

Data Path : Z:\HPCHEM1\BNA_F\Data\BF040616\
 Data File : BF086122.D
 Acq On : 7 Apr 2016 2:44
 Operator : UM/SJ
 Sample : H2104-02
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-1

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-BF040216.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	3.253	100	102	104	rVB2	36249	37657	1.67%	0.148%
2	3.824	148	152	154	rBV3	410211	942263	41.78%	3.702%
3	4.693	227	228	230	rVB	39599	27381	1.21%	0.108%
4	4.910	244	247	253	rBV	142311	208849	9.26%	0.820%
5	5.013	255	256	257	rVB	124089	85229	3.78%	0.335%
6	5.344	283	285	293	rBV	2056874	2255124	100.00%	8.859%
7	6.396	375	377	379	rBV	2102866	2200998	97.60%	8.646%
8	6.510	385	387	390	rBV	1836687	2076742	92.09%	8.158%
9	6.739	405	407	410	rBV	449624	557796	24.73%	2.191%
10	6.899	419	421	424	rBV	1985700	1726216	76.55%	6.781%
11	7.310	455	457	460	rBV	1165413	1180356	52.34%	4.637%
12	8.030	518	520	522	rBV	782113	663955	29.44%	2.608%
13	9.105	612	614	617	rBV	2053344	1874591	83.13%	7.364%
14	9.505	647	649	651	rBV	231493	195221	8.66%	0.767%
15	9.722	666	668	670	rBV	37927	42797	1.90%	0.168%
16	9.779	671	673	675	rVV	597249	544529	24.15%	2.139%
17	9.973	689	690	694	rVB2	20279	28959	1.28%	0.114%
18	10.213	710	711	713	rVB	63499	56521	2.51%	0.222%
19	10.305	718	719	720	rBV	31828	23166	1.03%	0.091%
20	10.431	725	730	732	rBV	23011	39850	1.77%	0.157%
21	10.476	732	734	736	rVB2	42805	50562	2.24%	0.199%
22	10.568	740	742	745	rBV2	892861	1068021	47.36%	4.196%
23	10.693	751	753	757	rVB	50148	82189	3.64%	0.323%
24	10.979	774	778	779	rBV4	12507	24376	1.08%	0.096%
25	11.139	788	792	793	rBV	21150	34055	1.51%	0.134%
26	11.265	801	803	805	rBV	665573	630070	27.94%	2.475%
27	11.345	807	810	811	rVB2	17994	25818	1.14%	0.101%
28	11.756	842	846	847	rBV3	31586	63657	2.82%	0.250%
29	11.802	847	850	852	rVV	87405	130516	5.79%	0.513%
30	11.882	855	857	861	rVB3	21157	44888	1.99%	0.176%
31	12.065	872	873	880	rVB5	13285	36375	1.61%	0.143%
32	12.385	899	901	905	rVB4	21673	33743	1.50%	0.133%
33	12.476	906	909	911	rBV	140156	131625	5.84%	0.517%
34	12.545	911	915	916	rVB3	17642	36452	1.62%	0.143%

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35	12.591	916	919	920	rBV2	31146	54968	2.44%	0.216%
36	12.705	925	929	932	rVB	119364	192701	8.55%	0.757%
37	12.854	939	942	944	rBV	1655401	2010290	89.14%	7.897%
38	12.899	944	946	951	rVB5	10302	23671	1.05%	0.093%
39	13.025	955	957	959	rVB2	22341	29329	1.30%	0.115%
40	13.082	959	962	965	rBV2	14910	40166	1.78%	0.158%
41	13.139	965	967	971	rVB3	22507	31856	1.41%	0.125%
42	13.414	989	991	993	rBV	22247	40463	1.79%	0.159%
43	13.539	1001	1002	1007	rVB3	13659	27560	1.22%	0.108%
44	13.654	1009	1012	1013	rBV2	23500	41928	1.86%	0.165%
45	13.699	1013	1016	1019	rVV	61311	95185	4.22%	0.374%
46	13.757	1019	1021	1025	rVV	121885	204859	9.08%	0.805%
47	13.905	1031	1034	1035	rBV	476895	639735	28.37%	2.513%
48	14.008	1039	1043	1044	rBV3	55870	140418	6.23%	0.552%
49	14.088	1049	1050	1052	rVB2	52379	40289	1.79%	0.158%
50	14.145	1052	1055	1056	rBV2	64031	89202	3.96%	0.350%
51	14.271	1064	1066	1067	rBV	71645	82676	3.67%	0.325%
52	14.362	1072	1074	1077	rBV3	28728	64722	2.87%	0.254%
53	14.419	1077	1079	1081	rBV2	38372	76711	3.40%	0.301%
54	14.545	1088	1090	1094	rBV5	27581	54830	2.43%	0.215%
55	14.602	1094	1095	1096	rBV	28308	33680	1.49%	0.132%
56	14.648	1096	1099	1103	rVV	483944	592253	26.26%	2.327%
57	14.797	1110	1112	1116	rVB	98763	117470	5.21%	0.461%
58	14.968	1126	1127	1129	rVB	70162	55858	2.48%	0.219%
59	15.071	1134	1136	1139	rBV3	40489	104231	4.62%	0.409%
60	15.254	1150	1152	1154	rBV2	53392	70255	3.12%	0.276%
61	15.300	1154	1156	1162	rVV	361763	529814	23.49%	2.081%
62	15.494	1172	1173	1177	rBV2	67664	155334	6.89%	0.610%
63	15.700	1189	1191	1194	rVB	93881	144606	6.41%	0.568%
64	15.780	1196	1198	1200	rBV3	48044	66541	2.95%	0.261%
65	15.814	1200	1201	1203	rVB2	78856	79603	3.53%	0.313%
66	16.763	1282	1284	1285	rBV2	59866	49354	2.19%	0.194%
67	16.980	1302	1303	1305	rVB2	85181	83442	3.70%	0.328%
68	17.037	1305	1308	1309	rBV3	65710	167901	7.45%	0.660%
69	17.311	1331	1332	1334	rBV2	36326	66427	2.95%	0.261%
70	17.723	1366	1368	1369	rBV2	39036	54146	2.40%	0.213%
71	17.837	1373	1378	1380	rVB6	43554	150390	6.67%	0.591%

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72	18.043	1393	1396	1399	rVB4	183173	419096	18.58%	1.646%
73	18.100	1399	1401	1404	rBV3	79264	175834	7.80%	0.691%
74	18.157	1404	1406	1409	rVV4	34251	87687	3.89%	0.344%
75	18.306	1417	1419	1422	rVB4	48391	60943	2.70%	0.239%
76	18.363	1422	1424	1428	rBV4	38068	70967	3.15%	0.279%
77	18.443	1430	1431	1432	rBV	67064	55728	2.47%	0.219%
78	18.614	1444	1446	1448	rBV3	33990	55883	2.48%	0.220%
79	18.740	1455	1457	1458	rBV2	29813	46466	2.06%	0.183%
80	18.763	1458	1459	1460	rVV	41899	53544	2.37%	0.210%
81	18.831	1463	1465	1466	rVB2	45142	49493	2.19%	0.194%
82	18.866	1466	1468	1469	rVB2	33300	41644	1.85%	0.164%
83	18.946	1469	1475	1477	rBV7	48002	171324	7.60%	0.673%
84	19.003	1477	1480	1483	rVV5	46799	117490	5.21%	0.462%
85	19.220	1495	1499	1501	rBV5	93522	253027	11.22%	0.994%
86	19.254	1501	1502	1503	rBV	50231	35586	1.58%	0.140%
87	19.334	1507	1509	1510	rBV2	44267	60948	2.70%	0.239%
88	19.517	1523	1525	1526	rVB2	33861	36353	1.61%	0.143%

