 15 Dec 2016 19:47
 Operator : UM/SJ
 Sample : H6057-03
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP20-COMP6

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

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

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.281	190	192	196	rBV	128653	213272	2.22%	0.217%
2	4.967	248	252	260	rVB	3110649	5072740	52.71%	5.167%
3	5.344	282	285	286	rBV	112227	138903	1.44%	0.141%
4	5.390	286	289	292	rVB	6314201	9624704	100.00%	9.804%
5	6.430	377	380	383	rBV	7082191	9066332	94.20%	9.236%
6	6.556	388	391	393	rBV	8471187	8708365	90.48%	8.871%
7	6.784	409	411	413	rBV	2265967	1983963	20.61%	2.021%
8	6.944	422	425	427	rBV	7040848	7537770	78.32%	7.678%
9	7.356	457	461	463	rBV	4974580	6733340	69.96%	6.859%
10	8.065	521	523	527	rBV2	2443962	2805366	29.15%	2.858%
11	9.150	615	618	621	rBV	8852431	8983966	93.34%	9.152%
12	9.242	623	626	628	rBV3	80837	141256	1.47%	0.144%
13	9.482	643	647	648	rBV	92508	105583	1.10%	0.108%
14	9.539	650	652	655	rVB	588998	536093	5.57%	0.546%
15	9.768	666	672	674	rVB3	130417	220506	2.29%	0.225%
16	9.825	674	677	678	rBV	2860037	2774913	28.83%	2.827%
17	9.848	678	679	682	rVB	202703	267936	2.78%	0.273%
18	10.019	693	694	696	rVB	72940	98931	1.03%	0.101%
19	10.259	714	715	717	rVB	168764	183935	1.91%	0.187%
20	10.362	723	724	728	rBV	183334	215188	2.24%	0.219%
21	10.488	733	735	737	rVB2	69772	109349	1.14%	0.111%
22	10.613	743	746	748	rBV	5125005	5411690	56.23%	5.513%
23	10.739	755	757	761	rVB	247960	346281	3.60%	0.353%
24	10.911	769	772	774	rBV3	58336	146023	1.52%	0.149%
25	11.025	781	782	784	rBV	101259	121393	1.26%	0.124%
26	11.162	791	794	795	rBV	66092	123789	1.29%	0.126%
27	11.185	795	796	800	rVV2	236850	319431	3.32%	0.325%
28	11.299	804	806	810	rVV2	2158989	3311737	34.41%	3.374%
29	11.379	810	813	814	rVB	274787	301502	3.13%	0.307%
30	11.528	824	826	830	rBV2	61800	138900	1.44%	0.141%
31	11.608	831	833	836	rVB2	135008	161599	1.68%	0.165%
32	11.802	848	850	851	rBV	146358	133924	1.39%	0.136%
33	11.836	851	853	855	rVV2	215301	367735	3.82%	0.375%
34	11.882	855	857	858	rVV	77024	107994	1.12%	0.110%

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Integration Parameters: rteint.p
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Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
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If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

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35	11.916	858	860	864	rVB2	328743	456651	4.74%	0.465%
36	12.099	874	876	880	rVB2	130720	171421	1.78%	0.175%
37	12.351	897	898	900	rBV	81222	101054	1.05%	0.103%
38	12.408	900	903	904	rVB3	68478	125391	1.30%	0.128%
39	12.511	910	912	914	rVB	1471653	1231890	12.80%	1.255%
40	12.556	914	916	919	rBV4	60298	96886	1.01%	0.099%
41	12.739	928	932	934	rBV	1252409	1474164	15.32%	1.502%
42	12.774	934	935	941	rVB4	87352	144631	1.50%	0.147%
43	12.888	943	945	948	rVB	5334840	6849524	71.17%	6.977%
44	12.957	948	951	953	rBV2	72082	115102	1.20%	0.117%
45	13.071	959	961	963	rVB	214586	220800	2.29%	0.225%
46	13.128	965	966	968	rVB2	156296	185429	1.93%	0.189%
47	13.174	968	970	971	rBV	140049	164349	1.71%	0.167%
48	13.379	985	988	992	rBV	1557205	1828424	19.00%	1.863%
49	13.448	992	994	997	rBV2	118324	230830	2.40%	0.235%
50	13.677	1012	1014	1016	rBV2	73747	136100	1.41%	0.139%
51	13.791	1022	1024	1027	rVB2	366252	445422	4.63%	0.454%
52	13.928	1033	1036	1041	rBV2	1492522	2807999	29.17%	2.860%
53	14.694	1101	1103	1104	rBV2	97739	105712	1.10%	0.108%
54	14.945	1122	1125	1130	rBV	466696	907260	9.43%	0.924%
55	15.220	1147	1149	1152	rVB	238665	380103	3.95%	0.387%
56	15.288	1152	1155	1157	rVV	340285	533276	5.54%	0.543%
57	15.345	1157	1160	1166	rVB	1365484	1869687	19.43%	1.905%
58	15.803	1197	1200	1201	rBV2	99547	185337	1.93%	0.189%
59	16.008	1215	1218	1224	rVB2	164892	355312	3.69%	0.362%
60	16.694	1275	1278	1283	rVB2	163303	333565	3.47%	0.340%
61	17.094	1311	1313	1317	rVB	106760	197098	2.05%	0.201%

