

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
 Data File : BF086107.D
 Acq On : 6 Apr 2016 19:35
 Operator : UM/SJ
 Sample : H2254-04
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-17

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	2.213	9	11	13	rVV2	75653	77732	2.65%	0.231%
2	2.270	13	16	20	rVV4	114590	280819	9.59%	0.836%
3	2.350	20	23	27	rVB3	191299	406935	13.90%	1.211%
4	2.407	27	28	31	rVB	29715	38276	1.31%	0.114%
5	2.498	33	36	39	rVB4	47422	88428	3.02%	0.263%
6	2.784	56	61	63	rVB3	75341	122076	4.17%	0.363%
7	2.956	73	76	78	rBV	116829	108048	3.69%	0.322%
8	3.036	82	83	86	rVB	20476	35392	1.21%	0.105%
9	3.138	90	92	96	rBV2	32187	52374	1.79%	0.156%
10	3.287	101	105	106	rBV3	40725	62738	2.14%	0.187%
11	3.344	106	110	111	rBV4	91066	105404	3.60%	0.314%
12	3.436	115	118	121	rVB2	105211	156099	5.33%	0.465%
13	4.739	230	232	236	rVB3	61348	112176	3.83%	0.334%
14	4.853	239	242	243	rBV	30484	46006	1.57%	0.137%
15	4.876	243	244	245	rVB	56001	38389	1.31%	0.114%
16	4.910	245	247	252	rBV	169381	303588	10.37%	0.904%
17	5.242	272	276	278	rVB2	27010	49790	1.70%	0.148%
