

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120216\
 Data File : BF091406.D
 Acq On : 2 Dec 2016 23:06
 Operator : UM/SJ
 Sample : H5829-03
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WESTWALL-120116

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-BF112916.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	2.458	12	15	25	rVB	188909	322010	7.04%	0.629%
2	3.144	73	75	90	rVB4	75804	197102	4.31%	0.385%
3	3.807	128	133	135	rBV	70563	122795	2.69%	0.240%
4	4.127	160	161	167	rVB3	49456	86331	1.89%	0.169%
5	4.310	175	177	182	rBV	32817	48232	1.06%	0.094%
6	4.447	185	189	192	rVB2	27203	62480	1.37%	0.122%
7	4.516	192	195	198	rBV2	77832	131810	2.88%	0.258%
8	4.687	209	210	216	rBV	28462	47168	1.03%	0.092%
9	5.041	238	241	244	rBV	1037079	1652288	36.14%	3.229%
10	5.453	274	277	280	rVB	2824790	3876178	84.79%	7.574%
11	5.693	296	298	302	rVB3	28843	55993	1.22%	0.109%
12	5.761	302	304	306	rBV	154796	142805	3.12%	0.279%
13	5.796	306	307	311	rVB	95615	81159	1.78%	0.159%
14	5.876	311	314	316	rBV2	39440	51958	1.14%	0.102%
15	6.013	323	326	328	rVB	44148	51869	1.13%	0.101%
16	6.104	333	334	338	rVB2	90195	138588	3.03%	0.271%
17	6.299	348	351	353	rBV	49771	72059	1.58%	0.141%
18	6.356	353	356	357	rVV2	53232	79934	1.75%	0.156%
19	6.390	357	359	362	rVV3	80394	127337	2.79%	0.249%
20	6.493	362	368	370	rVV	3702257	4571646	100.00%	8.933%
21	6.630	377	380	382	rVB	3428795	4052623	88.65%	7.919%
22	6.699	384	386	390	rVB	285792	325828	7.13%	0.637%
23	6.859	398	400	402	rBV	1723024	1463665	32.02%	2.860%
24	6.927	404	406	407	rBV2	71753	84057	1.84%	0.164%
25	7.019	411	414	416	rVV	3199843	3147790	68.85%	6.151%
26	7.144	418	425	426	rBV2	83137	132075	2.89%	0.258%
27	7.213	429	431	433	rVB	50314	54833	1.20%	0.107%
28	7.259	433	435	437	rVB	58569	58071	1.27%	0.113%
29	7.419	446	449	452	rBV	2327298	2268787	49.63%	4.433%
30	7.464	452	453	455	rVB	148893	113073	2.47%	0.221%
31	7.510	455	457	459	rVB2	68205	79340	1.74%	0.155%
32	8.139	510	512	515	rBV	1453431	1798400	39.34%	3.514%
33	8.859	572	575	576	rVB	61188	60825	1.33%	0.119%
34	8.950	581	583	588	rVB	44771	50293	1.10%	0.098%

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35	9.225	604	607	609 rBV	4928789	4472842	97.84%	8.740%
36	9.305	611	614	617 rVV2	42419	66624	1.46%	0.130%
37	9.556	634	636	637 rBV	60771	52155	1.14%	0.102%
38	9.613	639	641	645 rVB	784432	681822	14.91%	1.332%
39	9.762	652	654	656 rBV	72590	97056	2.12%	0.190%
40	9.842	659	661	663 rVB	68822	79817	1.75%	0.156%
41	9.899	663	666	668 rBV	2371621	2040010	44.62%	3.986%
42	10.093	682	683	685 rVB	38965	53120	1.16%	0.104%
43	10.310	697	702	703 rBV	59865	79330	1.74%	0.155%
44	10.333	703	704	707 rVB2	132695	128662	2.81%	0.251%
45	10.436	712	713	717 rVB3	37212	65760	1.44%	0.129%
46	10.562	720	724	725 rBV	54158	83659	1.83%	0.163%
47	10.688	732	735	737 rBV	2107621	2613888	57.18%	5.108%
48	10.813	744	746	749 rBV	163385	270465	5.92%	0.529%
49	11.145	770	775	777 rBV4	23993	73710	1.61%	0.144%
50	11.259	783	785	786 rBV	146118	128923	2.82%	0.252%
51	11.282	786	787	790 rVB	47020	58430	1.28%	0.114%
52	11.385	793	796	799 rBV	1390226	1955410	42.77%	3.821%
53	11.453	799	802	803 rVB	71739	78088	1.71%	0.153%
54	11.682	820	822	823 rBV	124251	96912	2.12%	0.189%
55	11.876	837	839	840 rBV	91177	89430	1.96%	0.175%
56	11.911	840	842	844 rVV	546882	572533	12.52%	1.119%
57	11.956	844	846	847 rVV2	52889	77076	1.69%	0.151%
58	11.991	847	849	854 rVB2	141897	256112	5.60%	0.500%
59	12.094	854	858	861 rBV2	61529	128351	2.81%	0.251%
60	12.174	863	865	870 rVB3	62984	124236	2.72%	0.243%
61	12.322	875	878	880 rBV3	44600	71662	1.57%	0.140%
62	12.356	880	881	884 rVB	63146	91182	1.99%	0.178%
63	12.425	885	887	889 rBV	116493	144711	3.17%	0.283%
64	12.482	889	892	893 rVB3	58180	98684	2.16%	0.193%
65	12.516	893	895	897 rVB3	49511	73584	1.61%	0.144%
66	12.585	899	901	903 rBV	391145	364986	7.98%	0.713%
67	12.631	903	905	908 rVB4	30846	66797	1.46%	0.131%
68	12.699	908	911	914 rBV3	177382	258889	5.66%	0.506%
69	12.814	918	921	923 rBV	476736	523224	11.44%	1.022%
70	12.871	925	926	928 rVB	50981	63160	1.38%	0.123%
71	12.962	932	934	937 rVB	2688679	3042944	66.56%	5.946%

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72	13.031	938	940	944	rBV3	74459	122085	2.67%	0.239%
73	13.145	946	950	951	rVB	92878	157672	3.45%	0.308%
74	13.179	951	953	954	rBV	64854	64592	1.41%	0.126%
75	13.214	954	956	957	rVB2	61586	71247	1.56%	0.139%
76	13.248	957	959	961	rBV	116632	120905	2.64%	0.236%
77	13.339	965	967	968	rBV	94558	74211	1.62%	0.145%
78	13.374	968	970	972	rVB	63352	64921	1.42%	0.127%
79	13.454	975	977	982	rBV	259435	563728	12.33%	1.102%
80	13.545	982	985	988	rVB3	52343	132085	2.89%	0.258%
81	13.648	992	994	997	rVB	60623	82742	1.81%	0.162%
82	13.762	1001	1004	1005	rBV2	60069	108082	2.36%	0.211%
83	13.865	1011	1013	1017	rVB2	350664	383297	8.38%	0.749%
84	14.014	1023	1026	1027	rBV	1644053	1745261	38.18%	3.410%
85	14.402	1058	1060	1061	rVB	60314	67167	1.47%	0.131%
86	14.437	1061	1063	1064	rVB	49709	55428	1.21%	0.108%
87	14.494	1066	1068	1069	rBV2	44708	49758	1.09%	0.097%
88	15.031	1112	1115	1119	rBV	170976	351950	7.70%	0.688%
89	15.134	1122	1124	1128	rVB2	65273	119331	2.61%	0.233%
90	15.328	1138	1141	1143	rVB	119361	171106	3.74%	0.334%
91	15.385	1143	1146	1148	rBV	123626	168931	3.70%	0.330%
92	15.442	1148	1151	1154	rBV	828025	1196214	26.17%	2.338%
93	15.968	1194	1197	1204	rVB7	94976	239751	5.24%	0.468%
94	16.837	1271	1273	1277	rVB2	57160	107560	2.35%	0.210%
95	17.260	1307	1310	1317	rVB	51569	123100	2.69%	0.241%