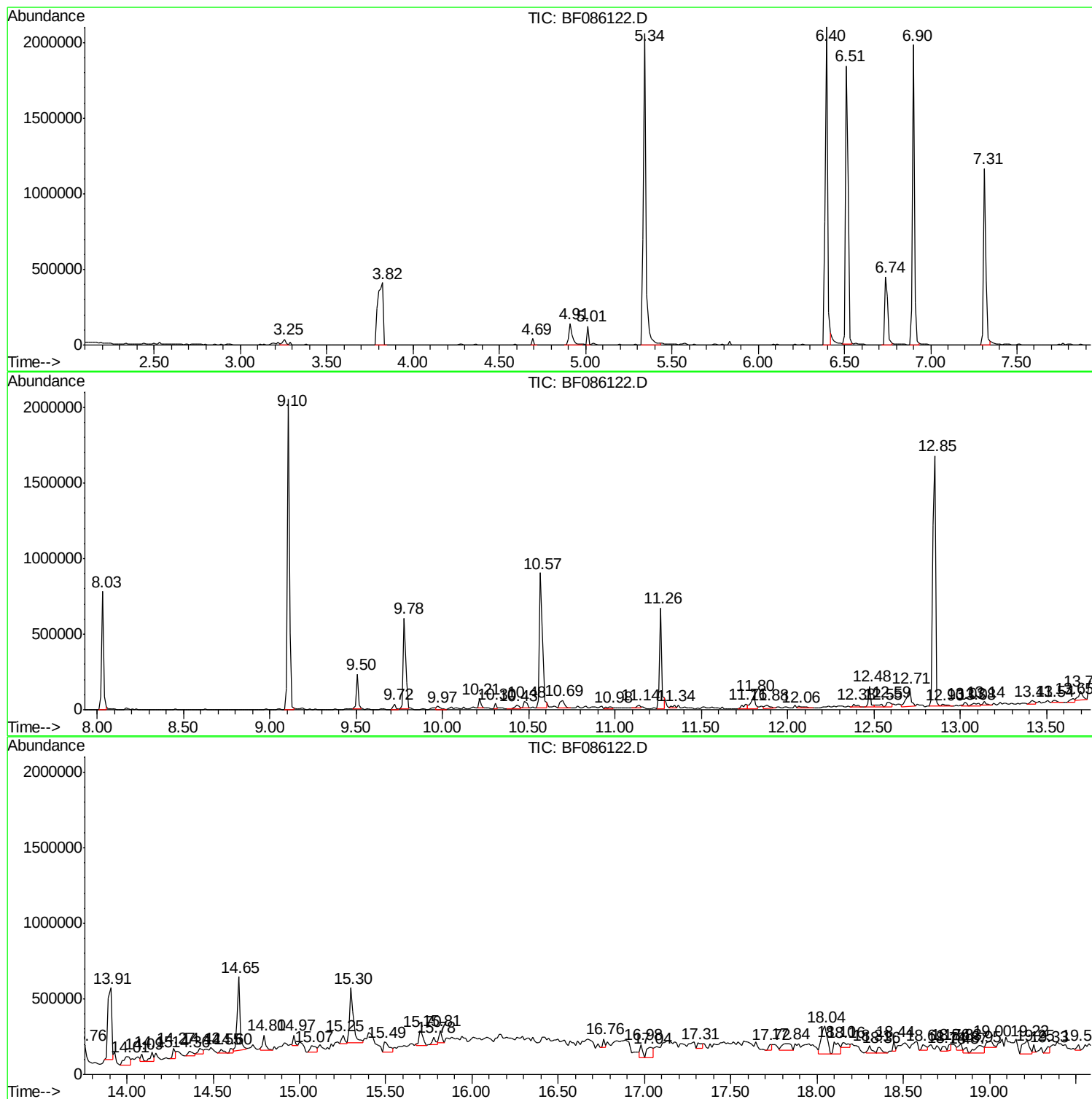
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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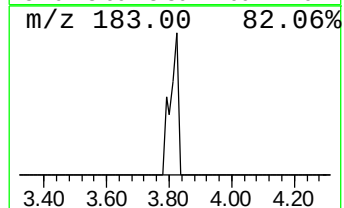
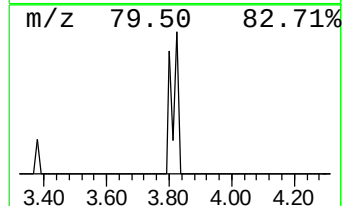
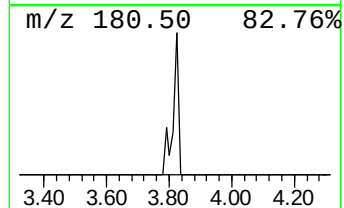
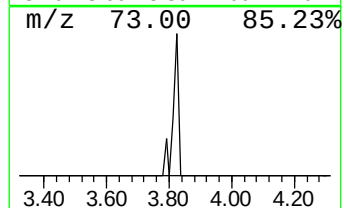
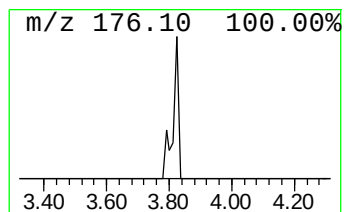
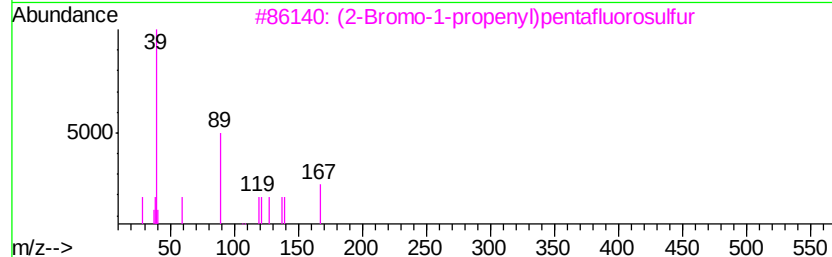
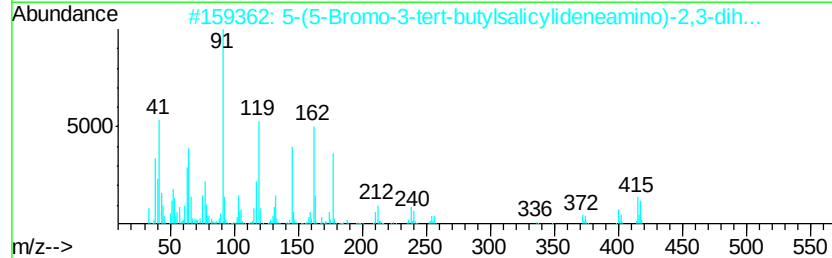
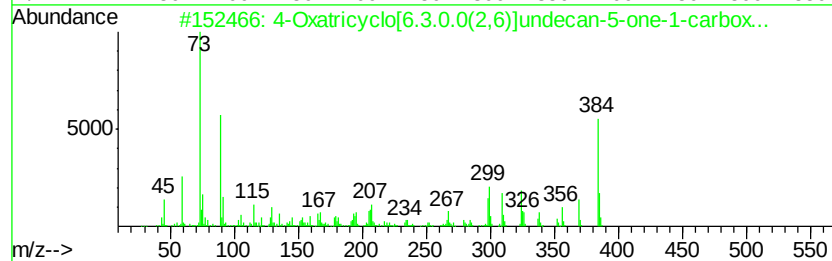
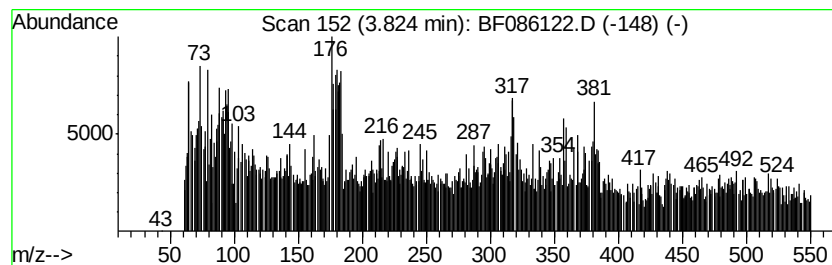
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown3.82 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.82	33.79 ng	942263	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Oxatricyclo[6.3.0.0(2,6)]undec...	384	C22H28O4Si	1000159-41-4	25
2		5-(5-Bromo-3-tert-butylsalicylid...	415	C19H18BrN3O3	330682-13-8	22
3		(2-Bromo-1-propenyl)pentafluoros...	246	C3H4BrF5S	070664-94-7	22
4		1H-Triazole-4-carboxamide, 1-[3,...	476	C23H14F6N4O	1000267-44-8	12
5		N,N'-(4-Methyl-1,3-phenylene)bis...	464	C27H20N4O4	1000223-72-7	12



