

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122015\
 Data File : BF083985.D
 Acq On : 20 Dec 2015 7:45
 Operator : UM/SJ
 Sample : G4851-04
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 UST-9

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-BF121815.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	2.773	50	60	64	rVB	140507	364215	1.86%	0.220%
2	3.767	139	147	150	rBV3	359910	1093095	5.59%	0.659%
3	3.904	155	159	162	rVV	215864	427389	2.19%	0.258%
4	4.064	168	173	175	rBV	371426	685479	3.51%	0.413%
5	4.110	175	177	183	rVB	158497	251371	1.29%	0.152%
6	4.224	183	187	189	rBV	143573	275051	1.41%	0.166%
7	4.281	189	192	196	rVB	190281	366534	1.87%	0.221%
8	4.430	201	205	210	rVV	963318	1789500	9.15%	1.079%
9	4.533	210	214	217	rVV	398062	574980	2.94%	0.347%
10	4.761	231	234	235	rBV	139555	222251	1.14%	0.134%
11	4.876	239	244	246	rBV	441465	790201	4.04%	0.476%
12	4.921	246	248	250	rBV3	220643	532576	2.72%	0.321%
13	4.967	250	252	255	rVB	1121153	1622937	8.30%	0.979%
14	5.024	255	257	260	rBV2	725878	1084712	5.55%	0.654%
15	5.093	262	263	266	rVB2	269314	299756	1.53%	0.181%
16	5.161	266	269	273	rVB	238864	352860	1.80%	0.213%
17	5.242	273	276	279	rBV	997678	1923084	9.84%	1.160%
18	5.310	279	282	283	rBV	1177073	1365257	6.98%	0.823%
19	5.344	283	285	286	rBV	344236	586423	3.00%	0.354%
20	5.367	286	287	290	rVB	773242	1166137	5.96%	0.703%
21	5.447	290	294	300	rVB2	3421944	8792051	44.97%	5.302%
22	5.539	300	302	304	rBV	547305	702515	3.59%	0.424%
23	5.584	304	306	307	rVB	164015	211254	1.08%	0.127%
24	5.619	307	309	311	rBV	1042274	1682588	8.61%	1.015%
25	5.699	311	316	320	rVB2	2552107	5157514	26.38%	3.110%
26	5.779	320	323	329	rVB	2743221	4561334	23.33%	2.750%
27	5.882	329	332	334	rBV	1296121	1522015	7.78%	0.918%
28	5.927	334	336	340	rBV3	329213	824874	4.22%	0.497%
29	6.019	341	344	346	rBV2	1379139	2306312	11.80%	1.391%
30	6.076	346	349	350	rBV	378092	625093	3.20%	0.377%
31	6.122	350	353	356	rBV2	1650338	3888872	19.89%	2.345%
32	6.167	356	357	366	rVB5	1025892	3891995	19.91%	2.347%
33	6.316	366	370	373	rBV2	1647415	4279648	21.89%	2.581%
34	6.407	373	378	379	rBV	3145351	8538049	43.67%	5.148%

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 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

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35	6.476	382	384	390	rVV4	2358904	5927439	30.32%	3.574%
36	6.567	390	392	394	rVB	1314285	1915414	9.80%	1.155%
37	6.716	401	405	415	rBV2	3950176	19552716	100.00%	11.790%
38	6.910	420	422	423	rBV	1198382	1828816	9.35%	1.103%
39	6.945	423	425	430	rVB2	1269831	3571542	18.27%	2.154%
40	7.059	430	435	440	rVB3	1320112	3838692	19.63%	2.315%
41	7.162	442	444	445	rBV	1285232	2120488	10.84%	1.279%
42	7.207	445	448	452	rBV2	1321224	3950179	20.20%	2.382%
43	7.459	464	470	471	rBV2	1385904	5096816	26.07%	3.073%
44	7.722	489	493	499	rVB3	1409106	5743399	29.37%	3.463%
45	7.859	499	505	507	rBV4	1823973	7307809	37.37%	4.407%
46	8.225	530	537	539	rBV2	2554281	11574747	59.20%	6.979%
47	12.145	877	880	881	rVB	2175854	3593400	18.38%	2.167%
48	12.396	900	902	906	rBV2	740451	1309219	6.70%	0.789%
49	12.511	908	912	915	rVV2	2667349	6880052	35.19%	4.149%
50	12.591	917	919	922	rVB4	738956	1031600	5.28%	0.622%
51	12.877	940	944	945	rVB2	2596546	4074328	20.84%	2.457%
52	12.911	945	947	949	rVV3	458240	800834	4.10%	0.483%
53	12.979	951	953	959	rVB2	1200108	2383293	12.19%	1.437%
54	13.219	971	974	975	rVB	2419582	3020390	15.45%	1.821%
55	13.254	975	977	978	rVB	196506	267287	1.37%	0.161%
56	13.551	1001	1003	1005	rVB	2125351	2330213	11.92%	1.405%
57	13.871	1026	1031	1036	rVB	1394005	2508107	12.83%	1.512%
58	13.951	1036	1038	1042	rVV3	325967	598124	3.06%	0.361%
59	14.020	1042	1044	1045	rVB	332482	297492	1.52%	0.179%
60	14.180	1055	1058	1059	rBV	613981	815397	4.17%	0.492%
61	14.477	1082	1084	1086	rVB	413494	372565	1.91%	0.225%
62	15.437	1166	1168	1173	rVB	273153	372167	1.90%	0.224%