18	5.310	278	282	283	rBV	29335	44384	1.52%	0.132%
19	5.344	283	285	287	rBV	2412177	2802185	95.70%	8.341%
20	5.744	318	320	324	rVB	30617	32416	1.11%	0.096%
21	6.396	374	377	379	rBV	2390836	2928195	100.00%	8.716%
22	6.510	384	387	390	rBV	2601858	2562753	87.52%	7.629%
23	6.739	405	407	410	rBV	999277	914865	31.24%	2.723%
24	6.899	419	421	423	rBV	2522047	2288155	78.14%	6.811%
25	7.310	455	457	460	rBV	1676573	1690154	57.72%	5.031%
26	8.030	517	520	522	rBV	1259600	1181121	40.34%	3.516%
27	9.105	611	614	617	rBV	2905521	2689808	91.86%	8.007%
28	9.505	647	649	651	rBV	444520	374887	12.80%	1.116%
29	9.722	666	668	671	rBV	22100	35254	1.20%	0.105%
30	9.779	671	673	675	rBV	1387294	1185248	40.48%	3.528%
31	10.213	710	711	713	rVB	67100	50591	1.73%	0.151%
32	10.431	727	730	734	rBV	20913	36726	1.25%	0.109%
33	10.568	739	742	745	rBV	1548026	1599789	54.63%	4.762%
34	10.682	751	752	757	rVB	61558	91416	3.12%	0.272%

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

35	11.128	789	791	797	rVB2	31708	65877	2.25%	0.196%
36	11.265	801	803	804	rBV	1100241	1058943	36.16%	3.152%
37	11.288	804	805	807	rVV	146764	114353	3.91%	0.340%