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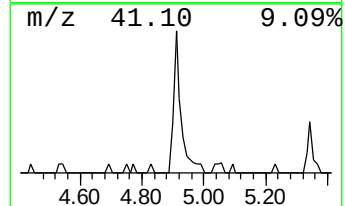
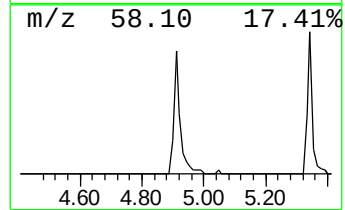
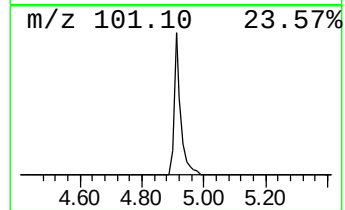
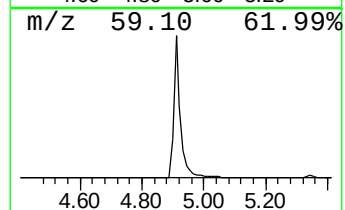
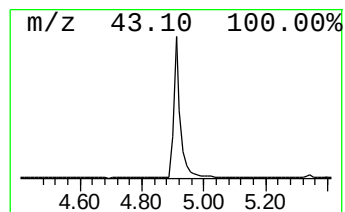
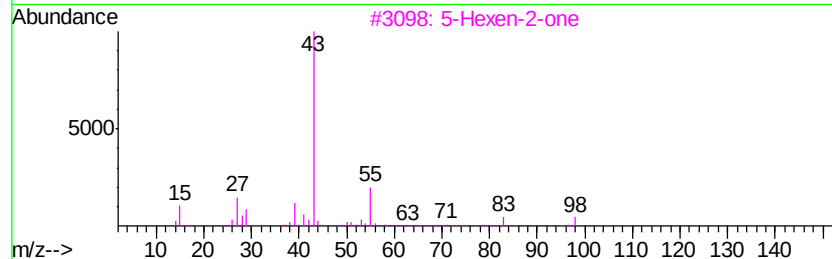
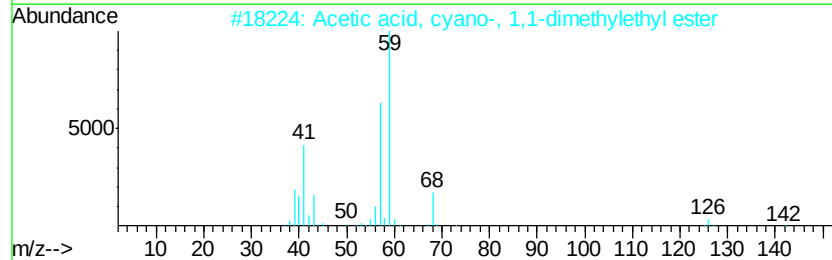
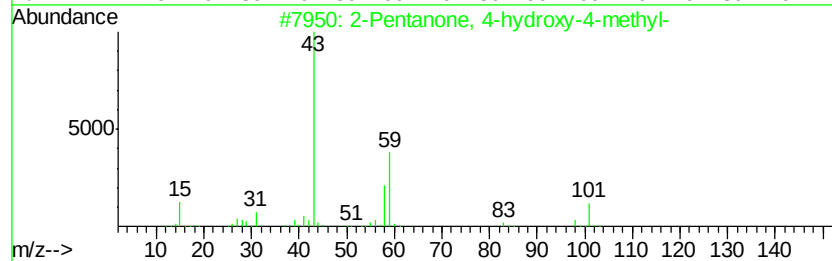
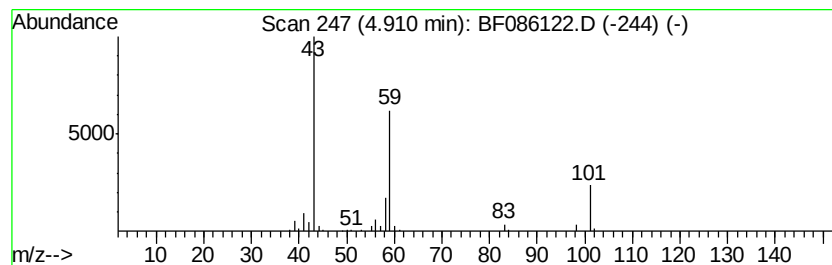
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.91	7.49 ng	208849	1,4-Dichlorobenzene-d4	6.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17		
3	5-Hexen-2-one	98	C6H10O	000109-49-9	9		
4	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9		
5	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9		



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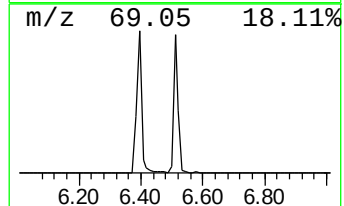
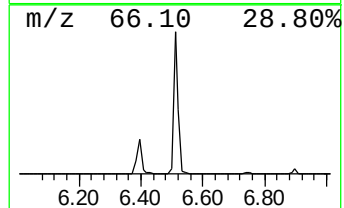
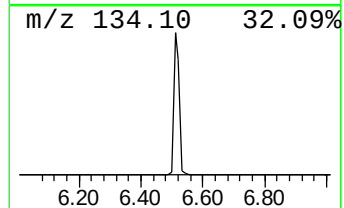
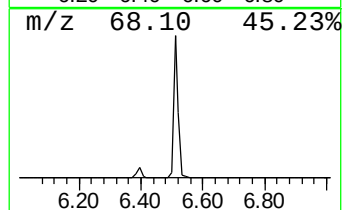
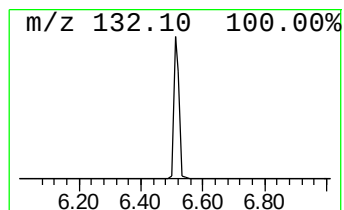
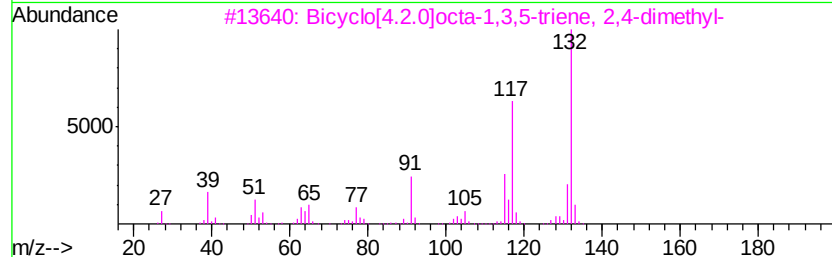
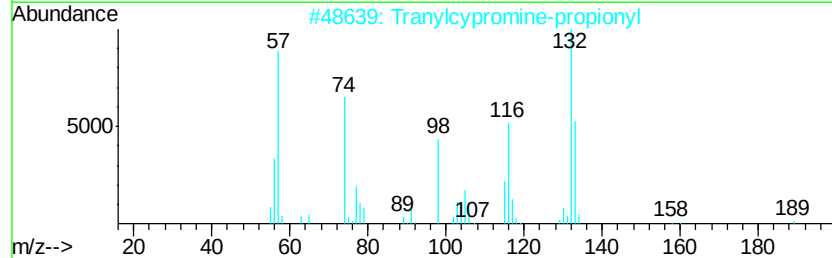
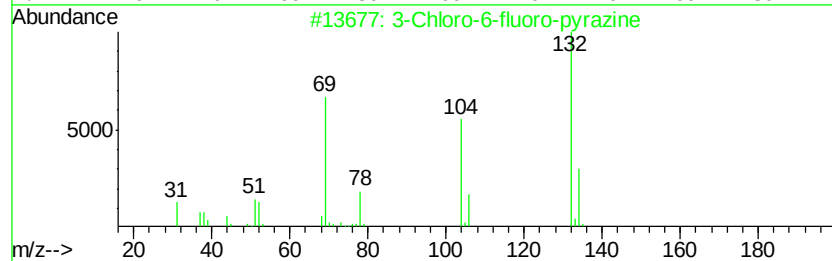
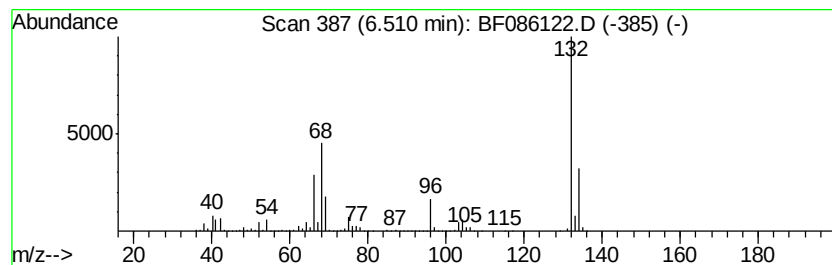
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.51 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	74.46 ng	2076740	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		1,3,4-Thiadiazole-2(3H)-thione,...	132	C3H4N2S2	029490-19-5	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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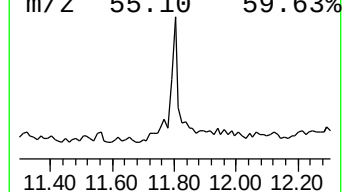
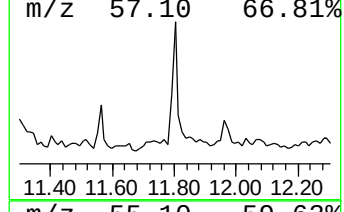
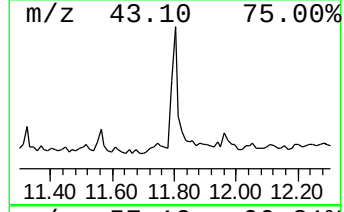
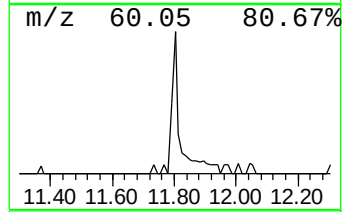
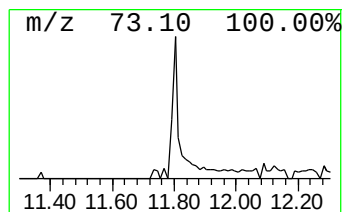
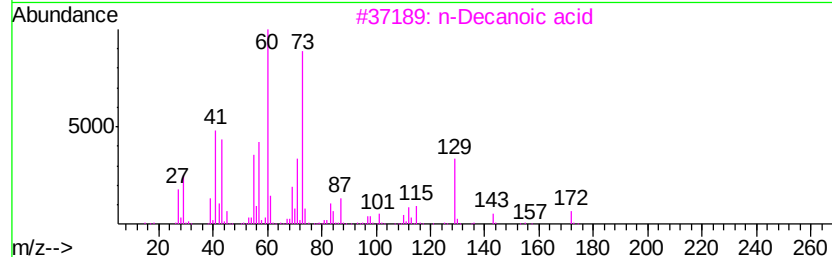
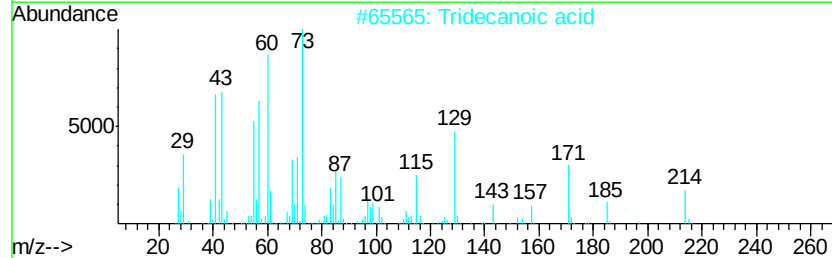
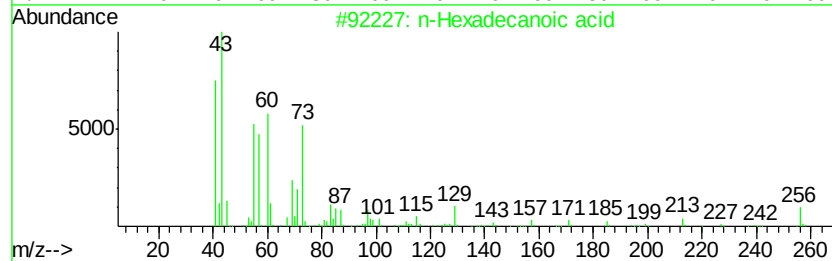
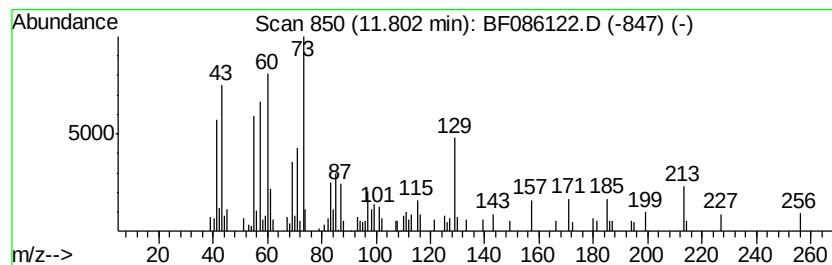
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.80	4.14 ng	130516	Phenanthrene-d10		11.26
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2	Tridecanoic acid	214	C13H26O2	000638-53-9	91
3	n-Decanoic acid	172	C10H20O2	000334-48-5	89
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	86



