

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032116\
 Data File : BF085629.D
 Acq On : 21 Mar 2016 18:34
 Operator : UM/SJ
 Sample : H1803-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-01

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-BF031816.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	5.030	238	240	248	rBV	46910	67266	2.33%	0.206%
2	5.442	273	276	278	rBV	2524984	2744779	95.09%	8.399%
3	6.470	363	366	369	rBV	1839035	2769337	95.94%	8.474%
4	6.607	375	378	381	rBV	2702530	2585730	89.58%	7.912%
5	6.836	396	398	401	rBV	752463	844274	29.25%	2.583%
6	6.996	410	412	415	rVB	2357510	2119315	73.42%	6.485%
7	7.407	445	448	450	rBV	1918380	1854322	64.24%	5.674%
8	7.499	453	456	459	rVV	45584	64862	2.25%	0.198%
9	7.865	484	488	491	rBV	18894	31535	1.09%	0.096%
10	8.128	509	511	513	rBV	1347516	1156700	40.07%	3.539%
11	9.202	602	605	608	rBV	3110169	2886537	100.00%	8.833%
12	9.591	638	639	641	rBV	231784	294981	10.22%	0.903%
13	9.739	650	652	654	rBV	40911	35635	1.23%	0.109%
14	9.819	654	659	661	rVV2	46966	87146	3.02%	0.267%
15	9.876	662	664	666	rVV	1163991	1092632	37.85%	3.343%
16	9.911	666	667	669	rVB	49056	38104	1.32%	0.117%
17	10.082	680	682	685	rBV3	25859	35029	1.21%	0.107%
18	10.288	697	700	701	rBV	32774	34999	1.21%	0.107%
19	10.311	701	702	704	rVB	116180	89233	3.09%	0.273%
20	10.425	710	712	715	rVB	47930	49513	1.72%	0.152%
21	10.528	715	721	724	rBV2	33783	79779	2.76%	0.244%
22	10.665	730	733	736	rBV	1531215	1489462	51.60%	4.558%
23	10.779	742	743	748	rBV	79547	144314	5.00%	0.442%
24	10.974	758	760	761	rBV2	24982	30909	1.07%	0.095%
25	11.168	775	777	779	rVB2	30996	30425	1.05%	0.093%
26	11.236	780	783	784	rBV2	47106	57446	1.99%	0.176%
27	11.362	792	794	795	rBV	1114837	988629	34.25%	3.025%
28	11.431	798	800	802	rVB2	65557	101956	3.53%	0.312%
29	11.591	811	814	817	rBV2	62520	105861	3.67%	0.324%
30	11.659	818	820	822	rVB	43774	38178	1.32%	0.117%
31	11.865	836	838	839	rBV	51794	57407	1.99%	0.176%
32	11.888	839	840	843	rVV	163653	177665	6.15%	0.544%
33	11.968	846	847	852	rVB	82616	151427	5.25%	0.463%
34	12.059	852	855	857	rBV2	26701	49331	1.71%	0.151%

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35	12.162	861	864	865	rBV2	21942	46407	1.61%	0.142%
36	12.334	878	879	882	rBV3	19402	31621	1.10%	0.097%
37	12.414	882	886	887	rVV2	46261	64918	2.25%	0.199%
38	12.448	887	889	892	rVV	33409	46486	1.61%	0.142%
39	12.494	892	893	895	rVB	80800	76244	2.64%	0.233%
40	12.574	898	900	902	rBV	506876	502821	17.42%	1.539%
41	12.677	906	909	911	rVV3	76652	117587	4.07%	0.360%
42	12.722	911	913	915	rVV2	29337	48466	1.68%	0.148%
43	12.768	915	917	918	rVV	251376	270880	9.38%	0.829%
44	12.802	918	920	921	rVV	362803	432754	14.99%	1.324%
45	12.848	921	924	927	rVB3	43410	76648	2.66%	0.235%
46	12.951	930	933	935	rVV	1958598	2334825	80.89%	7.144%
47	13.019	936	939	943	rBV3	41478	80632	2.79%	0.247%
48	13.122	945	948	950	rVB	78516	135374	4.69%	0.414%
49	13.191	950	954	956	rBV2	75024	127415	4.41%	0.390%
50	13.225	956	957	961	rVV3	39740	74803	2.59%	0.229%
51	13.294	961	963	967	rVB4	54524	116458	4.03%	0.356%
52	13.351	967	968	970	rBV	40779	42923	1.49%	0.131%
53	13.477	977	979	980	rBV	147962	141177	4.89%	0.432%
54	13.522	980	983	984	rVV2	20211	38787	1.34%	0.119%
55	13.637	991	993	995	rVB	39099	47632	1.65%	0.146%
56	13.740	1000	1002	1004	rBV	49287	78738	2.73%	0.241%
57	13.842	1009	1011	1016	rVB2	208494	320547	11.10%	0.981%
58	13.991	1021	1024	1030	rVV2	806143	1293628	44.82%	3.958%
59	14.105	1031	1034	1035	rBV2	33923	60163	2.08%	0.184%
60	14.380	1056	1058	1060	rBV	43132	47538	1.65%	0.145%
61	14.414	1060	1061	1063	rBV2	32921	47991	1.66%	0.147%
62	14.837	1096	1098	1101	rBV	83224	87358	3.03%	0.267%
63	15.020	1111	1114	1117	rBV	224846	506654	17.55%	1.550%
64	15.305	1136	1139	1142	rVB	135083	226517	7.85%	0.693%
65	15.363	1142	1144	1147	rBV	212232	315514	10.93%	0.965%
66	15.420	1147	1149	1156	rVB	571993	920853	31.90%	2.818%
67	15.557	1159	1161	1166	rVB3	71044	116120	4.02%	0.355%
68	15.637	1166	1168	1170	rBV2	72252	126155	4.37%	0.386%
69	15.923	1191	1193	1200	rVB4	85843	218839	7.58%	0.670%
70	16.083	1205	1207	1215	rVB2	162860	341771	11.84%	1.046%
71	16.494	1241	1243	1250	rVB2	121412	258321	8.95%	0.790%

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72	16.814	1268	1271	1275	rVB2	83626	172459	5.97%	0.528%
73	17.043	1289	1291	1295	rVB	83216	153999	5.34%	0.471%
74	17.237	1305	1308	1312	rBV	69230	147880	5.12%	0.453%

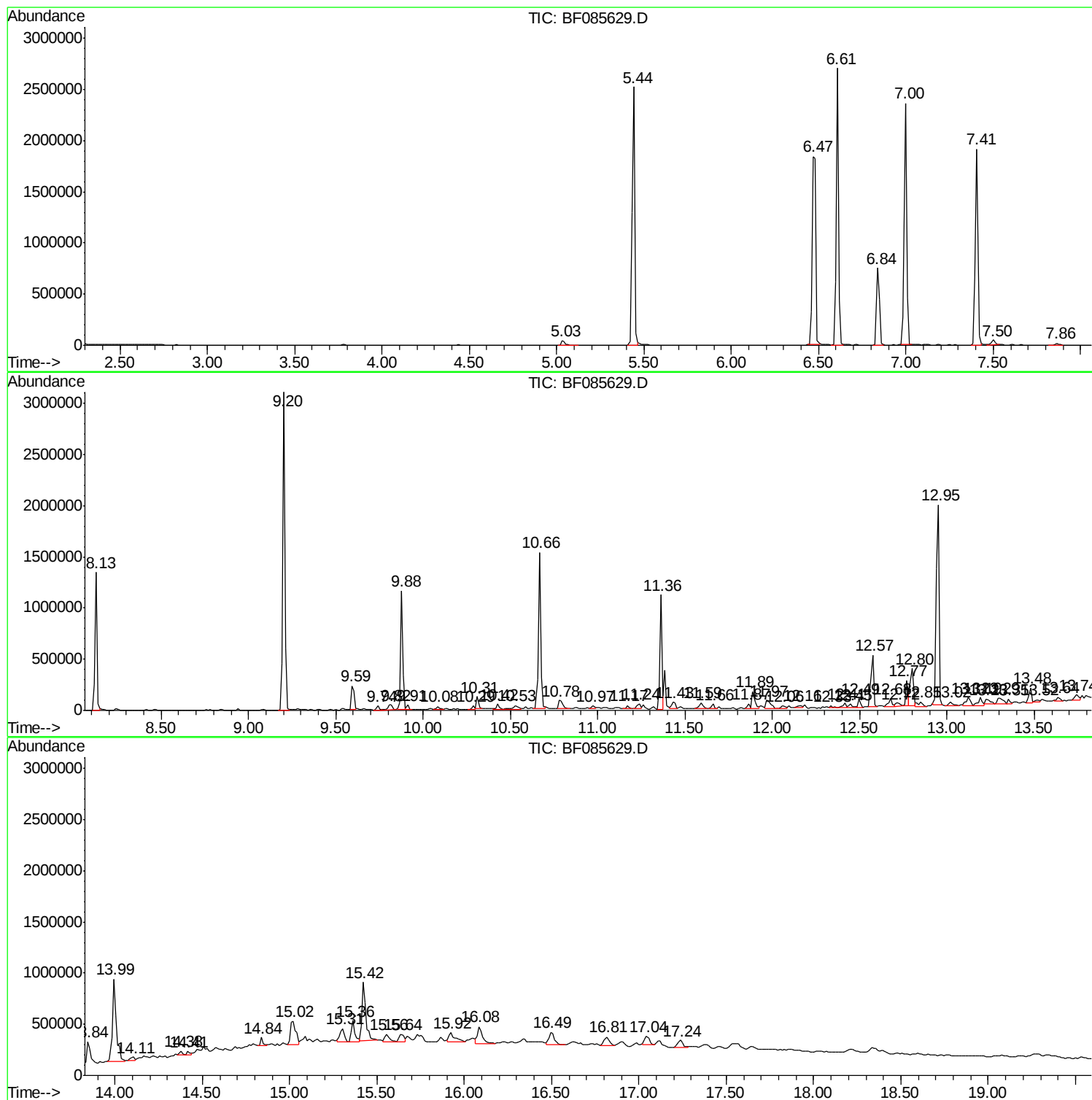
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Instrument :
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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ALS Vial : 14 Sample Multiplier: 1