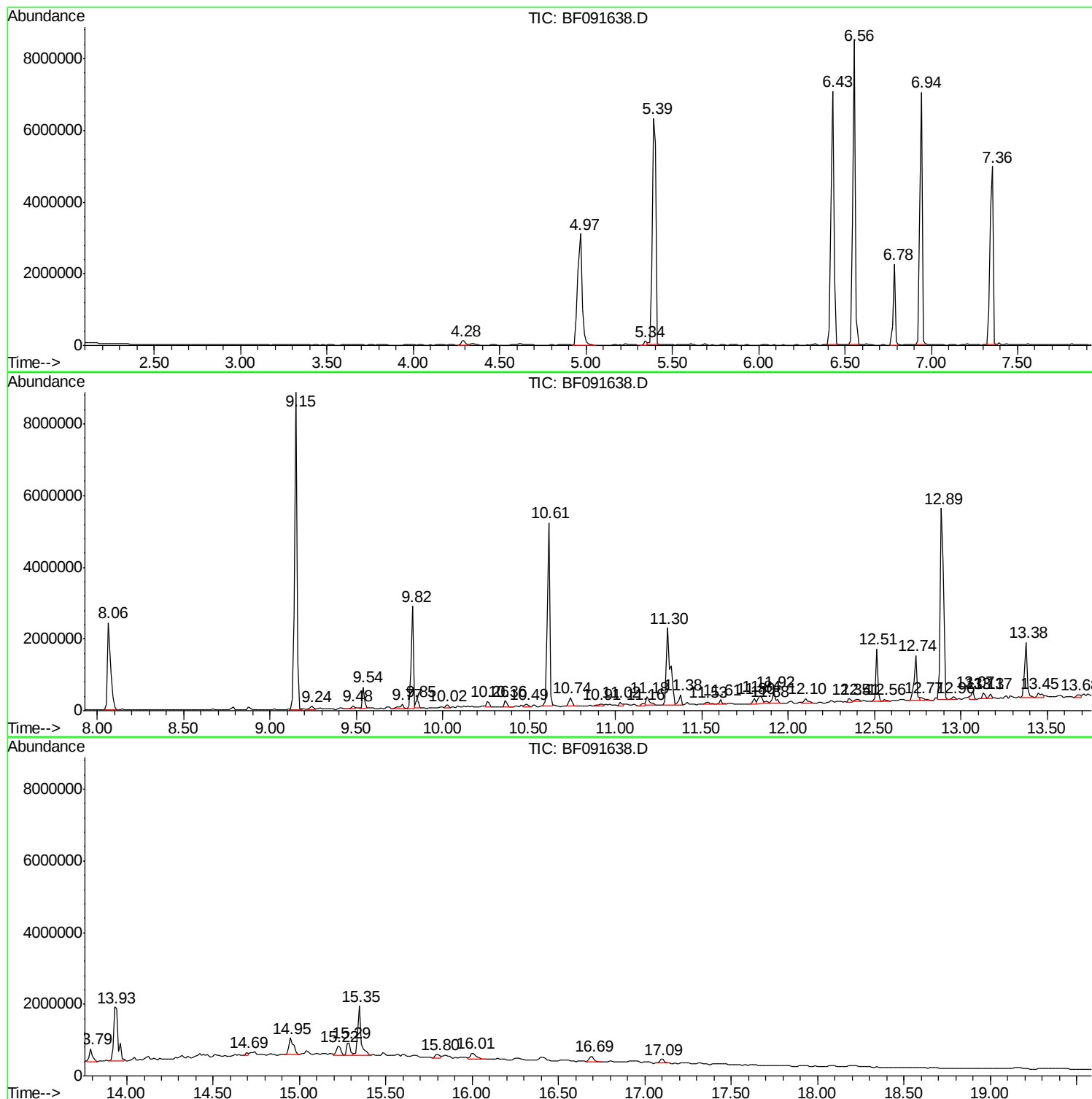
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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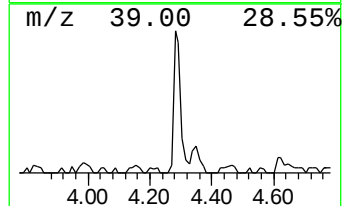
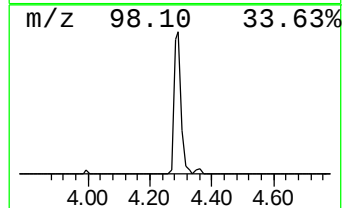
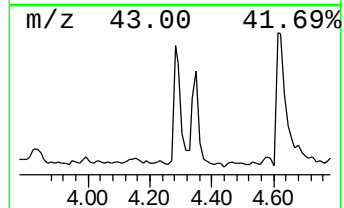
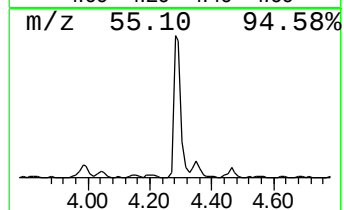
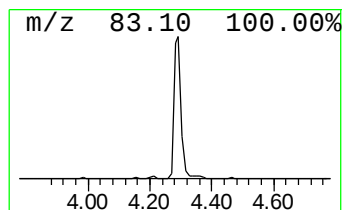
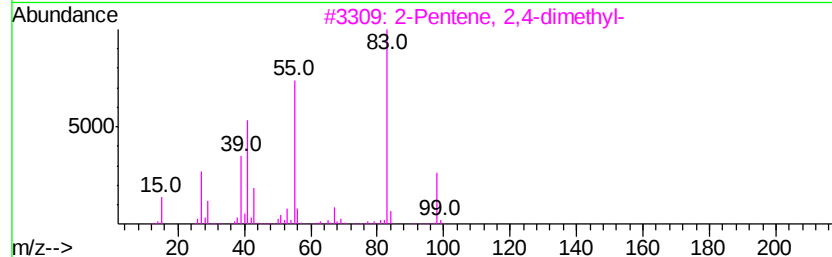
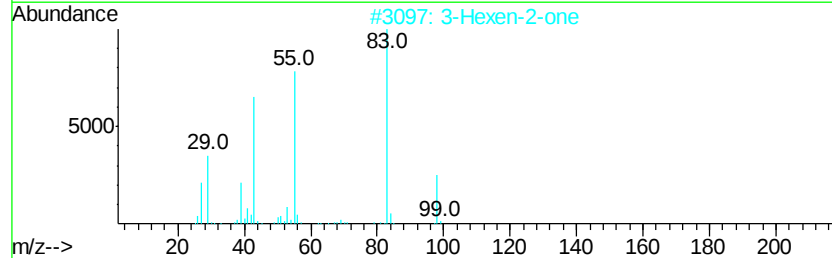
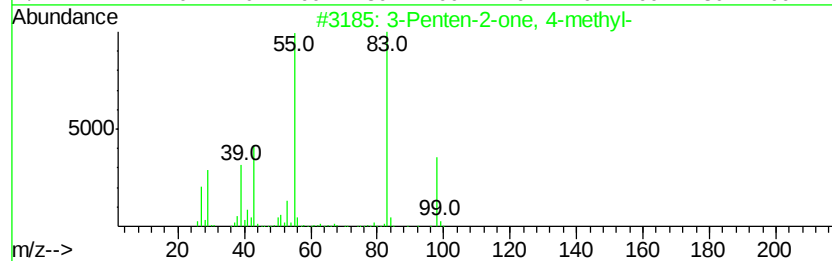
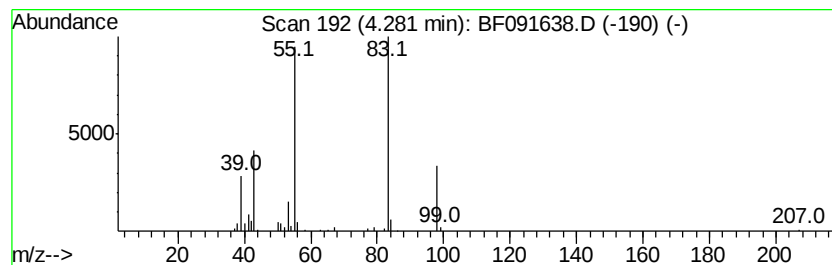
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.28	2.15 ng	213272	1,4-Dichlorobenzene-d4	6.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
2			3-Hexen-2-one	98	C6H10O	000763-93-9	87
3			2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	83
4			2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	80
5			2-Pentene, 4,4-dimethyl-, (Z)-	98	C7H14	000762-63-0	80