Data Path : Z:\HPCHEM1\BNA_F\Data\BF040616\
Data File : BF086122.D
Acq On : 7 Apr 2016 2:44
Operator : UM/SJ
Sample : H2104-02
Misc :
ALS Vial : 33 Sample Multiplier: 1

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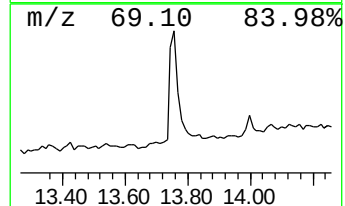
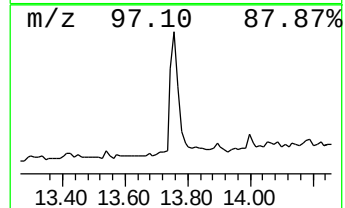
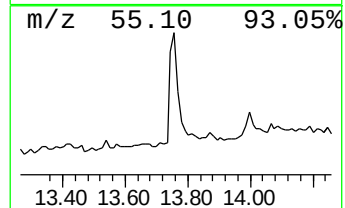
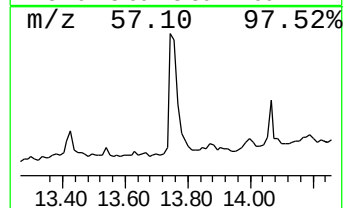
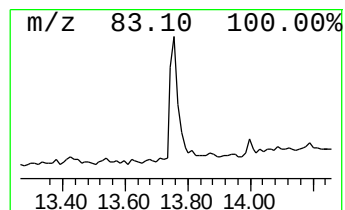
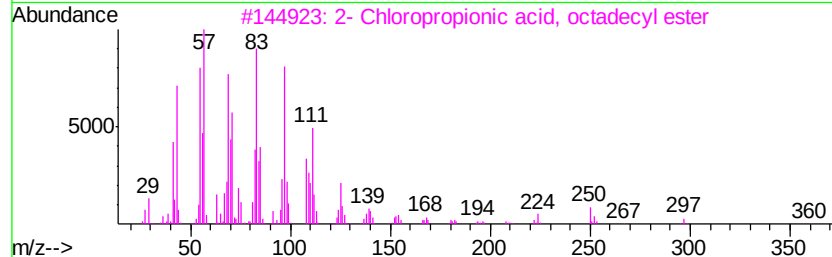
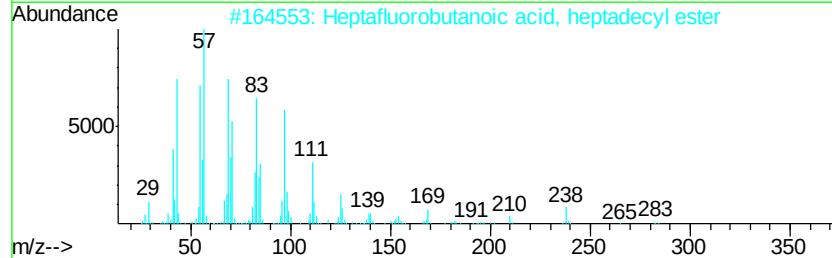
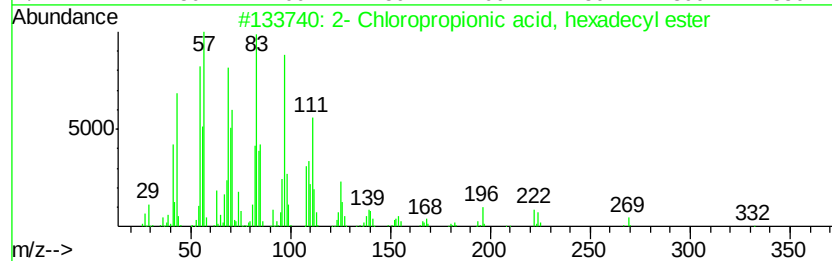
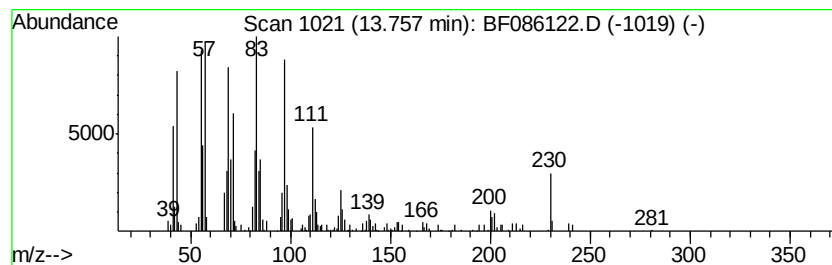
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2- Chloropropionic acid, he... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.76	6.40 ng	204859	Chrysene-d12	13.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2- Chloropropionic acid, hexadec...	332	C19H37ClO2	086711-81-1	94
2		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	94
3		2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94
4		1-Nonadecene	266	C19H38	018435-45-5	91
5		Trifluoroacetic acid, n-octadecy...	366	C20H37F3O2	079392-43-1	91