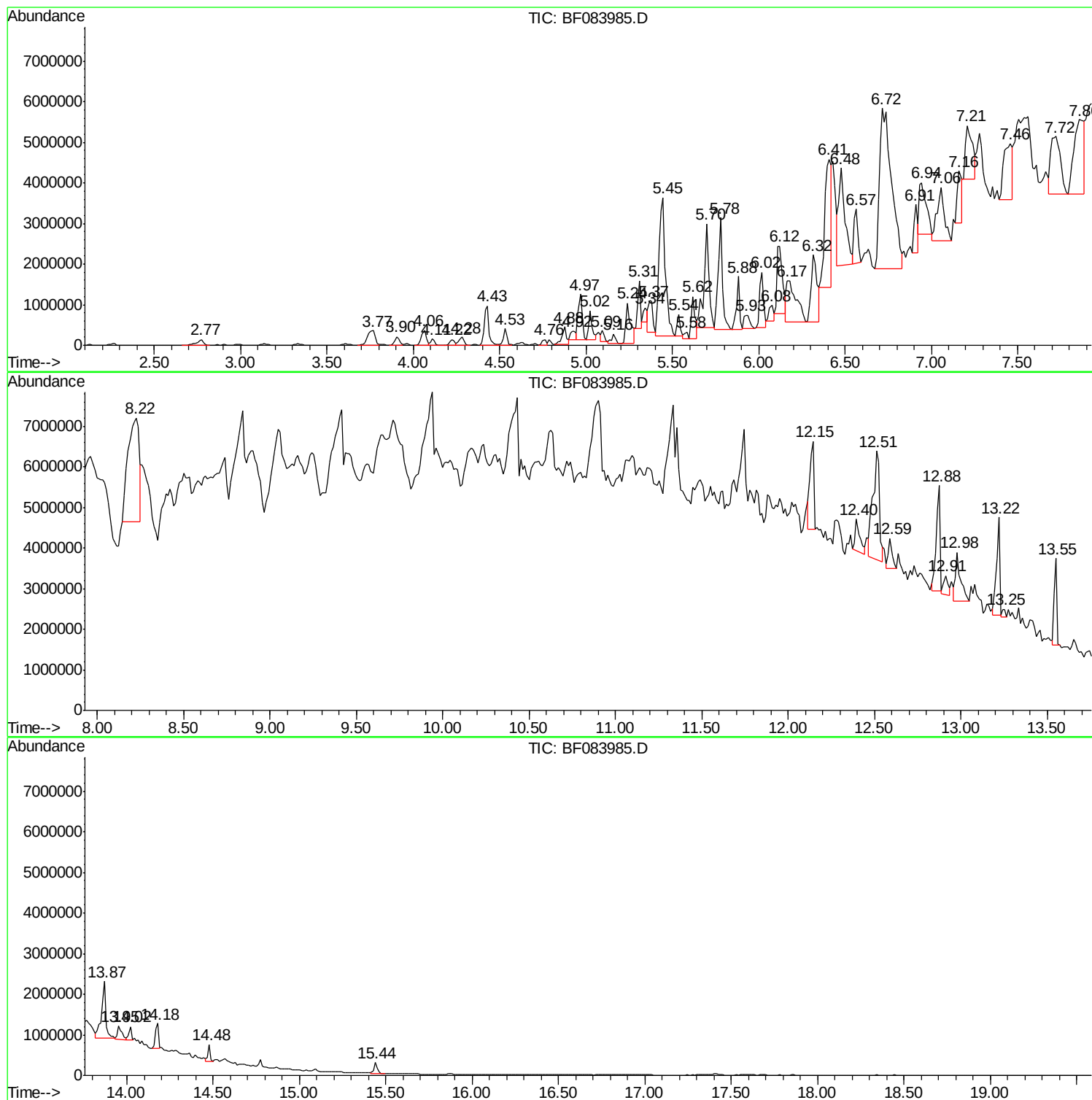
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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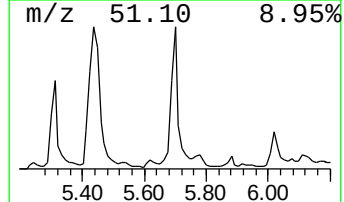
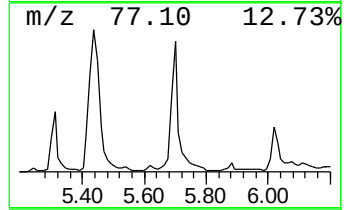
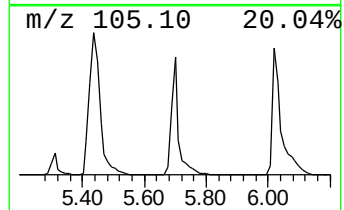
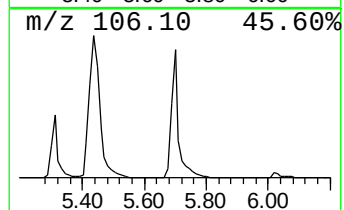
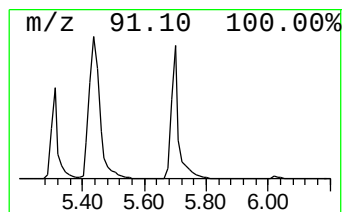
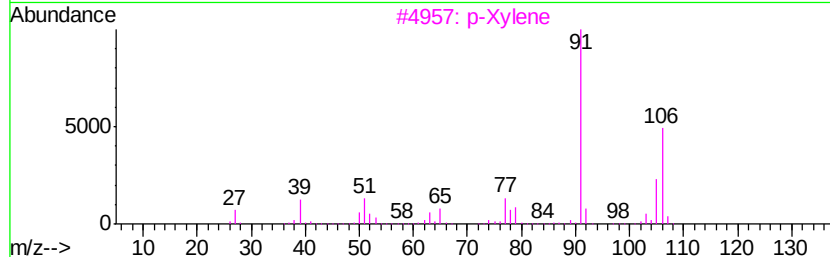
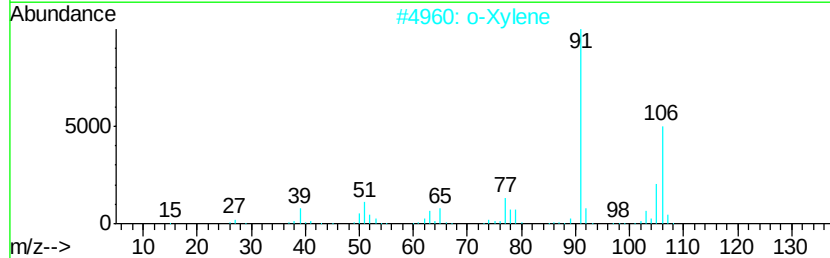
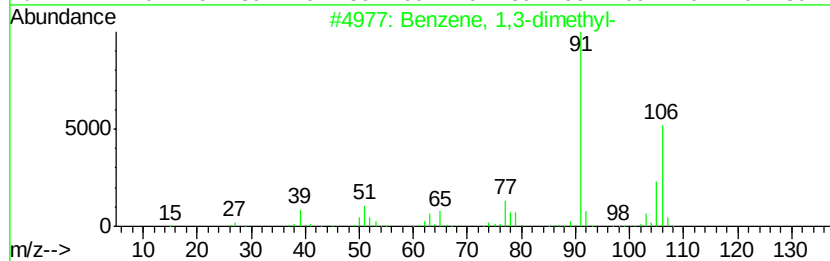
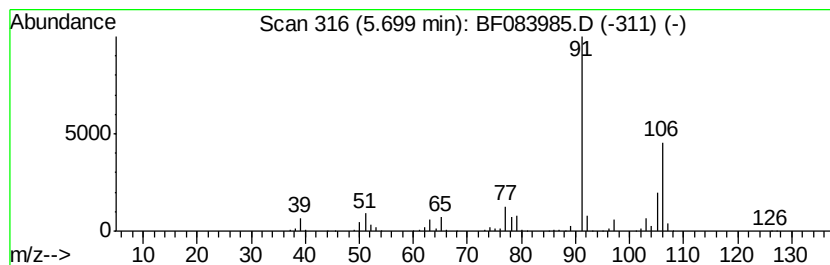
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Benzene, 1,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.70	56.40 ng	5157510	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2		o-Xylene	106	C8H10	000095-47-6	97
3		p-Xylene	106	C8H10	000106-42-3	97
4		1,3-Cyclopentadiene, 5-(1-methyl...	106	C8H10	002175-91-9	81
5		Ethylbenzene	106	C8H10	000100-41-4	81



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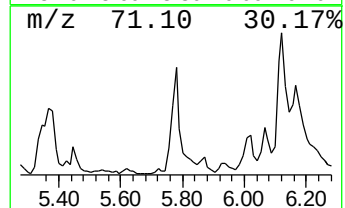
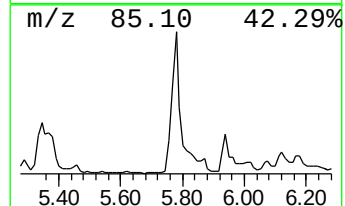
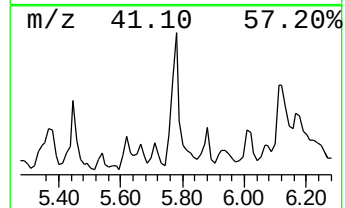
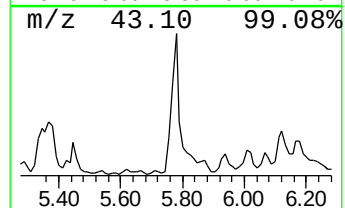
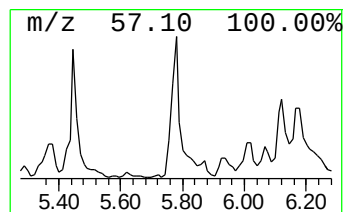
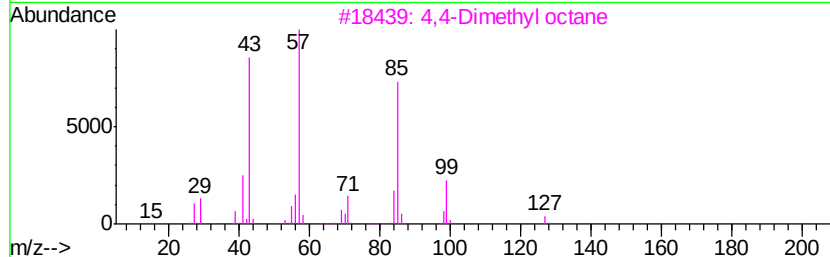
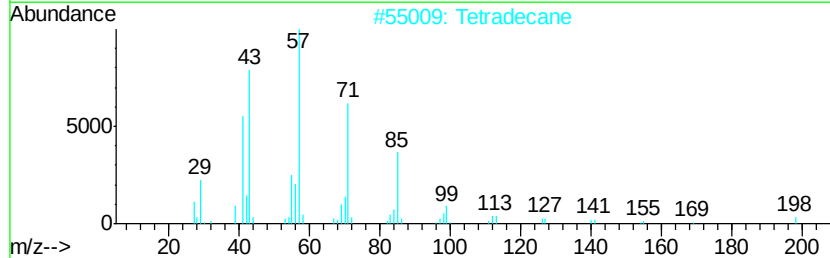
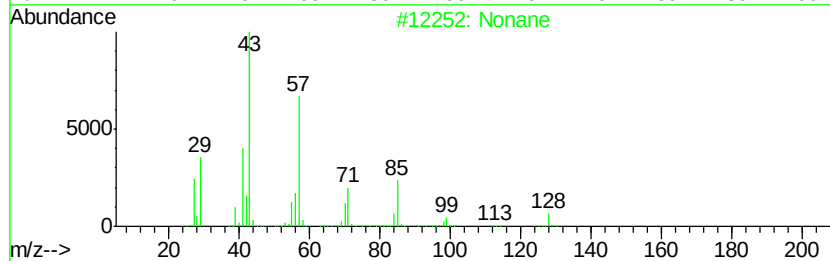
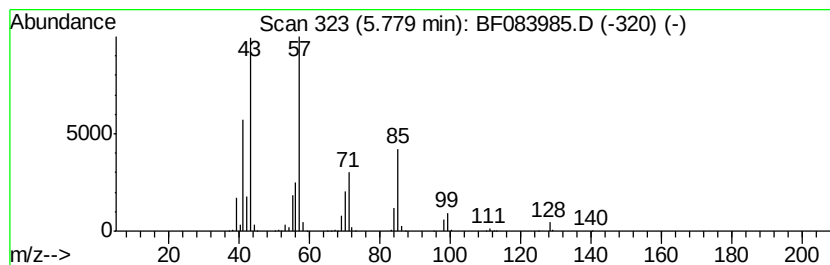
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121815.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Nonane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.78	49.88 ng	4561330	1,4-Dichlorobenzene-d4	6.88

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane		128	C9H20	000111-84-2	94
2	Tetradecane		198	C14H30	000629-59-4	72
3	4,4-Dimethyl octane		142	C10H22	015869-95-1	59
4	Octane, 4-methyl-		128	C9H20	002216-34-4	58
5	Heptane, 2,4-dimethyl-		128	C9H20	002213-23-2	53