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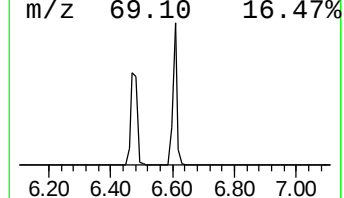
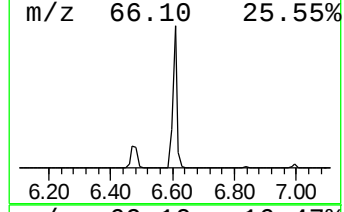
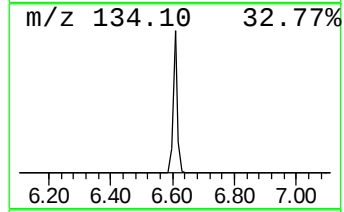
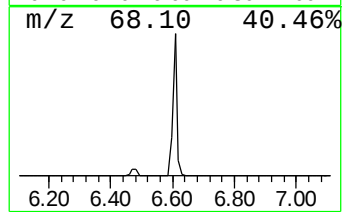
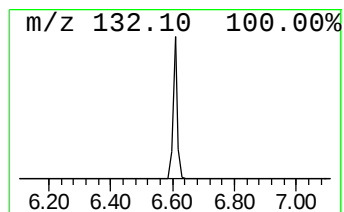
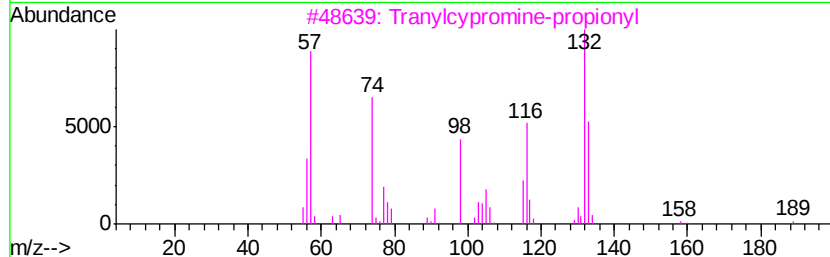
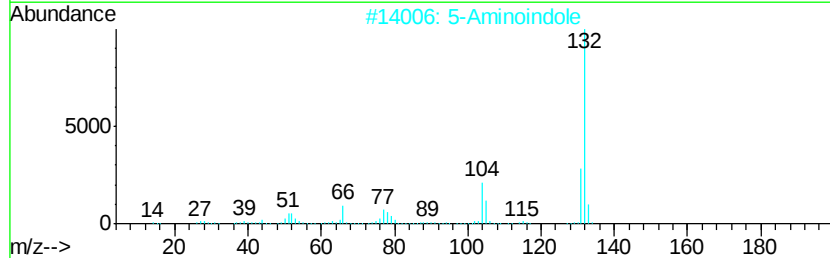
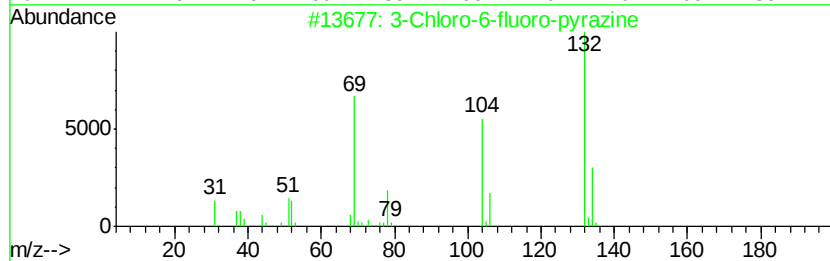
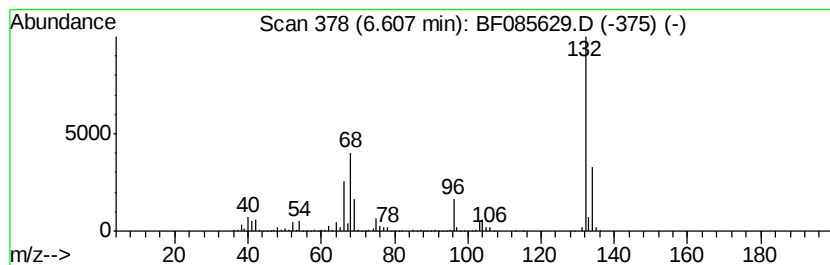
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.61	61.25 ng	2585730	1,4-Dichlorobenzene-d4	6.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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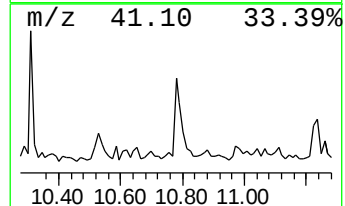
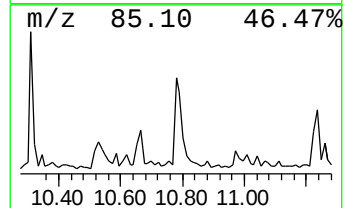
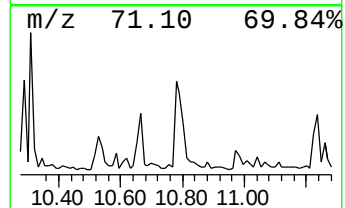
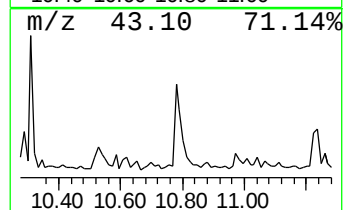
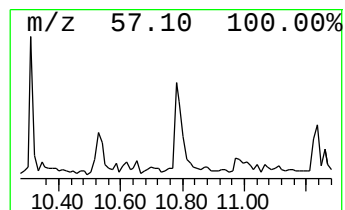
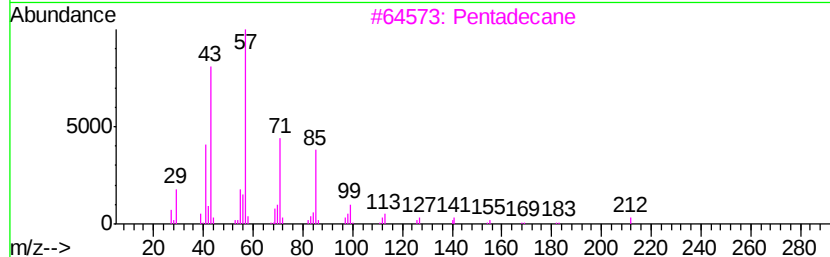
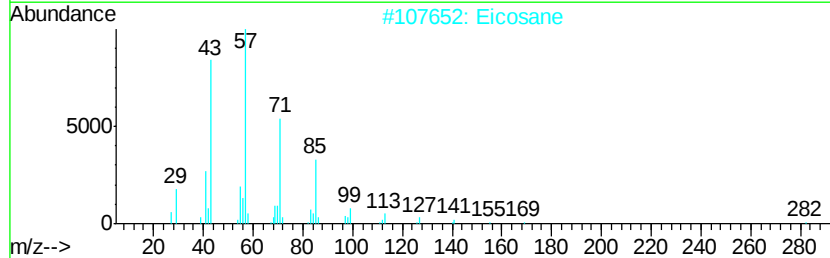
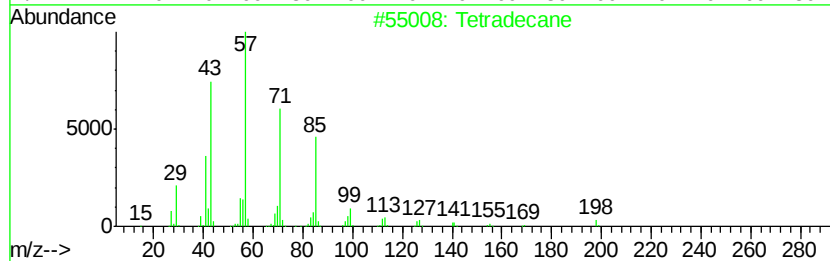
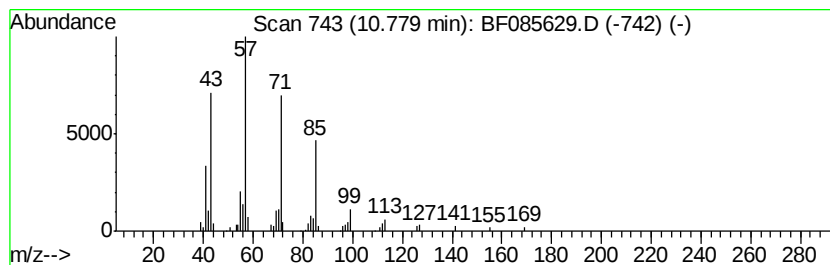
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Tetradecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.78	2.92 ng	144314	Phenanthrene-d10	11.36

