

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051115\
 Data File : BF079125.D
 Acq On : 11 May 2015 19:41
 Operator : TP/IZ
 Sample : G2146-08
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-4

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-BF043015.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	1.576	31	34	37	rVB	434955	618163	13.94%	1.500%
2	1.747	47	49	57	rVV2	760376	1168423	26.36%	2.836%
3	1.861	57	59	91	rVB	2439669	4433181	100.00%	10.759%
4	5.599	383	386	389	rBV	673161	683910	15.43%	1.660%
5	6.856	494	496	499	rBV	430256	434024	9.79%	1.053%
6	7.005	507	509	512	rBV	1048346	1336127	30.14%	3.243%
7	7.302	532	535	537	rBV	268370	341874	7.71%	0.830%
8	7.485	549	551	552	rBV	1472948	1338931	30.20%	3.250%
9	7.519	552	554	556	rVB	528015	599960	13.53%	1.456%
10	7.965	590	593	594	rBV2	173076	305756	6.90%	0.742%
11	7.988	594	595	598	rVV2	1268467	1255524	28.32%	3.047%
12	8.250	615	618	620	rBV	134659	125623	2.83%	0.305%
13	8.479	634	638	640	rBV	226939	305055	6.88%	0.740%
14	8.559	642	645	648	rBV	624861	812859	18.34%	1.973%
15	8.605	648	649	652	rVV2	87227	123825	2.79%	0.301%
16	8.673	652	655	657	rVB3	36825	64058	1.44%	0.155%
17	8.799	661	666	667	rBV	49505	75310	1.70%	0.183%
18	8.868	669	672	674	rVV2	544641	770918	17.39%	1.871%
19	8.902	674	675	677	rVV	108928	171182	3.86%	0.415%
20	8.936	677	678	681	rVB	262893	258182	5.82%	0.627%
21	8.993	681	683	686	rVB3	35810	55987	1.26%	0.136%
22	9.119	692	694	698	rVB3	25645	50928	1.15%	0.124%
23	9.222	701	703	705	rBV2	103527	165784	3.74%	0.402%
24	9.359	713	715	718	rVB	169375	162225	3.66%	0.394%
25	9.462	722	724	725	rVV	345309	333149	7.51%	0.809%
26	9.508	725	728	732	rVV	141333	237791	5.36%	0.577%
27	9.576	732	734	736	rVV2	66198	94312	2.13%	0.229%
28	9.611	736	737	740	rVB	205000	166264	3.75%	0.404%
29	9.702	743	745	748	rVV	226001	219765	4.96%	0.533%
30	9.759	748	750	753	rVV	1946916	2012833	45.40%	4.885%
31	9.828	753	756	757	rVV2	95979	110991	2.50%	0.269%
32	9.885	757	761	763	rVV	3078764	3082485	69.53%	7.481%
33	9.931	763	765	770	rVB2	93851	116441	2.63%	0.283%
34	10.011	770	772	774	rBV3	45646	81588	1.84%	0.198%

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35	10.056	774	776	779	rBV2	51462	84862	1.91%	0.206%
36	10.114	779	781	783	rVB3	51345	71053	1.60%	0.172%
37	10.216	788	790	792	rBV	1856854	2082930	46.98%	5.055%
38	10.274	792	795	797	rBV2	35689	79096	1.78%	0.192%
39	10.388	804	805	807	rVB2	59547	57640	1.30%	0.140%
40	10.456	809	811	815	rVB	444971	695949	15.70%	1.689%
41	10.548	815	819	821	rBV	747502	1257746	28.37%	3.053%
42	10.582	821	822	823	rBV	68810	55325	1.25%	0.134%
43	10.628	823	826	828	rVV	1117117	1399943	31.58%	3.398%
44	10.662	828	829	832	rVB	525848	556946	12.56%	1.352%
45	10.776	836	839	843	rBV	502055	732525	16.52%	1.778%
46	10.879	844	848	850	rVB	447916	514215	11.60%	1.248%
47	10.936	850	853	855	rVB3	54960	69075	1.56%	0.168%
48	11.005	858	859	861	rBV2	37729	57819	1.30%	0.140%
49	11.051	861	863	865	rBV2	677430	715062	16.13%	1.735%
50	11.085	865	866	870	rBV2	154643	246209	5.55%	0.598%
51	11.188	873	875	878	rBV	124861	215636	4.86%	0.523%
52	11.245	878	880	882	rBV2	83573	121017	2.73%	0.294%
53	11.302	882	885	888	rVV	173660	307797	6.94%	0.747%
54	11.359	888	890	892	rVB2	131157	147150	3.32%	0.357%
55	11.451	895	898	900	rBV	131886	182805	4.12%	0.444%
56	11.485	900	901	904	rVB2	96363	110710	2.50%	0.269%
57	11.565	904	908	912	rBV3	114148	338252	7.63%	0.821%
58	11.668	915	917	921	rVB5	55598	96765	2.18%	0.235%
59	11.725	921	922	924	rBV	245989	296877	6.70%	0.721%
60	11.817	927	930	934	rVV3	167487	341263	7.70%	0.828%
61	11.885	934	936	939	rVV4	45006	102729	2.32%	0.249%
62	11.954	939	942	946	rVB3	75803	136504	3.08%	0.331%
63	12.034	946	949	951	rBV	1434337	1451053	32.73%	3.522%
64	12.091	952	954	958	rVB3	51879	99191	2.24%	0.241%
65	12.159	958	960	961	rBV	42371	57104	1.29%	0.139%
66	12.194	961	963	965	rVB2	80799	114997	2.59%	0.279%
67	12.285	968	971	973	rBV3	44697	70099	1.58%	0.170%
68	12.331	973	975	979	rVB5	18207	45454	1.03%	0.110%
69	12.411	979	982	984	rBV	61766	109531	2.47%	0.266%
70	12.445	984	985	990	rBV4	77038	139464	3.15%	0.338%
71	12.891	1022	1024	1026	rBV	454310	562947	12.70%	1.366%

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72	12.982	1030	1032	1035	rVB2	54370	83505	1.88%	0.203%
73	13.040	1035	1037	1039	rBV	290354	290065	6.54%	0.704%
74	13.200	1049	1051	1052	rVV	61105	81964	1.85%	0.199%
75	13.520	1077	1079	1081	rBV2	64170	97391	2.20%	0.236%
76	13.565	1081	1083	1086	rVV	129055	159226	3.59%	0.386%
77	13.611	1086	1087	1089	rVV	56125	80492	1.82%	0.195%
78	13.657	1089	1091	1093	rVV	67317	93442	2.11%	0.227%
79	14.068	1124	1127	1130	rBV3	25356	66774	1.51%	0.162%
80	14.891	1197	1199	1202	rVB	2094440	2577878	58.15%	6.257%
81	16.068	1300	1302	1306	rBV	57801	77106	1.74%	0.187%
82	16.194	1310	1313	1315	rBV	688631	710384	16.02%	1.724%
83	16.846	1366	1370	1373	rBV	187282	341840	7.71%	0.830%
84	17.954	1464	1467	1470	rBV	590634	705749	15.92%	1.713%