Data Path : Z:\HPCHEM1\BNA_F\Data\BF040616\
Data File : BF086122.D
Acq On : 7 Apr 2016 2:44
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Sample : H2104-02
Misc :
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Instrument :
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ClientSampleId :
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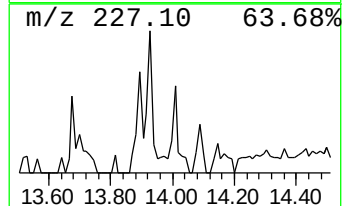
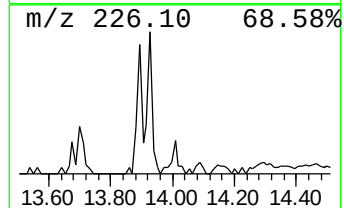
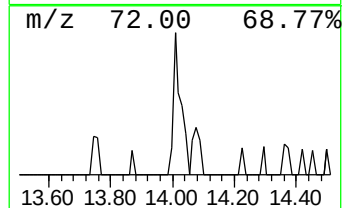
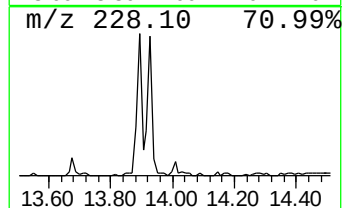
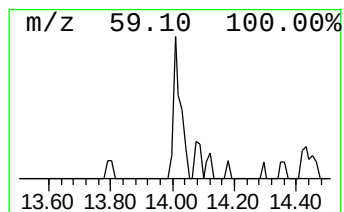
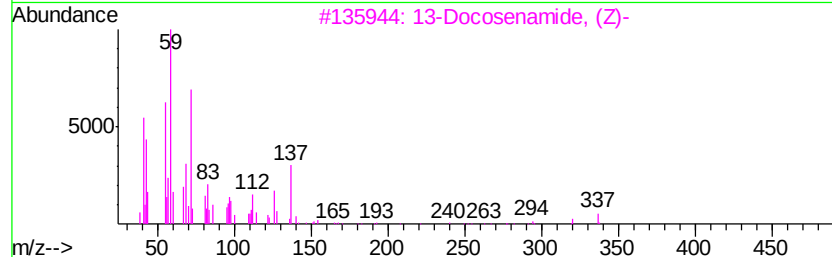
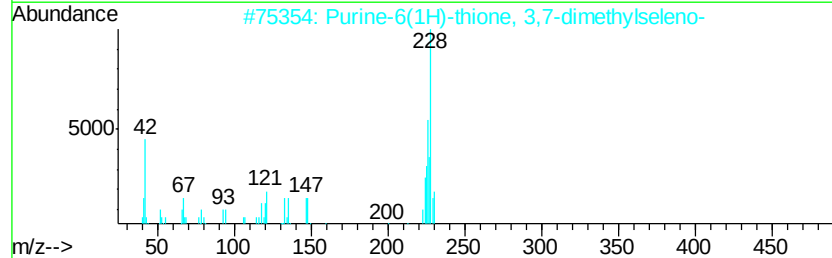
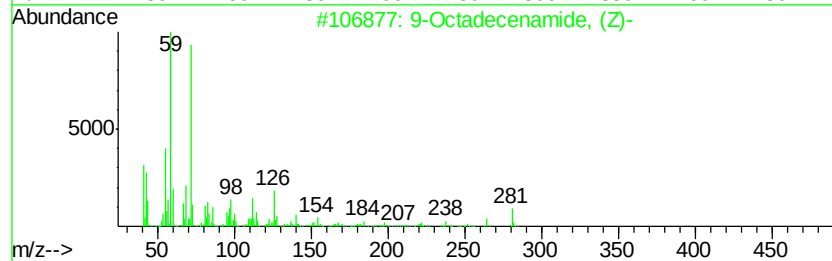
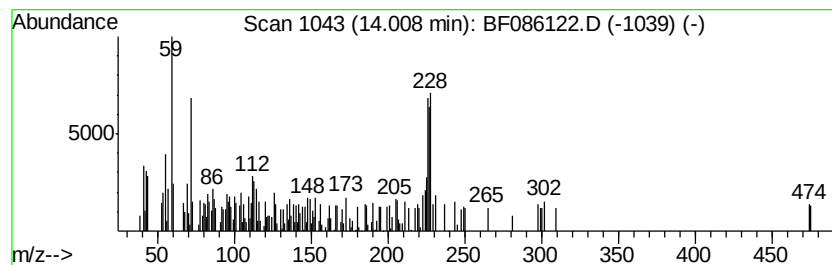
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown14.01 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.01	4.39 ng	140418	Chrysene-d12	13.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	30
2		Purine-6(1H)-thione, 3,7-dimethy...	228	C7H8N4Se	023663-58-3	30
3		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	27
4		8-Methyl-6-nonenamide	169	C10H19NO	1000293-20-9	27
5		Methyl N-acetyl-d-glucosamide	235	C9H17NO6	1000127-56-1	22



Data Path : Z:\HPCHEM1\BNA_F\Data\BF040616\
Data File : BF086122.D
Acq On : 7 Apr 2016 2:44
Operator : UM/SJ
Sample : H2104-02
Misc :
ALS Vial : 33 Sample Multiplier: 1

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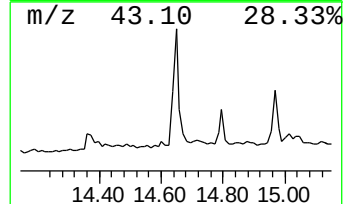
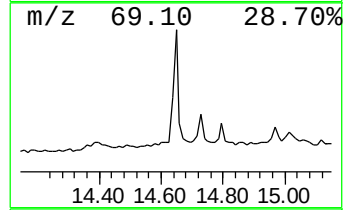
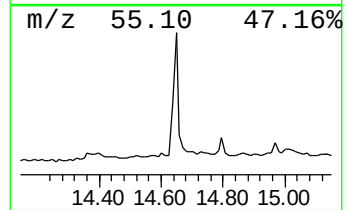
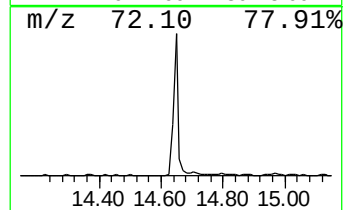
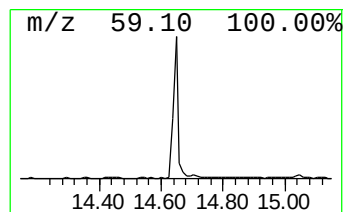
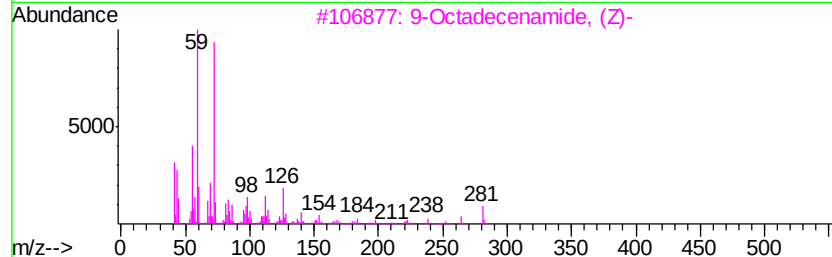
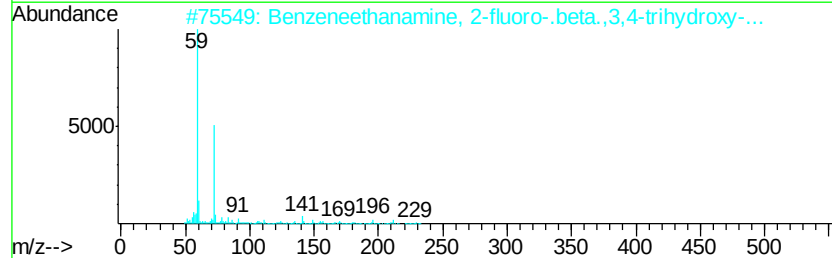
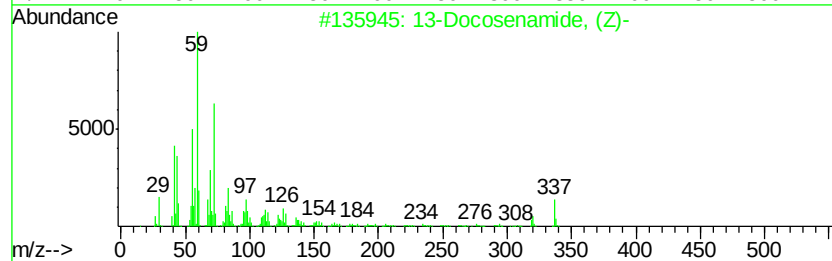
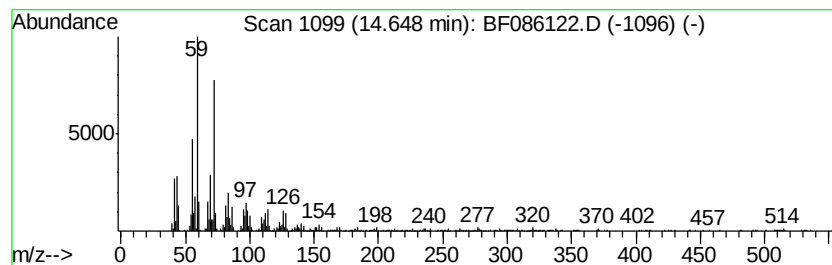
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 13-Docosenamide, (Z)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.65	22.36 ng	592253	Perylene-d12	15.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	93
2		Benzeneethanamine, 2-fluoro-.bet...	229	C11H16FN03	061338-98-5	72
3		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	72
4		Tetradecanamide	227	C14H29NO	000638-58-4	35
5		Benzo[b]dioxane, 3,3-dimethyl-2-...	235	C14H21N02	003523-36-2	30