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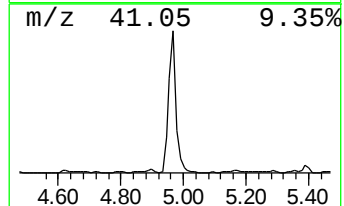
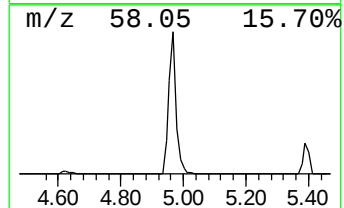
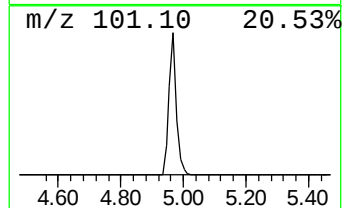
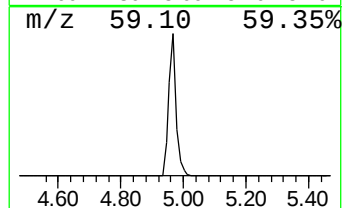
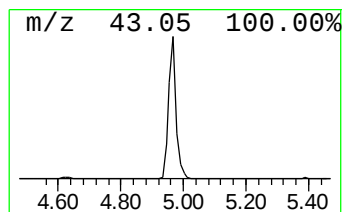
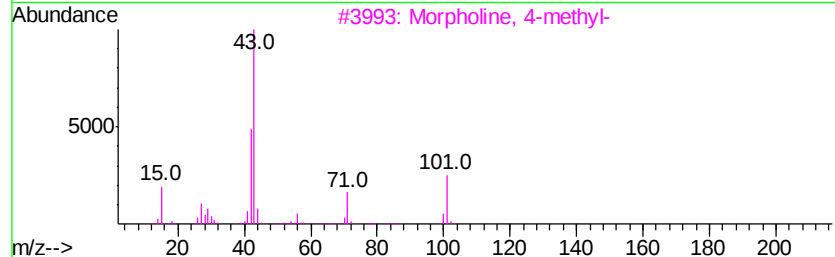
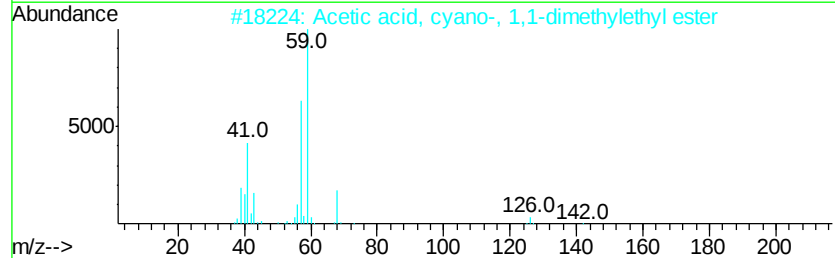
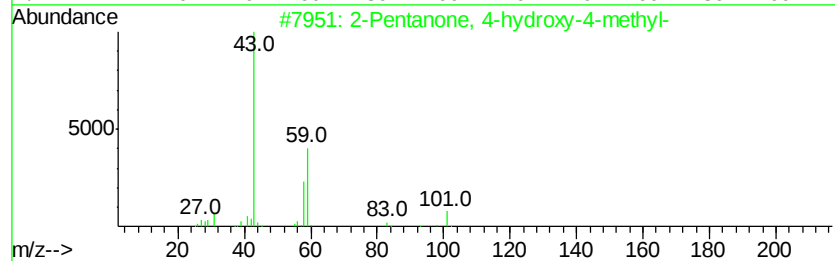
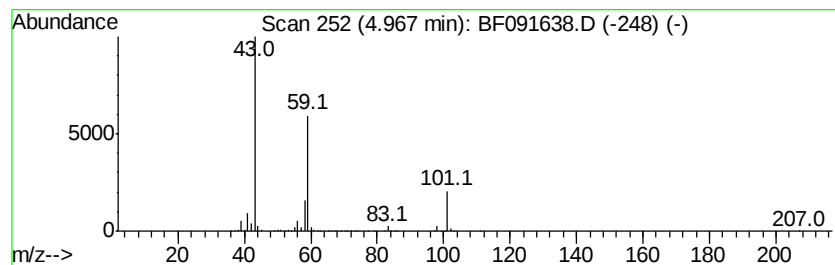
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.97	51.14 ng	5072740	1,4-Dichlorobenzene-d4	6.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	9