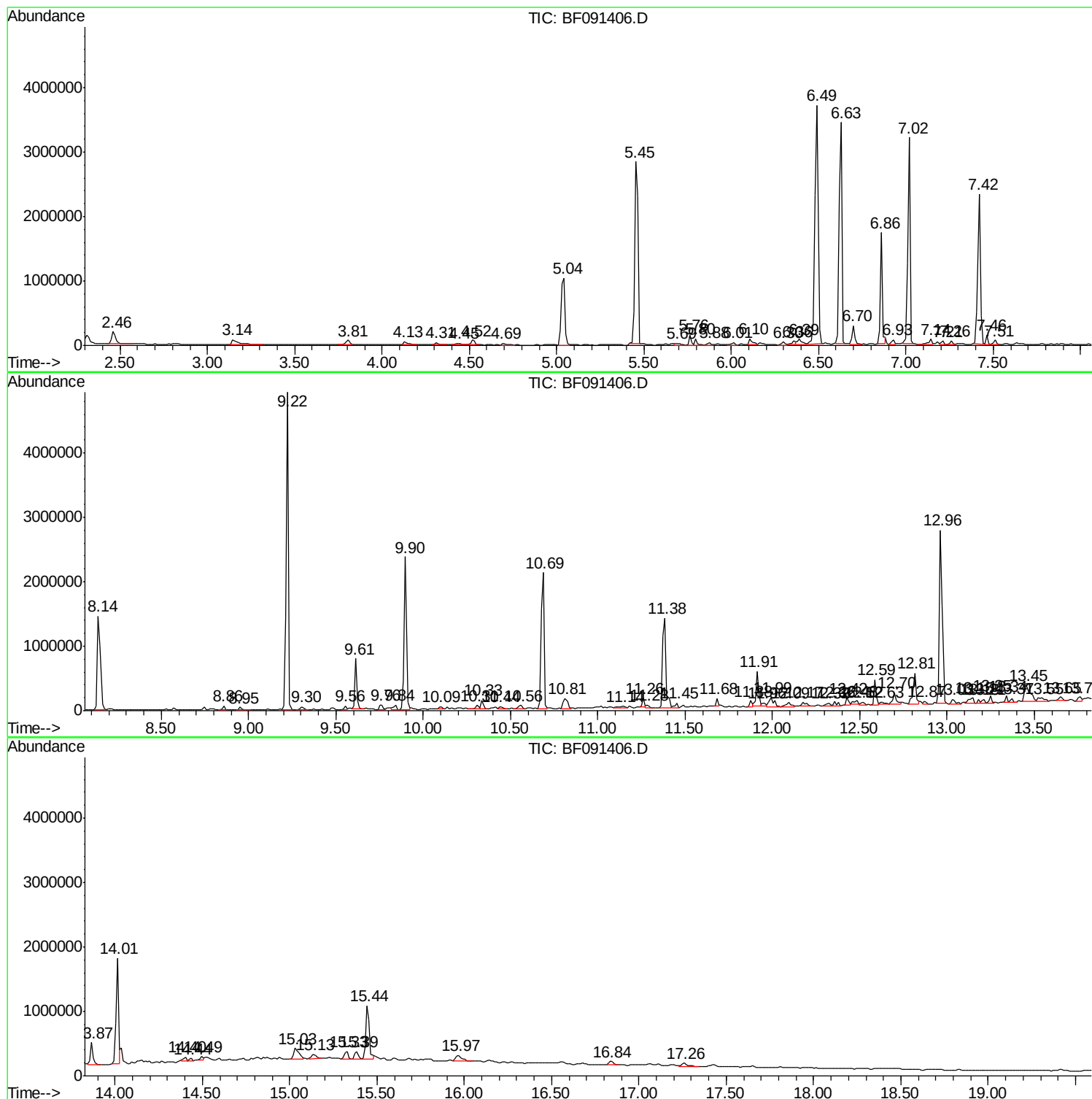
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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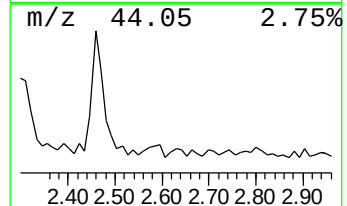
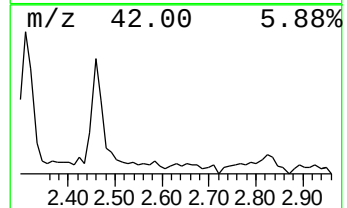
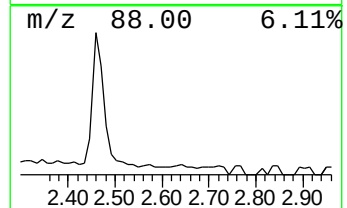
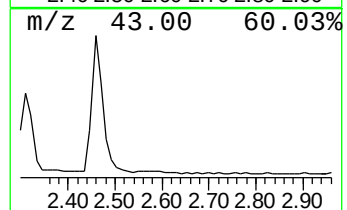
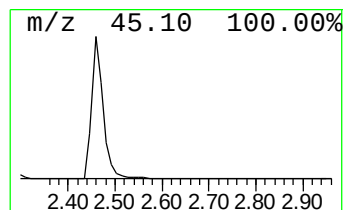
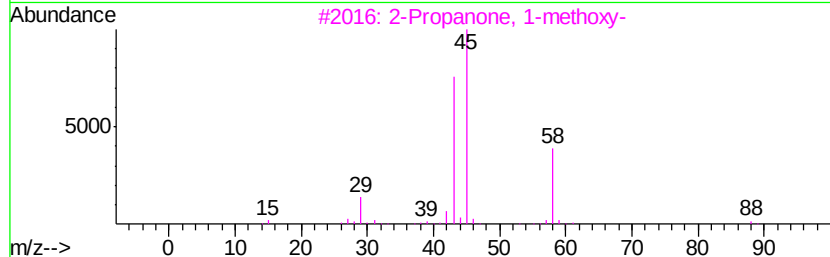
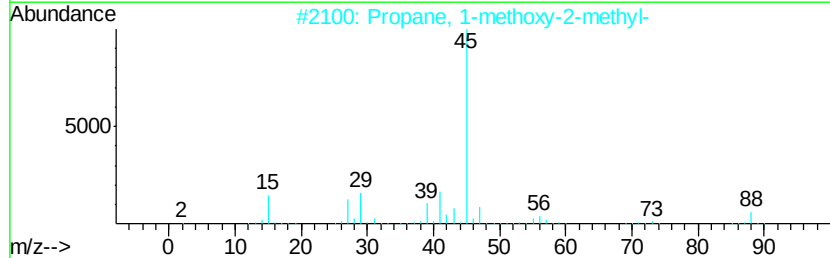
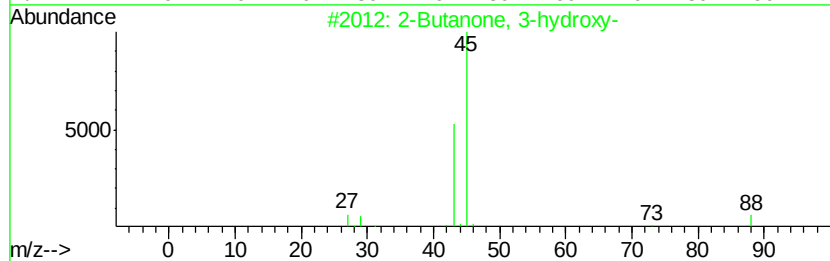
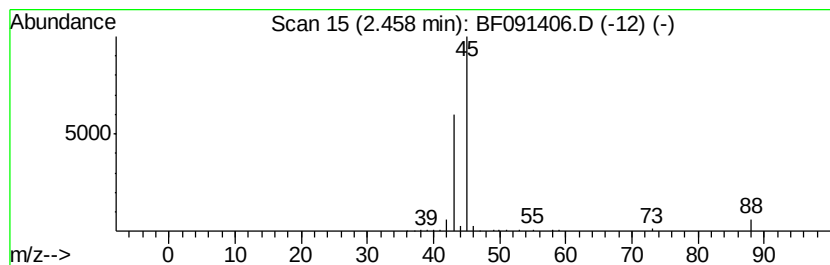
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Butanone, 3-hydroxy- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.46	4.40 ng	322010	1,4-Dichlorobenzene-d4	6.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Butanone, 3-hydroxy-	88	C4H8O2	000513-86-0	86
2			Propane, 1-methoxy-2-methyl-	88	C5H12O	000625-44-5	56
3			2-Propanone, 1-methoxy-	88	C4H8O2	005878-19-3	43
4			Propane, 2-ethoxy-	88	C5H12O	000625-54-7	9
5			Ethanol, 2-(vinylloxy)-	88	C4H8O2	000764-48-7	9