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	91
2		Eicosane	282	C20H42	000112-95-8	91
3		Pentadecane	212	C15H32	000629-62-9	90
4		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	90
5		Hexadecane	226	C16H34	000544-76-3	80



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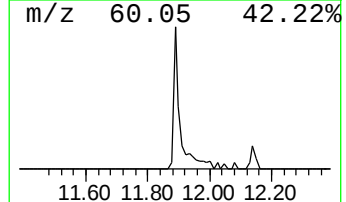
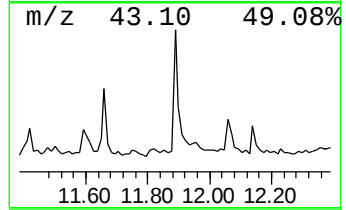
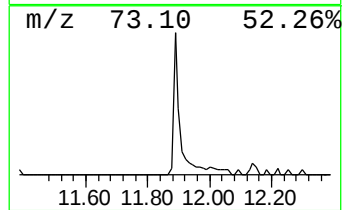
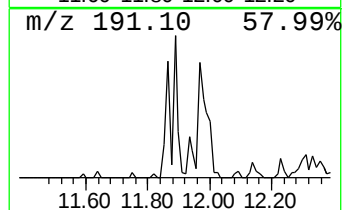
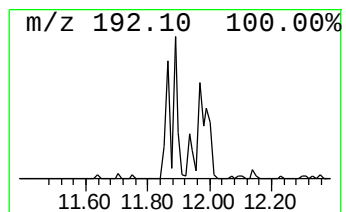
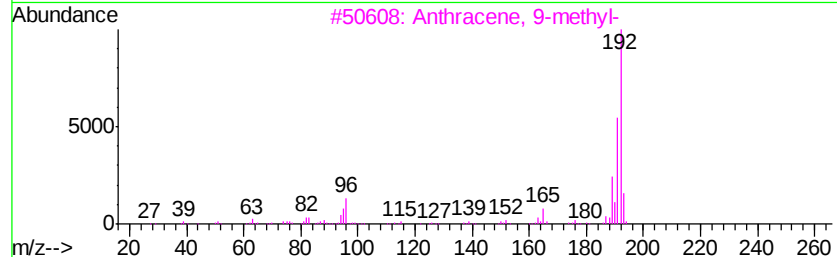
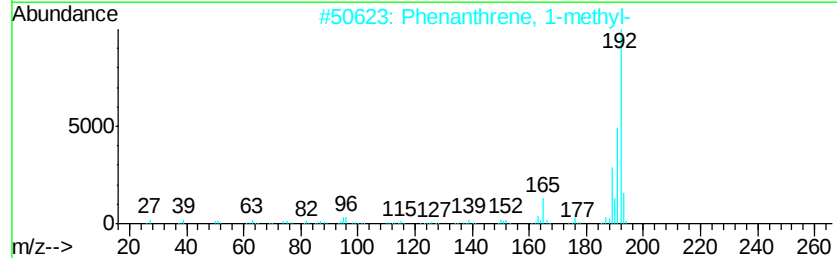
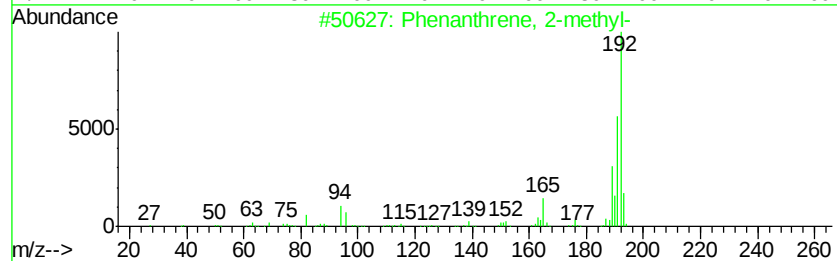
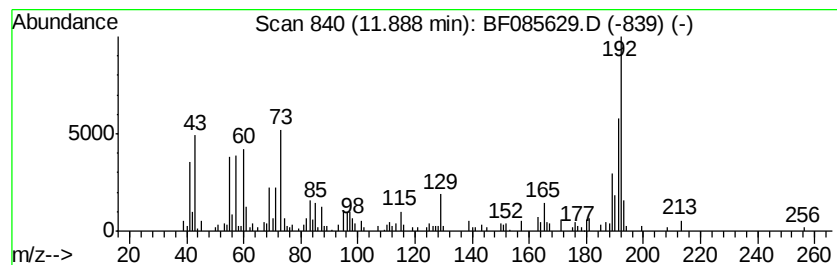
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.89	3.59 ng	177665	Phenanthrene-d10	11.36

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
3		Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	91
5		Anthracene, 1-methyl-	192	C15H12	000610-48-0	83