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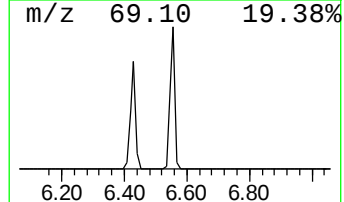
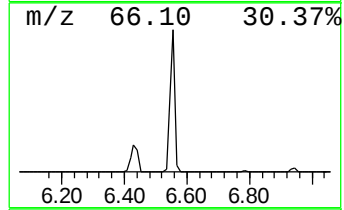
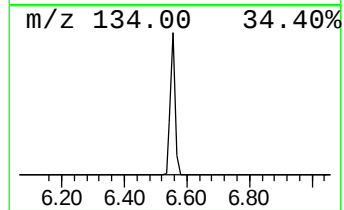
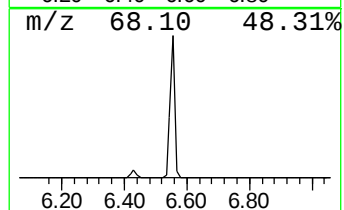
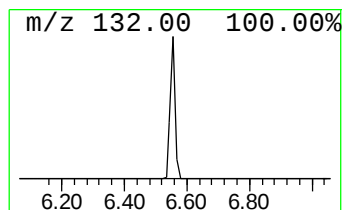
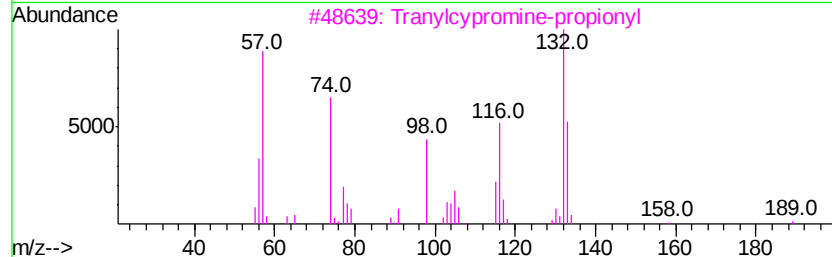
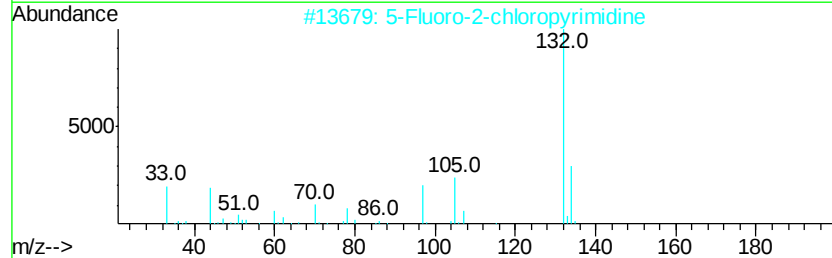
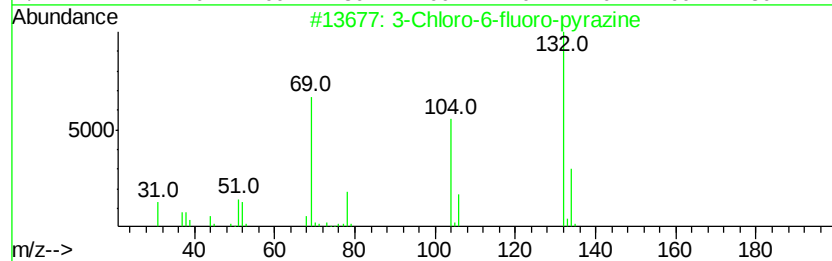
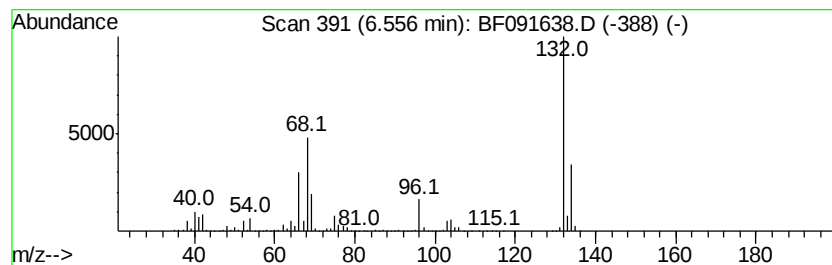
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	87.79 ng	8708370	1,4-Dichlorobenzene-d4	6.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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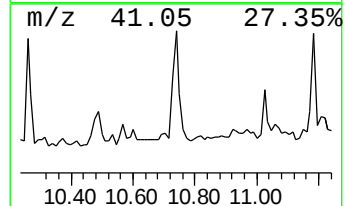
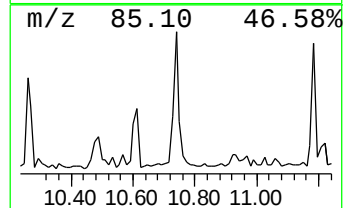
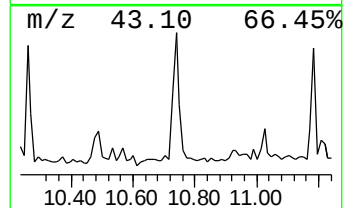
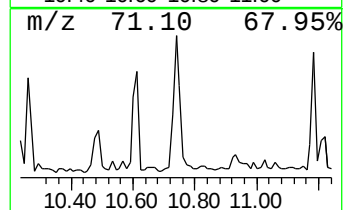
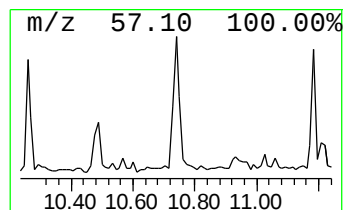
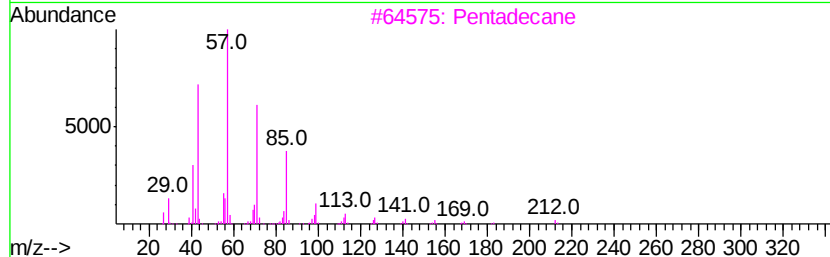
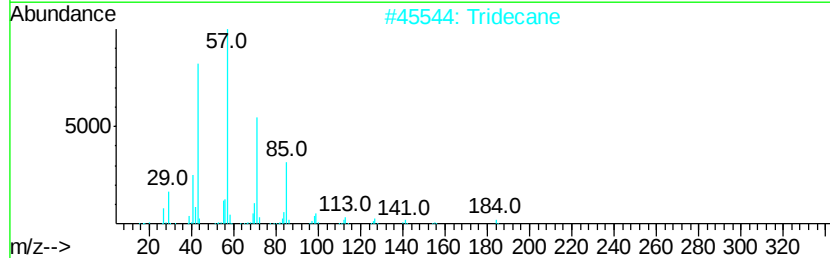
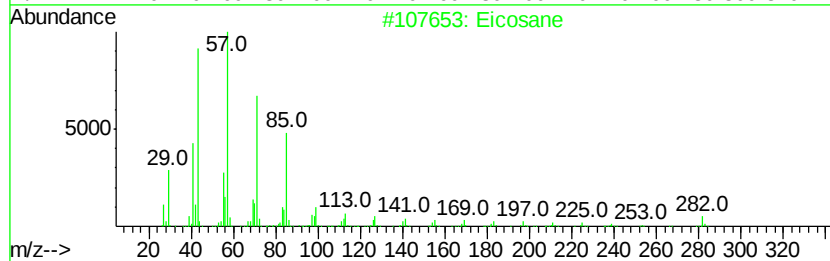
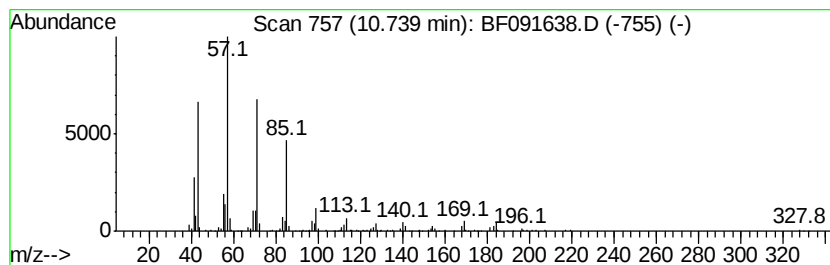
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Eicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.		
10.74	2.09 ng	346281	Phenanthrene-d10		11.30		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane			282	C20H42	000112-95-8	87
2	Tridecane			184	C13H28	000629-50-5	87
3	Pentadecane			212	C15H32	000629-62-9	86
4	Heptacosane, 1-chloro-			414	C27H55Cl	062016-79-9	86
5	Hexadecane			226	C16H34	000544-76-3	86