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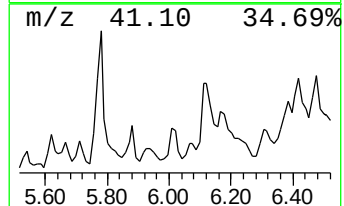
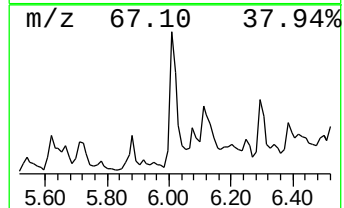
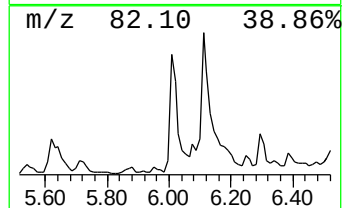
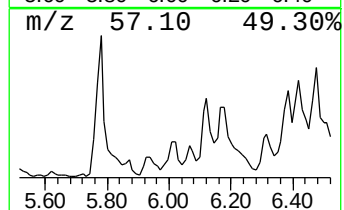
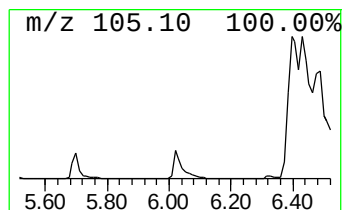
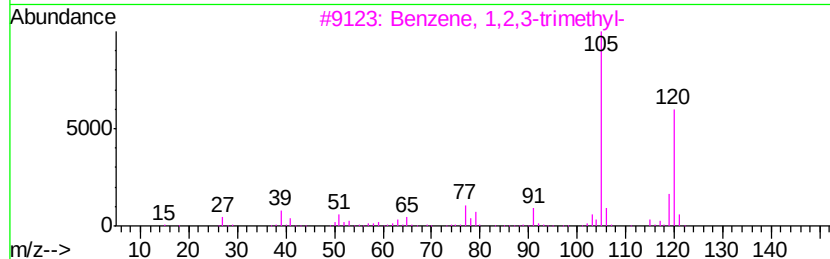
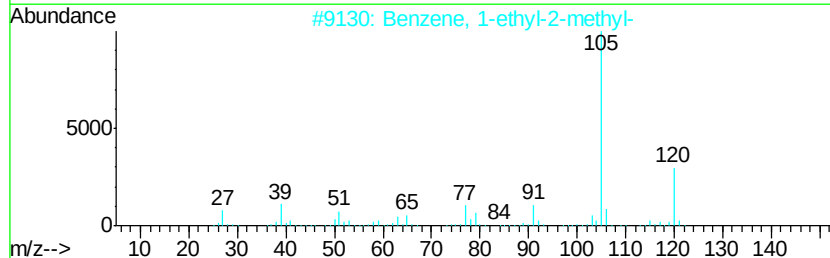
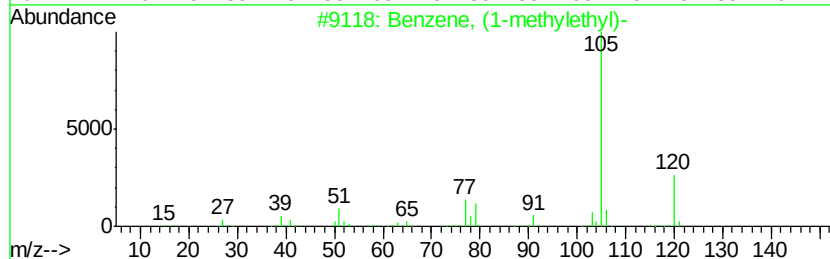
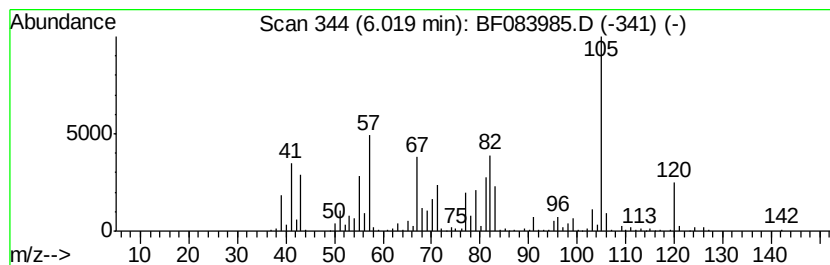
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Benzene, (1-methylethyl)- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.02	25.22 ng	2306310	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	42
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	27
3		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	27
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	27
5		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	27



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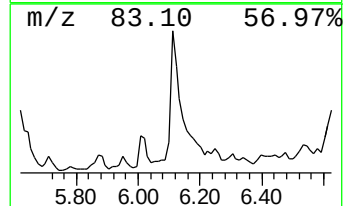
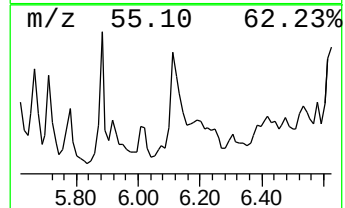
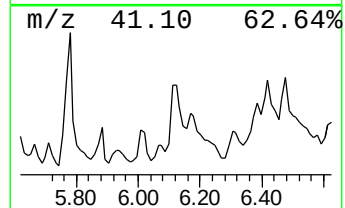
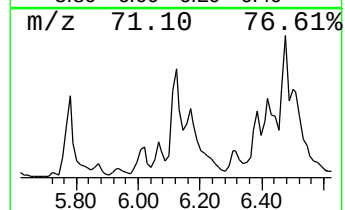
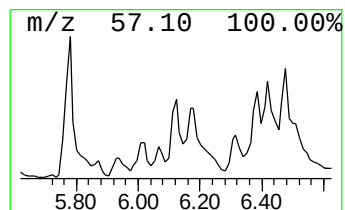
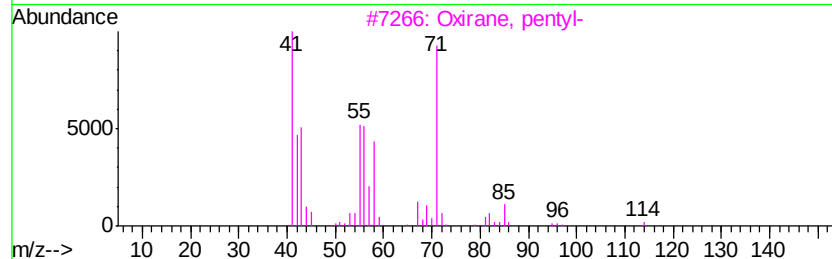
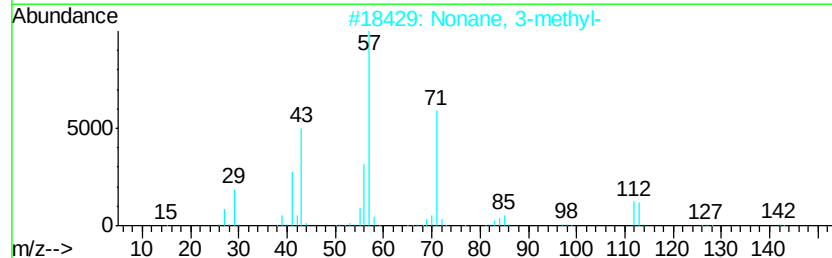
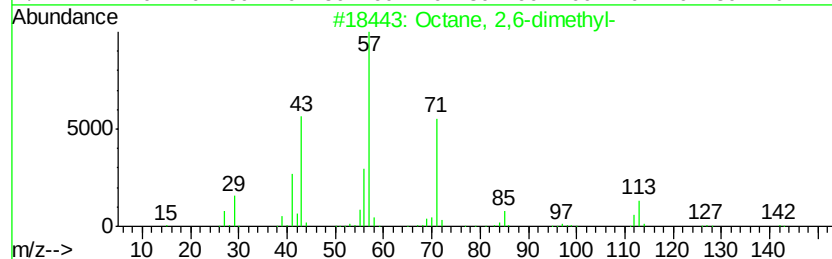
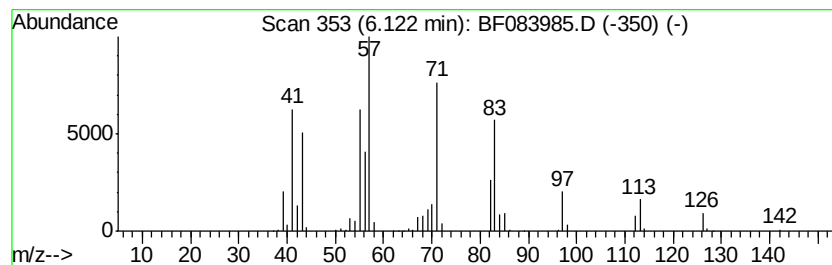
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Octane, 2,6-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.12	42.53 ng	3888870	1,4-Dichlorobenzene-d4	6.88

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	58
2		Nonane, 3-methyl-	142	C10H22	005911-04-6	58
3		Oxirane, pentyl-	114	C7H14O	005063-65-0	45
4		Cyclopentane, (2-methylpropyl)-	126	C9H18	003788-32-7	42
5		3-Bromooctane	192	C8H17Br	000999-64-4	38