38	11.345	807	810	812	rVB2	26456	44187	1.51%	0.132%
39	11.505	820	824	827	rBV	23133	35235	1.20%	0.105%
40	11.802	848	850	855	rBV2	88289	162534	5.55%	0.484%
41	11.871	855	856	860	rVB2	20783	35206	1.20%	0.105%
42	12.099	869	876	882	rBV5	22140	60023	2.05%	0.179%
43	12.477	906	909	910	rBV	173660	199917	6.83%	0.595%
44	12.511	910	912	914	rVV3	29694	41271	1.41%	0.123%
45	12.579	917	918	920	rVB	58707	51508	1.76%	0.153%
46	12.694	925	928	931	rBV2	120257	189587	6.47%	0.564%
47	12.842	939	941	944	rBV	1938619	2015454	68.83%	5.999%
48	13.037	954	958	959	rBV	18508	44554	1.52%	0.133%
49	13.094	959	963	965	rVB4	16116	41366	1.41%	0.123%
50	13.139	965	967	969	rBV	42492	48105	1.64%	0.143%
51	13.254	973	977	978	rBV3	18095	34521	1.18%	0.103%
52	13.425	989	992	993	rBV3	18256	30482	1.04%	0.091%
53	13.585	1004	1006	1009	rVB4	27670	63675	2.17%	0.190%
54	13.654	1009	1012	1013	rBV	19086	36433	1.24%	0.108%
55	13.700	1013	1016	1018	rBV	91872	119207	4.07%	0.355%
56	13.745	1018	1020	1023	rBV2	289338	400750	13.69%	1.193%
57	13.894	1030	1033	1038	rBV	913579	1119330	38.23%	3.332%
58	13.974	1038	1040	1043	rBV3	41346	65390	2.23%	0.195%