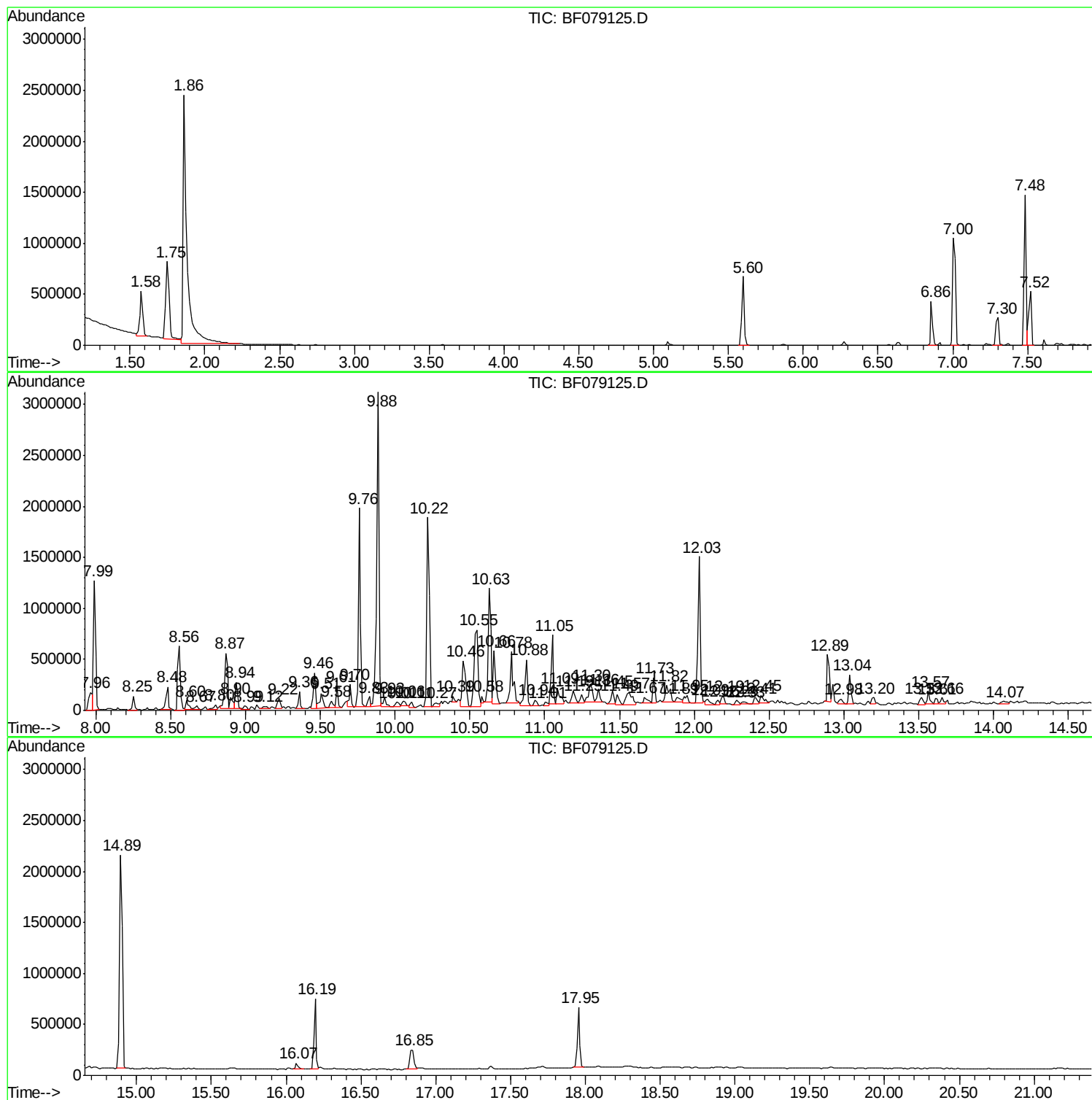
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :
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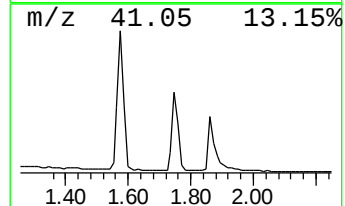
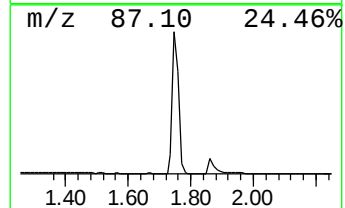
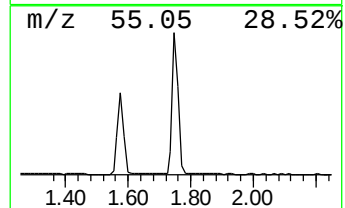
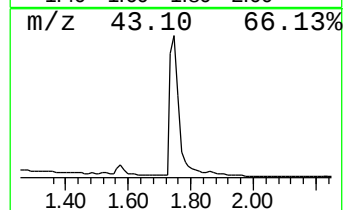
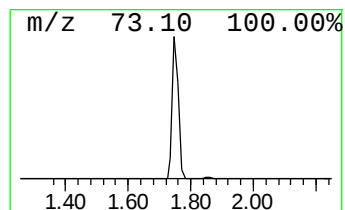
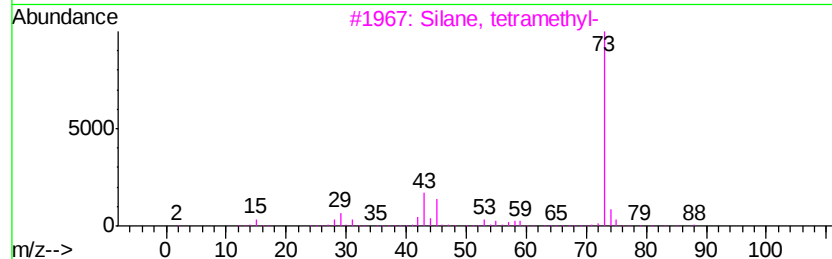
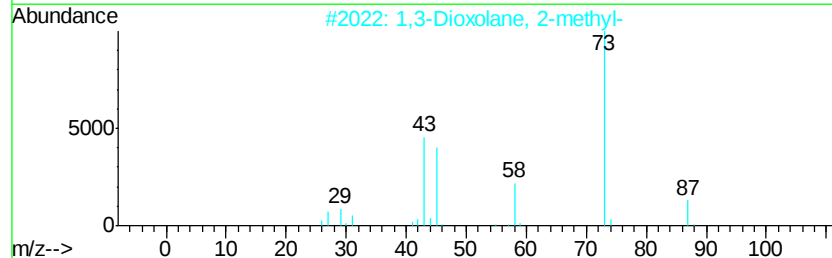
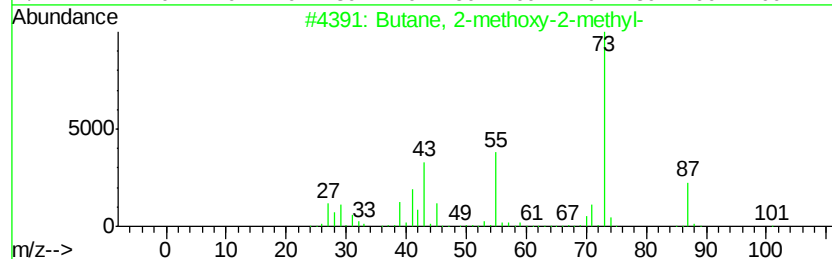
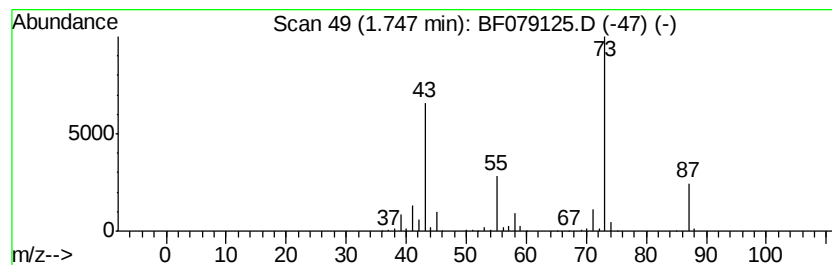
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.75	68.35 ng	1168420	1,4-Dichlorobenzene-d4	7.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
4		Acetamide, N-methyl-	73	C3H7NO	000079-16-3	12
5		Borane, dimethoxy-	74	C2H7BO2	004542-61-4	12



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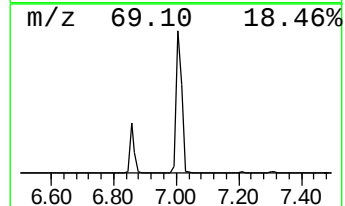
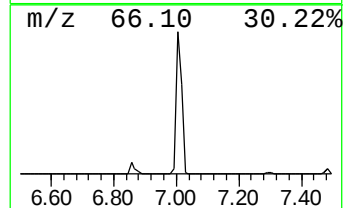
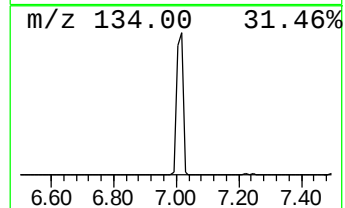
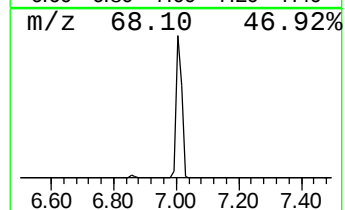
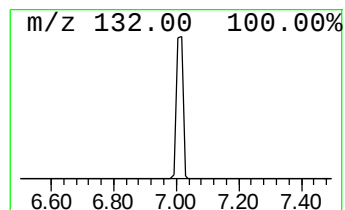
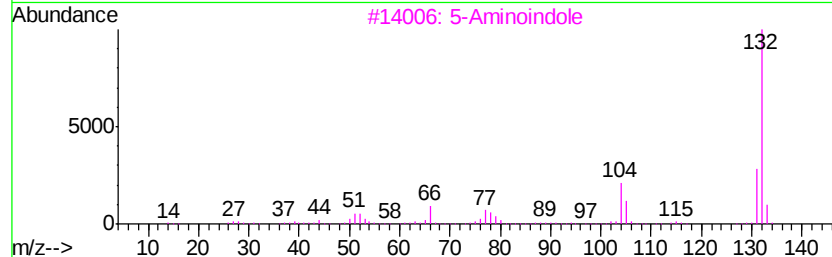
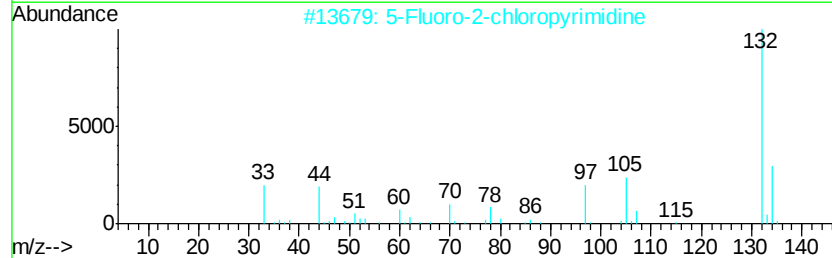
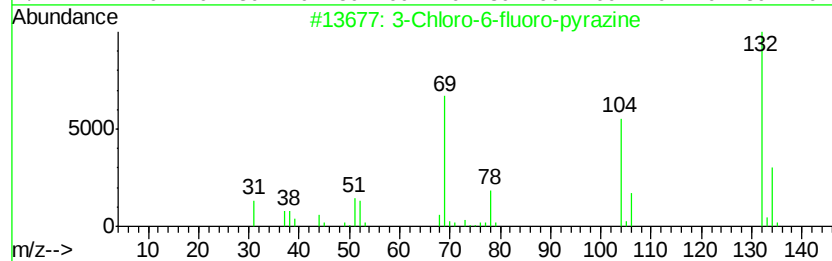
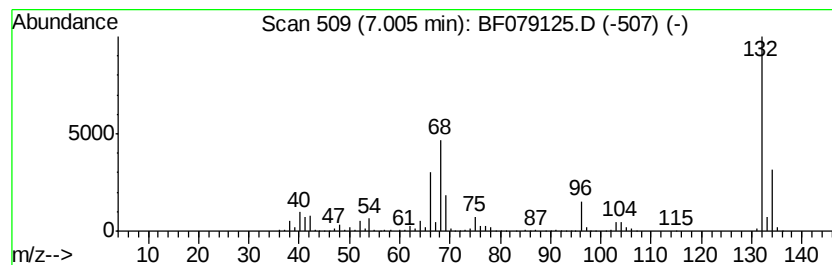
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown7.00 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.00	78.16 ng	1336130	1,4-Dichlorobenzene-d4	7.30