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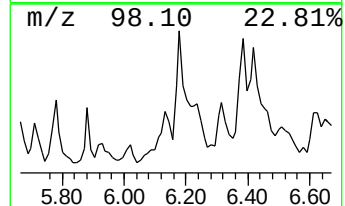
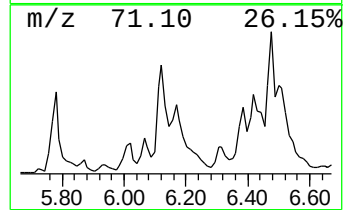
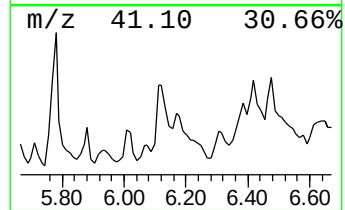
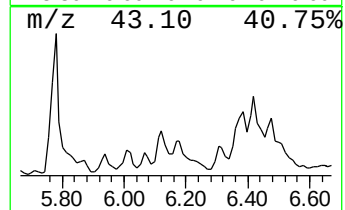
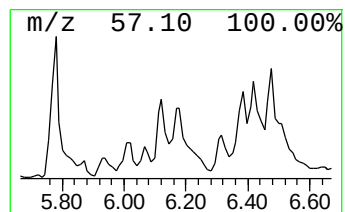
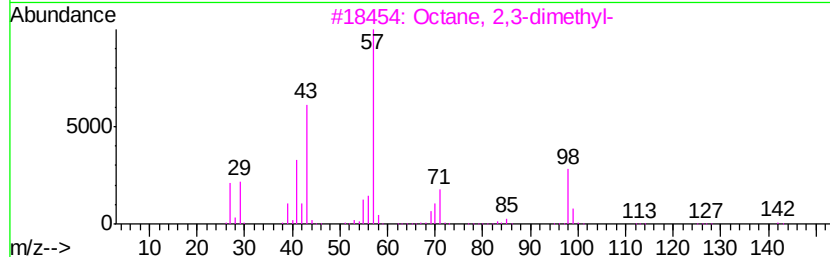
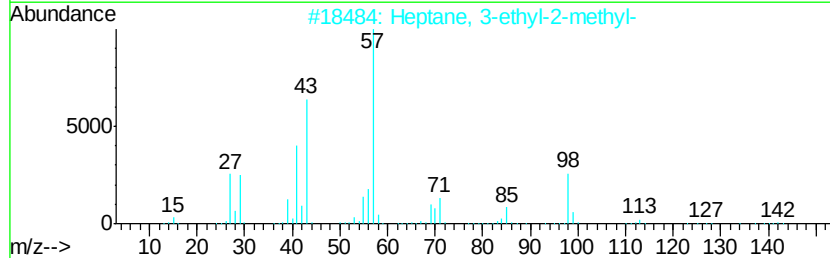
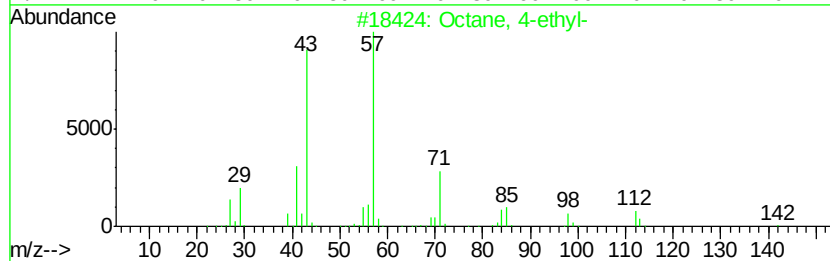
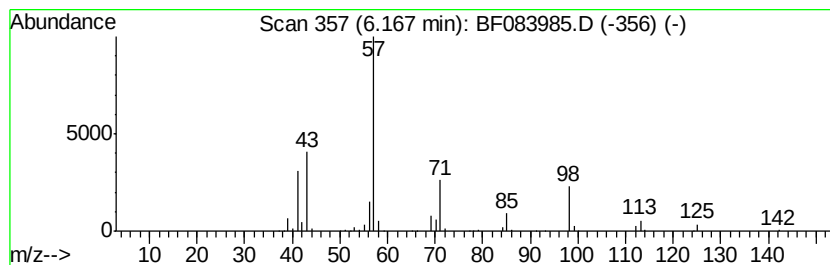
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Octane, 4-ethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.17	42.56 ng	3892000	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 4-ethyl-	142	C10H22	015869-86-0	64
2		Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	64
3		Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	59
4		Decane, 2,5-dimethyl-	170	C12H26	017312-50-4	53
5		Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122015\
Data File : BF083985.D
Acq On : 20 Dec 2015 7:45
Operator : UM/SJ
Sample : G4851-04
Misc :
ALS Vial : 36 Sample Multiplier: 1

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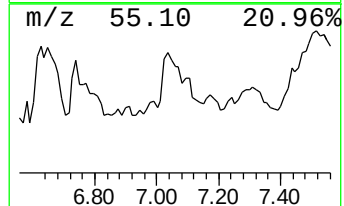
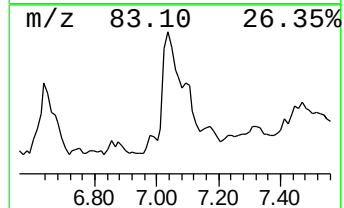
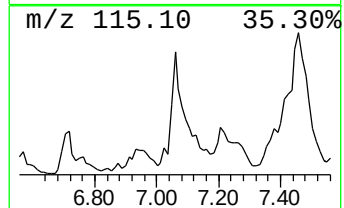
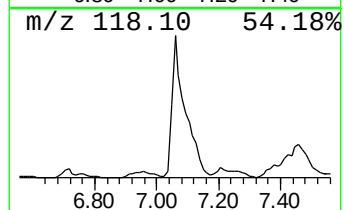
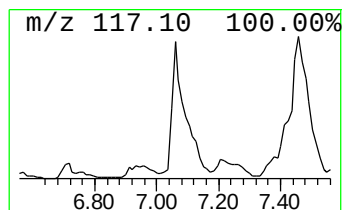
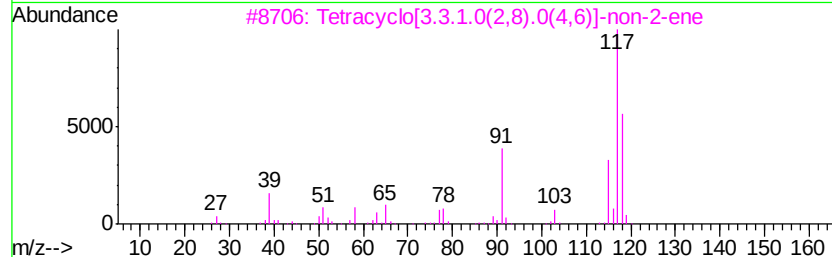
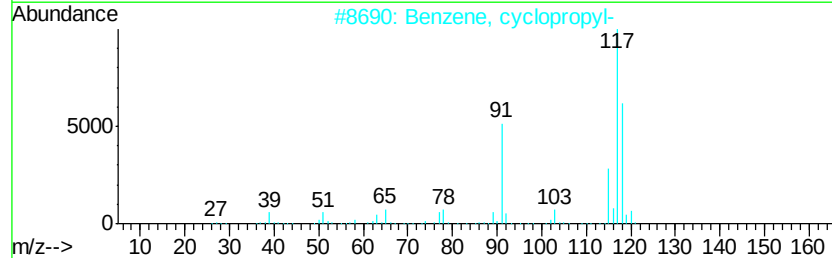
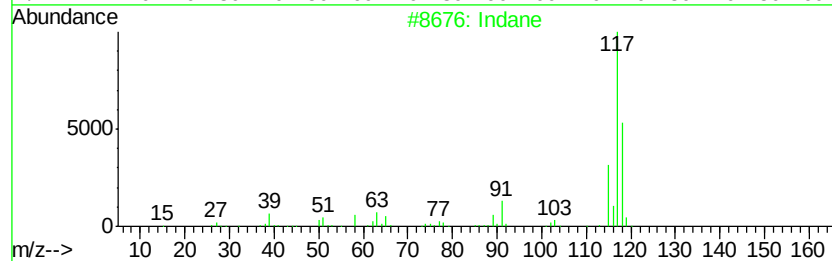
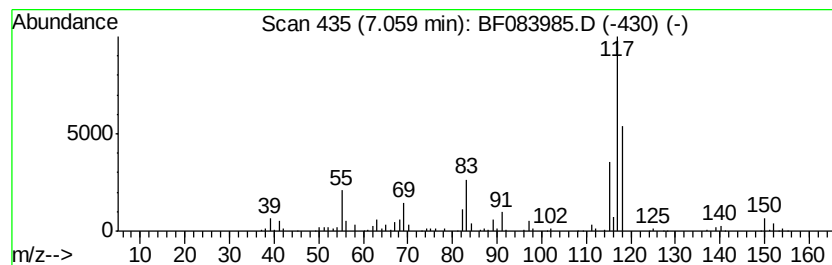
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Indane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.06	41.98 ng	3838690	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	70
2		Benzene, cyclopropyl-	118	C9H10	000873-49-4	58
3		Tetracyclo[3.3.1.0(2,8).0(4,6)]-	118	C9H10	1000191-13-7	53
4		Deltacyclene	118	C9H10	007785-10-6	53
5		Benzene, 1-propenyl-	118	C9H10	000637-50-3	42