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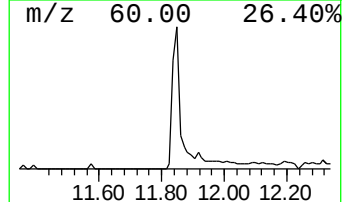
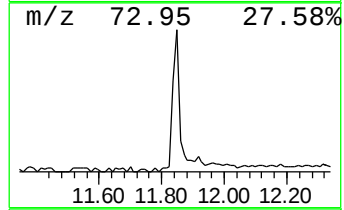
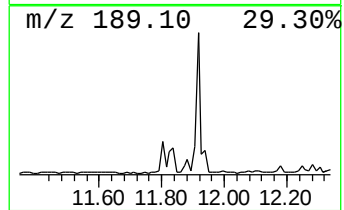
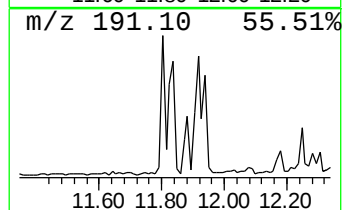
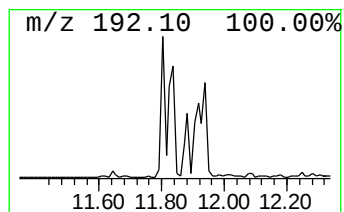
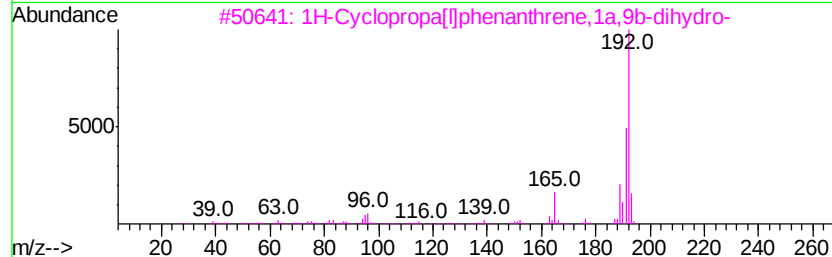
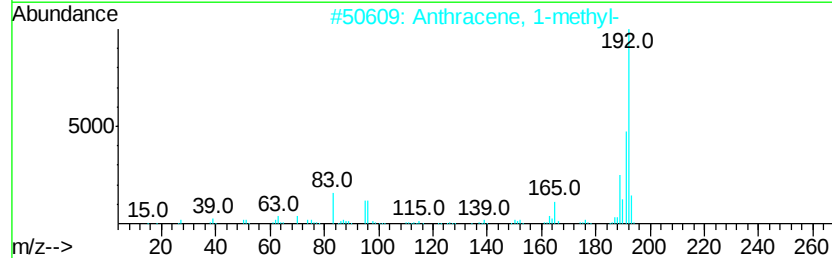
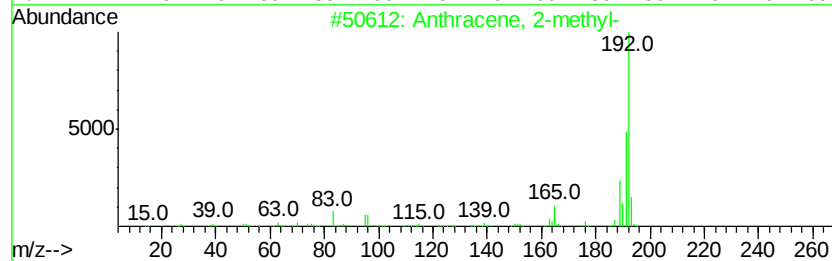
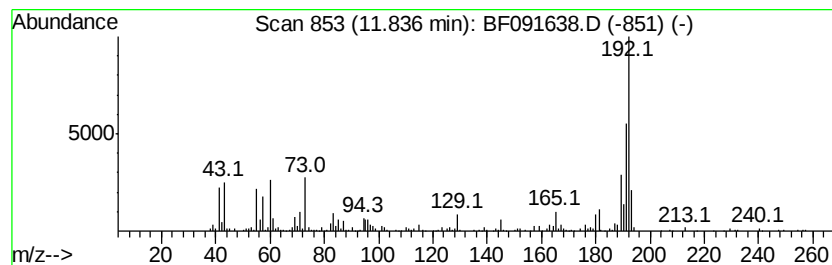
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Anthracene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.84	2.22 ng	367735	Phenanthrene-d10	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
2		Anthracene, 1-methyl-	192	C15H12	000610-48-0	92
3		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	91
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
5		Phenanthrene, 9-methyl-	192	C15H12	000883-20-5	83



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121516\
Data File : BF091638.D
Acq On : 15 Dec 2016 19:47
Operator : UM/SJ
Sample : H6057-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP20-COMP6

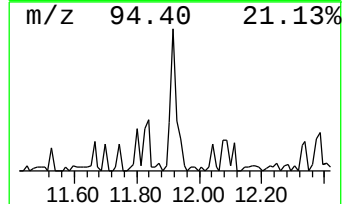
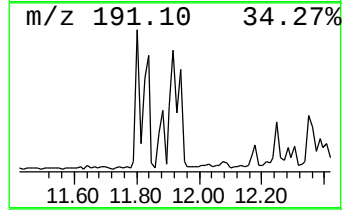
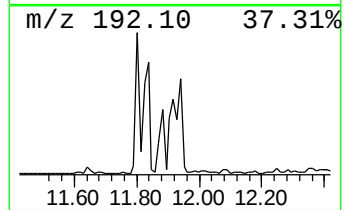
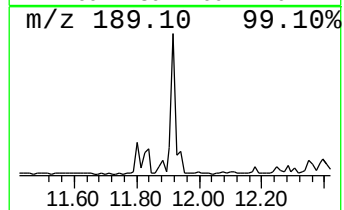
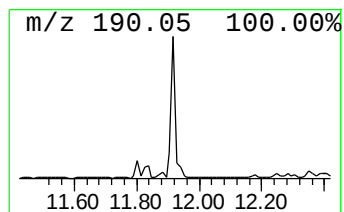
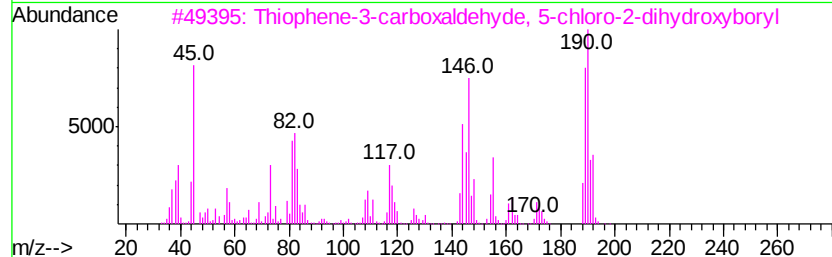
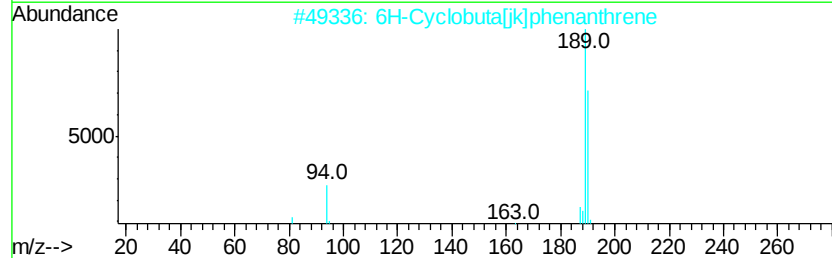
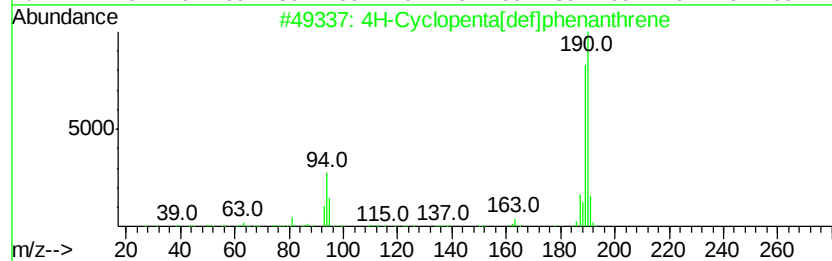
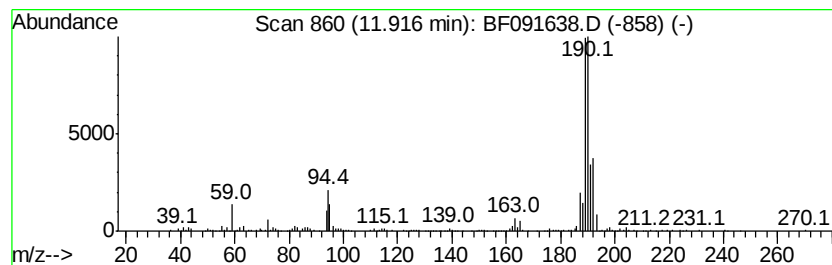
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.92	2.76 ng	456651	Phenanthrene-d10	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	95
2		6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	80
3		Thiophene-3-carboxaldehyde, 5-chloro-2-dihydroxyboryl	190	C5H4BClO3S	036155-87-0	49
4		2,3,5,6-Tetramethylterephthalaldehyde	190	C12H14O2	007072-01-7	47
5		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	42