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	12
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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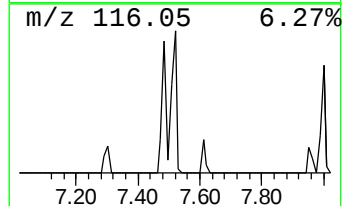
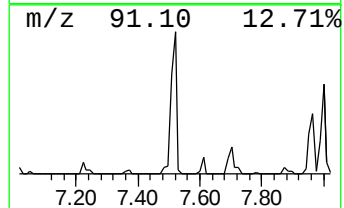
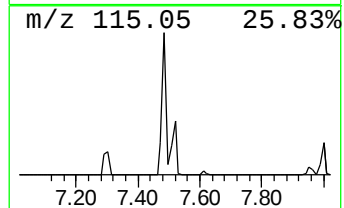
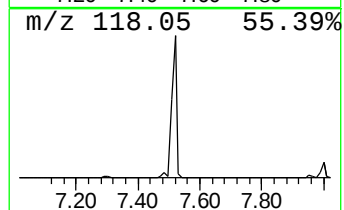
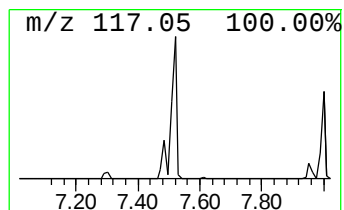
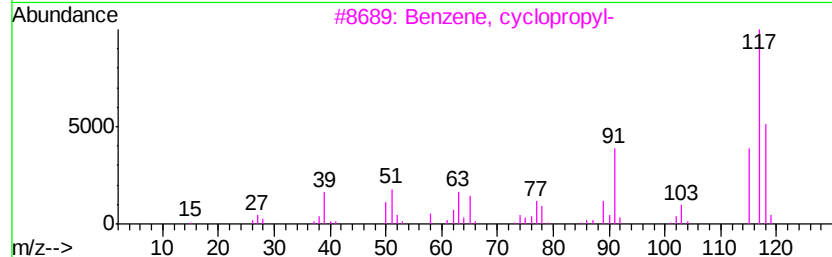
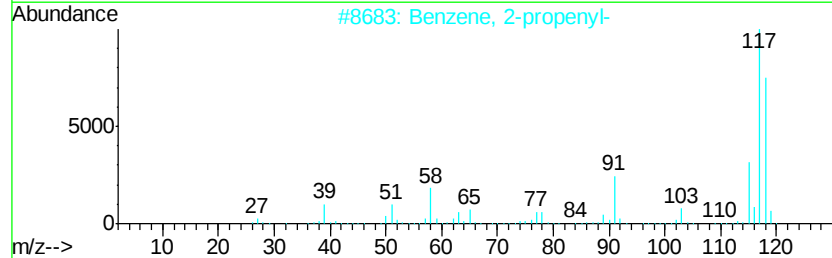
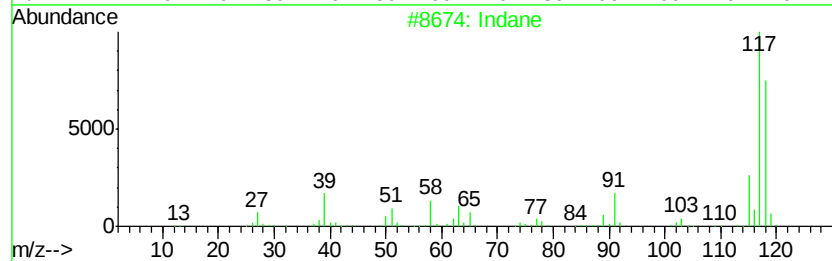
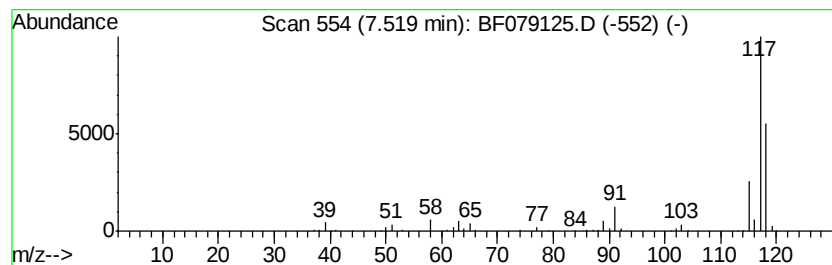
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Indane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.52	35.10 ng	599960	1,4-Dichlorobenzene-d4	7.30

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	87
2	Benzene, 2-propenyl-		118	C9H10	000300-57-2	64
3	Benzene, cyclopropyl-		118	C9H10	000873-49-4	64
4	Benzene, 1-propenyl-		118	C9H10	000637-50-3	50
5	m-Aminophenylacetylene		117	C8H7N	054060-30-9	43



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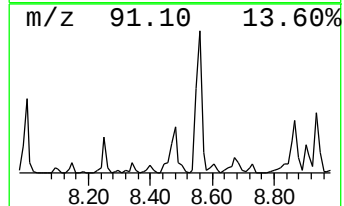
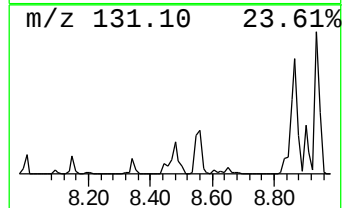
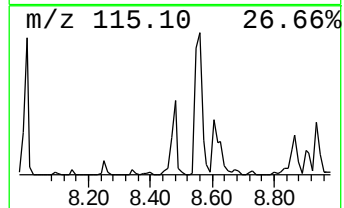
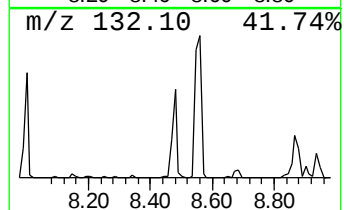
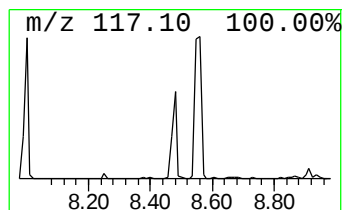
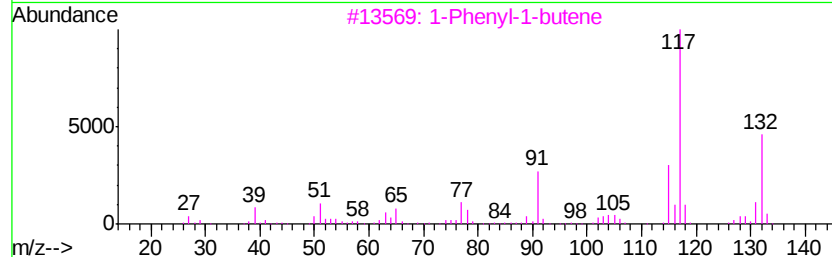
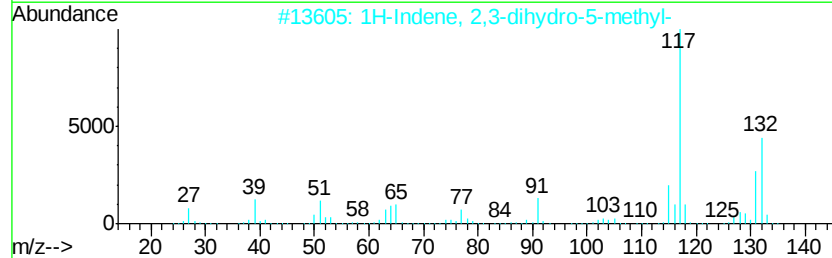
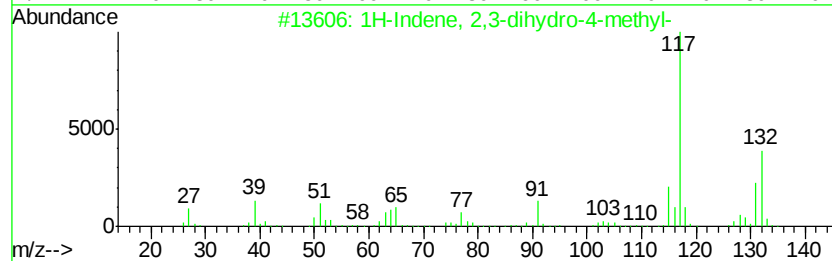
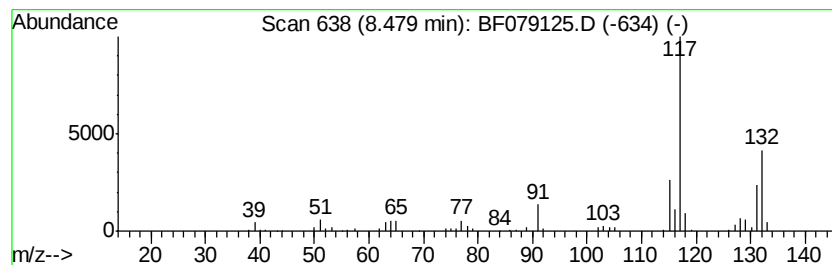
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.48	7.91 ng	305055	Napthalene-d8	8.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	87
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051115\
Data File : BF079125.D
Acq On : 11 May 2015 19:41
Operator : TP/IZ
Sample : G2146-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-4

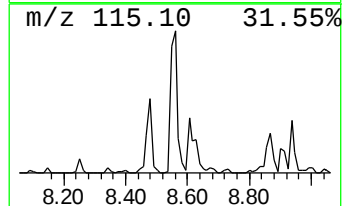
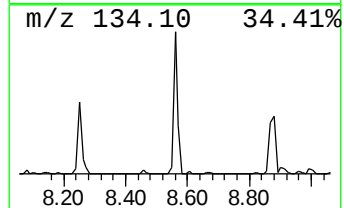
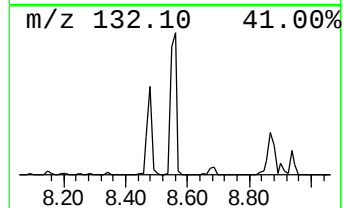
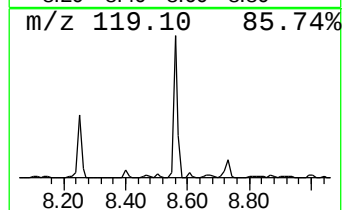
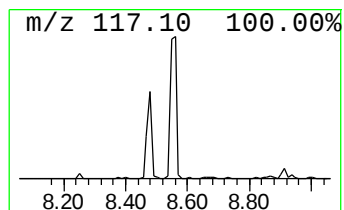
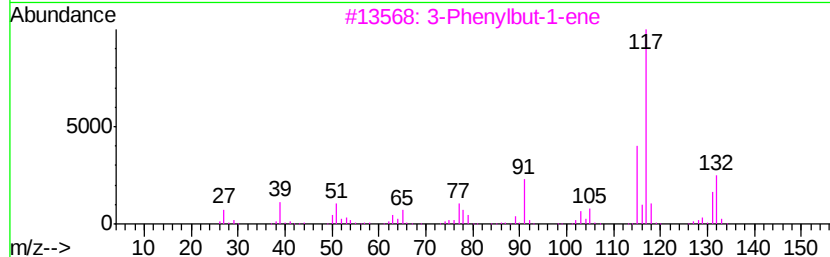
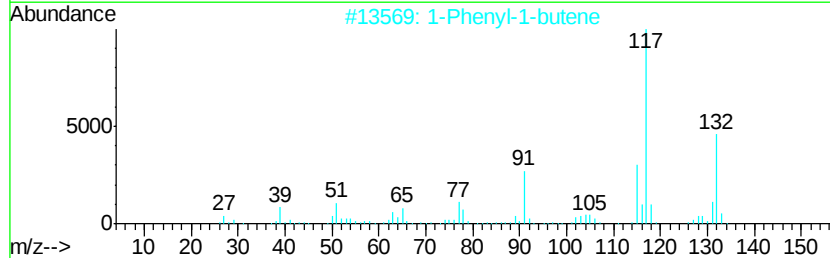
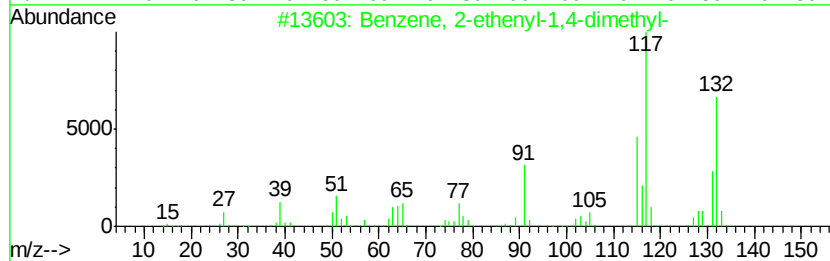
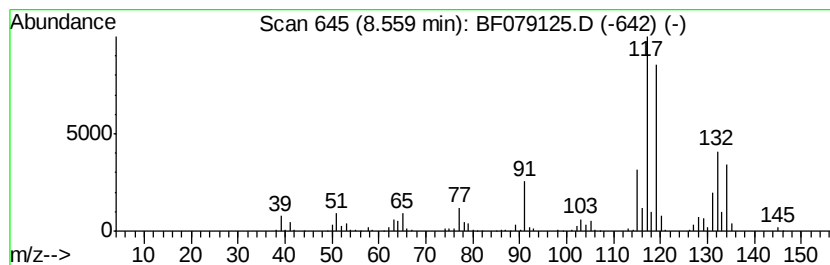
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.56	21.09 ng	812859	Naphthalene-d8	8.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	89
2		1-Phenyl-1-butene	132	C10H12	000824-90-8	87
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	60
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	55
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	55