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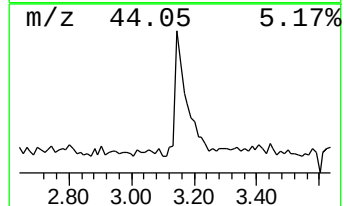
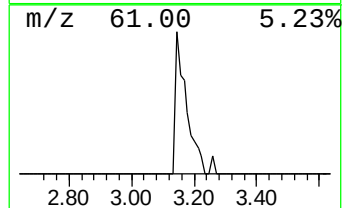
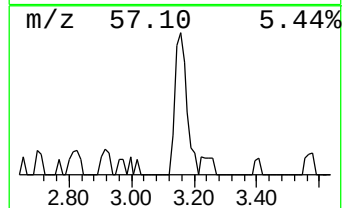
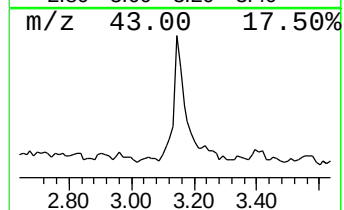
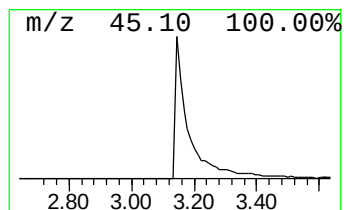
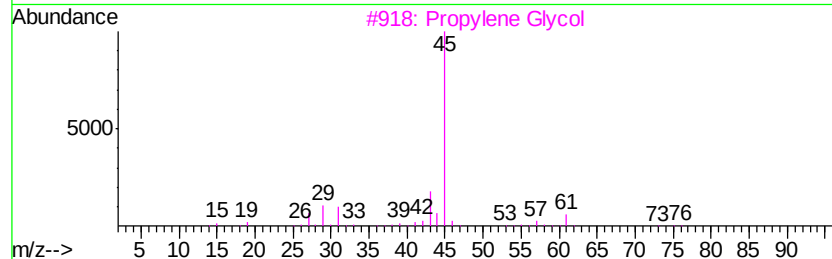
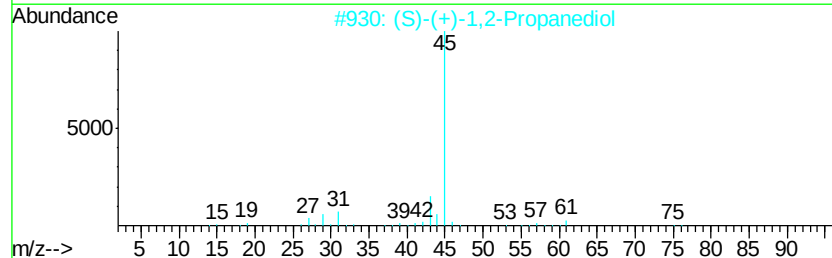
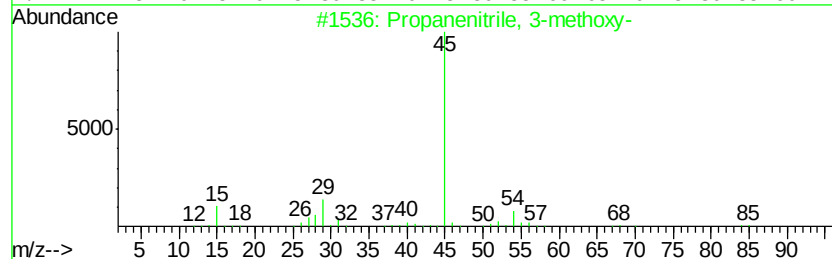
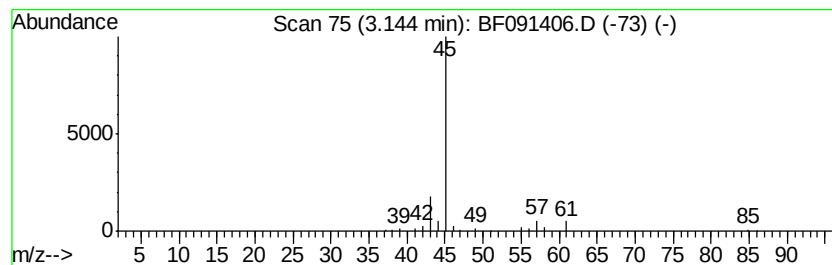
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Propanenitrile, 3-methoxy- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.14	2.69 ng	197102	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propanenitrile, 3-methoxy-	85	C4H7NO	000110-67-8	50
2		(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	40
3		Propylene Glycol	76	C3H8O2	000057-55-6	40
4		Propane, 2-ethoxy-	88	C5H12O	000625-54-7	9
5		Isopropyl Alcohol	60	C3H8O	000067-63-0	9



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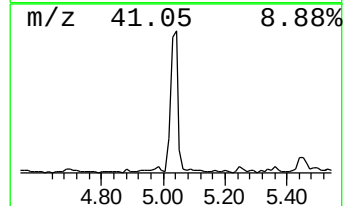
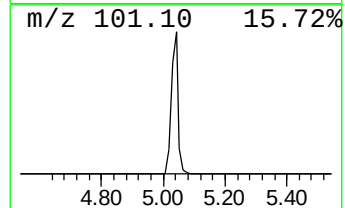
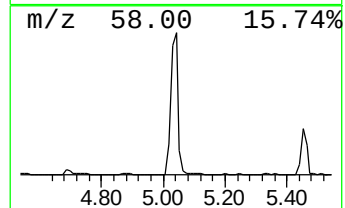
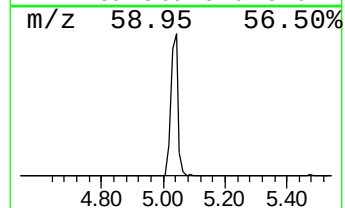
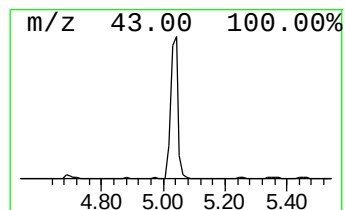
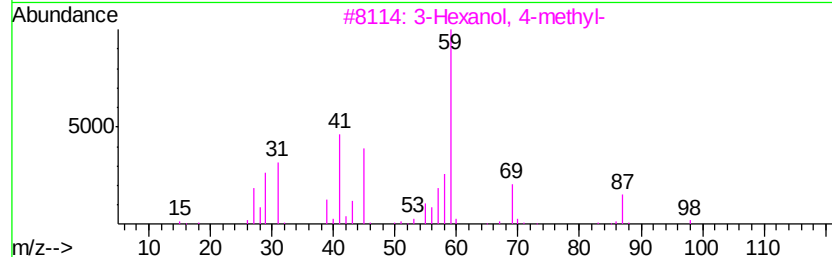
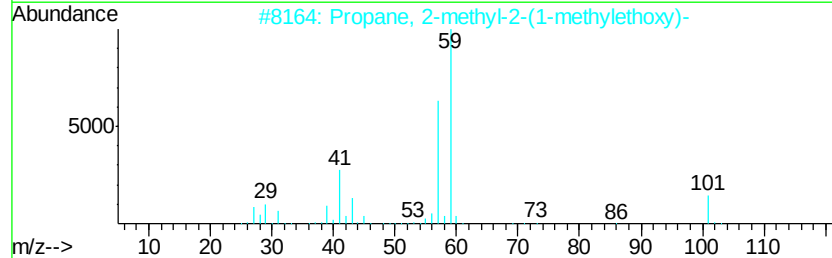
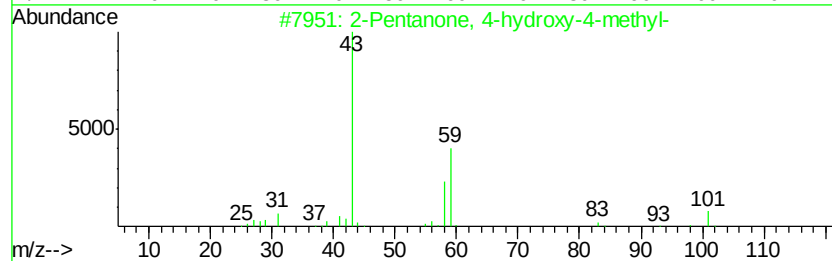
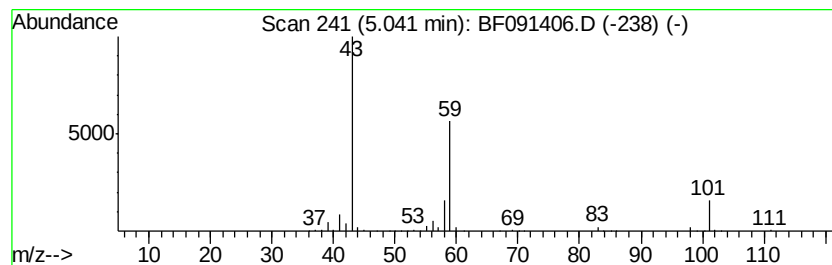
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.04	22.58 ng	1652290	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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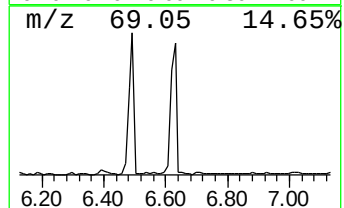
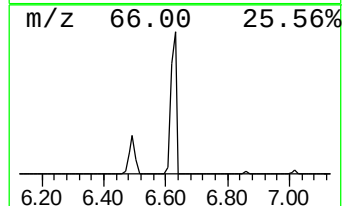
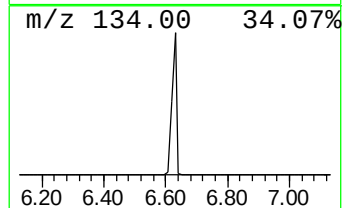
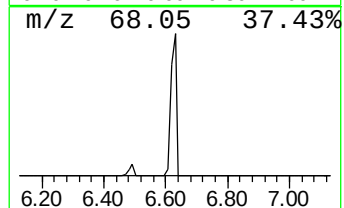
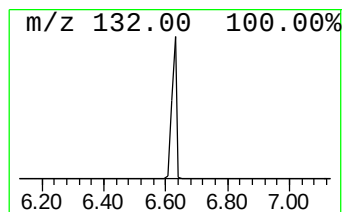
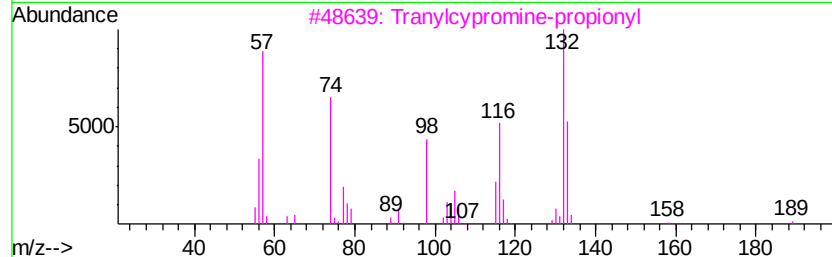
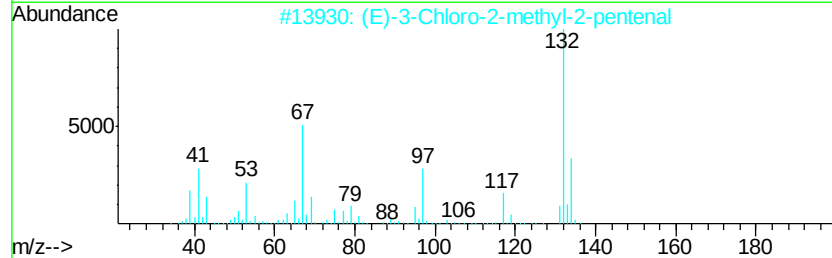
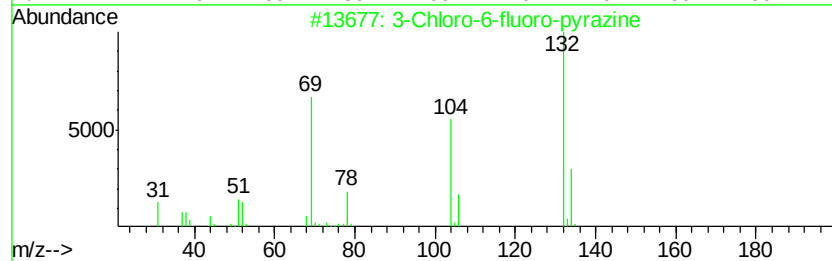
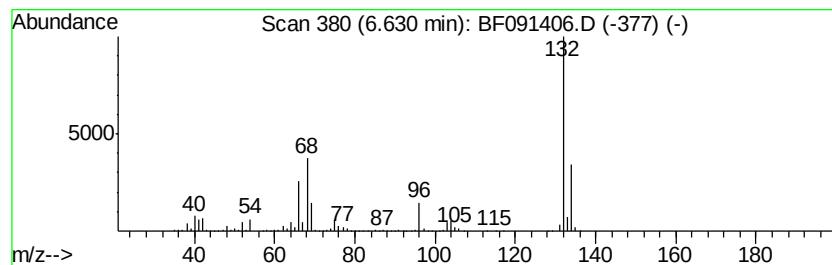
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 unknown6.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	55.38 ng	4052620	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120216\
Data File : BF091406.D
Acq On : 2 Dec 2016 23:06
Operator : UM/SJ
Sample : H5829-03
Misc :
ALS Vial : 21 Sample Multiplier: 1