Data Path : Z:\HPCHEM1\BNA_F\Data\BF040616\
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Acq On : 7 Apr 2016 2:44
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Sample : H2104-02
Misc :
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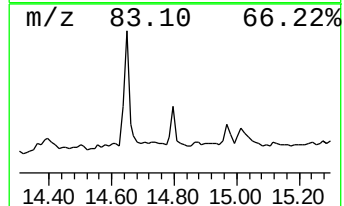
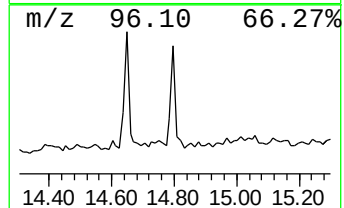
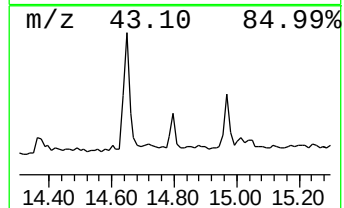
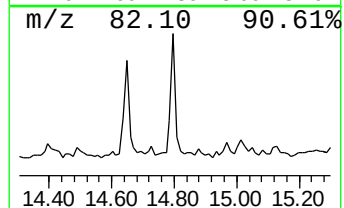
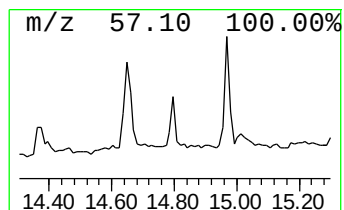
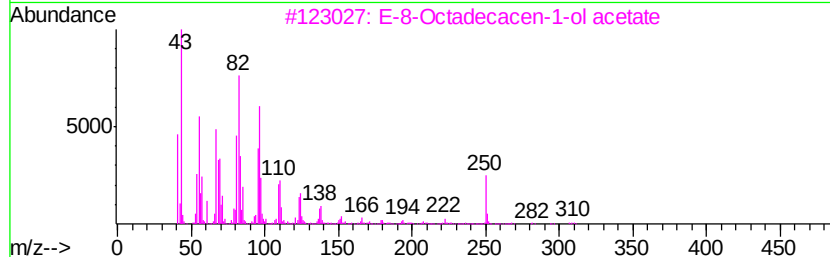
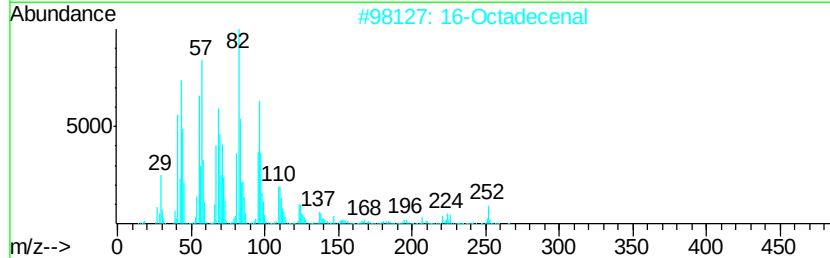
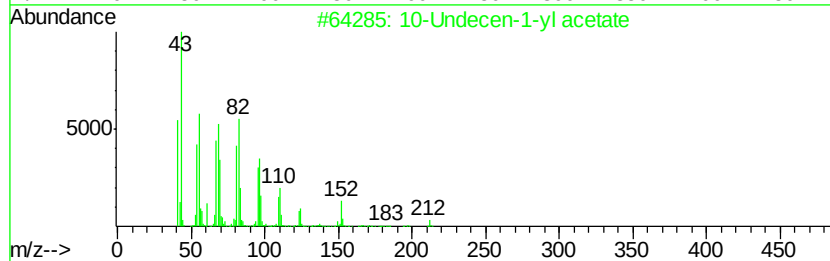
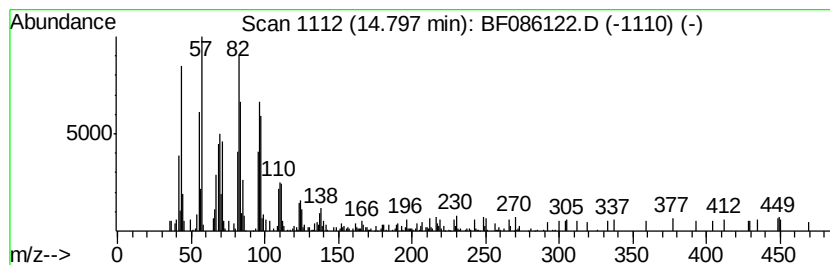
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 10-Undecen-1-yl acetate Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.80	4.43 ng	117470	Perylene-d12	15.30	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10-Undecen-1-yl acetate	212	C13H24O2	000112-19-6	68
2	16-Octadecenal	266	C18H34O	056554-87-1	64
3	E-8-Octadecacene-1-ol acetate	310	C20H38O2	002195-90-6	64
4	Oleyl Alcohol	268	C18H36O	000143-28-2	64
5	1,19-Eicosadiene	278	C20H38	014811-95-1	64



Data Path : Z:\HPCHEM1\BNA_F\Data\BF040616\
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Acq On : 7 Apr 2016 2:44
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Sample : H2104-02
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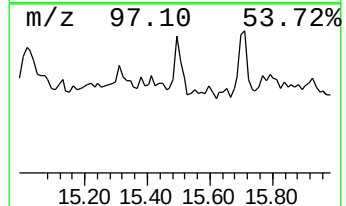
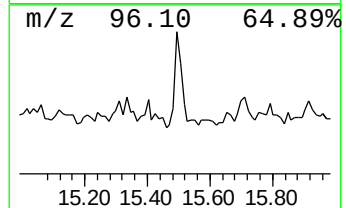
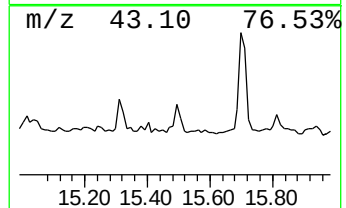
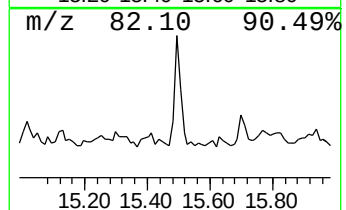
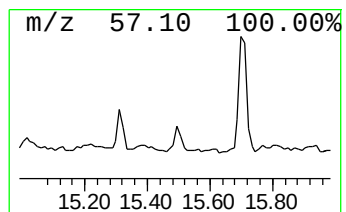
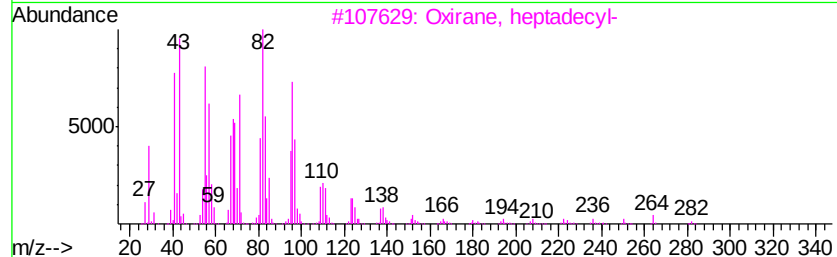
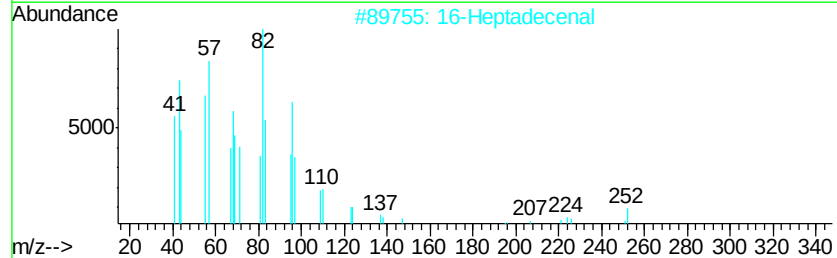
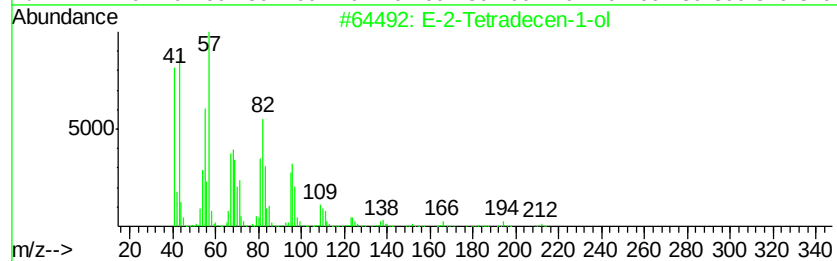
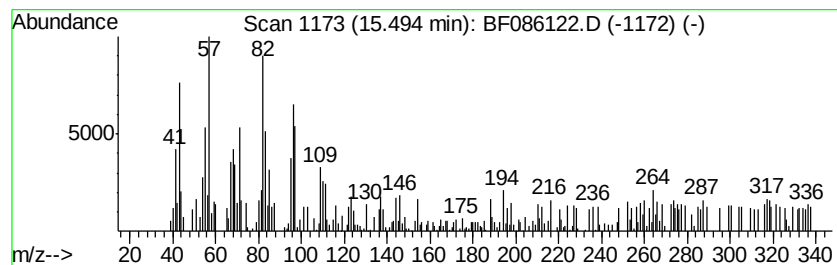
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 E-2-Tetradecen-1-ol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.49	5.86 ng	155334	Perylene-d12	15.30	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	E-2-Tetradecen-1-ol	212	C14H28O	1000130-83-7	86
2	16-Heptadecenal	252	C17H32O	1000144-57-9	70
3	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	70
4	Z-5-Methyl-6-tetradecen-1-ol ace...	268	C17H32O2	1000130-82-2	64
5	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	64