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051115\
Data File : BF079125.D
Acq On : 11 May 2015 19:41
Operator : TP/IZ
Sample : G2146-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-4

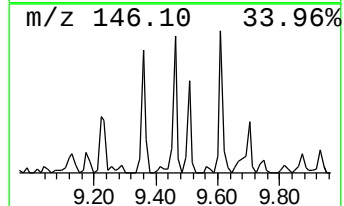
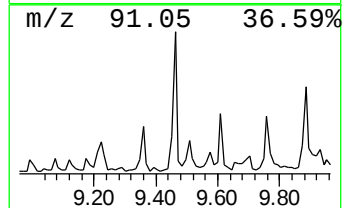
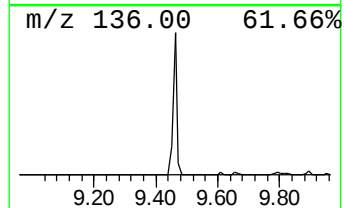
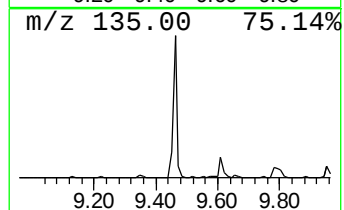
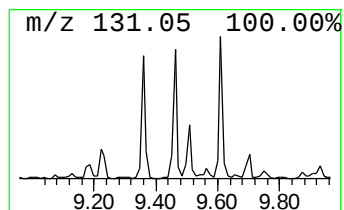
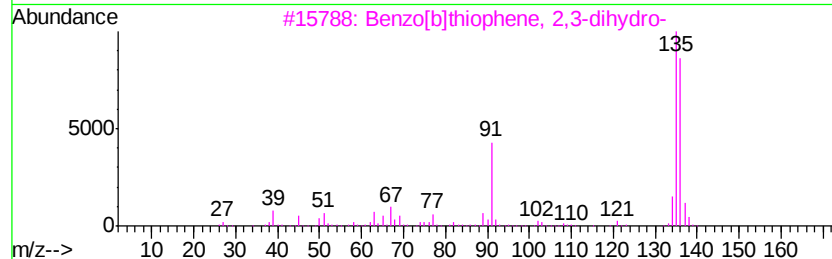
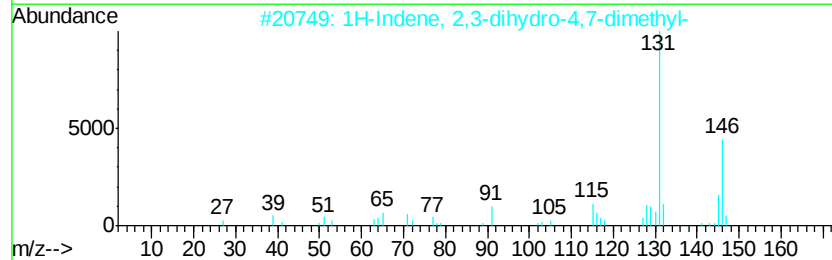
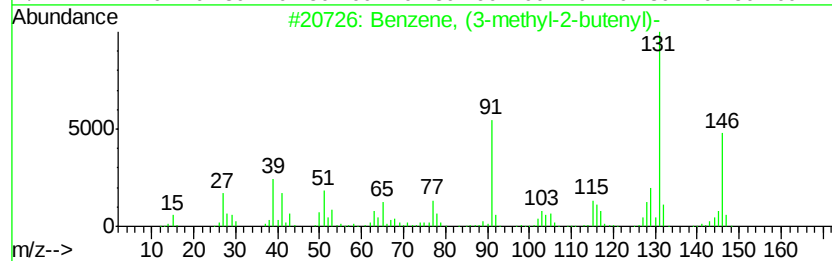
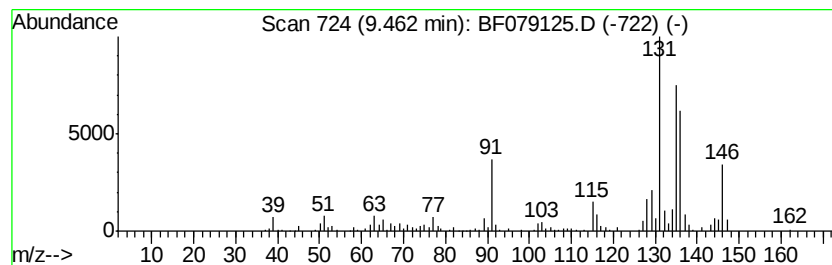
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzene, (3-methyl-2-butenyl)- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.46	8.64 ng	333149	Naphthalene-d8	8.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64
2			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	46
3			Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	46
4			Naphthalene, 1,2,3,4-tetrahydro...	146	C11H14	001559-81-5	46
5			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	46



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051115\
Data File : BF079125.D
Acq On : 11 May 2015 19:41
Operator : TP/IZ
Sample : G2146-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

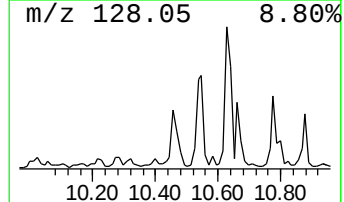
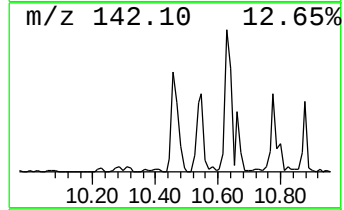
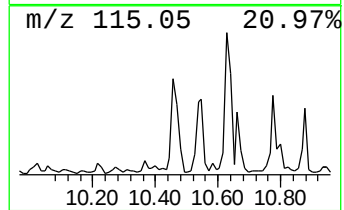
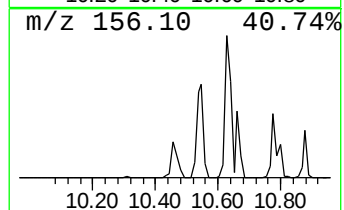
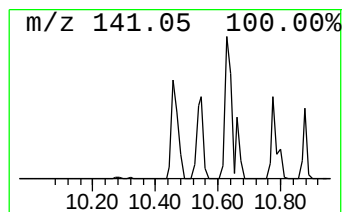
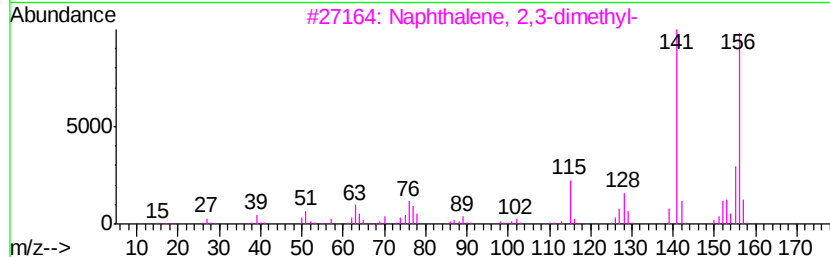
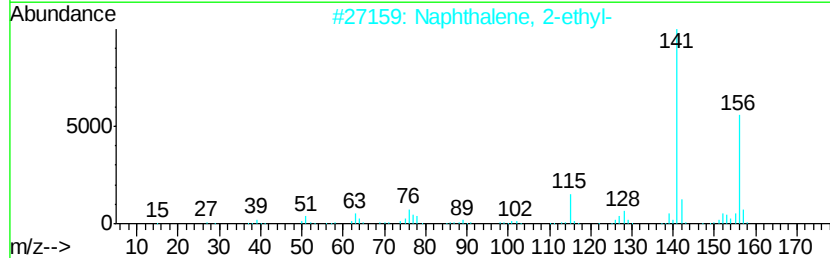
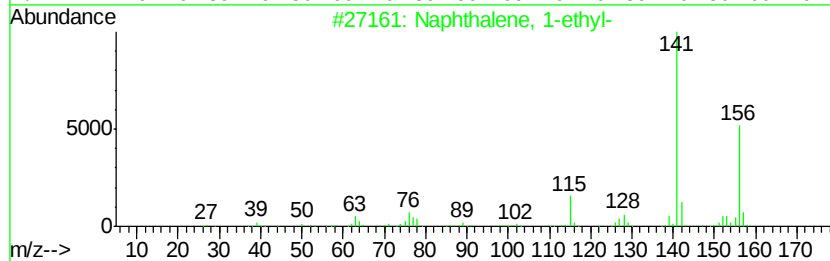
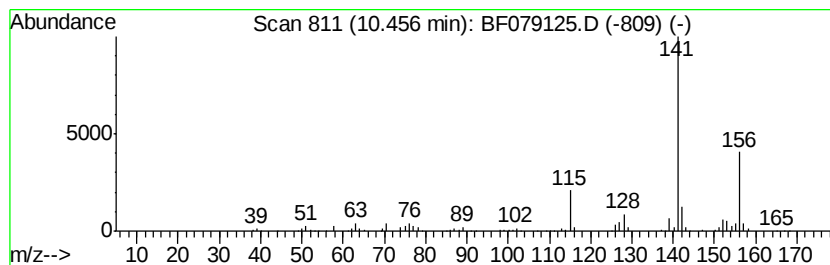
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Naphthalene, 2-ethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.46	19.47 ng	695949	Acenaphthene-d10	11.05
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1-ethyl-	156 C12H12	001127-76-0 95
2		Naphthalene, 2-ethyl-	156 C12H12	000939-27-5 94
3		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 78
4		2-Thiophenecarboxylic acid, 5-et...	156 C7H8O2S	023229-72-3 64
5		Methyl phenylphosphinic acid	156 C7H9O2P	004271-13-0 53



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051115\
Data File : BF079125.D
Acq On : 11 May 2015 19:41
Operator : TP/IZ
Sample : G2146-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-4