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122015\
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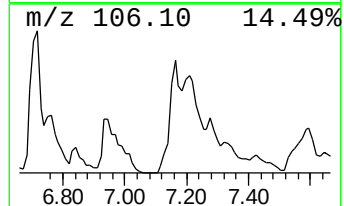
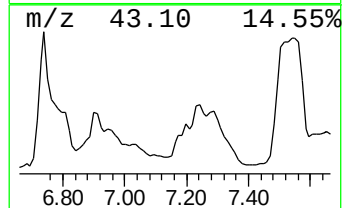
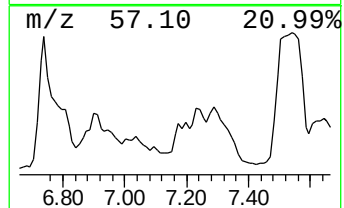
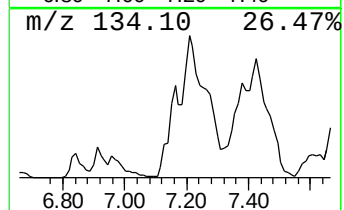
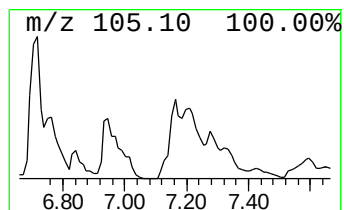
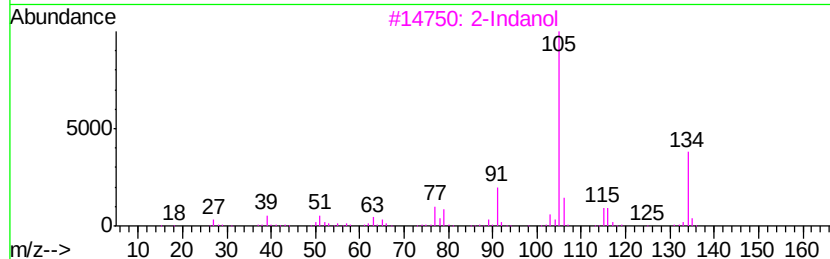
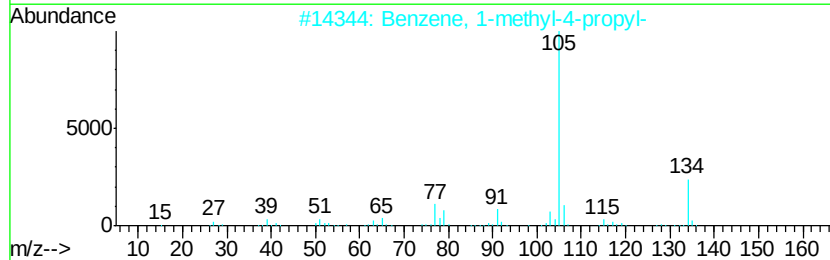
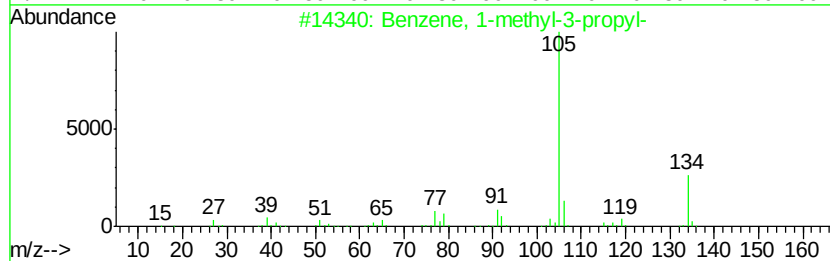
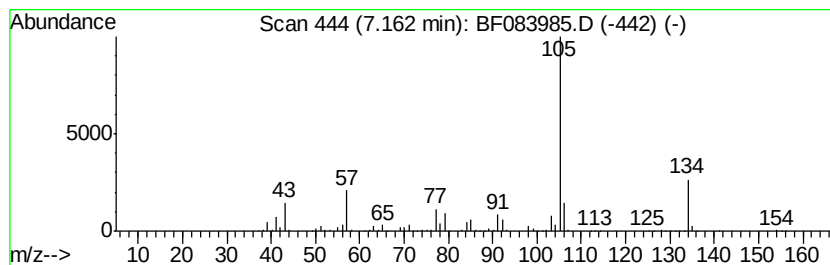
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzene, 1-methyl-3-propyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.16	23.19 ng	2120490	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	90
2		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	87
3		2-Indanol	134	C9H10O	004254-29-9	80
4		Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	80
5		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	80



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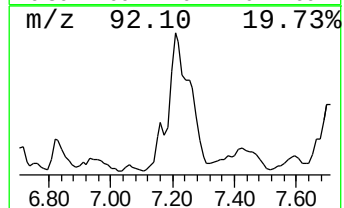
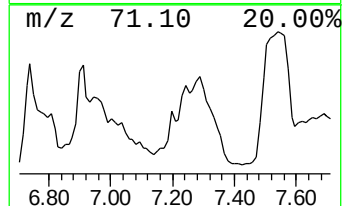
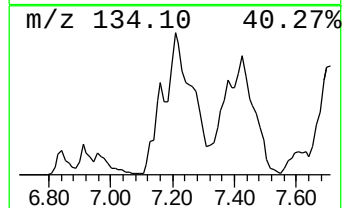
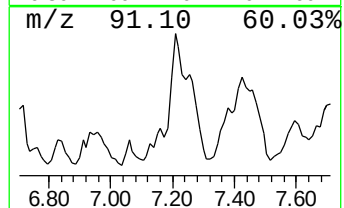
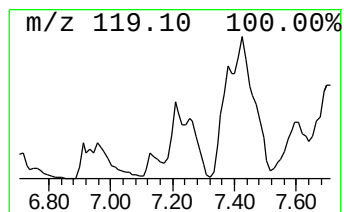
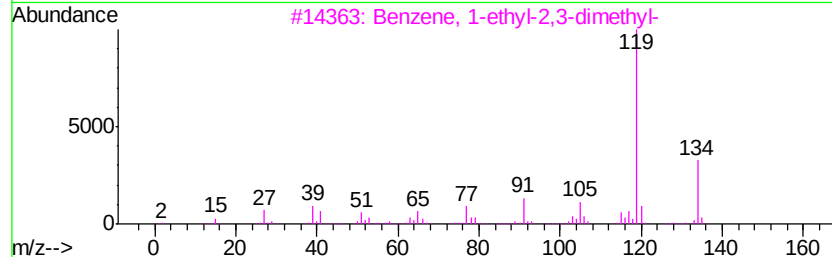
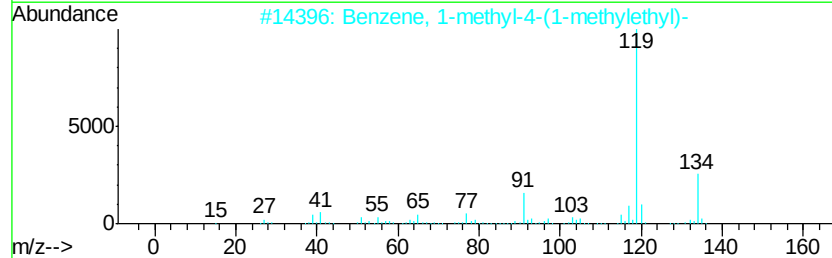
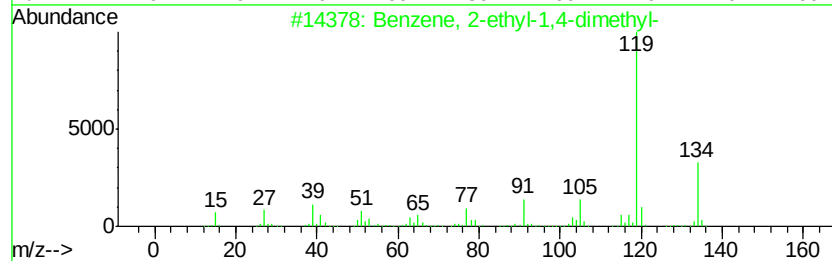
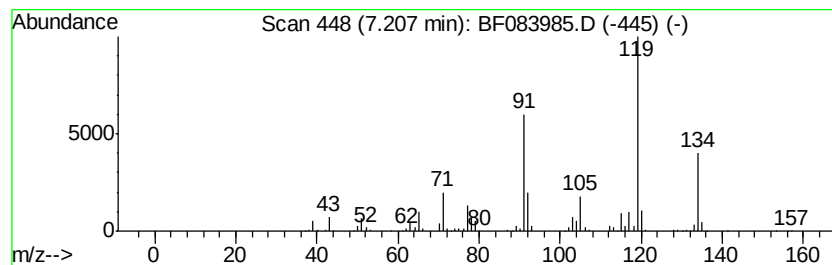
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.21	43.20 ng	3950180	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
2		Benzene, 1-methyl-4-(1-methyleth...	134	C10H14	000099-87-6	93
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	90
4		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	81
5		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	81