59	14.088	1048	1050	1051	rBV2	26913	34980	1.19%	0.104%
60	14.202	1057	1060	1061	rBV3	28920	51815	1.77%	0.154%
61	14.282	1065	1067	1069	rVV	36805	59987	2.05%	0.179%
62	14.385	1071	1076	1082	rVB8	69211	295240	10.08%	0.879%
63	14.522	1084	1088	1091	rBV5	40305	143846	4.91%	0.428%
64	14.728	1104	1106	1108	rBV2	27740	34831	1.19%	0.104%
65	14.911	1119	1122	1126	rBV	109084	228044	7.79%	0.679%
66	15.185	1144	1146	1149	rBV	84326	101127	3.45%	0.301%
67	15.243	1149	1151	1153	rVV	95554	109892	3.75%	0.327%
68	15.300	1153	1156	1161	rVB	768480	926344	31.64%	2.757%
69	15.563	1177	1179	1183	rBV2	54700	119860	4.09%	0.357%
70	16.088	1222	1225	1228	rVB2	69982	127843	4.37%	0.381%
71	16.157	1228	1231	1233	rBV4	30782	88409	3.02%	0.263%

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Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

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72	16.294	1241	1243	1246	rVB2	37068	49769	1.70%	0.148%
73	16.385	1250	1251	1253	rBV2	26094	46101	1.57%	0.137%
74	16.637	1271	1273	1277	rBV	41172	108281	3.70%	0.322%
75	16.808	1286	1288	1290	rBV2	24982	43224	1.48%	0.129%