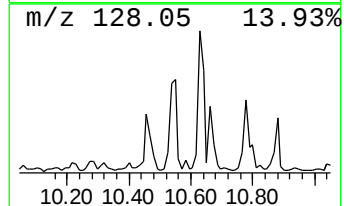
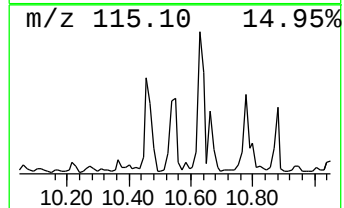
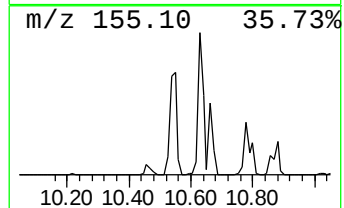
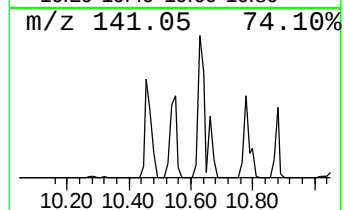
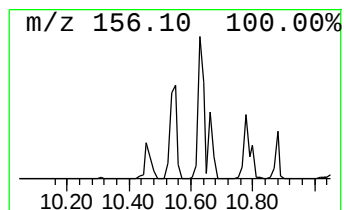
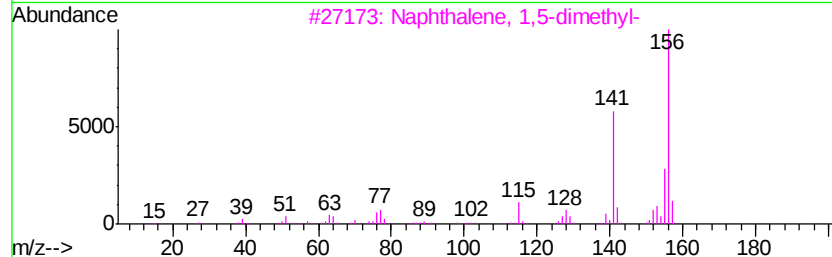
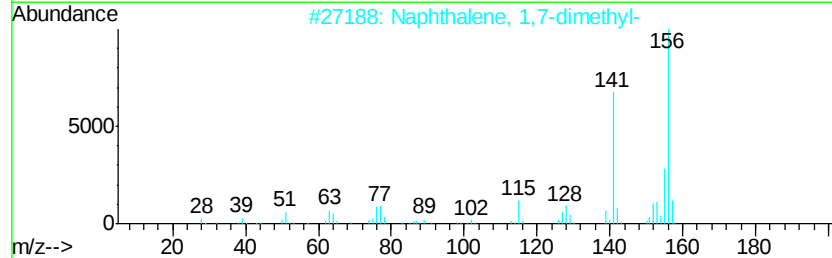
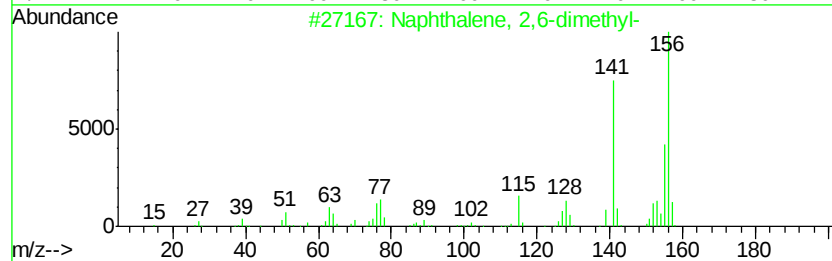
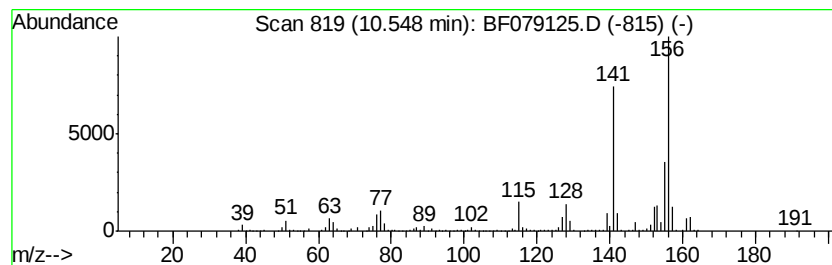
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Naphthalene, 2,6-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.55	35.18 ng	1257750	Acenaphthene-d10	11.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
2		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
4		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
5		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051115\
Data File : BF079125.D
Acq On : 11 May 2015 19:41
Operator : TP/IZ
Sample : G2146-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-4

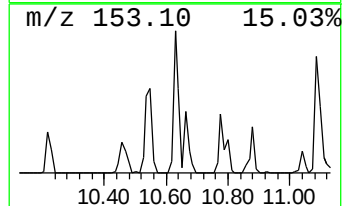
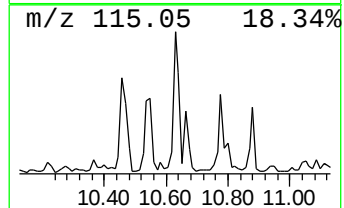
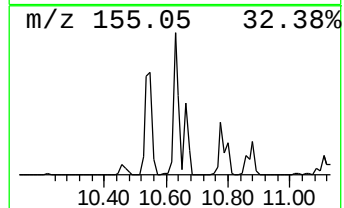
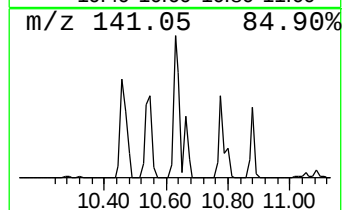
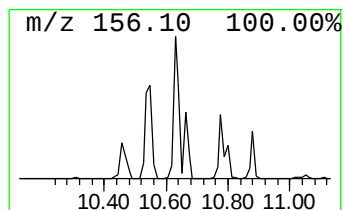
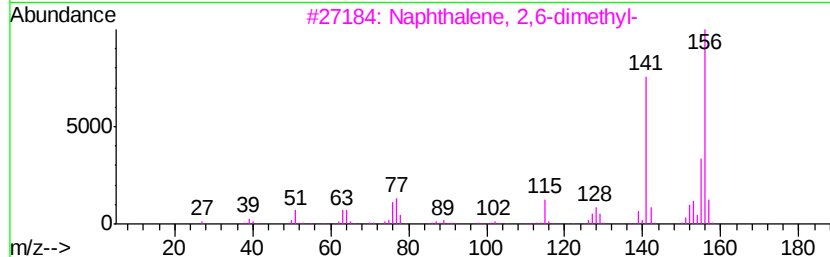
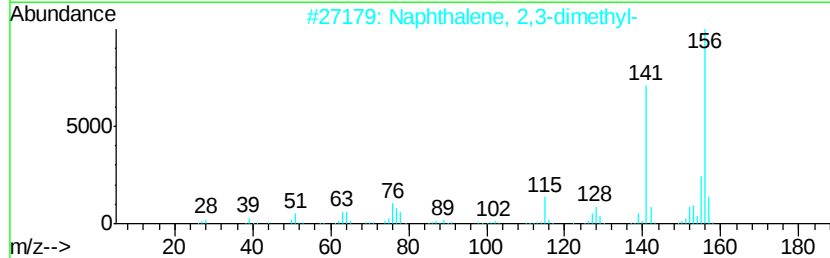
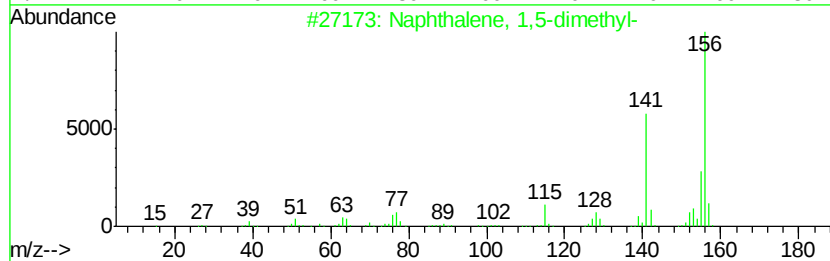
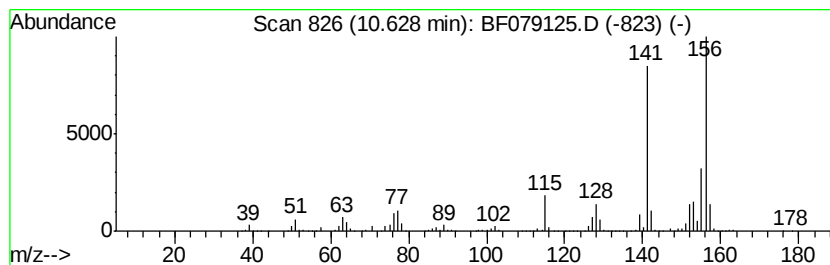
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Naphthalene, 1,5-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.63	39.16 ng	1399940	Acenaphthene-d10	11.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
5			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051115\
Data File : BF079125.D
Acq On : 11 May 2015 19:41
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Sample : G2146-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-4