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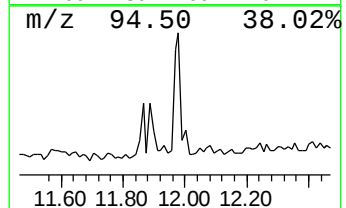
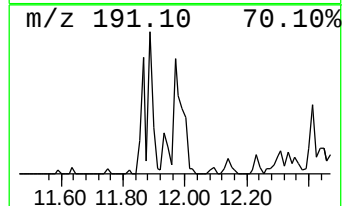
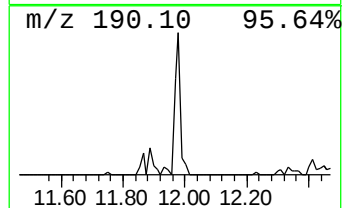
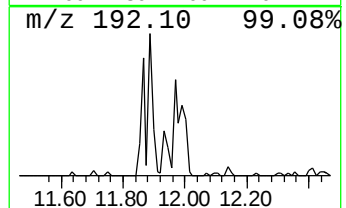
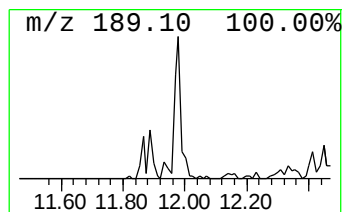
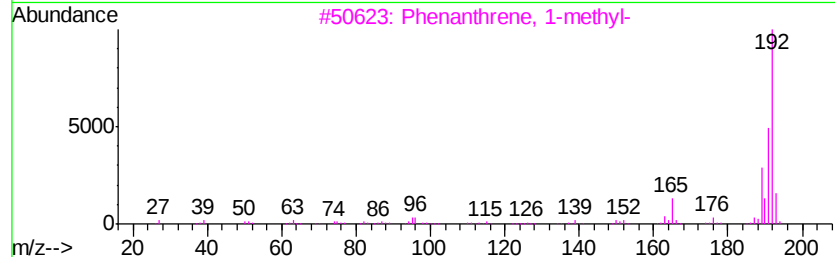
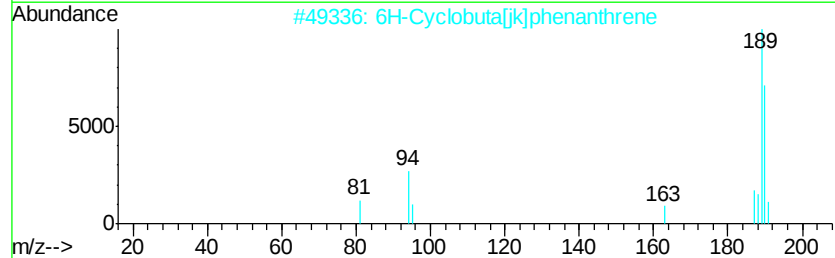
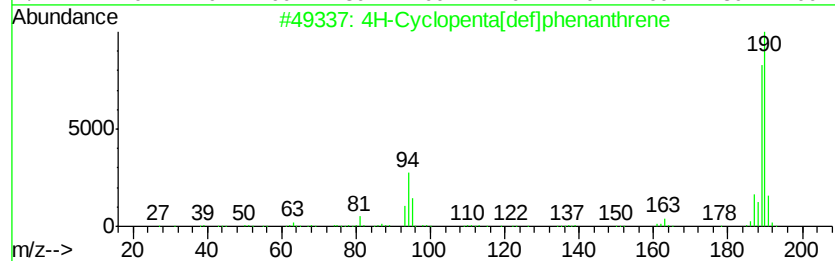
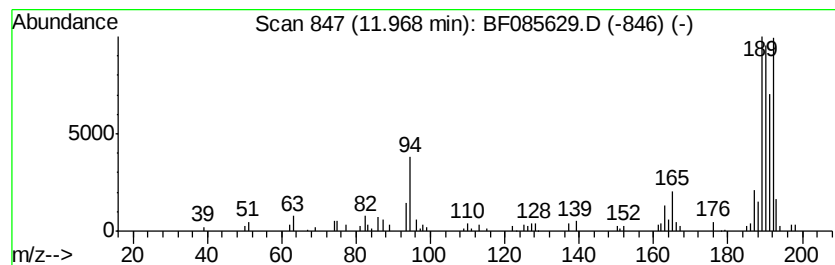
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.97	3.06 ng	151427	Phenanthrene-d10	11.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	53
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	49
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	42
4	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	42
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	42



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032116\
Data File : BF085629.D
Acq On : 21 Mar 2016 18:34
Operator : UM/SJ
Sample : H1803-01
Misc :
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ClientSampleId :
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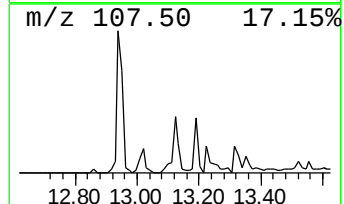
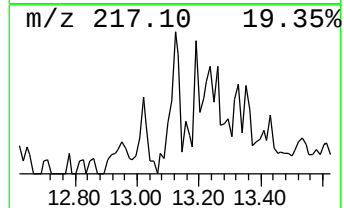
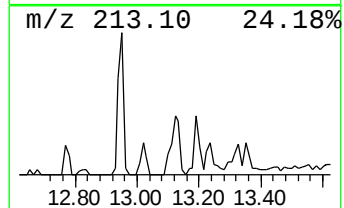
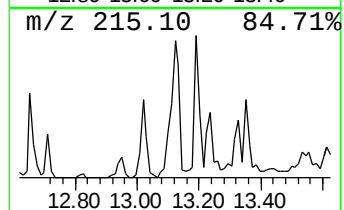
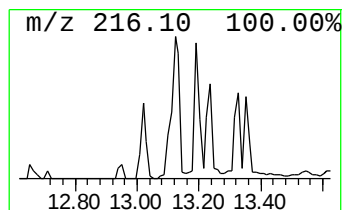
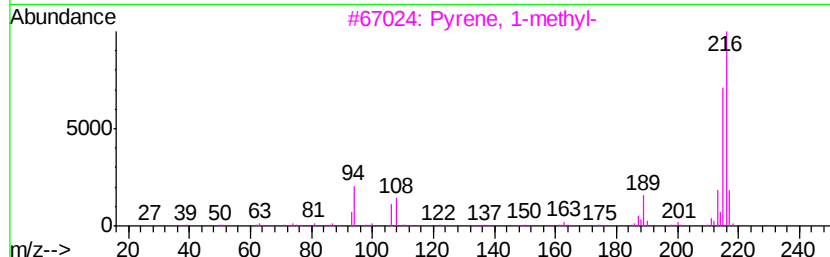
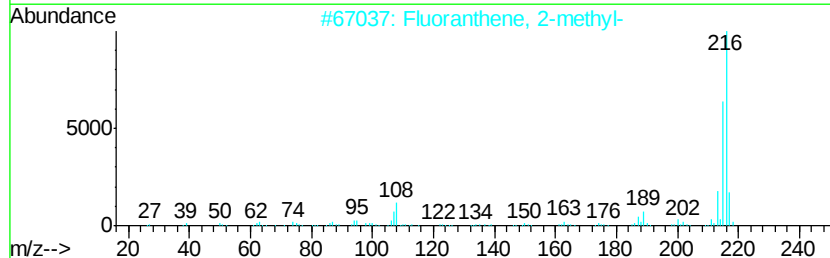
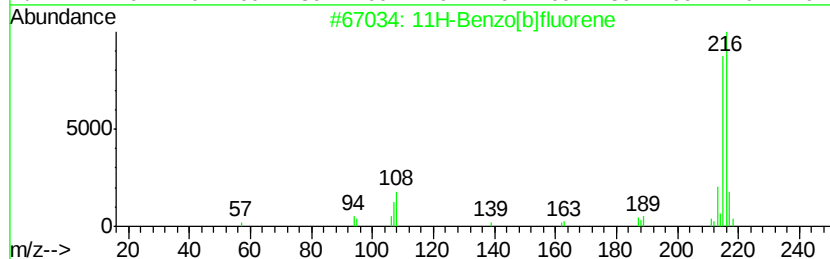
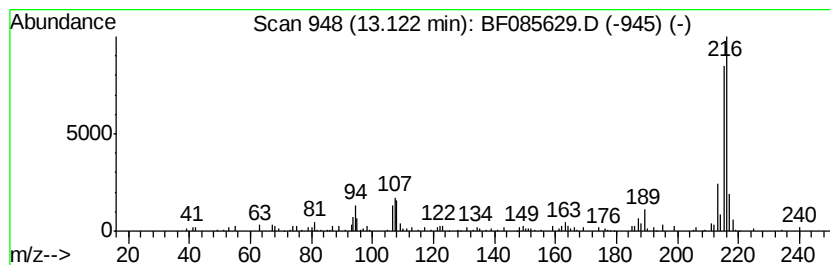
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 11H-Benzo[b]fluorene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	2.09 ng	135374	Chrysene-d12	13.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	95
2		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	94
3		Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
4		Pyrene, 2-methyl-	216	C17H12	003442-78-2	93
5		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	83