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122015\
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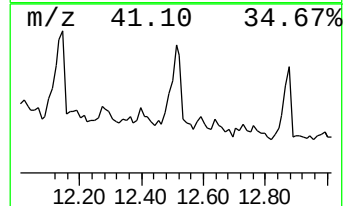
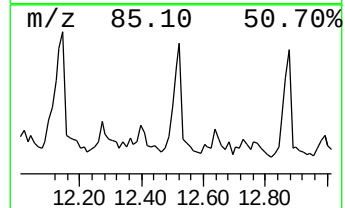
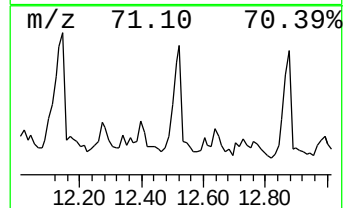
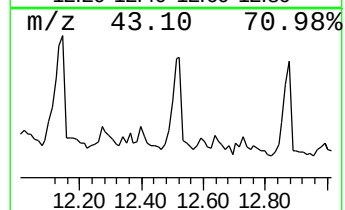
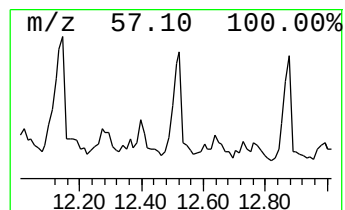
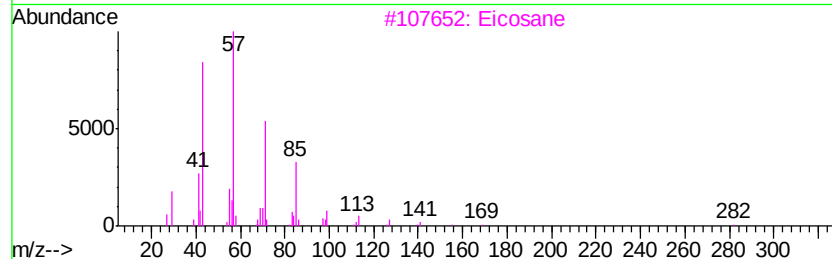
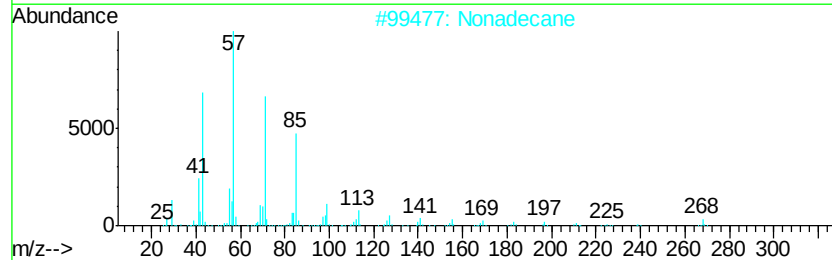
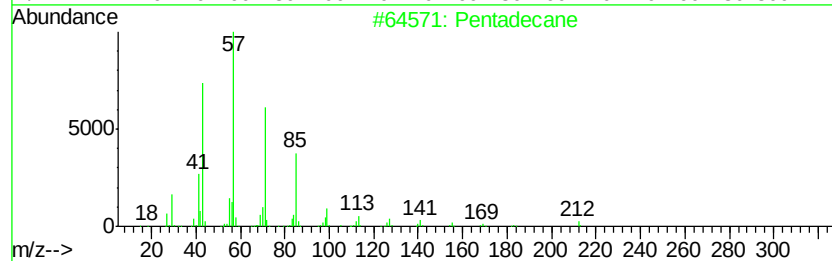
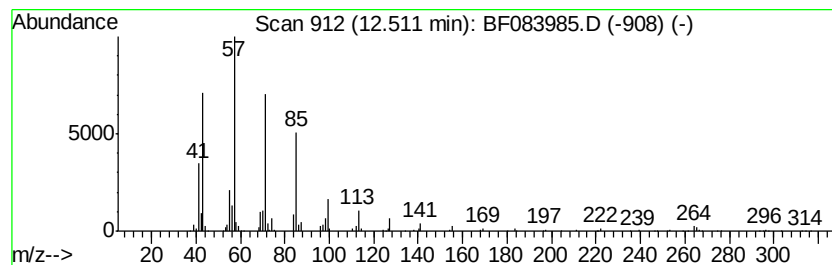
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Pentadecane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.51	38.29 ng	6880050	Phenanthrene-d10	11.44

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane		212	C15H32	000629-62-9	91
2	Nonadecane		268	C19H40	000629-92-5	91
3	Eicosane		282	C20H42	000112-95-8	90
4	Tetradecane		198	C14H30	000629-59-4	90
5	Octacosane		394	C28H58	000630-02-4	90



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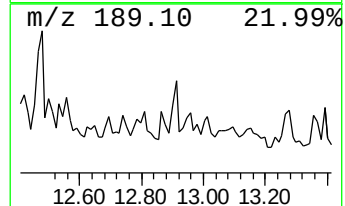
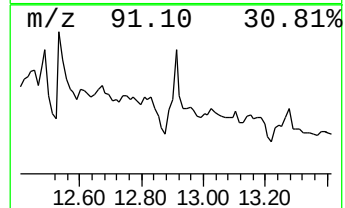
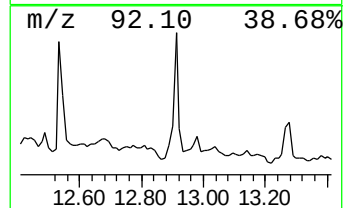
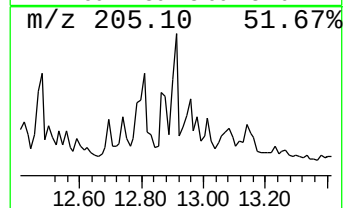
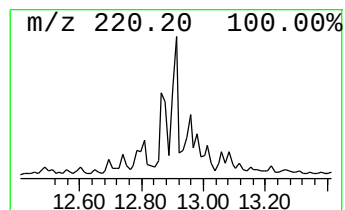
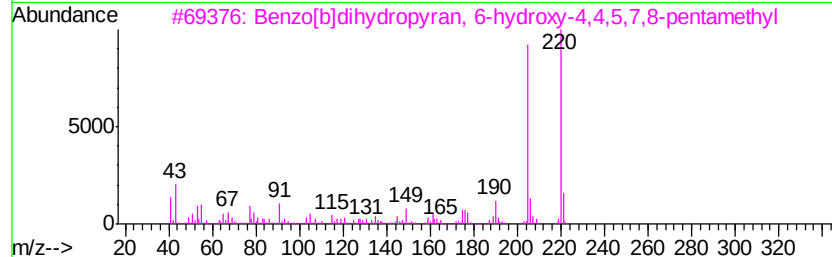
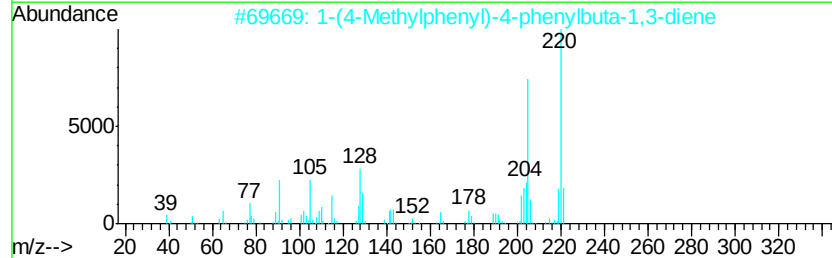
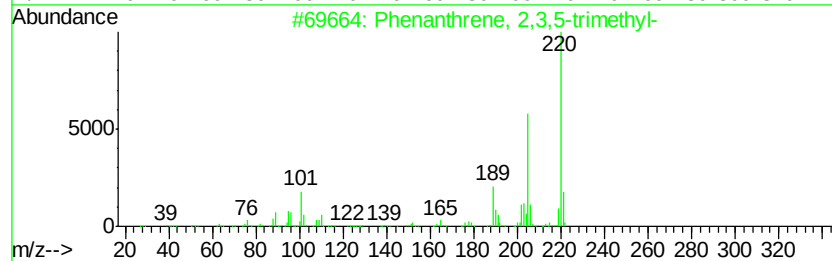
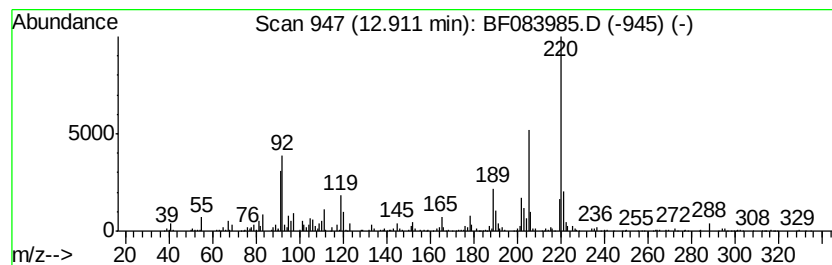
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Phenanthrene, 2,3,5-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.91	53.84 ng	800834	Chrysene-d12	14.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,3,5-trimethyl-	220	C17H16	003674-73-5	70
2		1-(4-Methylphenyl)-4-phenylbuta-...	220	C17H16	037985-11-8	62
3		Benzo[b]dihydropyran, 6-hydroxyv-...	220	C14H20O2	050442-70-1	50
4		Pyrrolo[1,2-a]thieno[2,3-d]pyrim...	220	C11H12N2OS	056929-61-4	49
5		6,7,8,9-Tetrahydro-5-methylisoth...	220	C11H12N2OS	339156-87-5	47