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121516\
Data File : BF091638.D
Acq On : 15 Dec 2016 19:47
Operator : UM/SJ
Sample : H6057-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP20-COMP6

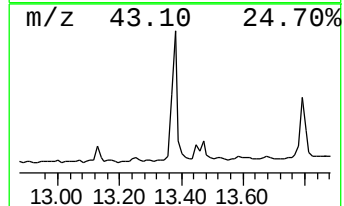
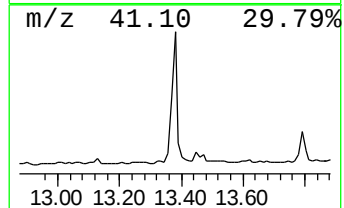
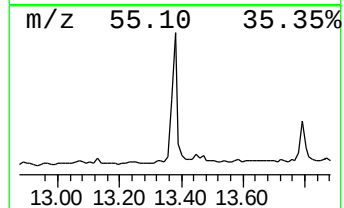
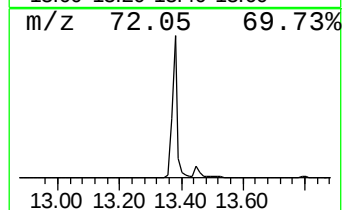
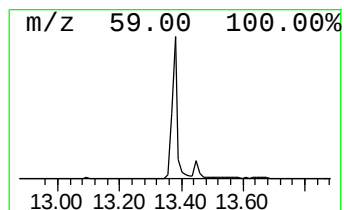
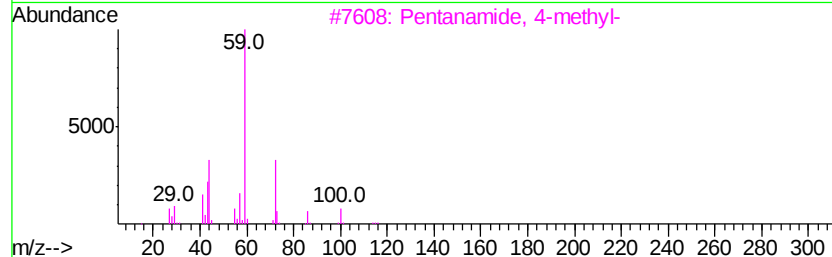
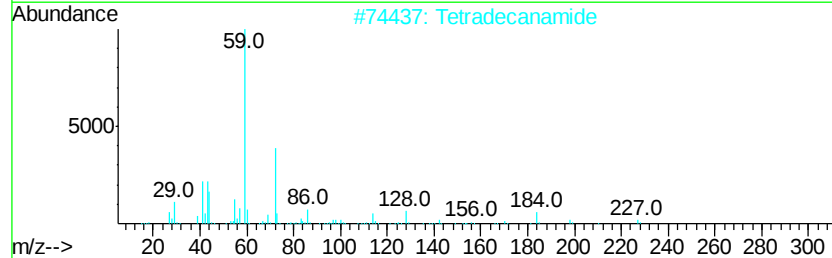
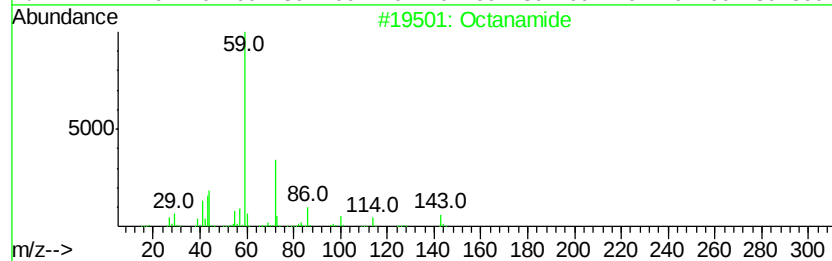
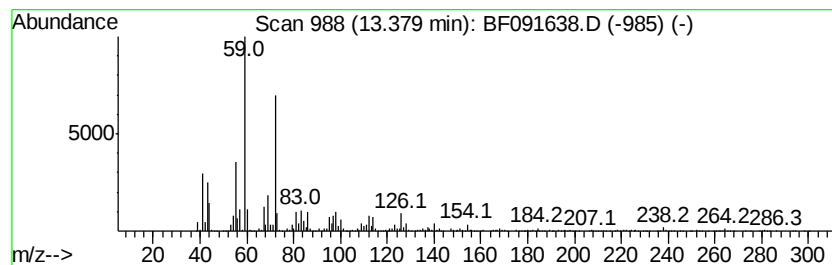
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octanamide Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.38	13.02 ng	1828420	Chrysene-d12	13.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octanamide	143	C8H17NO	000629-01-6	59
2		Tetradecanamide	227	C14H29NO	000638-58-4	50
3		Pentanamide, 4-methyl-	115	C6H13NO	001119-29-5	50
4		Hexadecanamide	255	C16H33NO	000629-54-9	45
5		Dodecanamide	199	C12H25NO	001120-16-7	45



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121516\
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Acq On : 15 Dec 2016 19:47
Operator : UM/SJ
Sample : H6057-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