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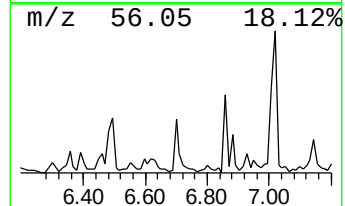
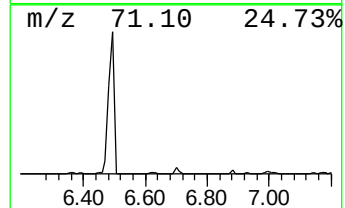
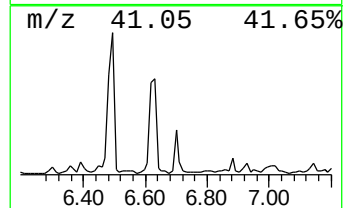
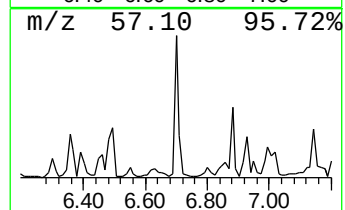
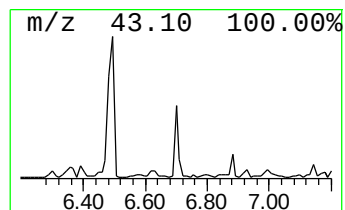
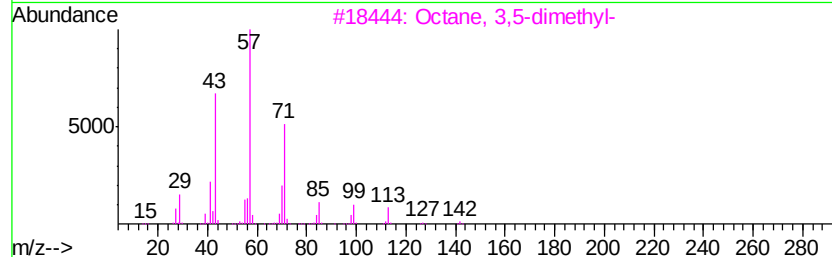
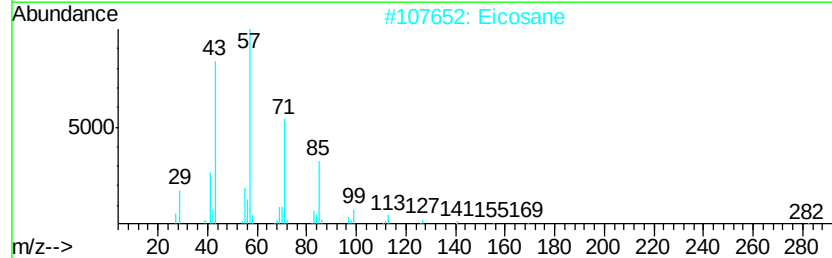
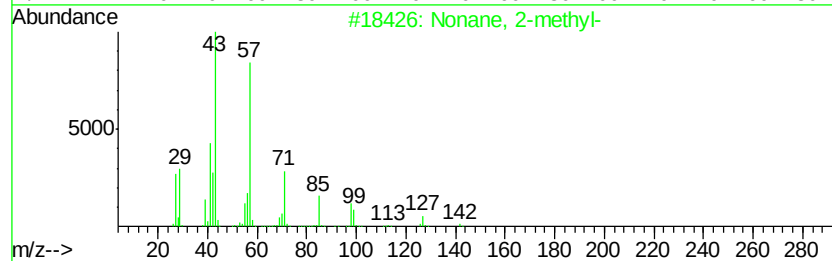
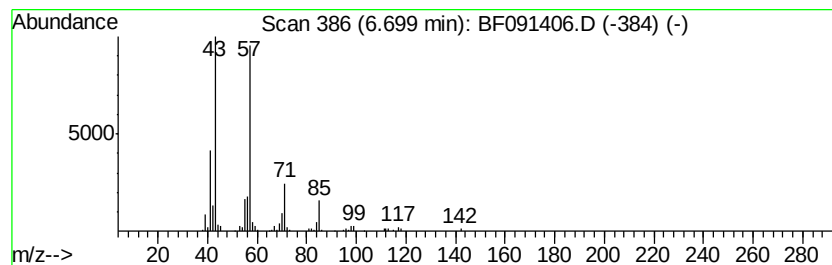
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Nonane, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.70	4.45 ng	325828	1,4-Dichlorobenzene-d4	6.86