76	17.060	1307	1310	1311	rBV	92048	146656	5.01%	0.437%
77	17.083	1311	1312	1315	rVV3	108278	166521	5.69%	0.496%
78	17.163	1315	1319	1320	rVV3	278191	397816	13.59%	1.184%
79	17.220	1323	1324	1326	rVV2	40164	58476	2.00%	0.174%
80	17.346	1326	1335	1336	rVV8	279405	643465	21.97%	1.915%
81	17.380	1336	1338	1341	rVB4	110618	217207	7.42%	0.647%
82	17.768	1368	1372	1379	rVB9	49464	264510	9.03%	0.787%
83	18.043	1393	1396	1401	rVB7	24669	58404	1.99%	0.174%
84	18.157	1404	1406	1407	rVB2	51025	50918	1.74%	0.152%
85	18.191	1407	1409	1412	rVB4	22379	40652	1.39%	0.121%
86	18.249	1412	1414	1415	rBV	30799	32148	1.10%	0.096%
87	18.969	1474	1477	1483	rVB3	82083	246503	8.42%	0.734%
88	19.289	1503	1505	1508	rVB4	16352	30761	1.05%	0.092%

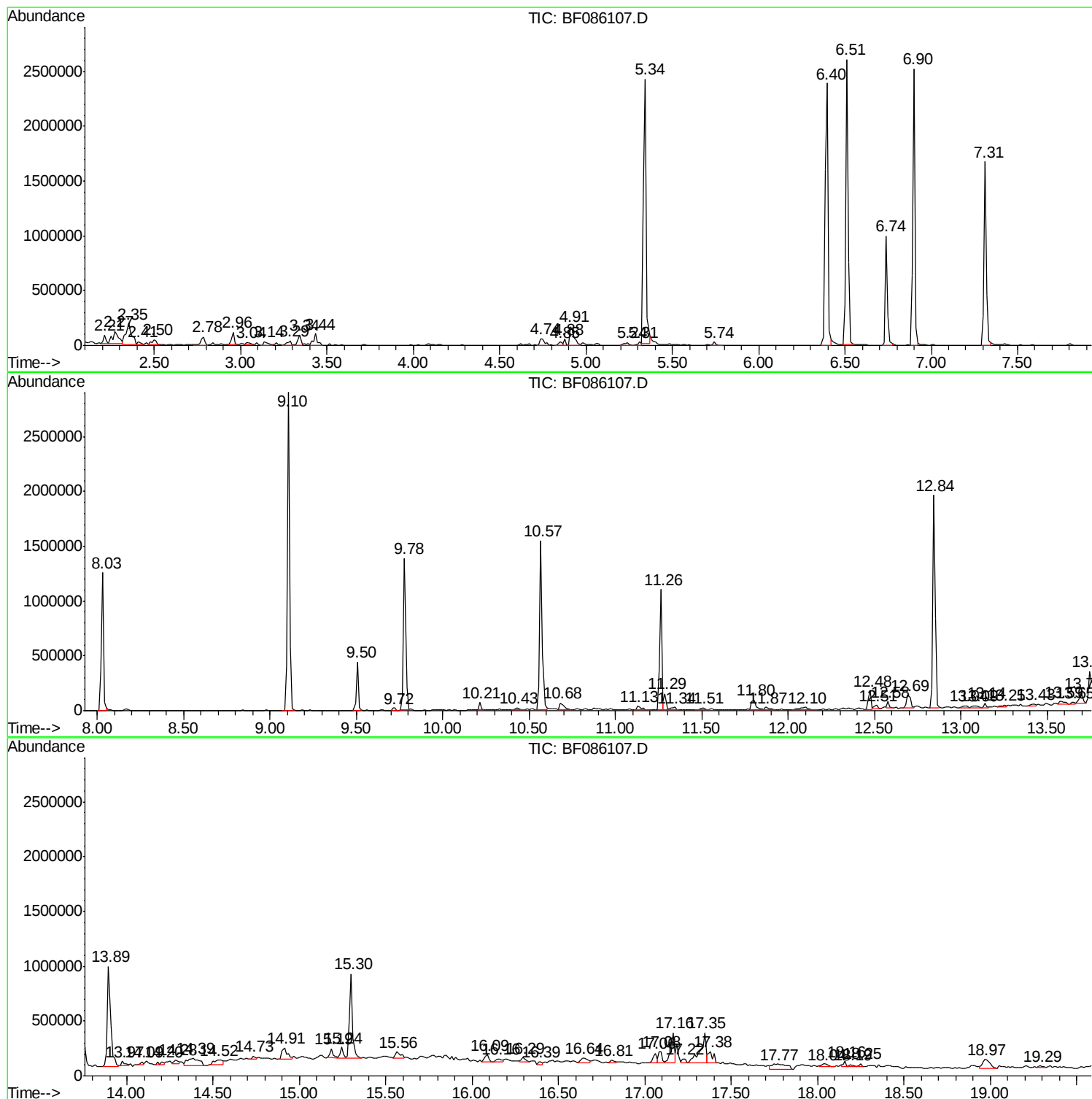
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Instrument :
BNA_F
ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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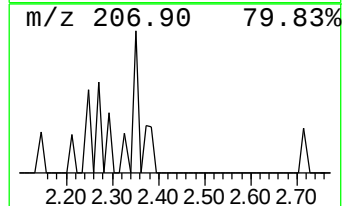
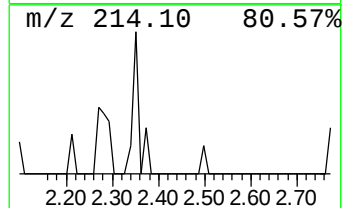
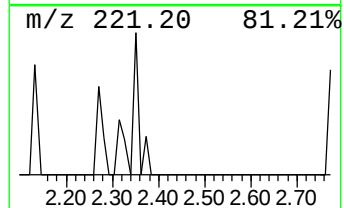
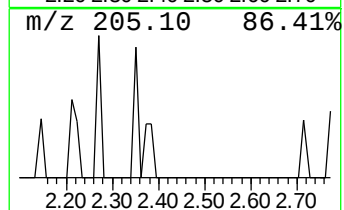