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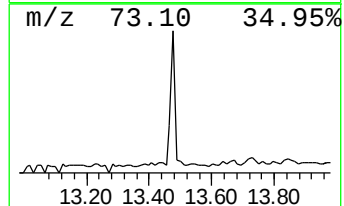
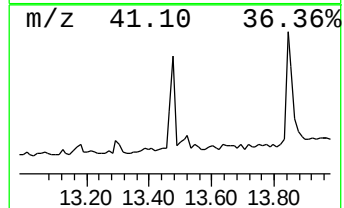
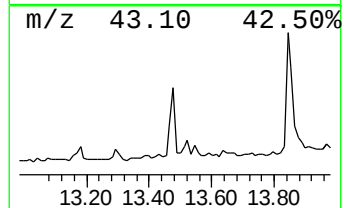
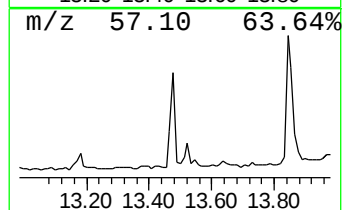
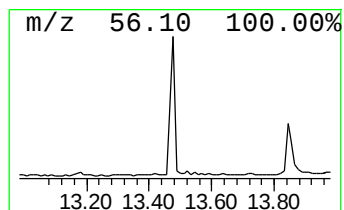
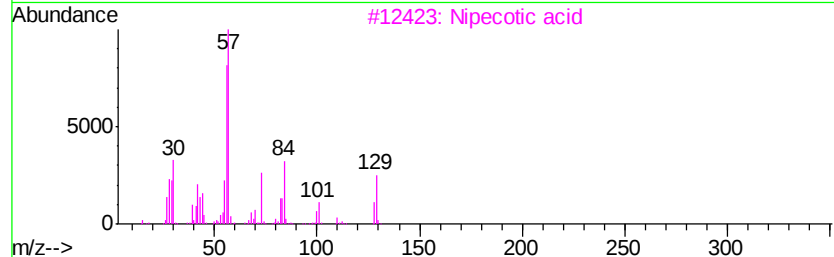
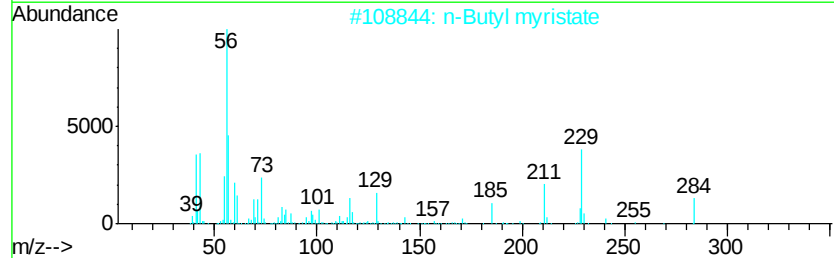
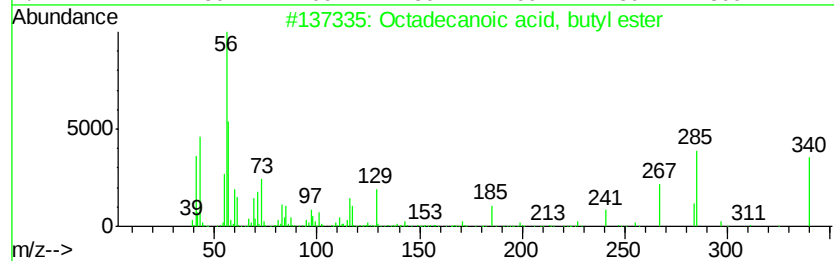
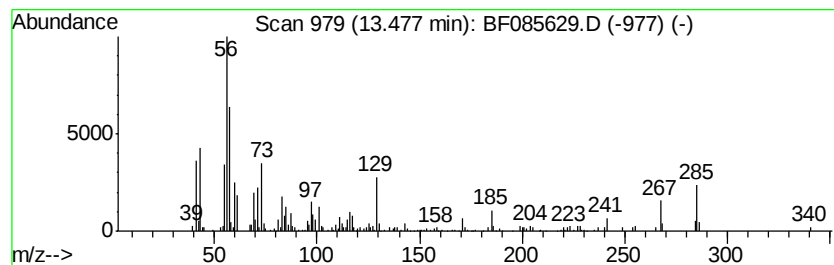
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Octadecanoic acid, butyl ester Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.48	2.18 ng	141177	Chrysene-d12	13.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	87
2		n-Butyl myristate	284	C18H36O2	000110-36-1	58
3		Nipecotic acid	129	C6H11NO2	000498-95-3	47
4		Octadecanoic acid, 2-methylpropy...	340	C22H44O2	000646-13-9	43
5		1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	32



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032116\
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Misc :
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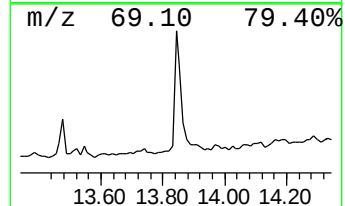
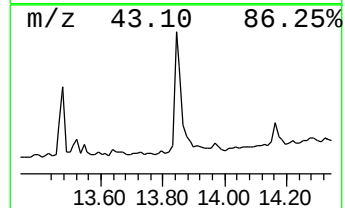
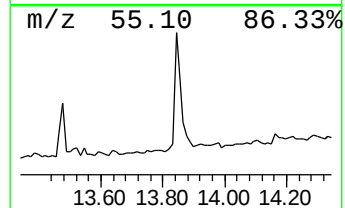
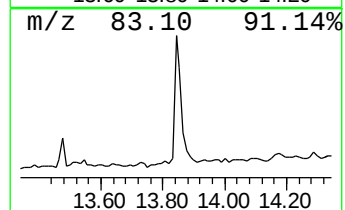
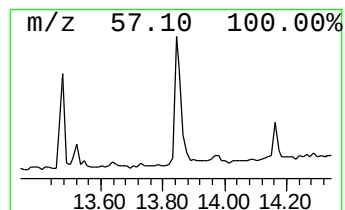
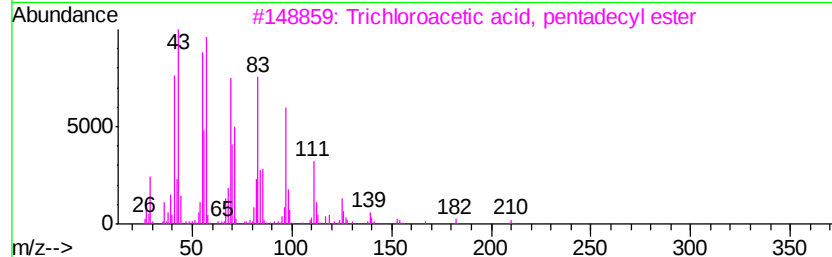
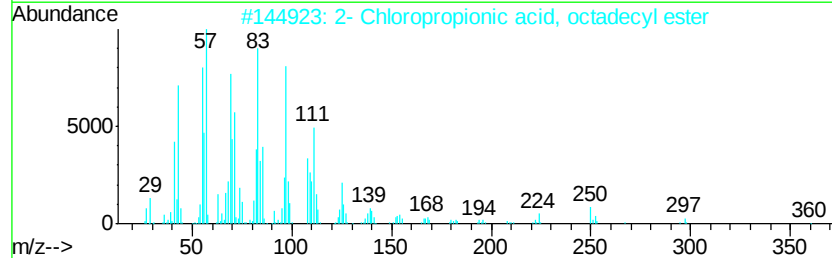
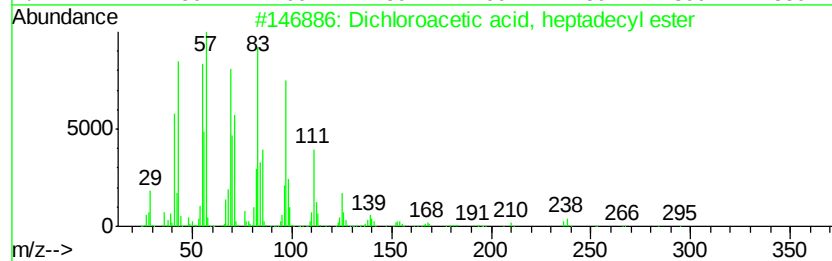
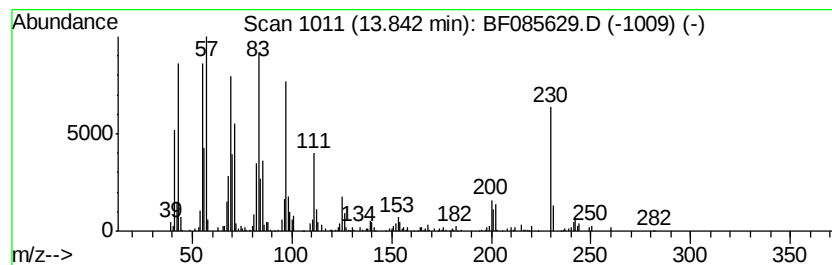
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Dichloroacetic acid, heptad... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.84	4.96 ng	320547	Chrysene-d12	13.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	93
2		2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	93
3		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	93
4		1-Pentadecanethiol	244	C15H32S	025276-70-4	91
5		Trifluoroacetic acid, n-octadecy...	366	C20H37F3O2	079392-43-1	90