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane, 2-methyl-	142	C10H22	000871-83-0	80	
2	Eicosane	282	C20H42	000112-95-8	72	
3	Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	64	
4	Nonane	128	C9H20	000111-84-2	59	
5	Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	59	



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120216\
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Sample : H5829-03
Misc :
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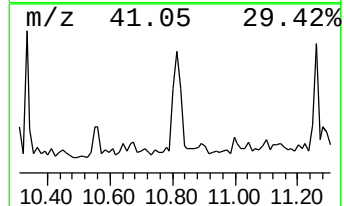
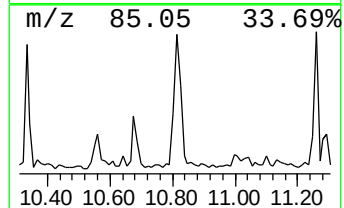
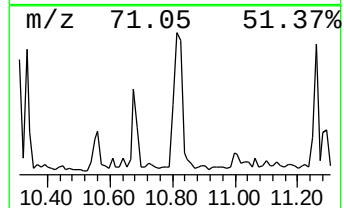
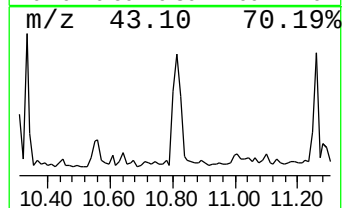
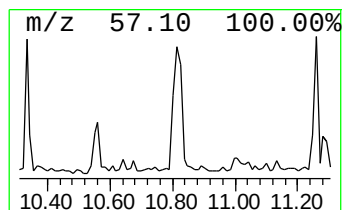
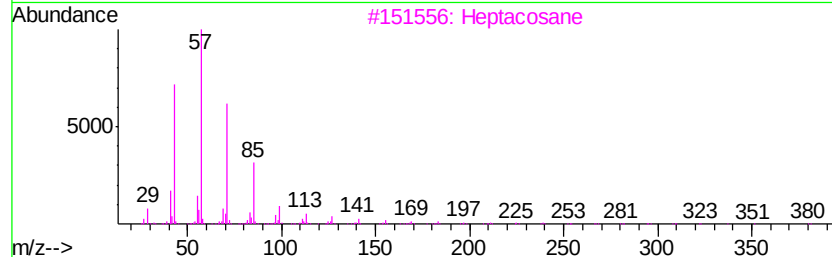
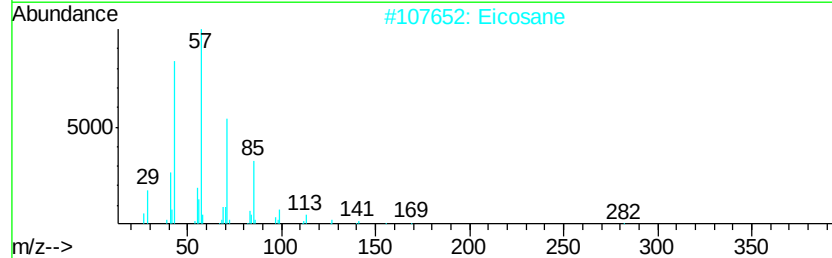
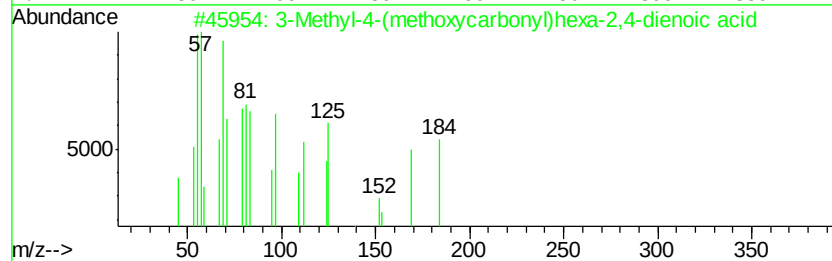
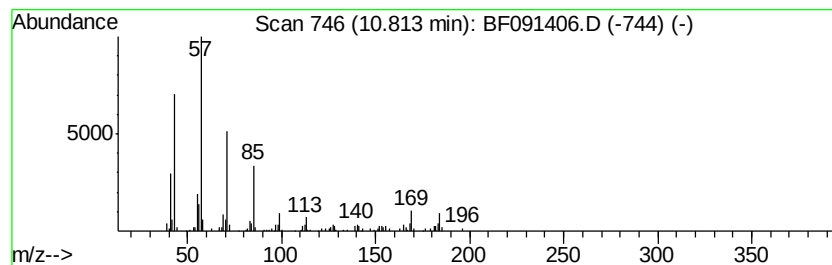
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 3-Methyl-4-(methoxycarbonyl... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.81	2.77 ng	270465	Phenanthrene-d10	11.38		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Methyl-4-(methoxycarbonyl)hexa...	184	C9H12O4	1000104-10-8	87	
2	Eicosane	282	C20H42	000112-95-8	70	
3	Heptacosane	380	C27H56	000593-49-7	62	
4	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	58	
5	5-Ethyldecane	170	C12H26	017302-36-2	58	