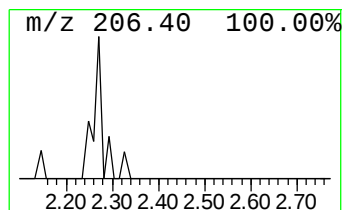
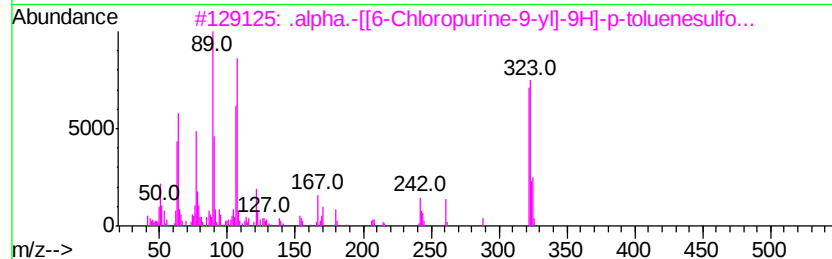
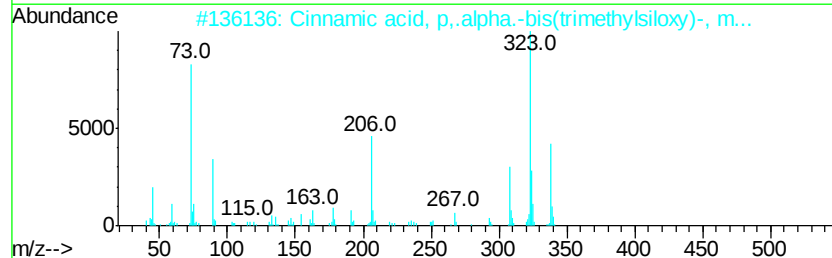
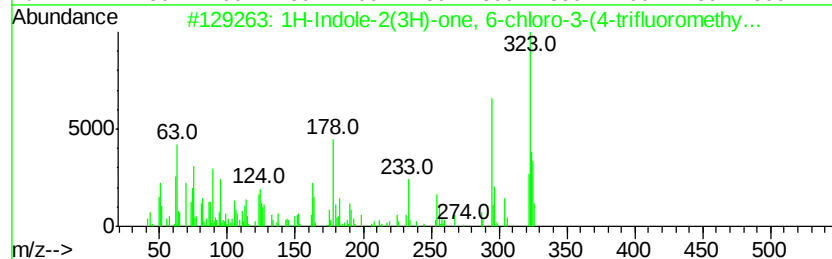
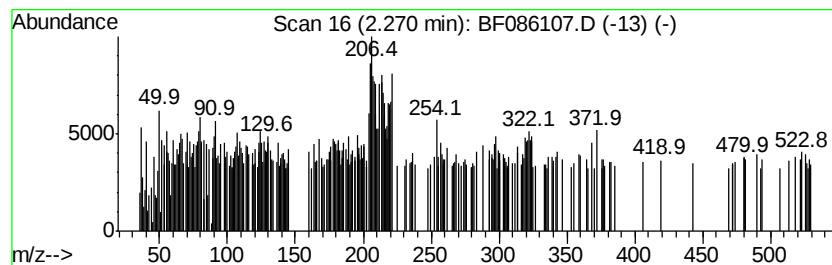
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown2.27 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.27	6.14 ng	280819	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indole-2(3H)-one, 6-chloro-3-...	323	C16H9ClF3NO	035315-59-4	32
2		Cinnamic acid, p,.alpha.-bis(tri...	338	C16H26O4Si2	029372-28-9	23
3		.alpha.-[[6-Chloropurine-9-yl]-9...	323	C12H10ClN5O2S	021267-95-8	23
4		Quinoline, 1,2-dihydro-7-methoxy...	338	C20H22N2O3	300589-50-8	9
5		.eta.-Pentamethylcyclopentadieny...	527	C19H32IMoN3	1000288-06-1	7



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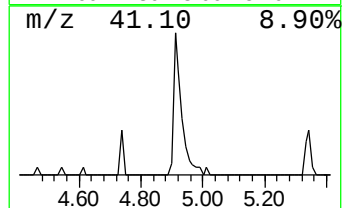
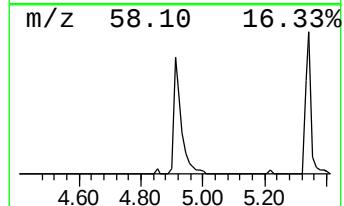
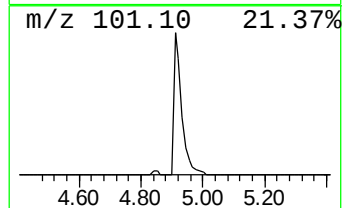
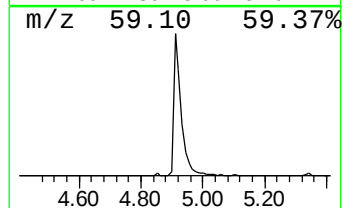
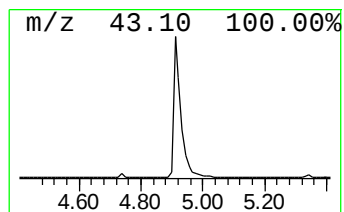
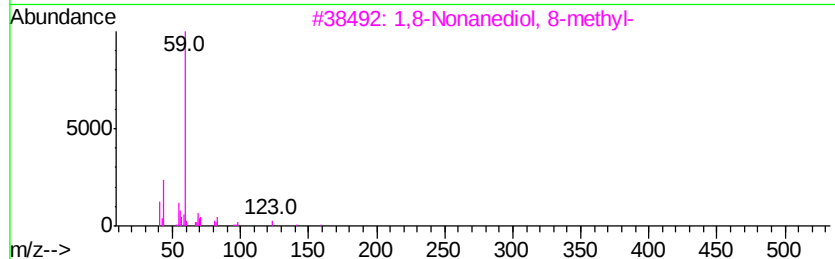
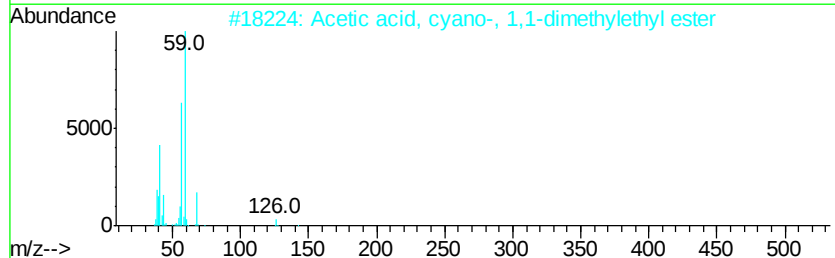
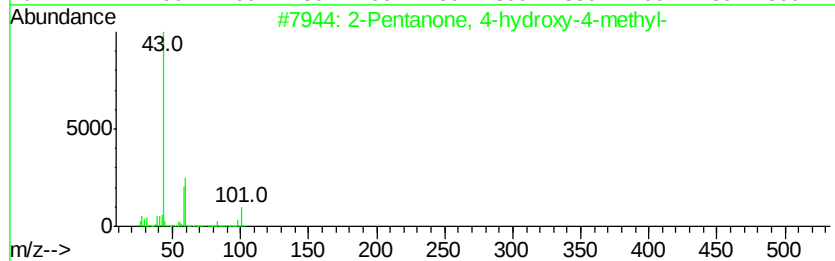