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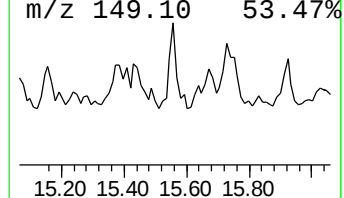
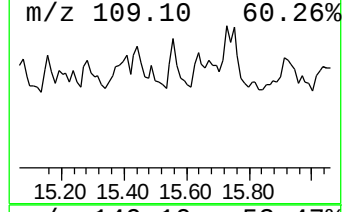
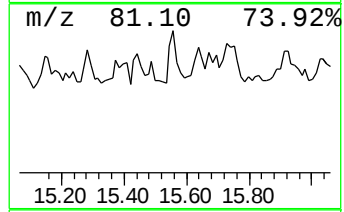
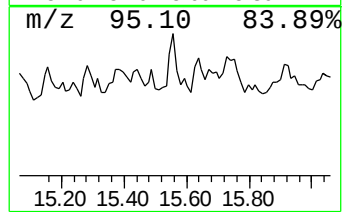
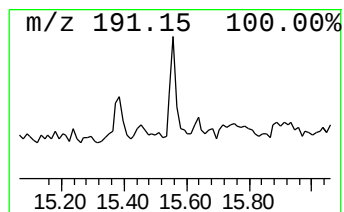
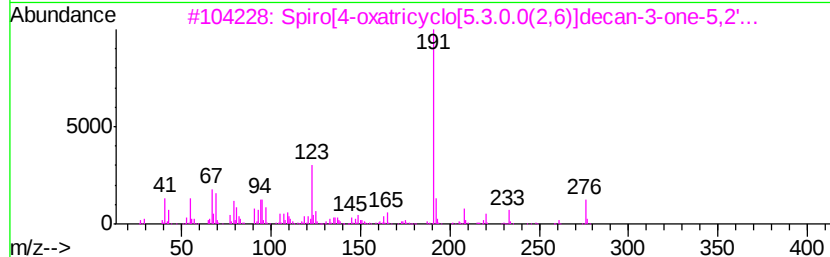
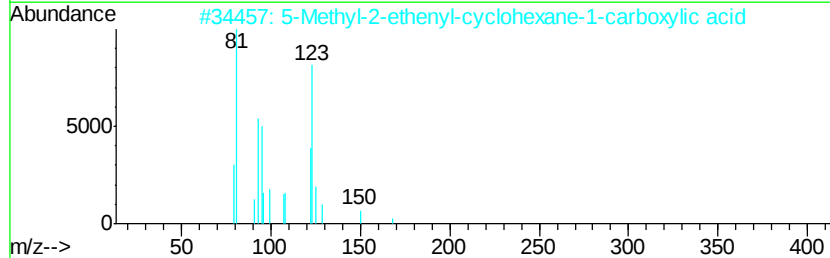
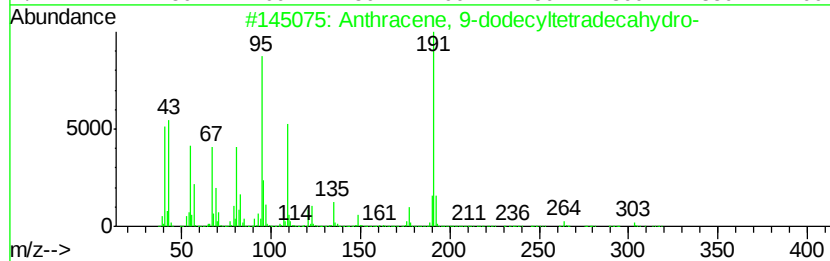
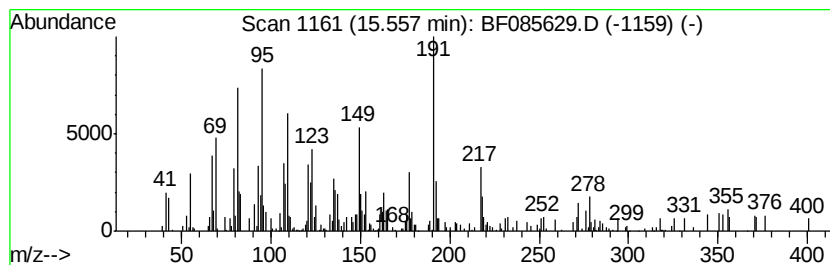
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown15.56 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.56	2.52 ng	116120	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 9-dodecyltetradecahy...	360	C26H48	055401-75-7	43
2		5-Methyl-2-ethenyl-cyclohexane-1...	168	C10H16O2	1000144-53-6	38
3		Spiro[4-oxatricyclo[5.3.0.0(2,6)...	276	C18H28O2	1000156-82-9	25
4		Naphthalene, decahydro-1,6-dimet...	208	C15H28	034315-85-0	25
5		2-Cyanomethylcyclohexanecarboxyl...	181	C10H15NO2	1000186-80-0	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032116\
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Misc :
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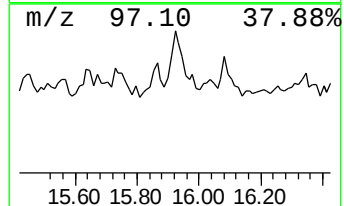
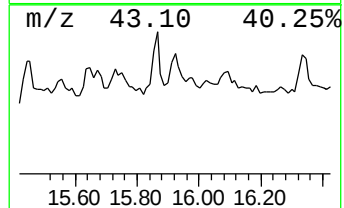
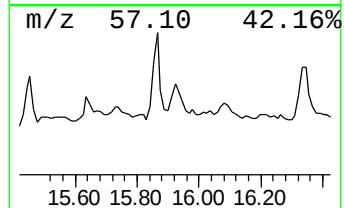
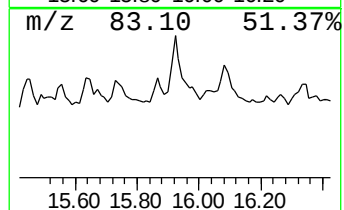
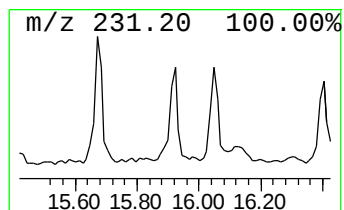
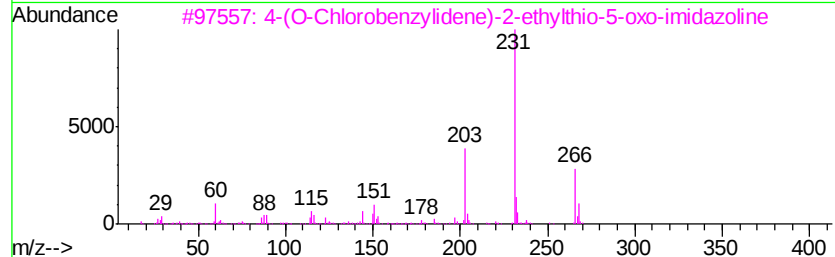
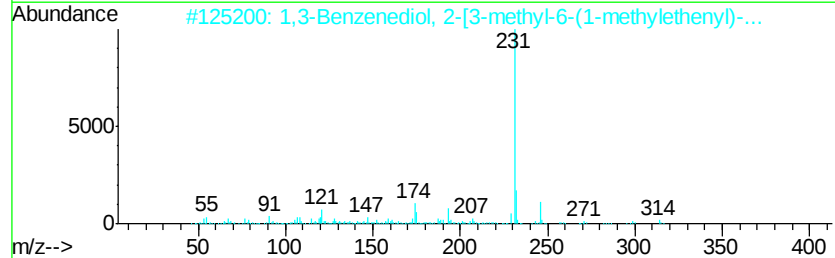
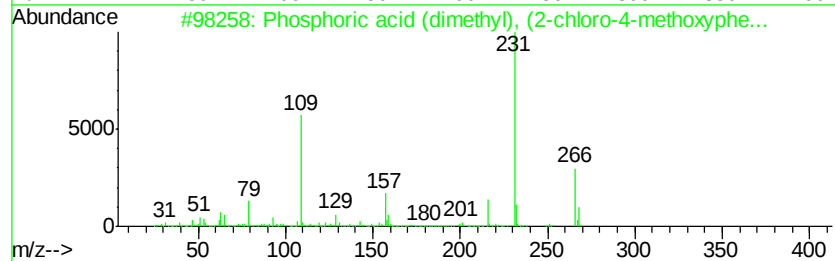
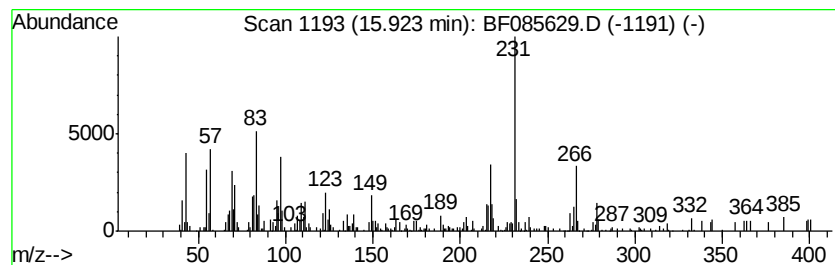
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Phosphoric acid (dimethyl),... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.92	4.75 ng	218839	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phosphoric acid (dimethyl), (2-c...	266	C9H12ClO5P	094420-73-2	50
2		1,3-Benzenediol, 2-[3-methyl-6-(...	314	C21H30O2	013956-29-1	50
3		4-(O-Chlorobenzylidene)-2-ethylt...	266	C12H11ClN2OS	033449-97-7	49
4		1,3-Dimethyl-6-(trifluoromethyl)...	231	C9H8F3N3O	332867-83-1	49
5		5-Trimethylsilyloxy-2-indolecarb...	321	C15H23N03Si2	1000292-72-0	47