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120216\
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Operator : UM/SJ
Sample : H5829-03
Misc :
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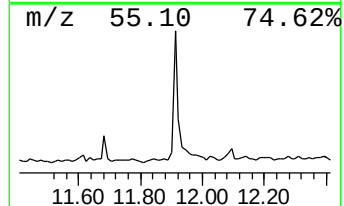
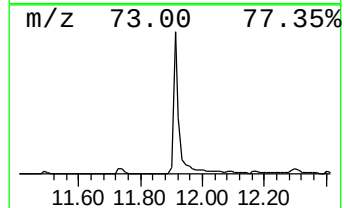
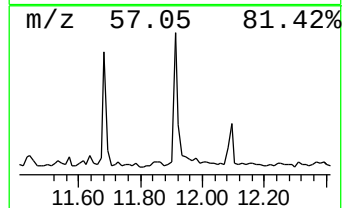
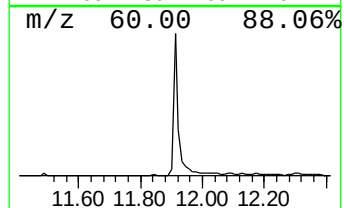
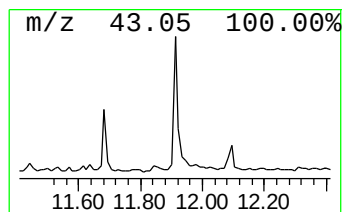
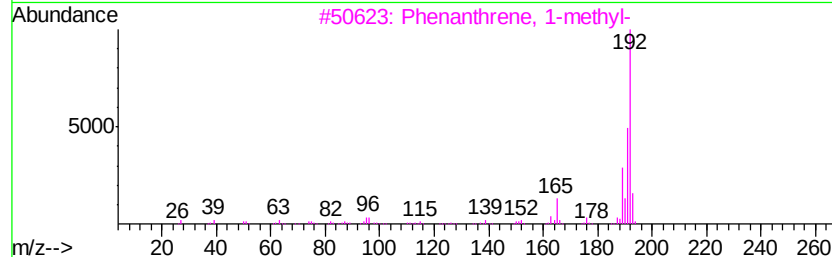
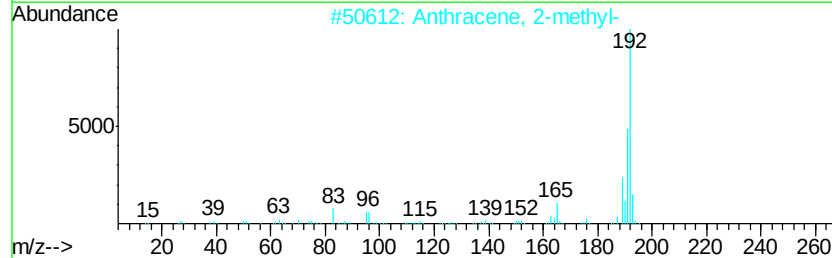
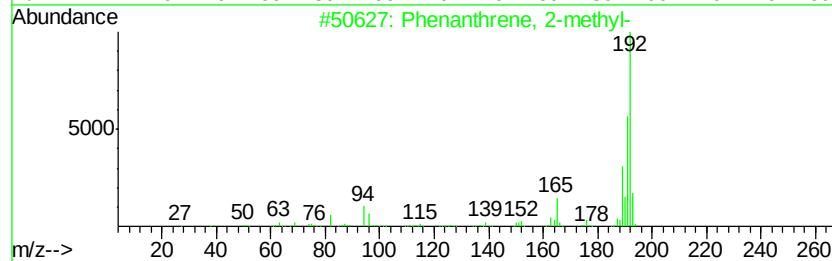
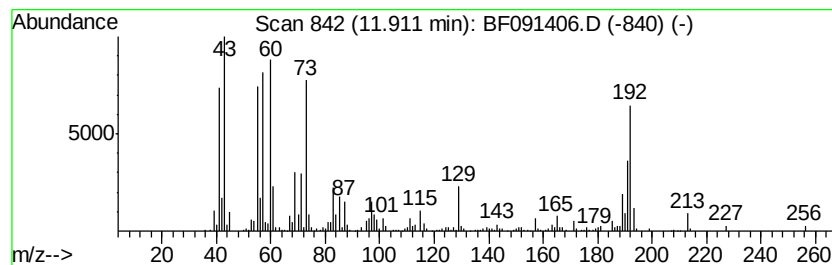
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.91	5.86 ng	572533	Phenanthrene-d10	11.38	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	91
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	89
5	1H-Cyclopropa[l]phenanthrene, 1a, ...	192	C15H12	000949-41-7	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120216\
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Sample : H5829-03
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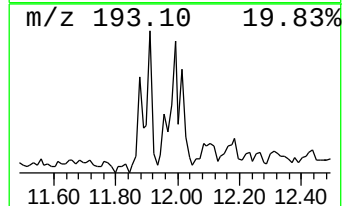
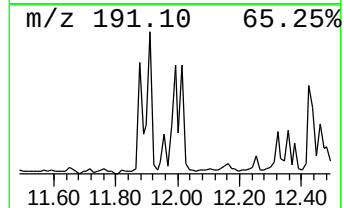
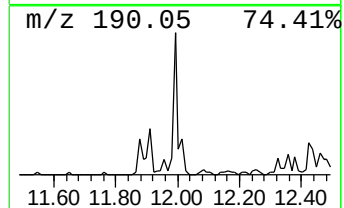
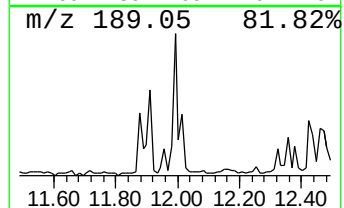
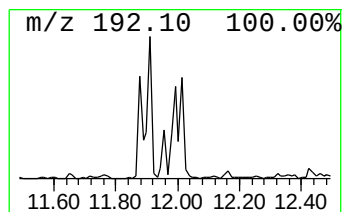
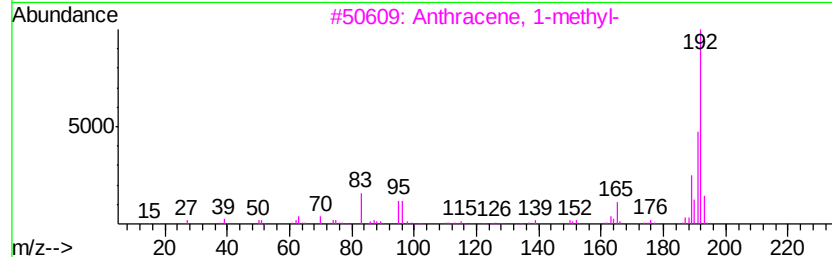
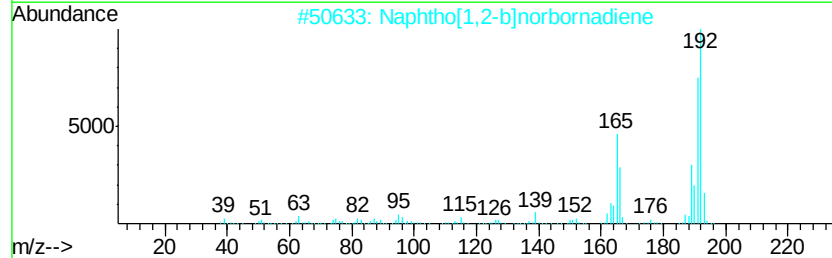
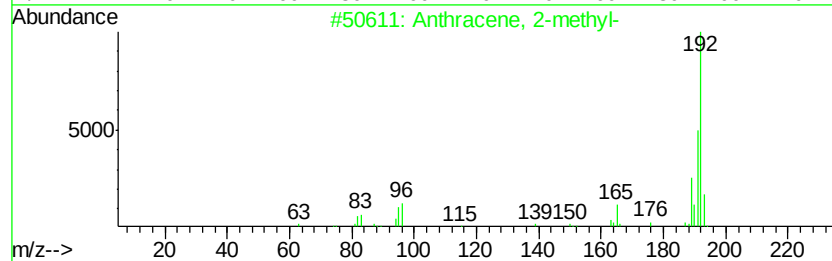
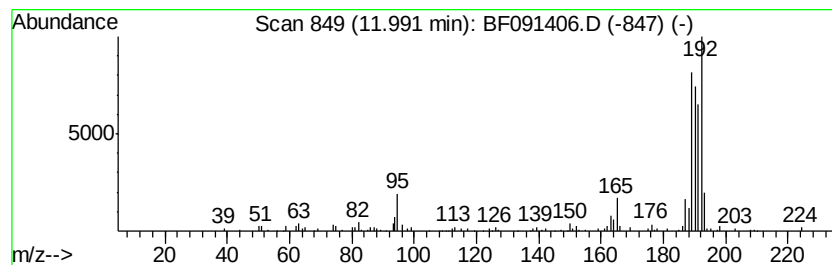
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Anthracene, 2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.99	2.62 ng	256112	Phenanthrene-d10		11.38	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 2-methyl-	192	C15H12	000613-12-7	64
2		Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	50
3		Anthracene, 1-methyl-	192	C15H12	000610-48-0	50
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	46
5		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	46