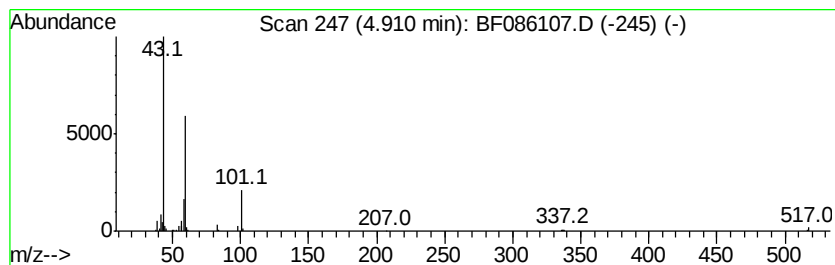
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.91	6.64 ng	303588	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	59
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	12
3			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	12
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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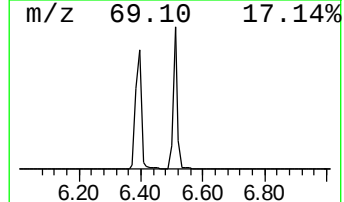
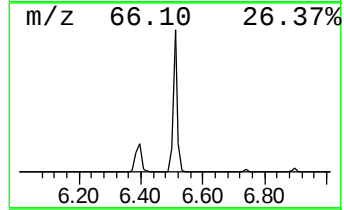
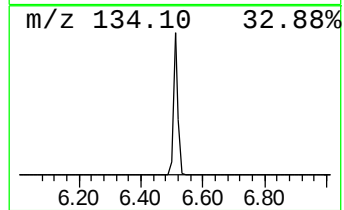
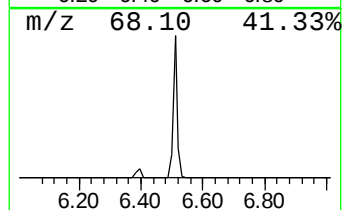
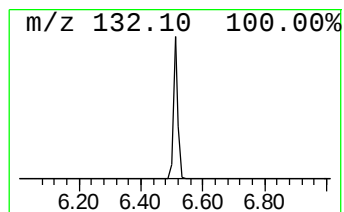
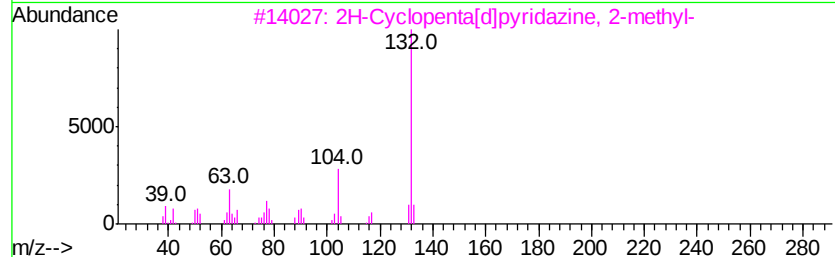
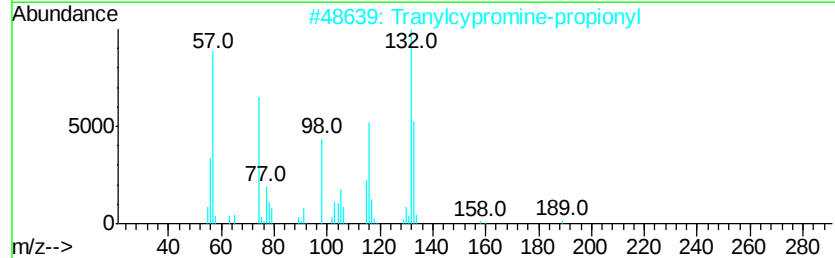
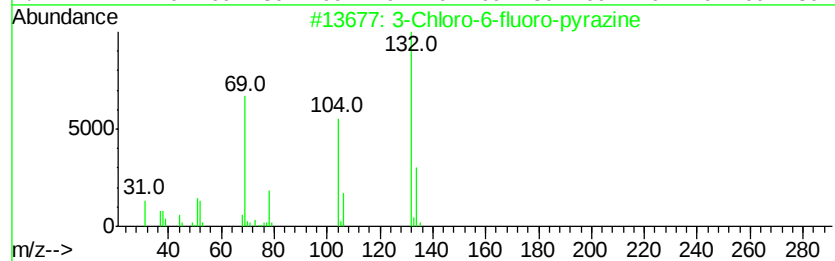
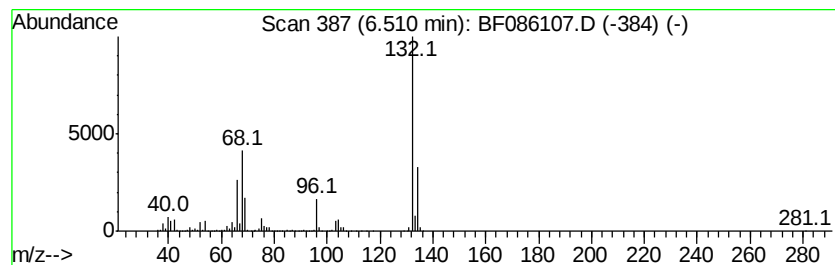
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 unknown6.51 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	56.02 ng	2562750	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		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
3		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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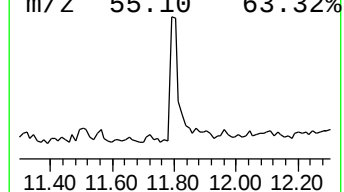