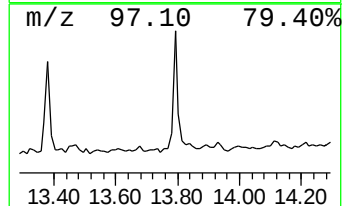
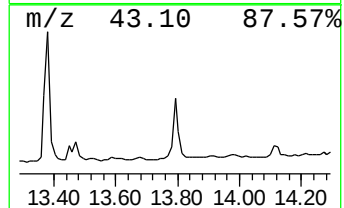
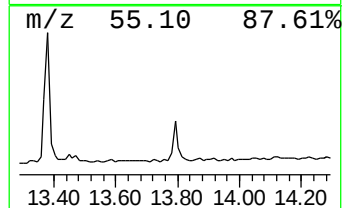
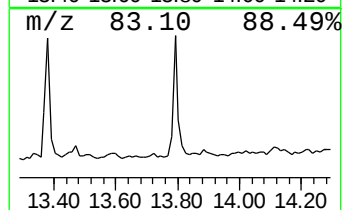
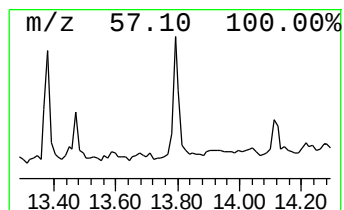
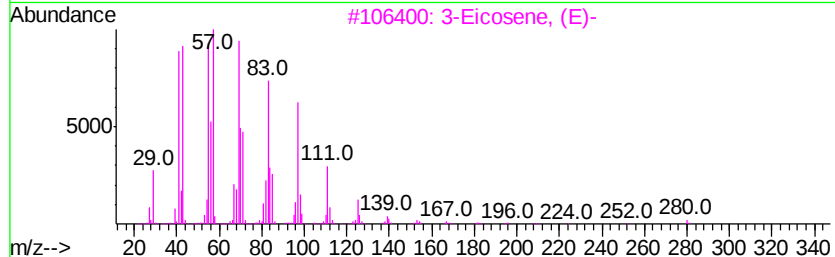
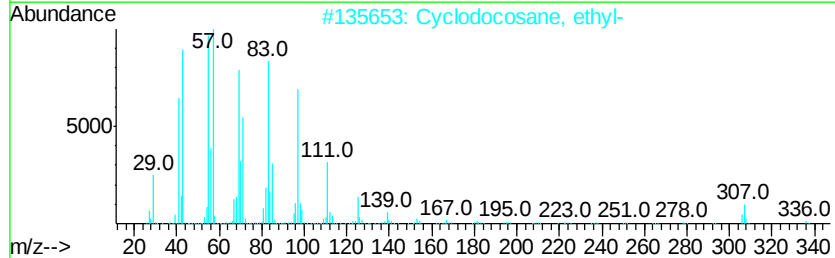
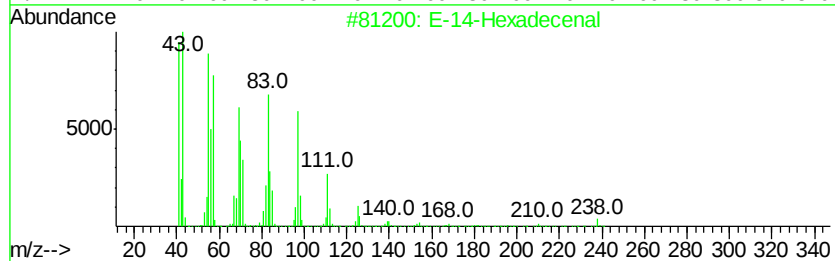
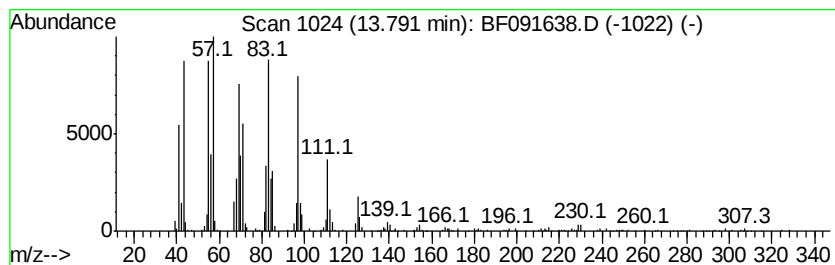
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 E-14-Hexadecenal Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.79	3.17 ng	445422	Chrysene-d12	13.94
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	E-14-Hexadecenal	238 C16H30O	330207-53-9	93
2	Cyclodocosane, ethyl-	336 C24H48	1000151-22-6	91
3	3-Eicosene, (E)-	280 C20H40	074685-33-9	91
4	Dichloroacetic acid, heptadecyl ...	366 C19H36Cl2O2	1000282-98-2	91
5	1-Nonadecene	266 C19H38	018435-45-5	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121516\
Data File : BF091638.D
Acq On : 15 Dec 2016 19:47
Operator : UM/SJ
Sample : H6057-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

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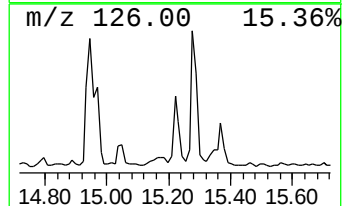
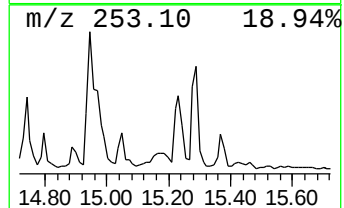
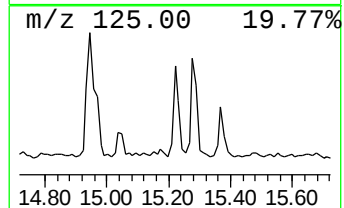
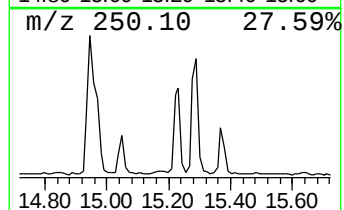
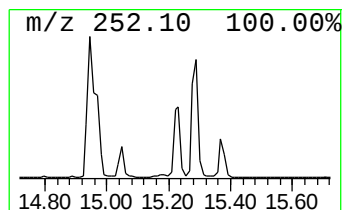
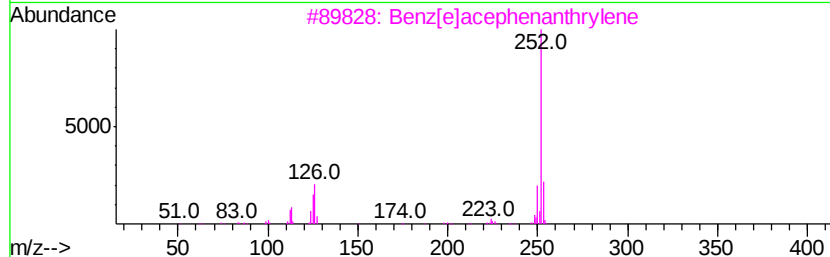
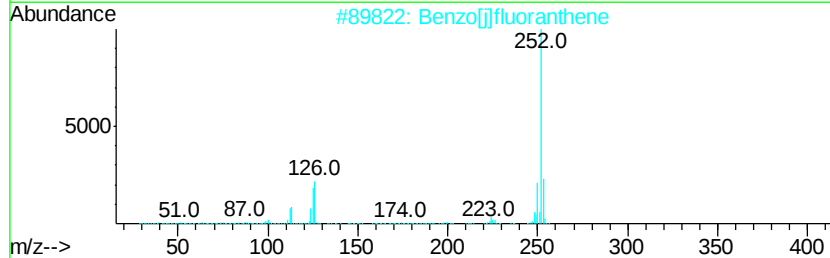
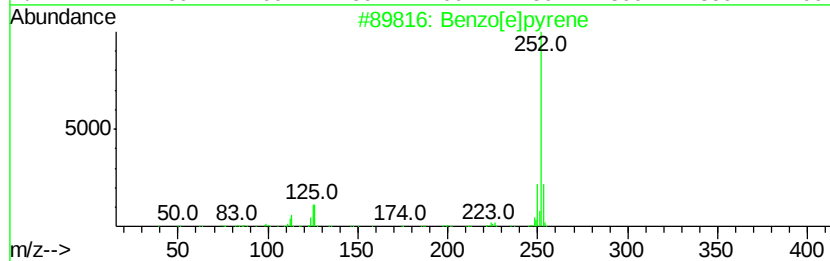
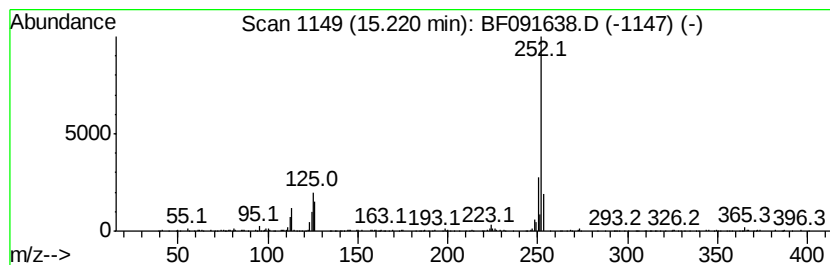
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzo[e]pyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.22	4.07 ng	380103	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2		Benzo[il]fluoranthene	252	C20H12	000205-82-3	97
3		Benz[el]acephenanthrvlene	252	C20H12	000205-99-2	97
4		Perylene	252	C20H12	000198-55-0	96
5		Benzo[a]pyrene	252	C20H12	000050-32-8	95