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120216\
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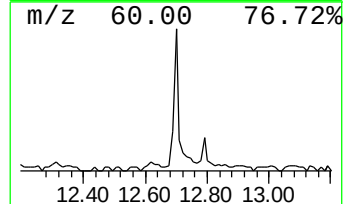
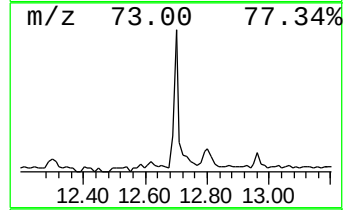
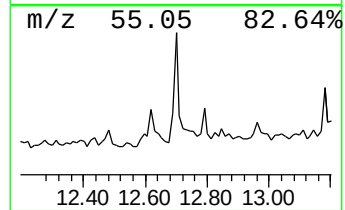
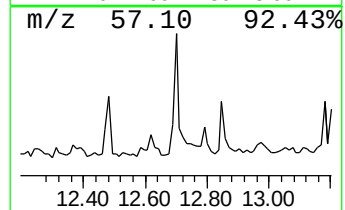
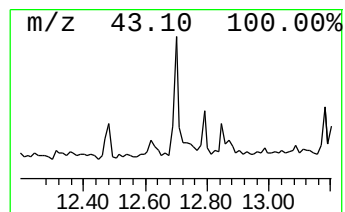
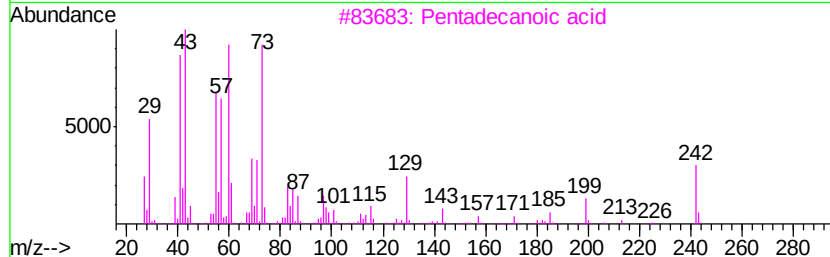
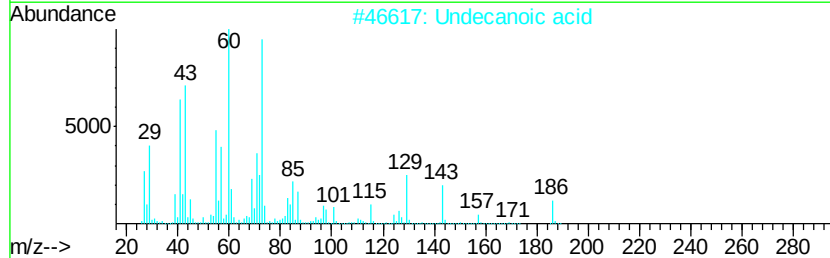
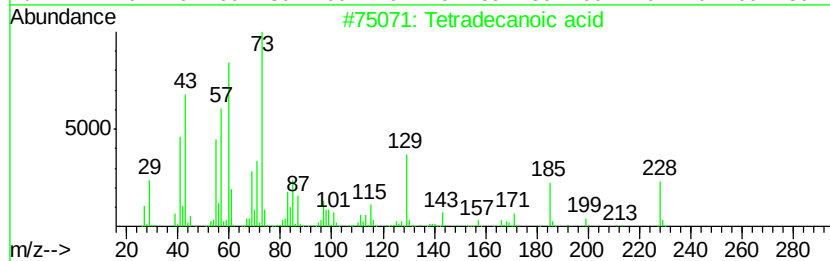
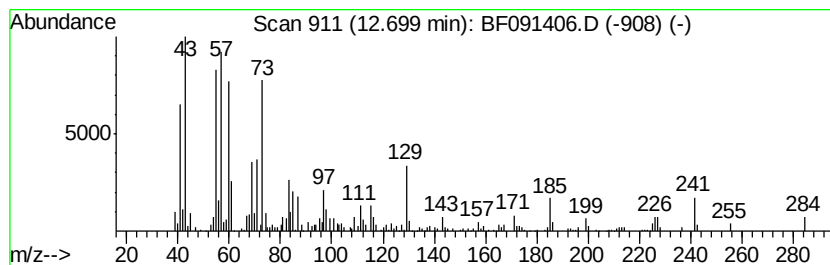
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tetradecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.70	2.97 ng	258889	Chrysene-d12	14.01	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecanoic acid	228	C14H28O2	000544-63-8	74
2	Undecanoic acid	186	C11H22O2	000112-37-8	64
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	64
4	Tridecanoic acid	214	C13H26O2	000638-53-9	59
5	n-Decanoic acid	172	C10H20O2	000334-48-5	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120216\
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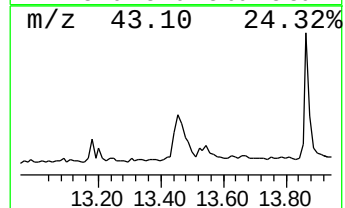
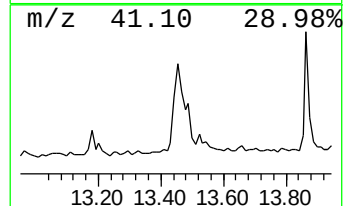
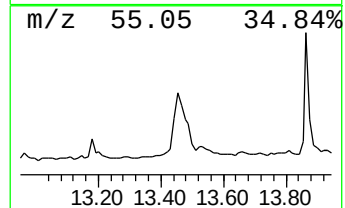
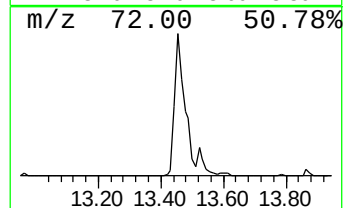
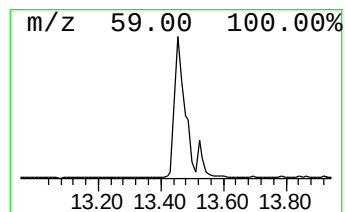
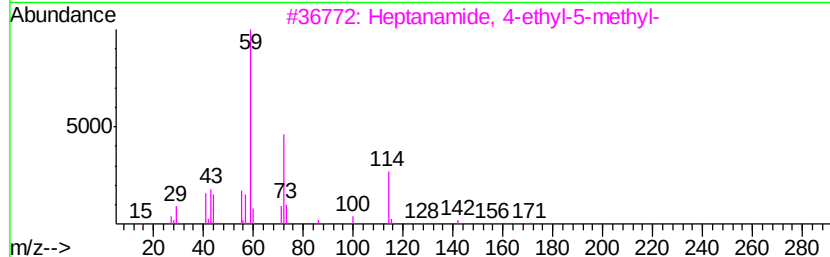
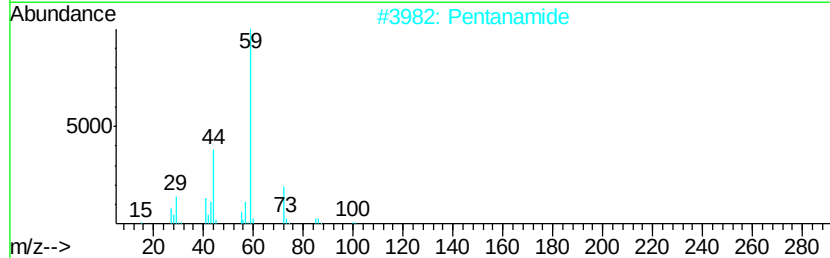
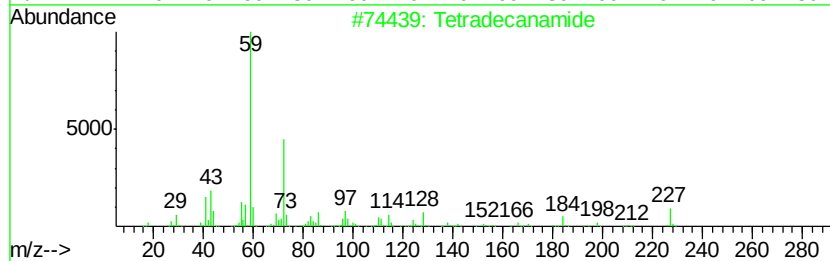
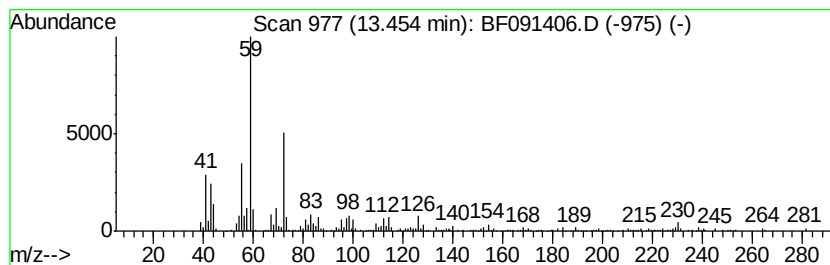
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tetradecanamide Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.45	6.46 ng	563728	Chrysene-d12	14.01
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tetradecanamide	227 C14H29NO	000638-58-4 72
2		Pentanamide	101 C5H11NO	000626-97-1 64
3		Heptanamide, 4-ethyl-5-methyl-	171 C10H21NO	054789-40-1 58
4		Octanamide	143 C8H17NO	000629-01-6 56
5		Nonanamide	157 C9H19NO	001120-07-6 53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120216\
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Sample : H5829-03
Misc :
ALS Vial : 21 Sample Multiplier: 1

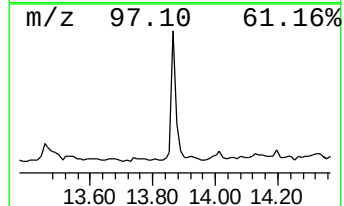
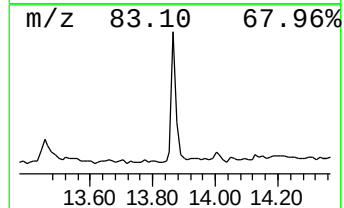
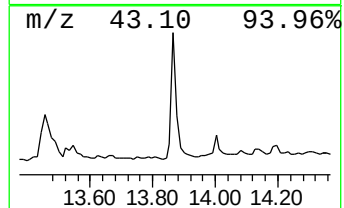
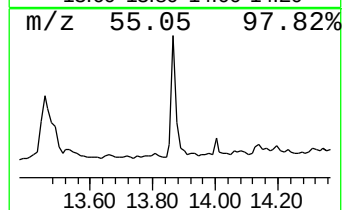
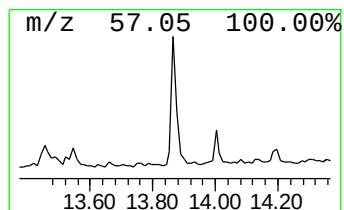
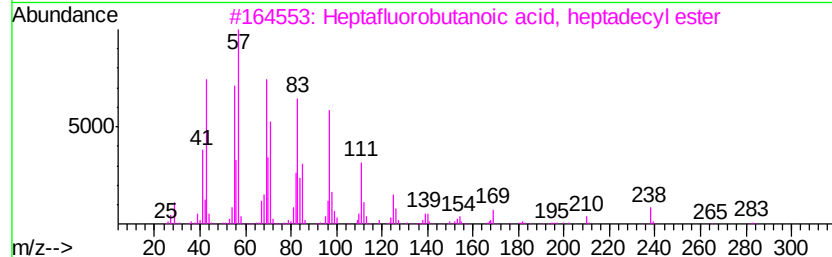
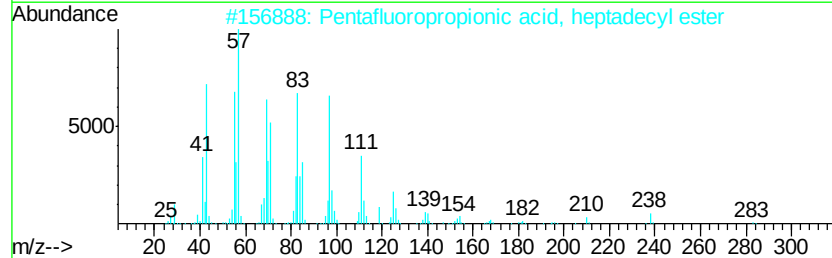
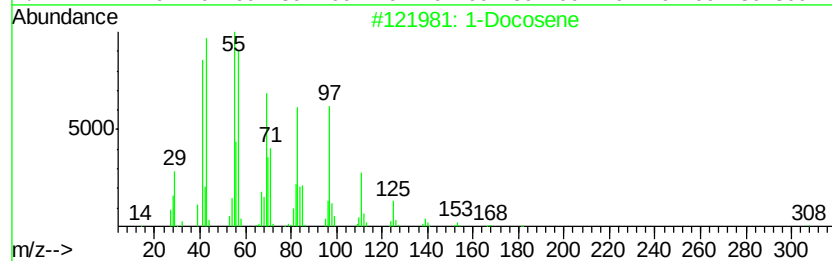
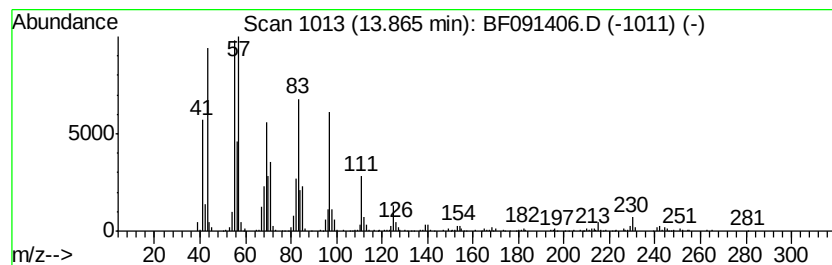
Instrument :
BNA_F
ClientSampleId :
WESTWALL-120116