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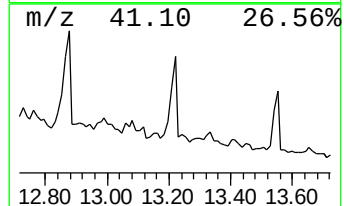
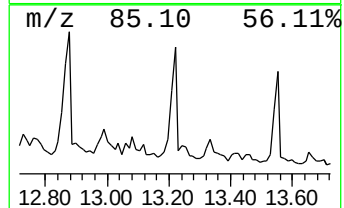
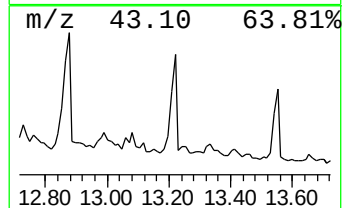
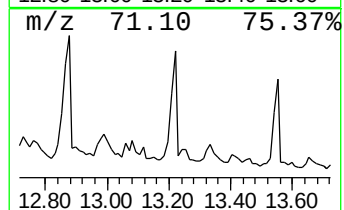
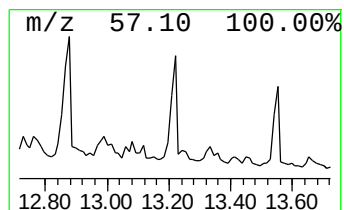
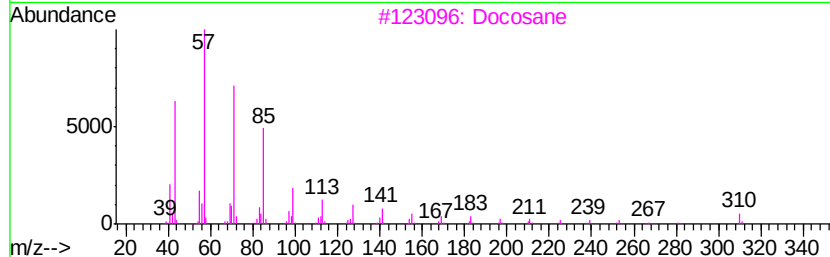
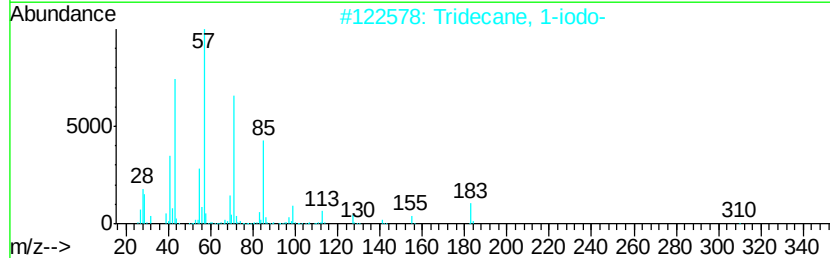
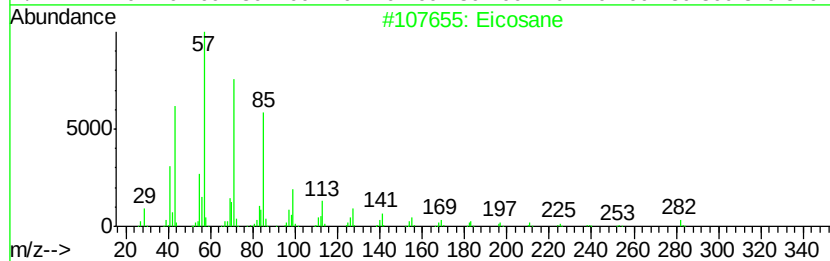
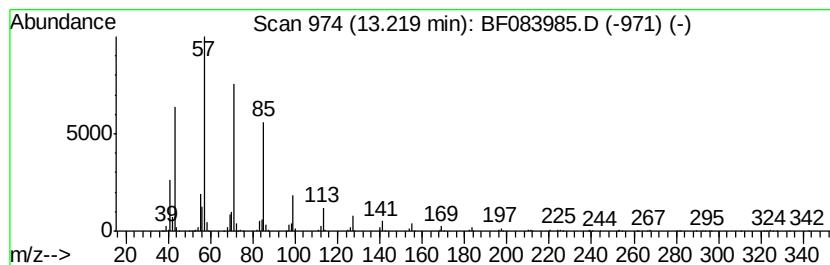
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Eicosane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	203.06 ng	3020390	Chrysene-d12	14.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	94
2		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	91
3		Docosane	310	C22H46	000629-97-0	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Tricosane	324	C23H48	000638-67-5	90



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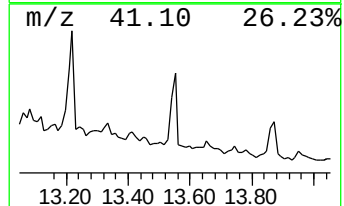
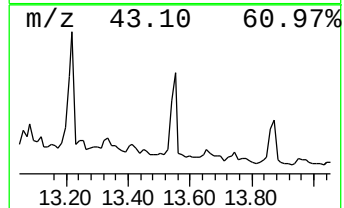
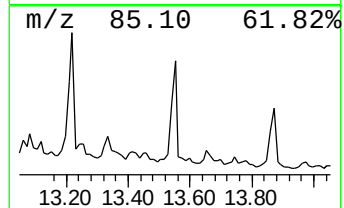
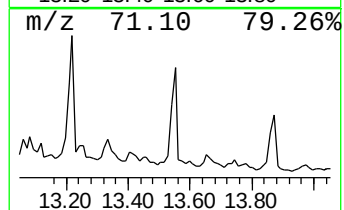
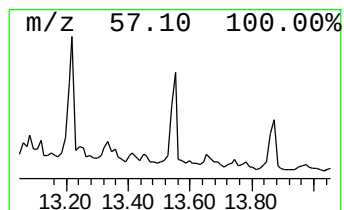
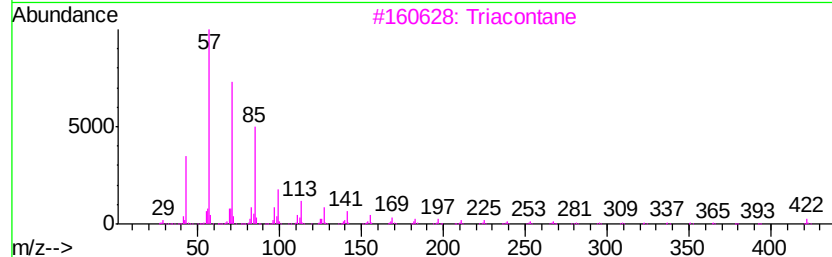
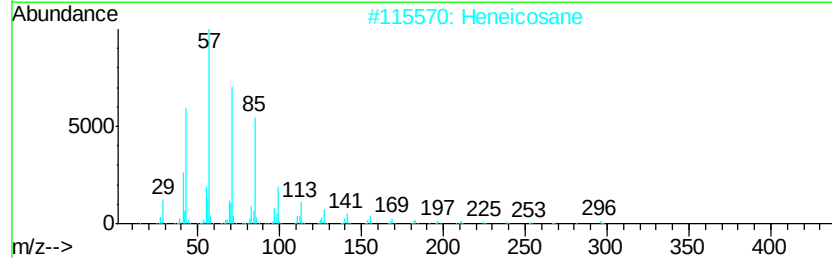
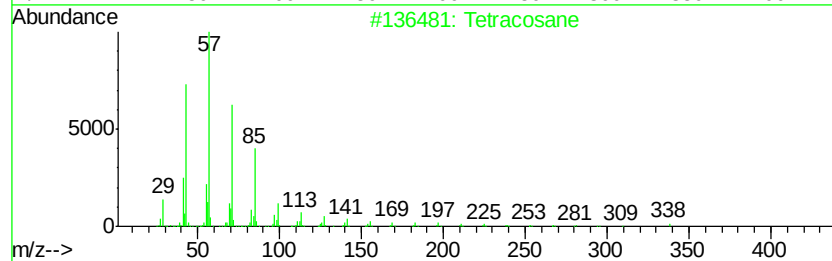
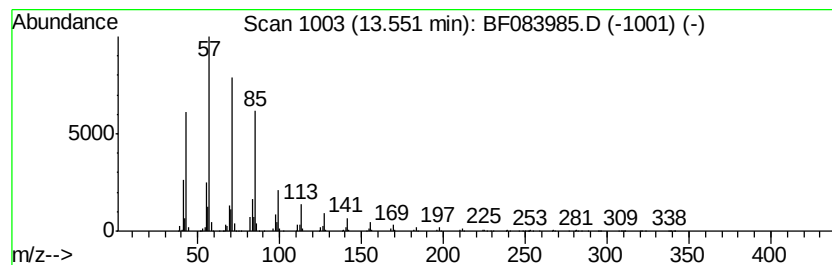
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Tetracosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.55	156.66 ng	2330210	Chrysene-d12	14.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	93
2		Heneicosane	296	C21H44	000629-94-7	91
3		Triacontane	422	C30H62	000638-68-6	90
4		Octadecane	254	C18H38	000593-45-3	87
5		Nonadecane	268	C19H40	000629-92-5	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122015\
Data File : BF083985.D
Acq On : 20 Dec 2015 7:45
Operator : UM/SJ
Sample : G4851-04
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
UST-9