Data Path : Z:\HPCHEM1\BNA_F\Data\BF040616\
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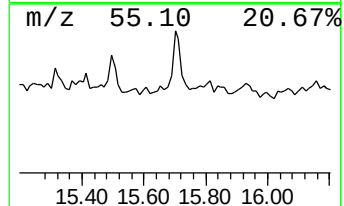
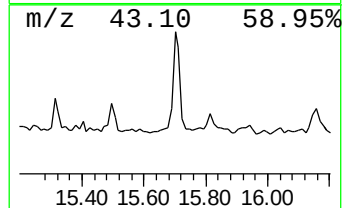
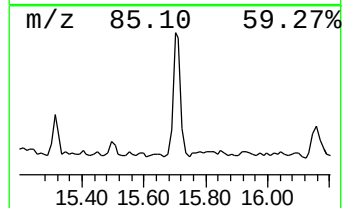
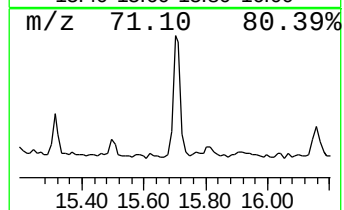
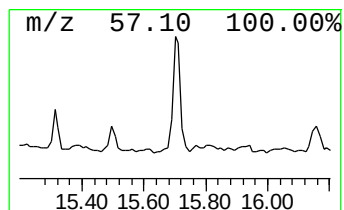
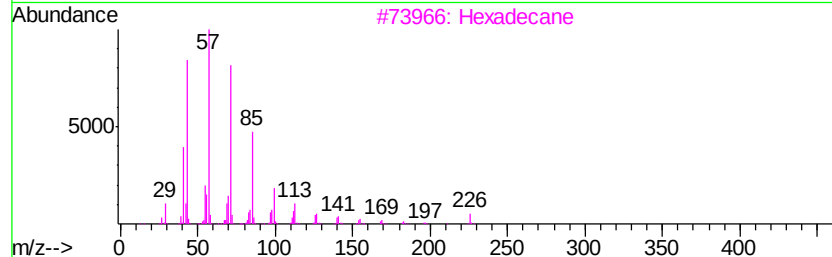
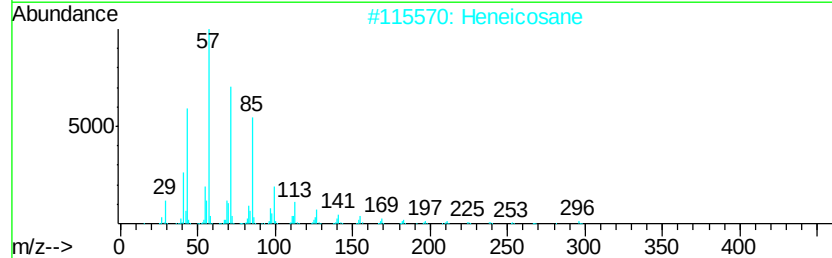
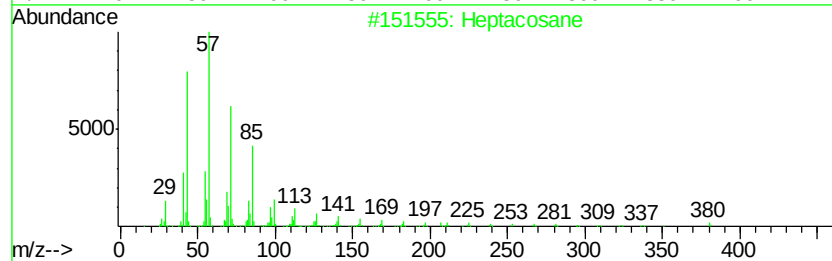
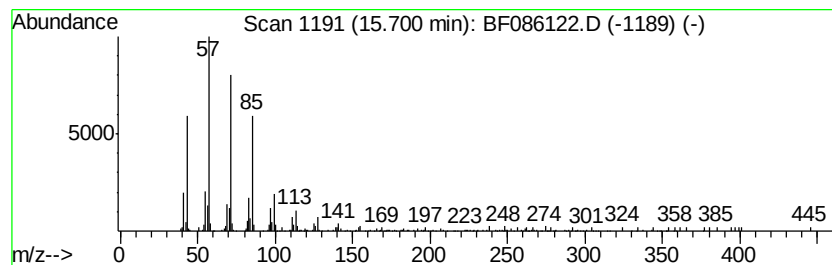
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heptacosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.70	5.46 ng	144606	Perylene-d12	15.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptacosane	380 C27H56	000593-49-7	94
2	Heneicosane	296 C21H44	000629-94-7	91
3	Hexadecane	226 C16H34	000544-76-3	87
4	Pentadecane, 8-heptyl-	310 C22H46	071005-15-7	86
5	Tridecane, 6-propyl-	226 C16H34	055045-10-8	83



Data Path : Z:\HPCHEM1\BNA_F\Data\BF040616\
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ALS Vial : 33 Sample Multiplier: 1

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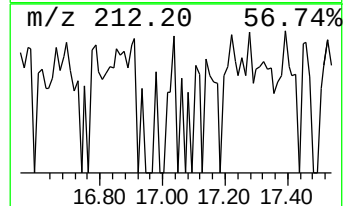
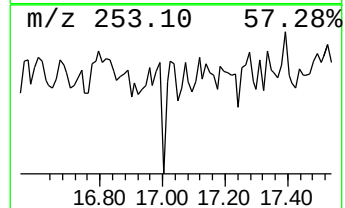
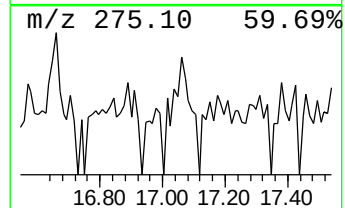
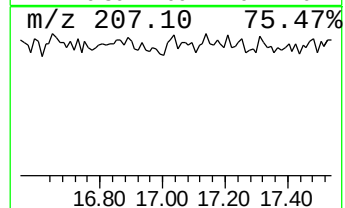
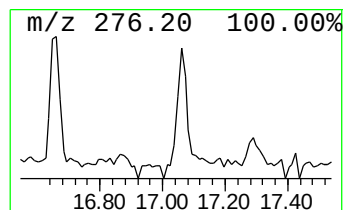
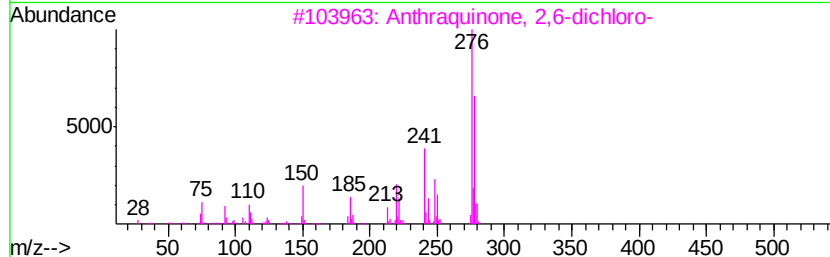
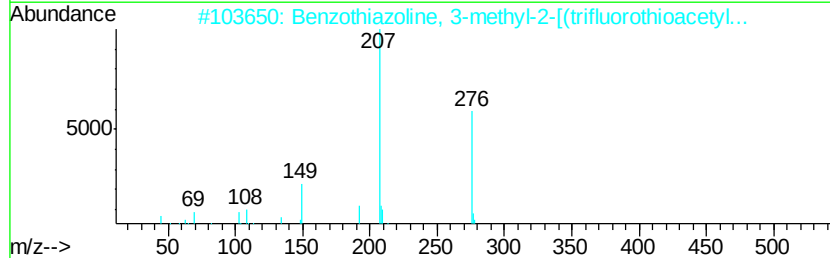
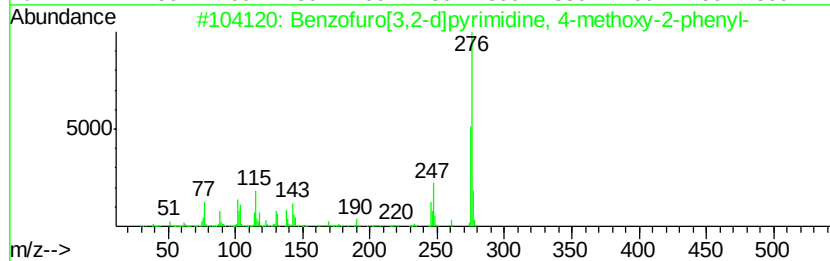
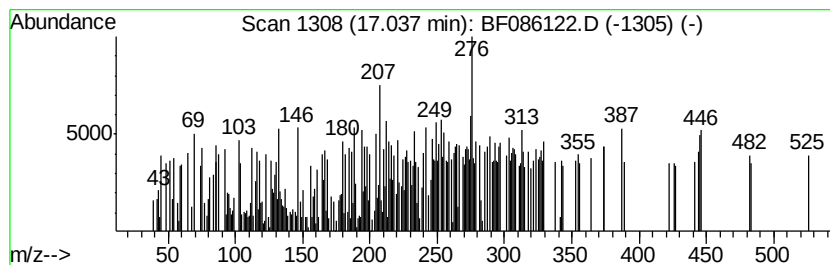
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown17.04 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.04	6.34 ng	167901	Perylene-d12	15.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzofuro[3,2-d]pyrimidine, 4-me...	276	C17H12N2O2	312265-36-4	25
2		Benzothiazoline, 3-methyl-2-[(tr...	276	C10H7F3N2S2	118215-08-0	15
3		Anthraquinone, 2,6-dichloro-	276	C14H6Cl2O2	000605-40-3	15
4		1,3,5-Cyclohexanetrione, tris(di...	291	C15H21N3O3	1000162-62-2	11
5		Benzene, 1,2,3,4-tetrachloro-5,6...	274	C8H6Cl4O2	000944-61-6	11