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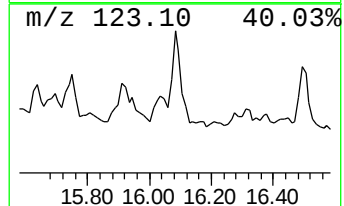
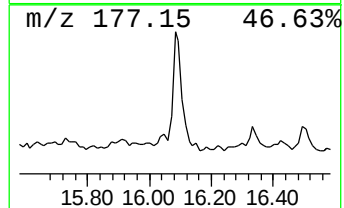
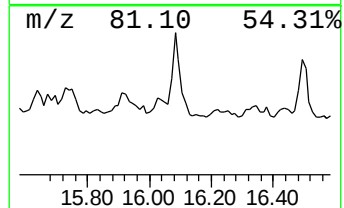
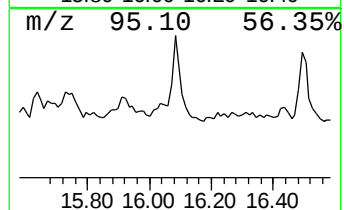
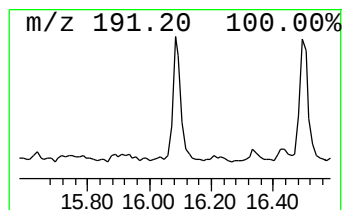
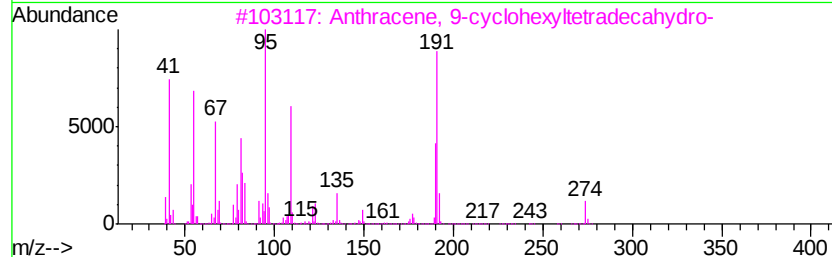
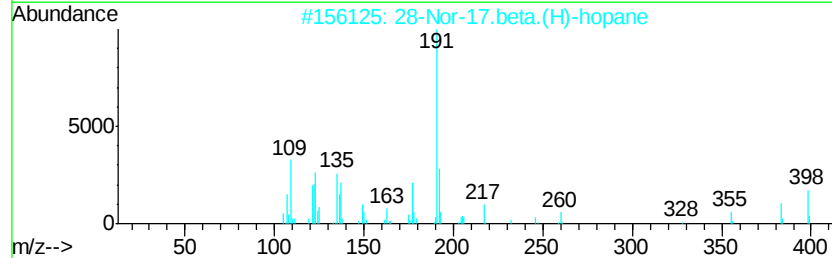
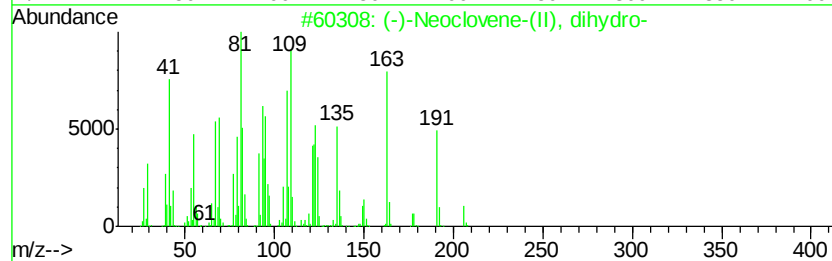
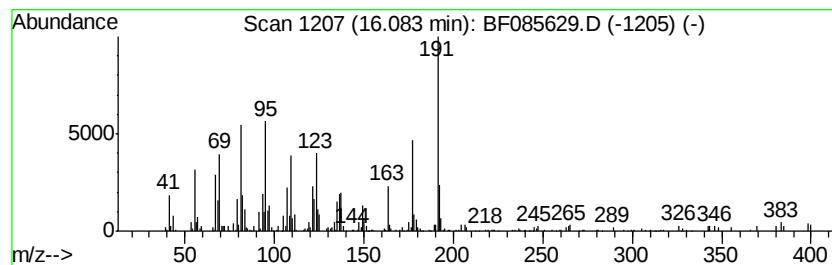
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 (-)-Neoclovene-(II), dihydro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.08	7.42 ng	341771	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(-)-Neoclovene-(II), dihydro-	206	C15H26	1000152-83-6	90
2		28-Nor-17.beta.(H)-hopane	398	C29H50	036728-72-0	58
3		Anthracene, 9-cyclohexyltetradec...	274	C20H34	055255-70-4	43
4		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	35
5		2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	32



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032116\
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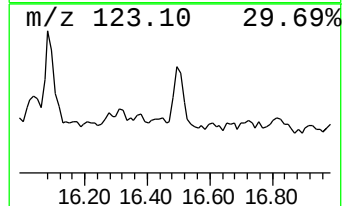
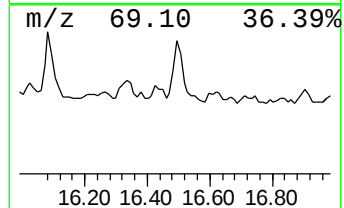
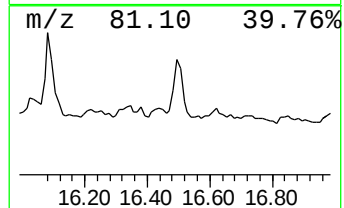
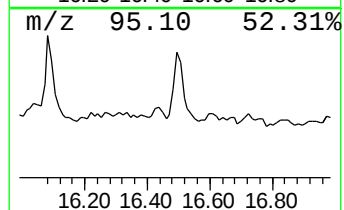
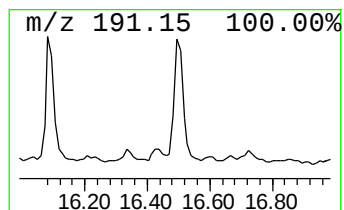
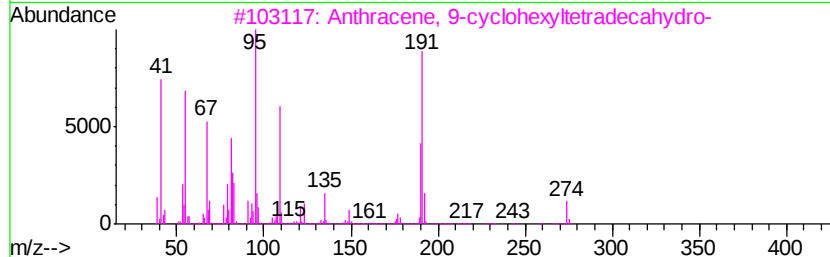
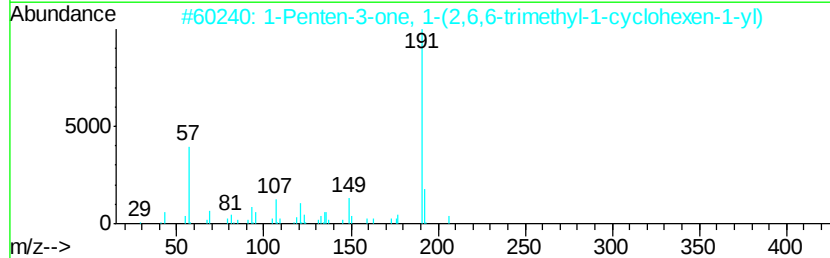
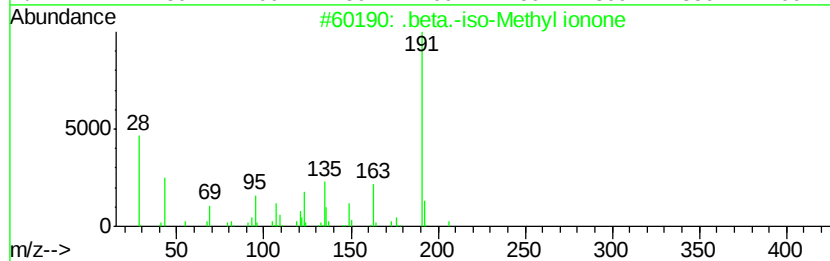
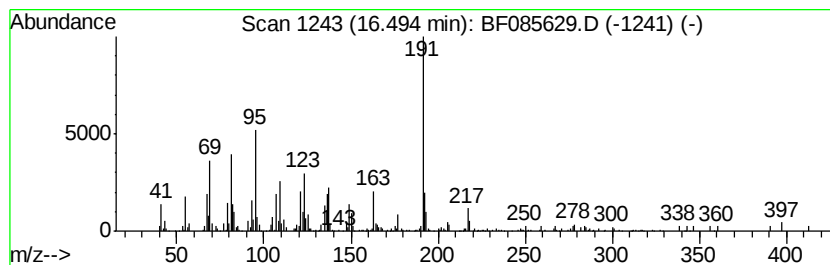
Instrument :
BNA_F
ClientSampleId :
SP-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 .beta.-iso-Methyl ionone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.49	5.61 ng	258321	Perylene-d12	15.42
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		.beta.-iso-Methyl ionone	206 C14H22O	1000285-40-2 59
2		1-Penten-3-one, 1-(2,6,6-trimeth...	206 C14H22O	000127-43-5 35
3		Anthracene, 9-cyclohexyltetradec...	274 C20H34	055255-70-4 35
4		2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206 C14H22O	1000195-40-9 32
5		5-Hexen-1-one, 1-(1H-imidazol-4-...	192 C11H16N2O	069393-41-5 30