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-Docosene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	4.39 ng	383297	Chrysene-d12	14.01
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Docosene		308 C22H44	001599-67-3 91
2	Pentafluoropropionic acid, hepta...		402 C20H35F5O2	1000283-04-2 87
3	Heptafluorobutanoic acid, heptad...		452 C21H35F7O2	1000282-97-3 74
4	5-Eicosene, (E)-		280 C20H40	074685-30-6 74
5	11-Tricosene		322 C23H46	052078-56-5 74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120216\
Data File : BF091406.D
Acq On : 2 Dec 2016 23:06
Operator : UM/SJ
Sample : H5829-03
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WESTWALL-120116

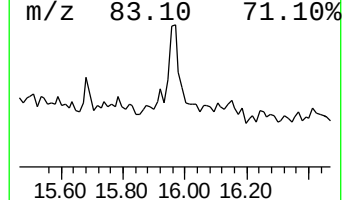
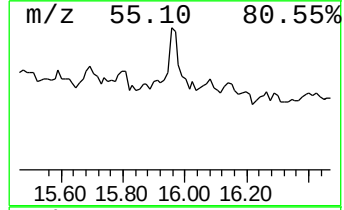
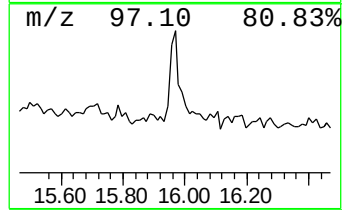
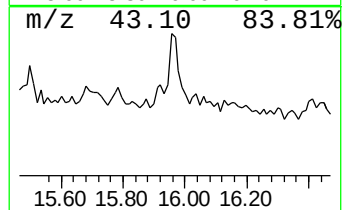
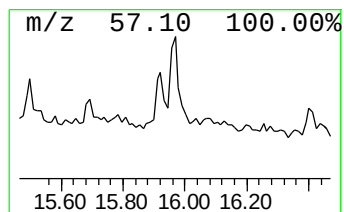
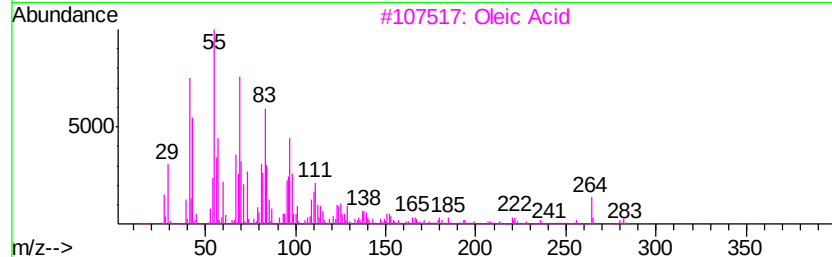
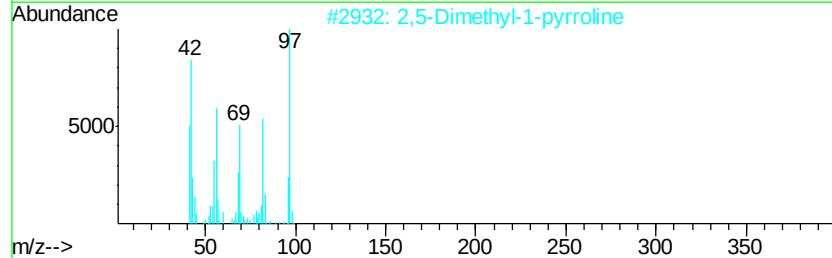
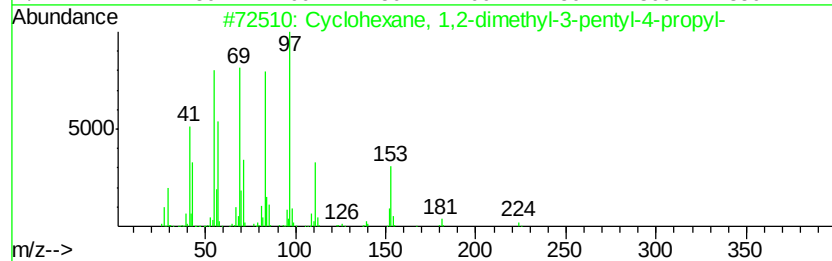
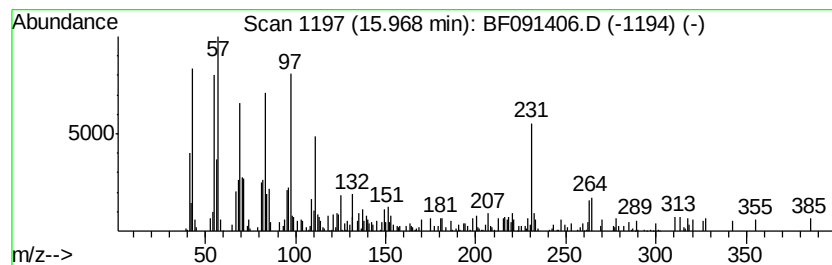
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown15.97 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.97	4.01 ng	239751	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,2-dimethyl-3-pent...	224	C16H32	062376-17-4	30
2		2,5-Dimethyl-1-pyrroline	97	C6H11N	000822-64-0	25
3		Oleic Acid	282	C18H34O2	000112-80-1	25
4		Heptane, 1,7-dibromo-	256	C7H14Br2	004549-31-9	18
5		Cyclopentane, 1-methyl-3-(2-meth...	140	C10H20	029053-04-1	18



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120216\
Data File : BF091406.D
Acq On : 2 Dec 2016 23:06
Operator : UM/SJ
Sample : H5829-03
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WESTWALL-120116

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.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
2-Butanone, 3-hyd...	2.46	4.4	ng	322010	1	6.86	1463670	20.0
Propanenitrile, 3...	3.14	2.7	ng	197102	1	6.86	1463670	20.0
2-Pentanone, 4-hy...	5.04	22.6	ng	1652290	1	6.86	1463670	20.0
unknown6.63	6.63	55.4	ng	4052620	1	6.86	1463670	20.0
Nonane, 2-methyl-	6.70	4.5	ng	325828	1	6.86	1463670	20.0
3-Methyl-4-(metho...	10.81	2.8	ng	270465	4	11.38	1955410	20.0
Phenanthrene, 2-m...	11.91	5.9	ng	572533	4	11.38	1955410	20.0
Anthracene, 2-met...	11.99	2.6	ng	256112	4	11.38	1955410	20.0
Tetradecanoic acid	12.70	3.0	ng	258889	5	14.01	1745260	20.0
Tetradecanamide	13.45	6.5	ng	563728	5	14.01	1745260	20.0
1-Docosene	13.87	4.4	ng	383297	5	14.01	1745260	20.0
unknown15.97	15.97	4.0	ng	239751	6	15.44	1196210	20.0