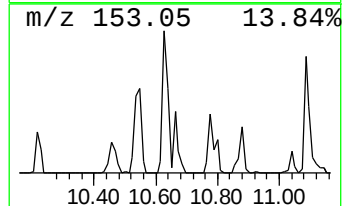
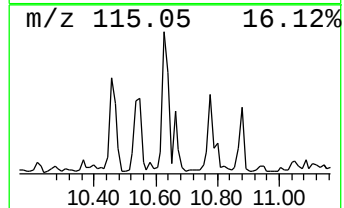
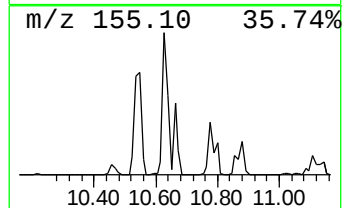
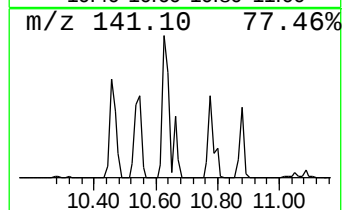
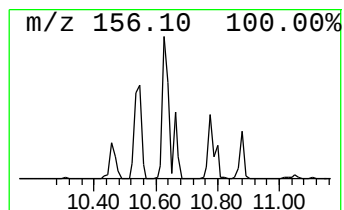
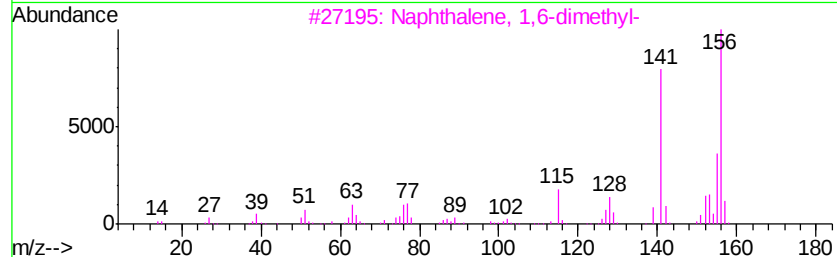
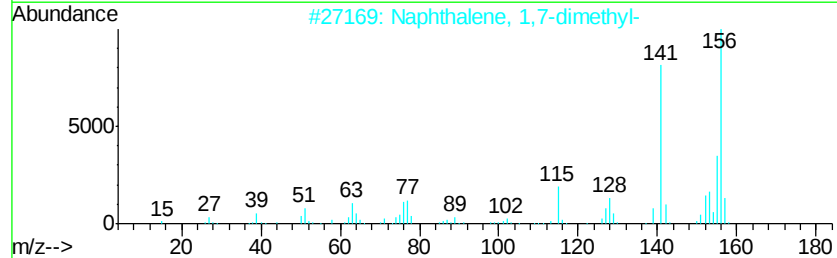
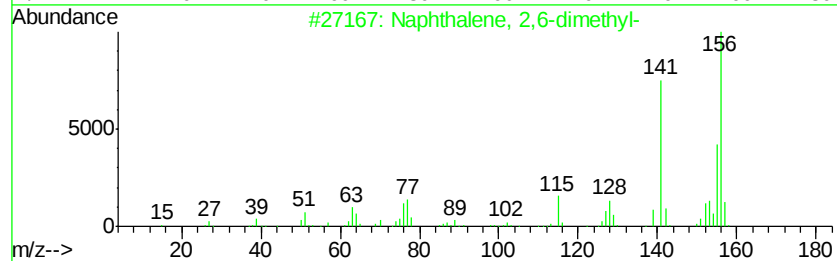
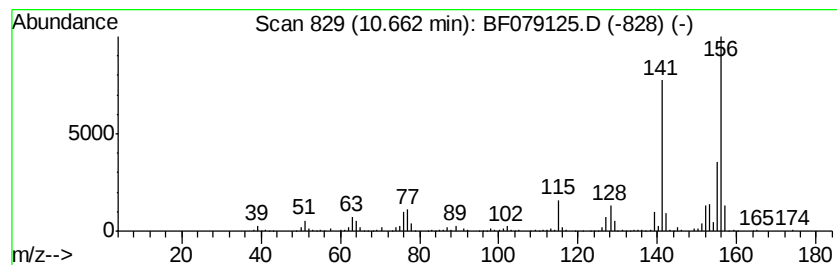
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Naphthalene, 1,7-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.66	15.58 ng	556946	Acenaphthene-d10	11.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
2		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
3		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
5		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051115\
Data File : BF079125.D
Acq On : 11 May 2015 19:41
Operator : TP/IZ
Sample : G2146-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-4

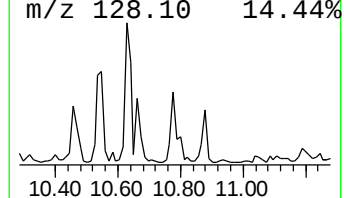
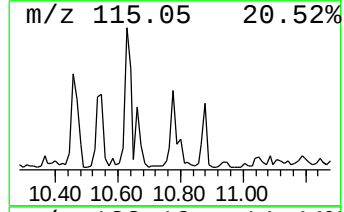
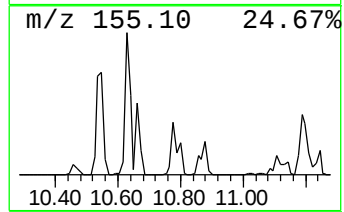
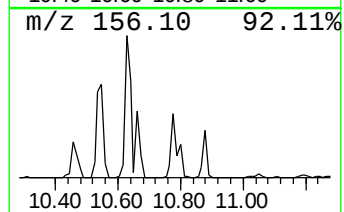
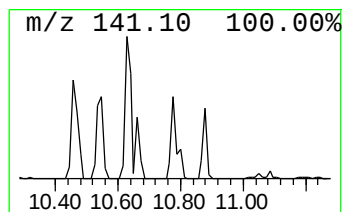
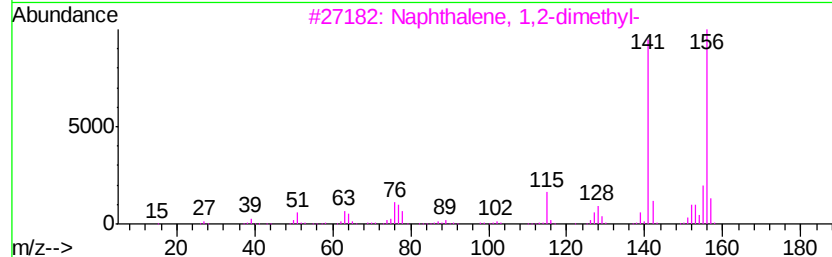
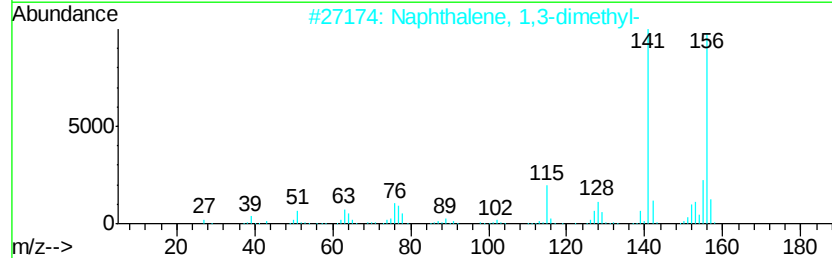
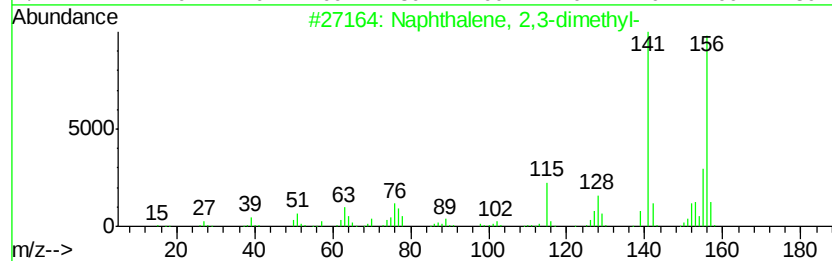
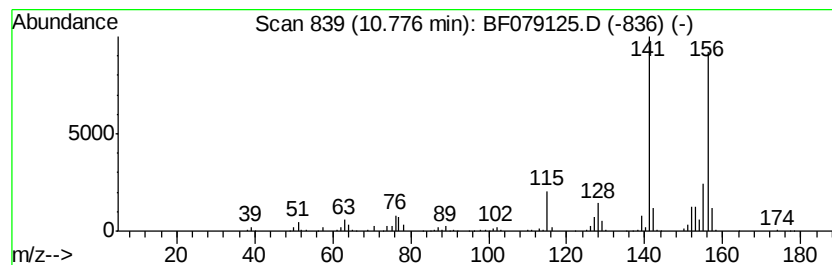
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Naphthalene, 2,3-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.78	20.49 ng	732525	Acenaphthene-d10	11.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
3		Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	96
4		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5		Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	96



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051115\
Data File : BF079125.D
Acq On : 11 May 2015 19:41
Operator : TP/IZ
Sample : G2146-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-4

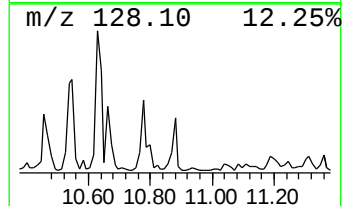
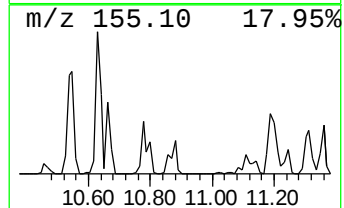
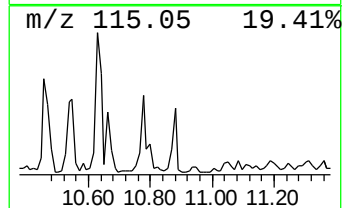
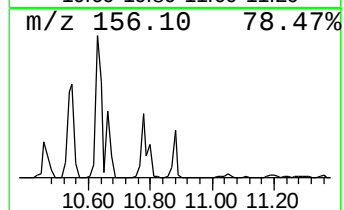
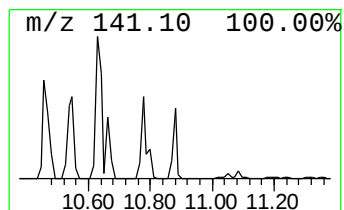
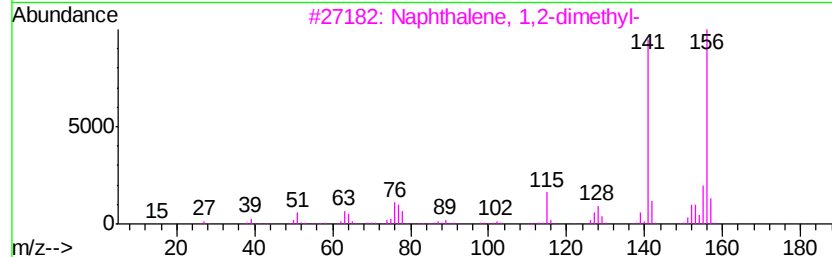
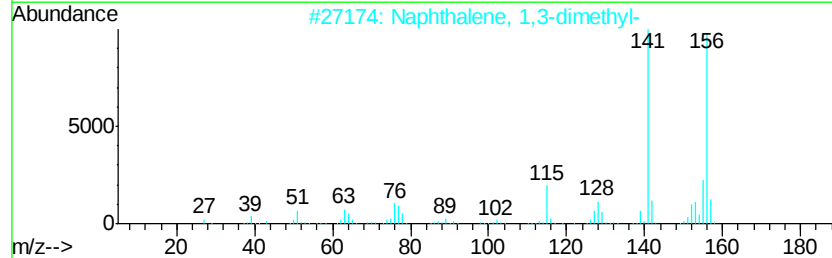
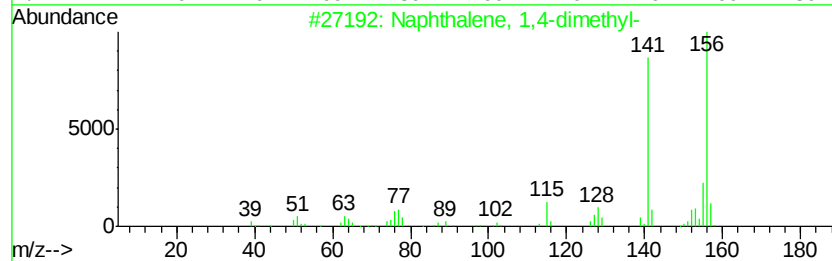
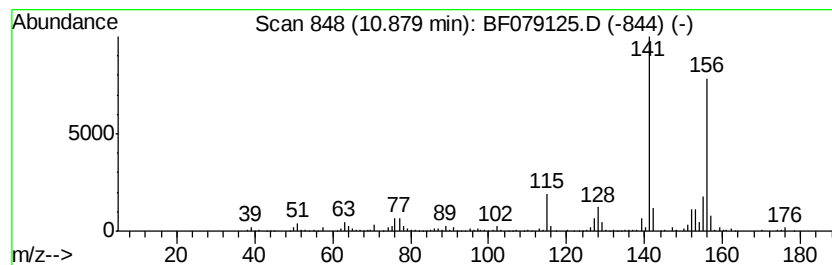
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Naphthalene, 1,4-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.88	14.38 ng	514215	Acenaphthene-d10	11.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
2		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
3		Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	96
4		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
5		Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	94