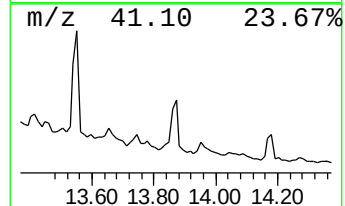
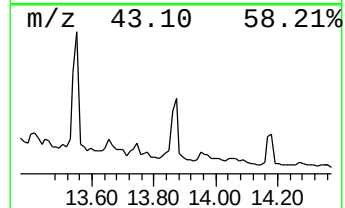
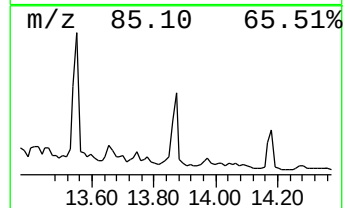
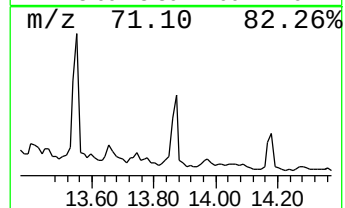
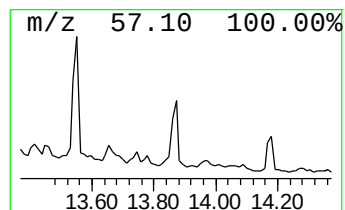
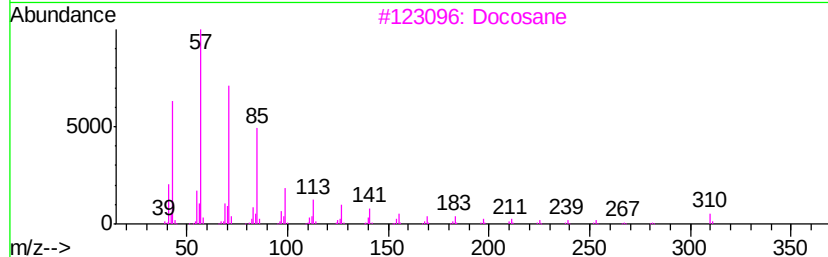
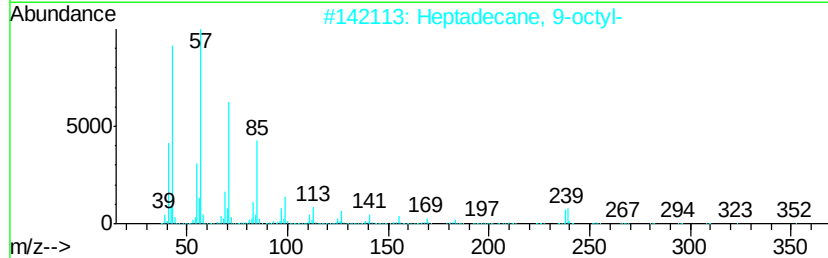
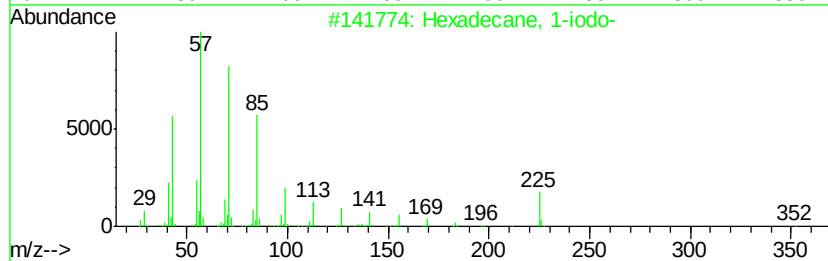
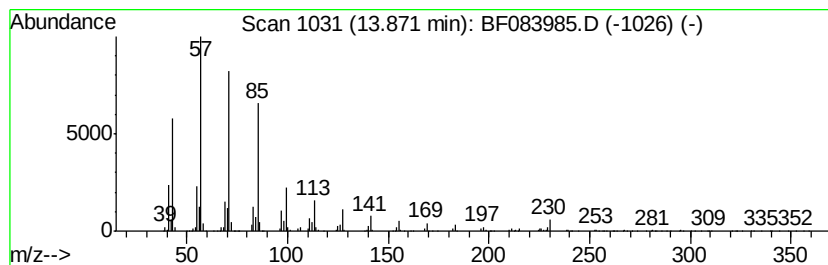
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Hexadecane, 1-iodo- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	168.62 ng	2508110	Chrysene-d12	14.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	97
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3		Docosane	310	C22H46	000629-97-0	91
4		Heneicosane	296	C21H44	000629-94-7	90
5		Nonadecane, 2-methyl-	282	C20H42	001560-86-7	90



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122015\
Data File : BF083985.D
Acq On : 20 Dec 2015 7:45
Operator : UM/SJ
Sample : G4851-04
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
UST-9

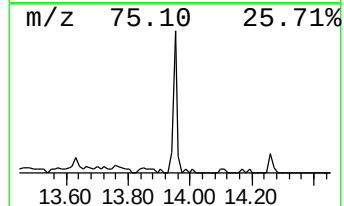
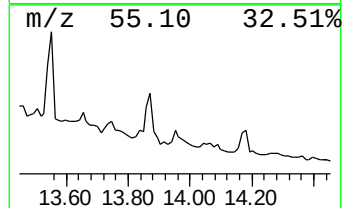
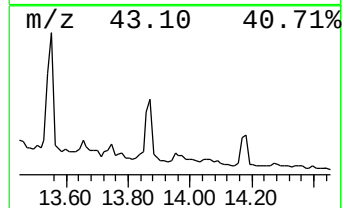
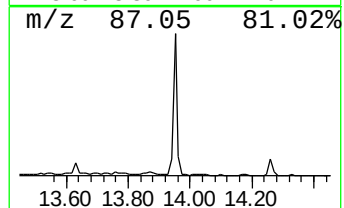
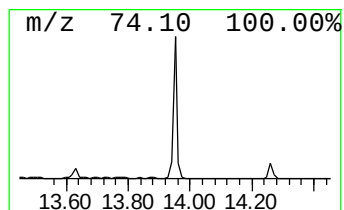
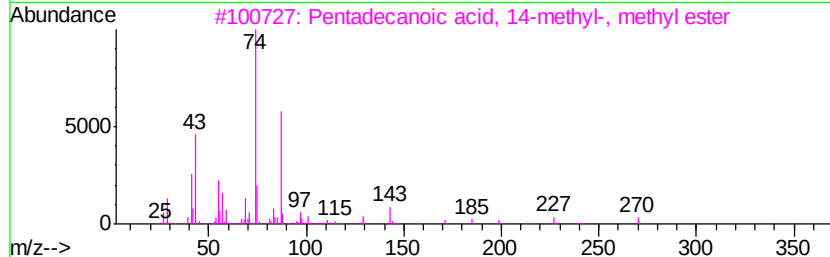
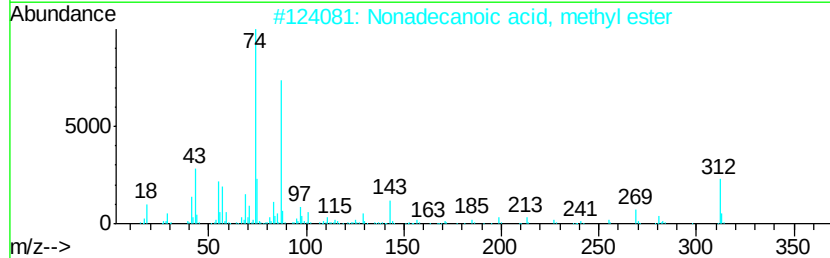
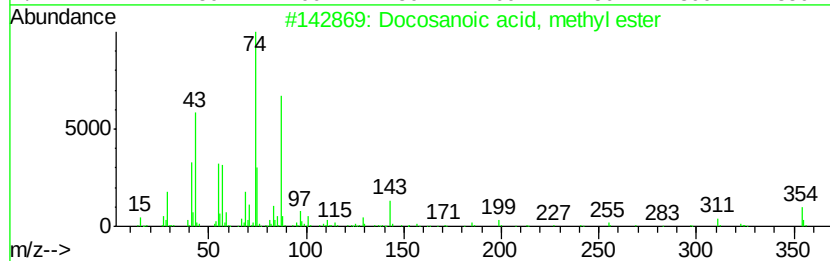
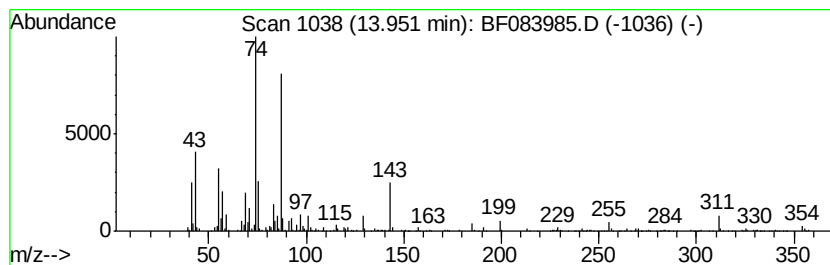
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Docosanoic acid, methyl ester Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.95	40.21 ng	598124	Chrysene-d12	14.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosanoic acid, methyl ester	354	C23H46O2	000929-77-1	97
2		Nonadecanoic acid, methyl ester	312	C20H40O2	001731-94-8	91
3		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	87
4		Octadecanoic acid, methyl ester	298	C19H38O2	000112-61-8	86
5		Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122015\
 Data File : BF083985.D
 Acq On : 20 Dec 2015 7:45
 Operator : UM/SJ
 Sample : G4851-04
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 UST-9

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121815.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
Benzene, 1,3-dime...	5.70	56.4	ng	5157510	1	6.88	1828820	20.0
Nonane	5.78	49.9	ng	4561330	1	6.88	1828820	20.0
Benzene, (1-methy...	6.02	25.2	ng	2306310	1	6.88	1828820	20.0
Octane, 2,6-dimet...	6.12	42.5	ng	3888870	1	6.88	1828820	20.0
Octane, 4-ethyl-	6.17	42.6	ng	3892000	1	6.88	1828820	20.0
Indane	7.06	42.0	ng	3838690	1	6.88	1828820	20.0
Benzene, 1-methyl...	7.16	23.2	ng	2120490	1	6.88	1828820	20.0
Benzene, 2-ethyl-...	7.21	43.2	ng	3950180	1	6.88	1828820	20.0
Pentadecane	12.51	38.3	ng	6880050	4	11.44	3593400	20.0
Phenanthrene, 2,3...	12.91	53.8	ng	800834	5	14.02	297492	20.0
Eicosane	13.22	203.1	ng	3020390	5	14.02	297492	20.0
Tetracosane	13.55	156.7	ng	2330210	5	14.02	297492	20.0
Hexadecane, 1-iodo-	13.87	168.6	ng	2508110	5	14.02	297492	20.0
Docosanoic acid, ...	13.95	40.2	ng	598124	5	14.02	297492	20.0