Data Path : Z:\HPCHEM1\BNA_F\Data\BF040616\
Data File : BF086122.D
Acq On : 7 Apr 2016 2:44
Operator : UM/SJ
Sample : H2104-02
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-1

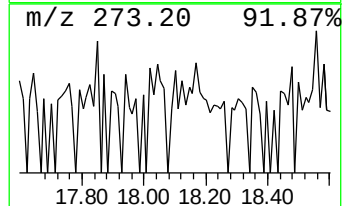
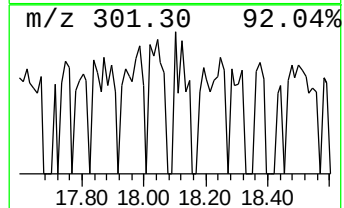
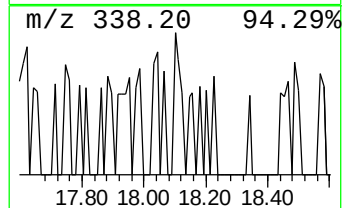
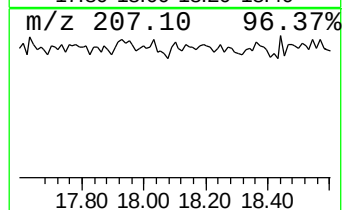
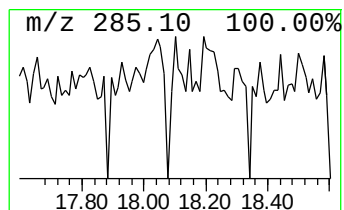
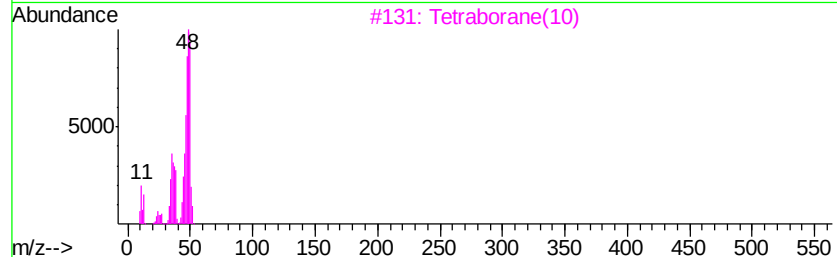
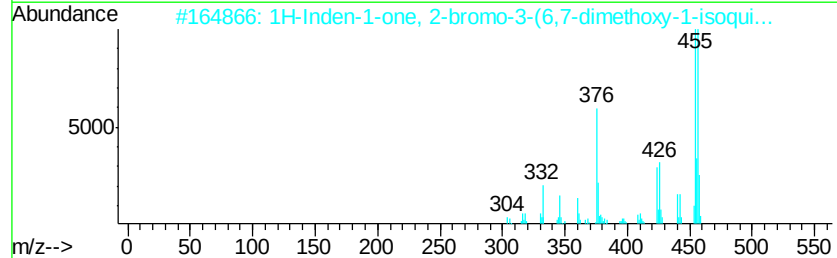
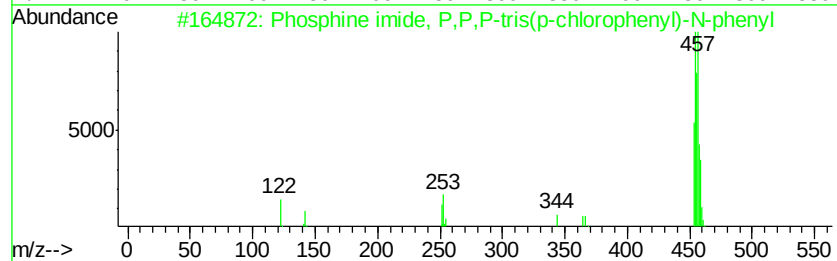
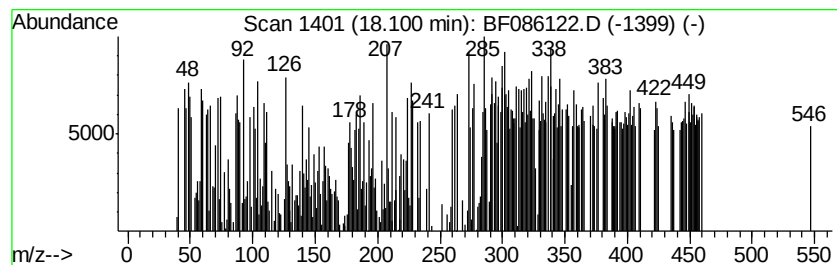
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown18.10 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.10	6.64 ng	175834	Perylene-d12	15.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phosphine imide, P,P,P-tris(p-ch...	455	C24H17Cl3NP	014796-92-0	46
2		1H-Inden-1-one, 2-bromo-3-(6,7-d...	455	C22H18BrN05	023434-75-5	38
3		Tetraborane(10)	54	B4H10	018283-93-7	14
4		1,3,5,7,9,11-Hexaethyl-1-chloroc...	478	C12H35Cl06Si6	073420-32-3	11
5		Epi-epoxy-buphanamine	317	C17H19N05	006835-78-5	10



Data Path : Z:\HPCHEM1\BNA_F\Data\BF040616\
Data File : BF086122.D
Acq On : 7 Apr 2016 2:44
Operator : UM/SJ
Sample : H2104-02
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.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
unknown3.82	3.82	33.8	ng	942263	1	6.74	557796	20.0
2-Pentanone, 4-hy...	4.91	7.5	ng	208849	1	6.74	557796	20.0
unknown6.51	6.51	74.5	ng	2076740	1	6.74	557796	20.0
n-Hexadecanoic acid	11.80	4.1	ng	130516	4	11.26	630070	20.0
2-Chloropropioni...	13.76	6.4	ng	204859	5	13.91	639735	20.0
unknown14.01	14.01	4.4	ng	140418	5	13.91	639735	20.0
13-Docosenamide, ...	14.65	22.4	ng	592253	6	15.30	529814	20.0
10-Undecen-1-yl a...	14.80	4.4	ng	117470	6	15.30	529814	20.0
E-2-Tetradecen-1-ol	15.49	5.9	ng	155334	6	15.30	529814	20.0
Heptacosane	15.70	5.5	ng	144606	6	15.30	529814	20.0
unknown17.04	17.04	6.3	ng	167901	6	15.30	529814	20.0
unknown18.10	18.10	6.6	ng	175834	6	15.30	529814	20.0