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051115\
Data File : BF079125.D
Acq On : 11 May 2015 19:41
Operator : TP/IZ
Sample : G2146-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-4

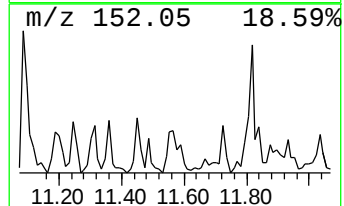
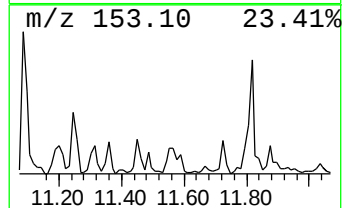
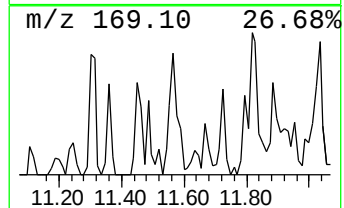
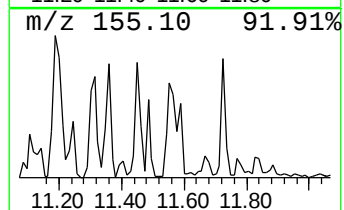
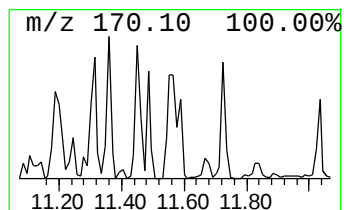
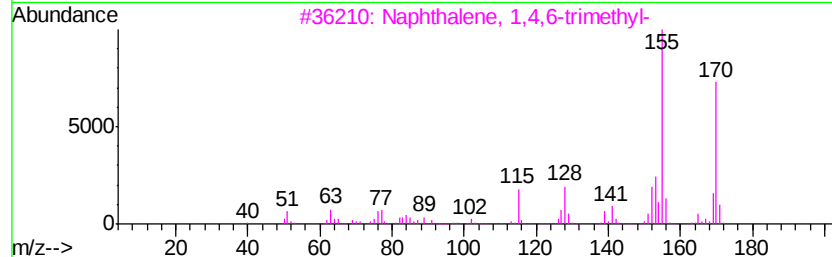
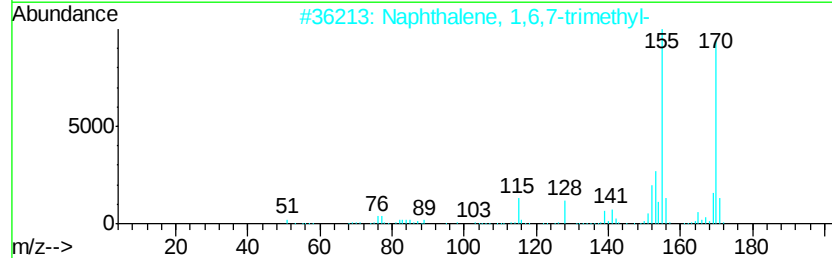
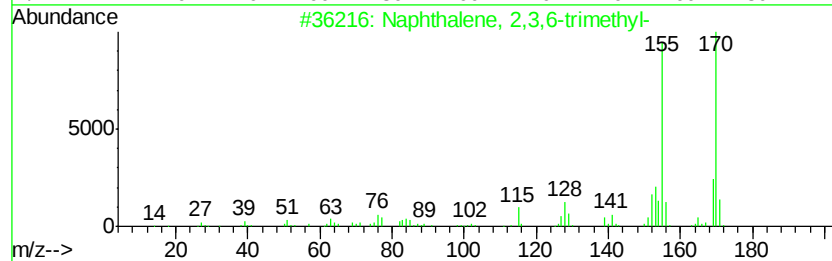
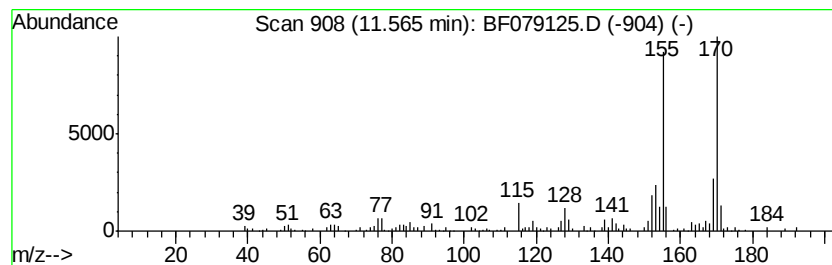
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Naphthalene, 2,3,6-trimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	9.46 ng	338252	Acenaphthene-d10	11.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	98
2			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	98
3			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	96
4			Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	96
5			Azulene, 4,6,8-trimethyl-	170	C13H14	000941-81-1	94



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051115\
Data File : BF079125.D
Acq On : 11 May 2015 19:41
Operator : TP/IZ
Sample : G2146-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-4

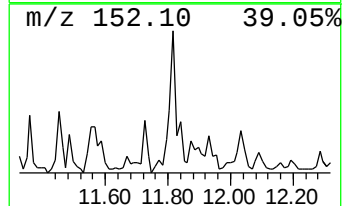
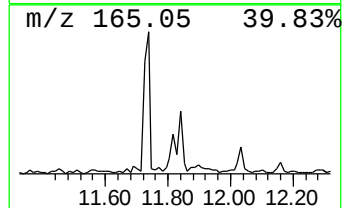
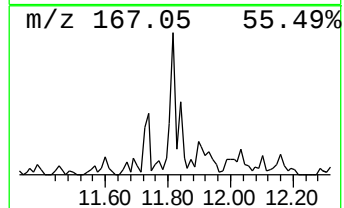
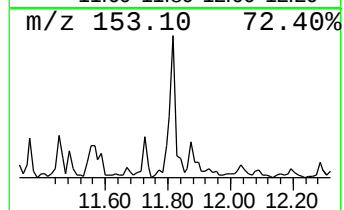
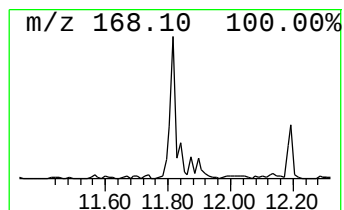
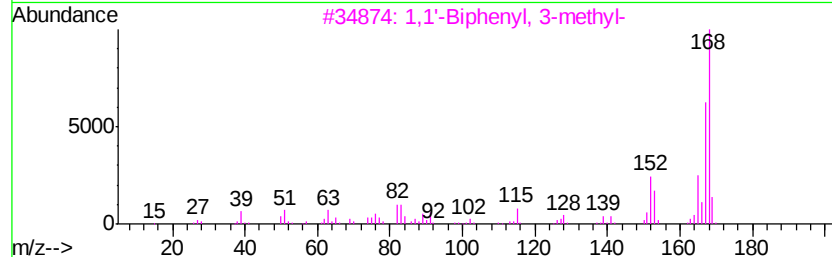
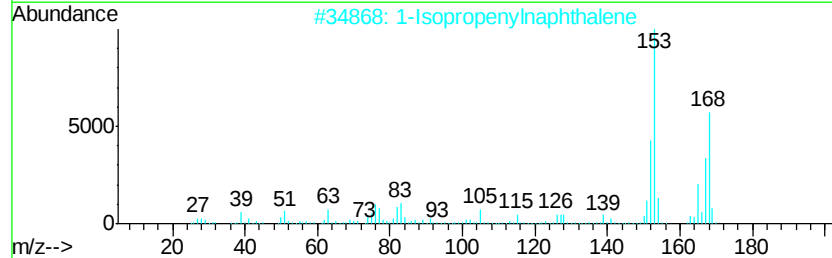
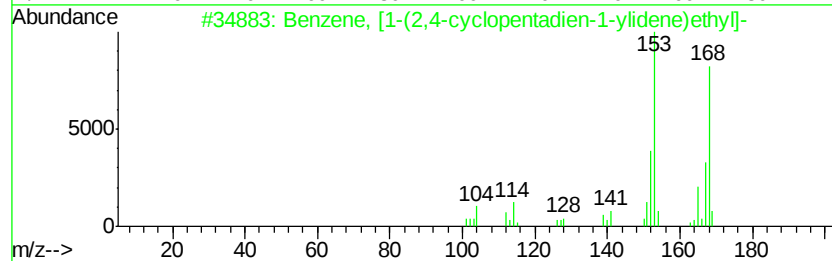
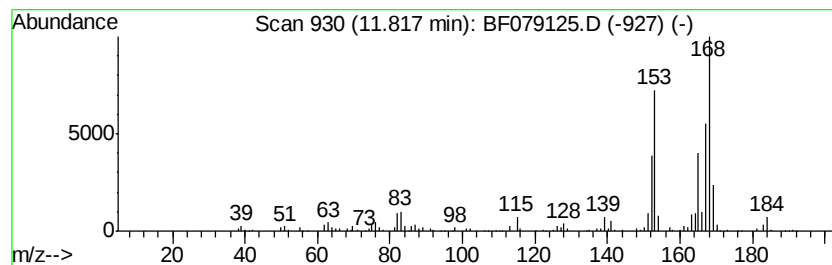
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Benzene, [1-(2,4-cyclopenta... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.82	9.54 ng	341263	Acenaphthene-d10	11.05

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, [1-(2,4-cyclopentadien-...	168	C13H12	002320-32-3	64
2		1-Isopropenylnaphthalene	168	C13H12	001855-47-6	58
3		1,1'-Biphenyl, 3-methyl-	168	C13H12	000643-93-6	49
4		1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	49
5		Naphthalene, 2-(1-methylethenyl)-	168	C13H12	003710-23-4	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051115\
Data File : BF079125.D
Acq On : 11 May 2015 19:41
Operator : TP/IZ
Sample : G2146-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-4