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032116\
Data File : BF085629.D
Acq On : 21 Mar 2016 18:34
Operator : UM/SJ
Sample : H1803-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-01

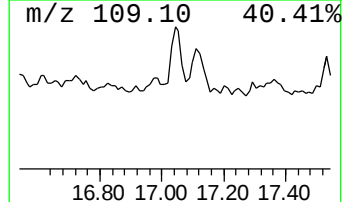
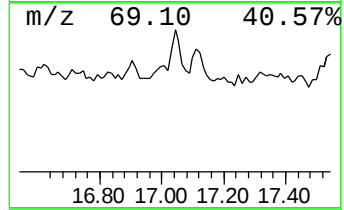
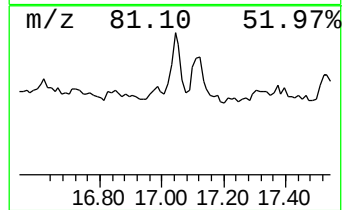
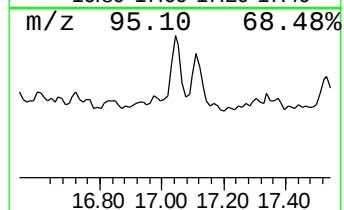
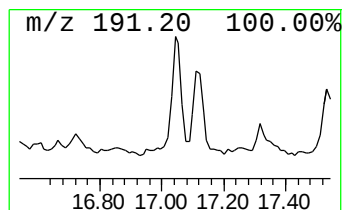
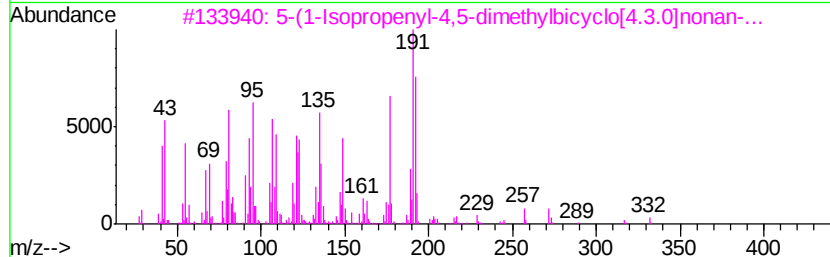
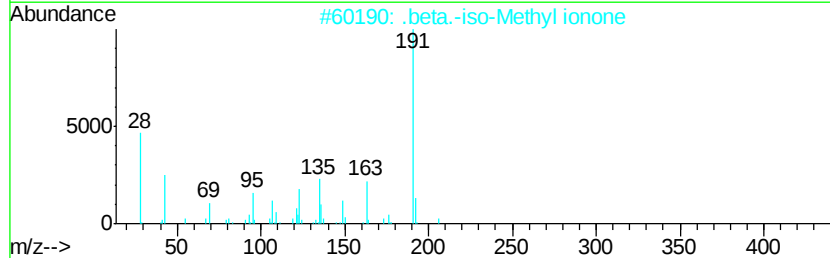
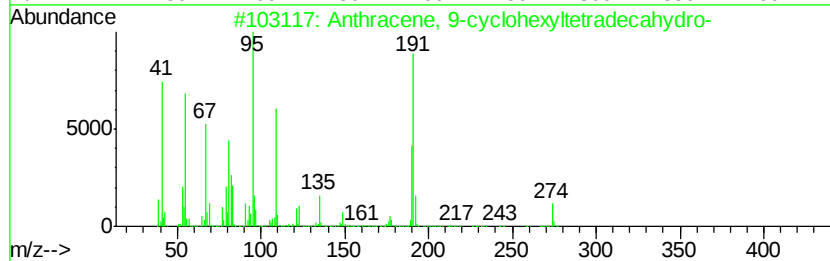
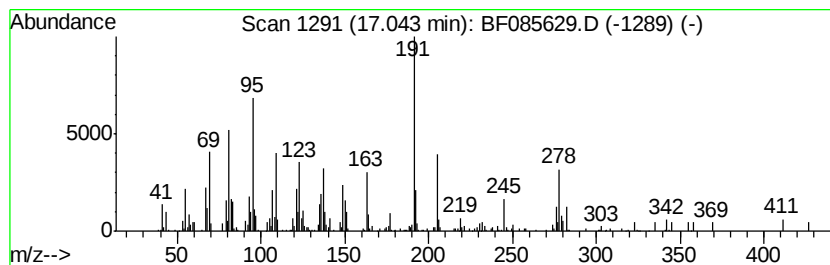
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Anthracene, 9-cyclohexyltet... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.04	3.34 ng	153999	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 9-cyclohexyltetradec...	274	C20H34	055255-70-4	60
2		.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	38
3		5-(1-Isopropenyl-4,5-dimethylbic...	332	C22H36O2	1000195-69-8	38
4		3-Octyne, 2,2,7-trimethyl-	152	C11H20	055402-13-6	35
5		Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	020536-40-7	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032116\
Data File : BF085629.D
Acq On : 21 Mar 2016 18:34
Operator : UM/SJ
Sample : H1803-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.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
unknown6.61	6.61	61.3	ng	2585730	1	6.84	844274	20.0
Tetradecane	10.78	2.9	ng	144314	4	11.36	988629	20.0
Phenanthrene, 2-m...	11.89	3.6	ng	177665	4	11.36	988629	20.0
4H-Cyclopenta[def...	11.97	3.1	ng	151427	4	11.36	988629	20.0
11H-Benzo[b]fluorene	13.12	2.1	ng	135374	5	13.99	1293630	20.0
Octadecanoic acid...	13.48	2.2	ng	141177	5	13.99	1293630	20.0
Dichloroacetic ac...	13.84	5.0	ng	320547	5	13.99	1293630	20.0
unknown15.56	15.56	2.5	ng	116120	6	15.42	920853	20.0
Phosphoric acid (...)	15.92	4.8	ng	218839	6	15.42	920853	20.0
(-)-Neoclovene-(I...	16.08	7.4	ng	341771	6	15.42	920853	20.0
.beta.-iso-Methyl...	16.49	5.6	ng	258321	6	15.42	920853	20.0
Anthracene, 9-cyc...	17.04	3.3	ng	153999	6	15.42	920853	20.0