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121516\
Data File : BF091638.D
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Operator : UM/SJ
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Misc :
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Instrument :
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ClientSampleId :
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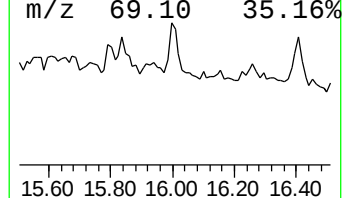
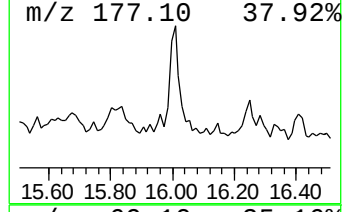
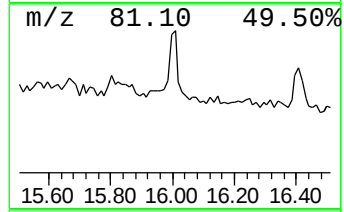
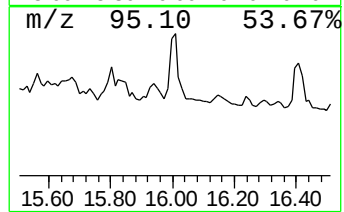
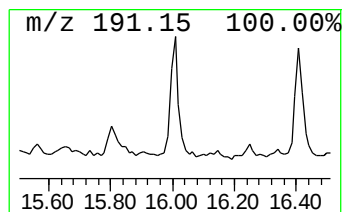
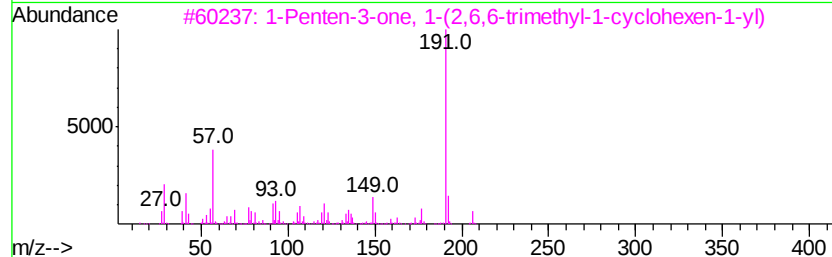
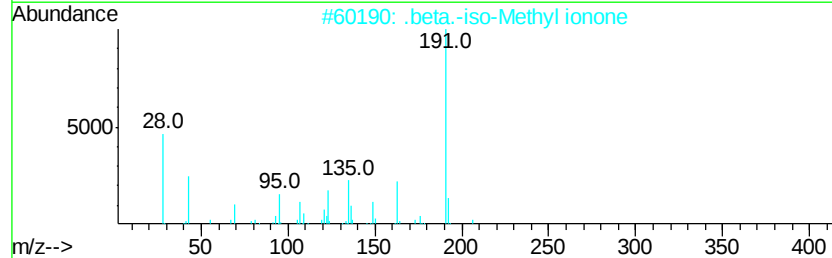
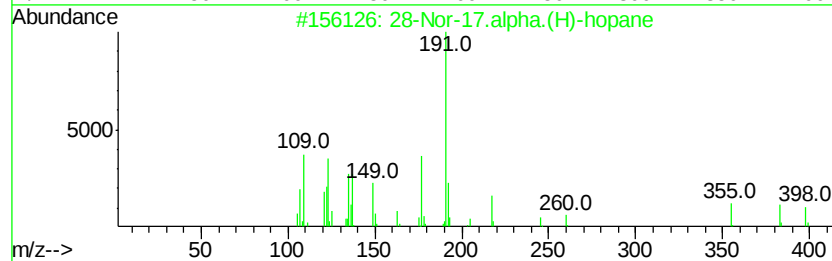
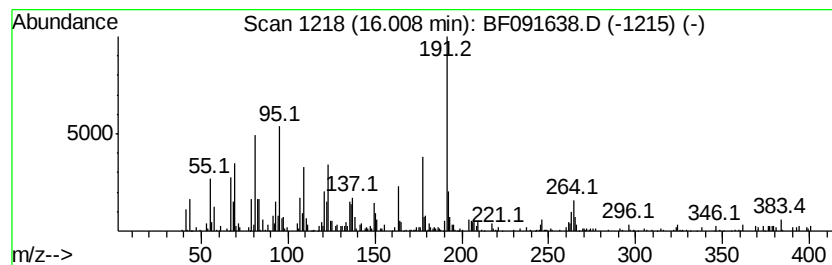
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 28-Nor-17.alpha.(H)-hopane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.01	3.80 ng	355312	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	28-Nor-17.alpha.(H)-hopane	398	C29H50	053584-60-4	72
2		.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	68
3		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	38
4		Pyrido[3,4-d]pyridazin-1(2H)-one...	191	C8H9N5O	1000263-69-7	38
5		5-(p-Aminophenyl)-2-thiazolamine	191	C9H9N3S	090349-87-4	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121516\
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Acq On : 15 Dec 2016 19:47
Operator : UM/SJ
Sample : H6057-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Penten-2-one, 4...	4.28	2.1 ng		213272	1	6.78	1983960	20.0
2-Pentanone, 4-hy...	4.97	51.1 ng		5072740	1	6.78	1983960	20.0
unknown6.56	6.56	87.8 ng		8708370	1	6.78	1983960	20.0
Eicosane	10.74	2.1 ng		346281	4	11.30	3311740	20.0
Anthracene, 2-met...	11.84	2.2 ng		367735	4	11.30	3311740	20.0
4H-Cyclopenta[def...	11.92	2.8 ng		456651	4	11.30	3311740	20.0
Octanamide	13.38	13.0 ng		1828420	5	13.94	2808000	20.0
E-14-Hexadecenal	13.79	3.2 ng		445422	5	13.94	2808000	20.0
Benzo[e]pyrene	15.22	4.1 ng		380103	6	15.35	1869690	20.0
28-Nor-17.alpha.(...	16.01	3.8 ng		355312	6	15.35	1869690	20.0