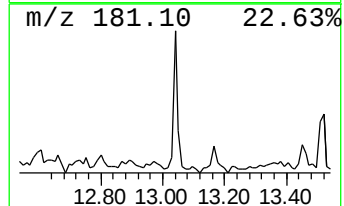
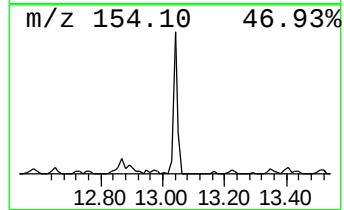
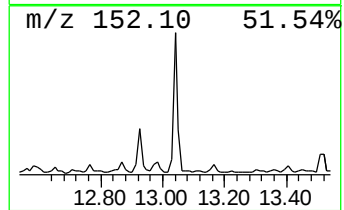
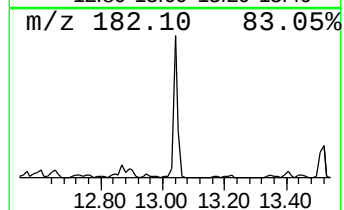
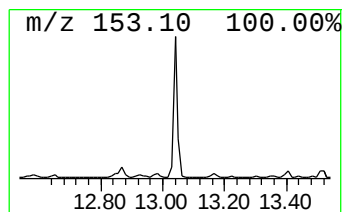
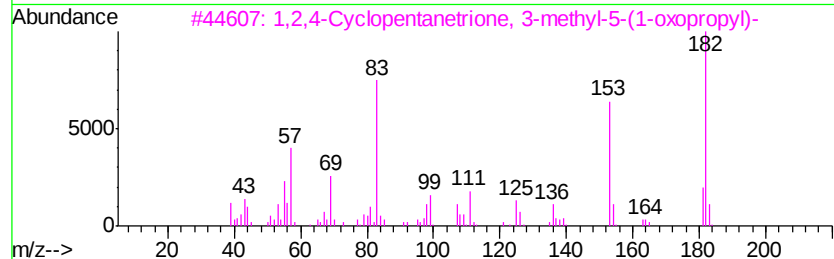
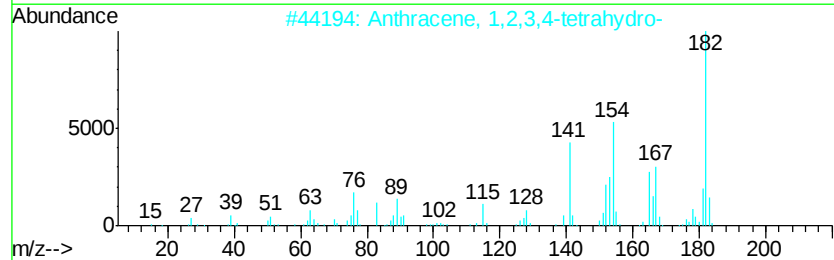
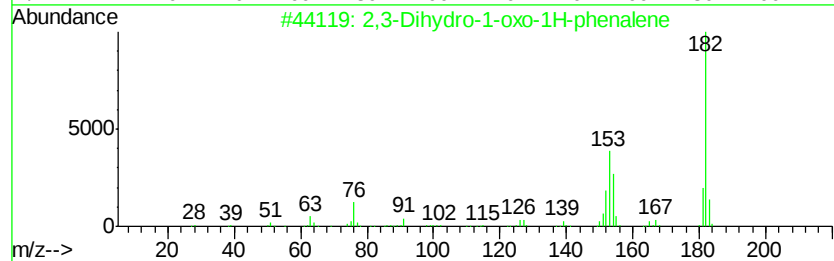
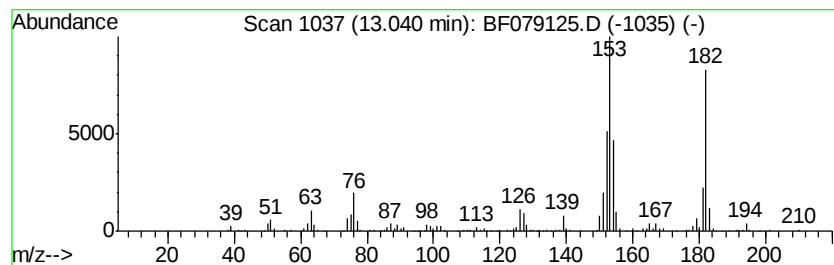
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 2,3-Dihydro-1-oxo-1H-phenalene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.04	10.31 ng	290065	Phenanthrene-d10	12.89

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,3-Dihydro-1-oxo-1H-phenalene	182	C13H10O	000518-85-4	70
2			Anthracene, 1,2,3,4-tetrahydro-	182	C14H14	002141-42-6	58
3			1,2,4-Cyclopentanetrione, 3-meth...	182	C9H10O4	057174-14-8	47
4			4,4'-Dimethylbiphenyl	182	C14H14	000613-33-2	30
5			9H-Pyrido[3,4-b]indole, 1-methyl-	182	C12H10N2	000486-84-0	27



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051115\
Data File : BF079125.D
Acq On : 11 May 2015 19:41
Operator : TP/IZ
Sample : G2146-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-4

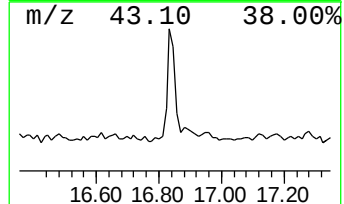
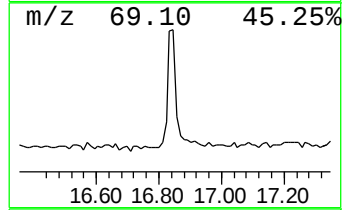
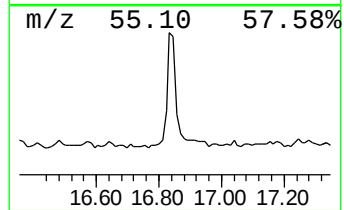
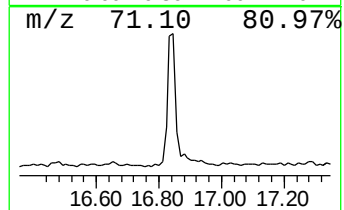
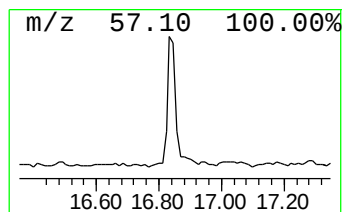
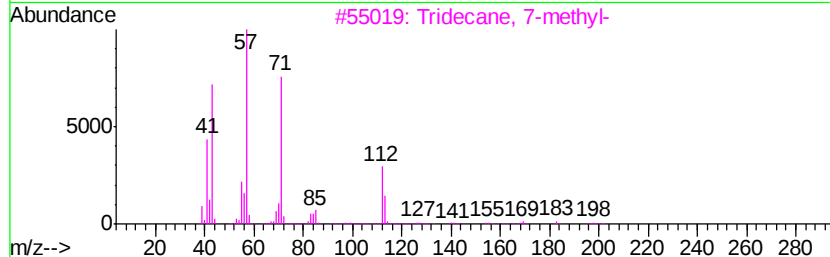
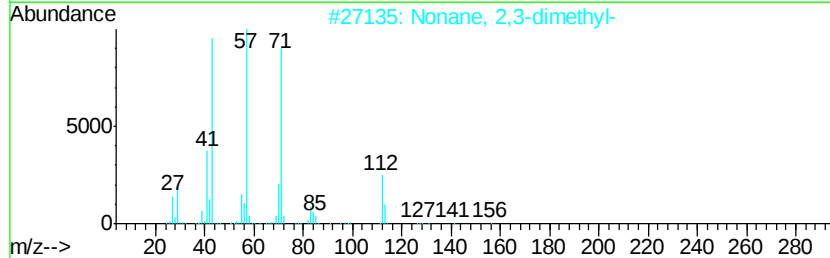
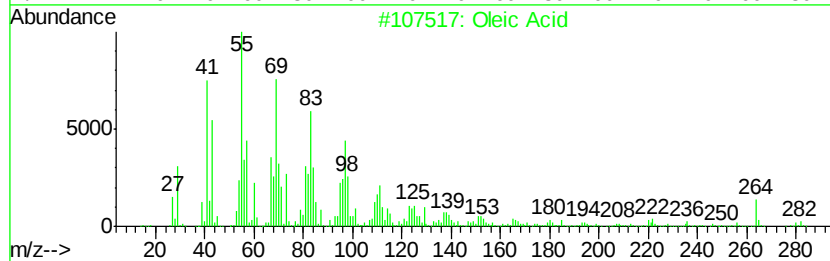
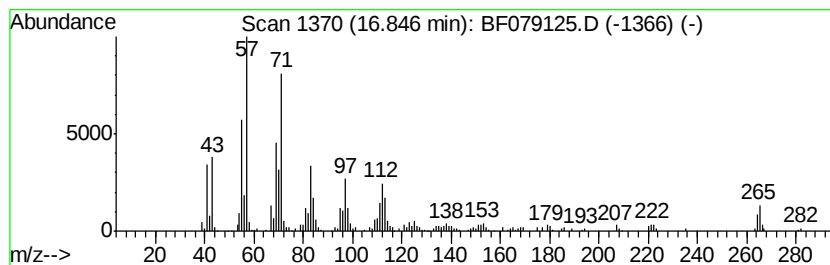
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Oleic Acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.85	9.62 ng	341840	Chrysene-d12	16.19

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oleic Acid	282	C18H34O2	000112-80-1	72
2			Nonane, 2,3-dimethyl-	156	C11H24	002884-06-2	47
3			Tridecane, 7-methyl-	198	C14H30	026730-14-3	47
4			Heptane, 3-ethyl-5-methyl-	142	C10H22	052896-90-9	43
5			Octane, 2-bromo-	192	C8H17Br	000557-35-7	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051115\
 Data File : BF079125.D
 Acq On : 11 May 2015 19:41
 Operator : TP/IZ
 Sample : G2146-08
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.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
Butane, 2-methoxy...	1.75	68.3	ng	1168420	1	7.30	341874	20.0
unknown7.00	7.00	78.2	ng	1336130	1	7.30	341874	20.0
Indane	7.52	35.1	ng	599960	1	7.30	341874	20.0
1H-Indene, 2,3-di...	8.48	7.9	ng	305055	2	8.88	770918	20.0
Benzene, 2-etheny...	8.56	21.1	ng	812859	2	8.88	770918	20.0
Benzene, (3-methy...	9.46	8.6	ng	333149	2	8.88	770918	20.0
Naphthalene, 2-et...	10.46	19.5	ng	695949	3	11.05	715062	20.0
Naphthalene, 2,6-...	10.55	35.2	ng	1257750	3	11.05	715062	20.0
Naphthalene, 1,5-...	10.63	39.2	ng	1399940	3	11.05	715062	20.0
Naphthalene, 1,7-...	10.66	15.6	ng	556946	3	11.05	715062	20.0
Naphthalene, 2,3-...	10.78	20.5	ng	732525	3	11.05	715062	20.0
Naphthalene, 1,4-...	10.88	14.4	ng	514215	3	11.05	715062	20.0
Naphthalene, 2,3,...	11.57	9.5	ng	338252	3	11.05	715062	20.0
Benzene, [1-(2,4-...	11.82	9.5	ng	341263	3	11.05	715062	20.0
2,3-Dihydro-1-oxo...	13.04	10.3	ng	290065	4	12.89	562947	20.0
Oleic Acid	16.85	9.6	ng	341840	5	16.19	710384	20.0