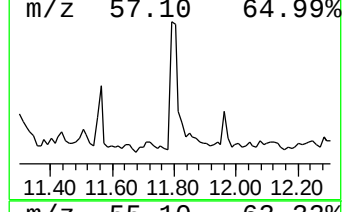
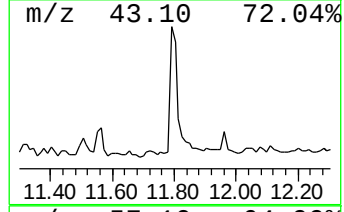
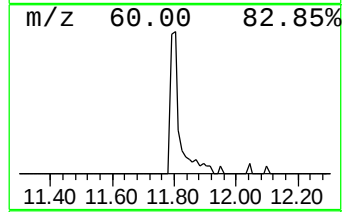
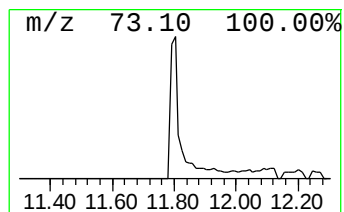
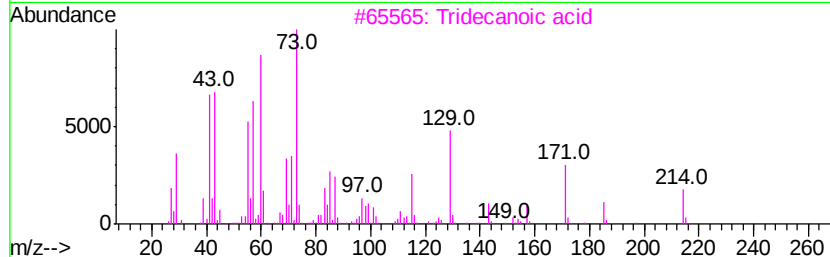
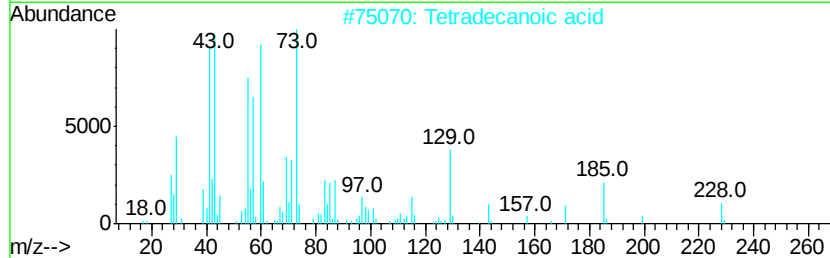
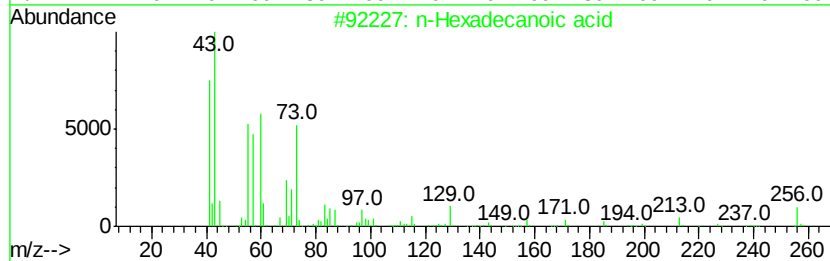
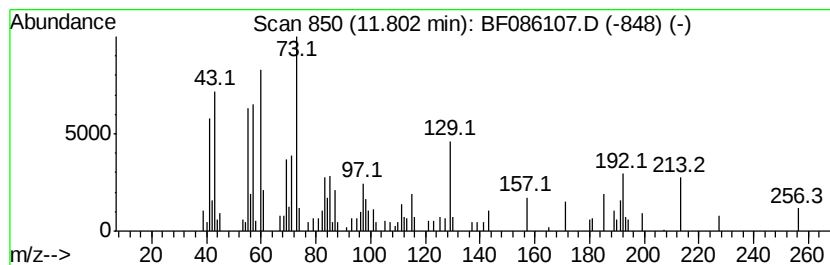
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.80	3.07 ng	162534	Phenanthrene-d10	11.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	74
3			Tridecanoic acid	214	C13H26O2	000638-53-9	64
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	52
5			Dodecanoic acid	200	C12H24O2	000143-07-7	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
Data File : BF086107.D
Acq On : 6 Apr 2016 19:35
Operator : UM/SJ
Sample : H2254-04
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-17

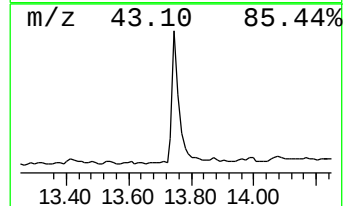
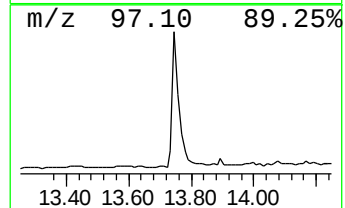
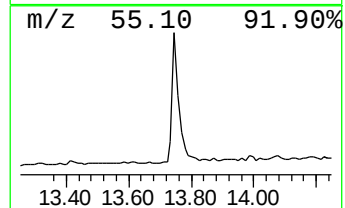
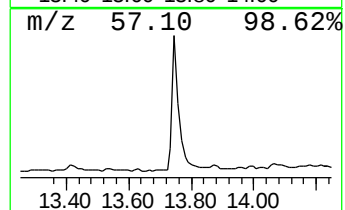
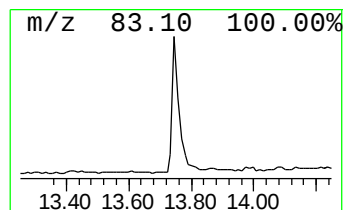
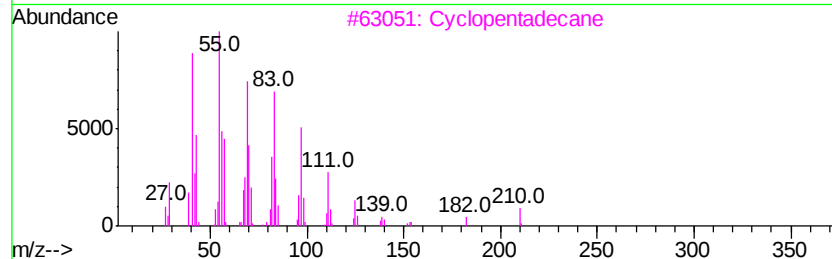
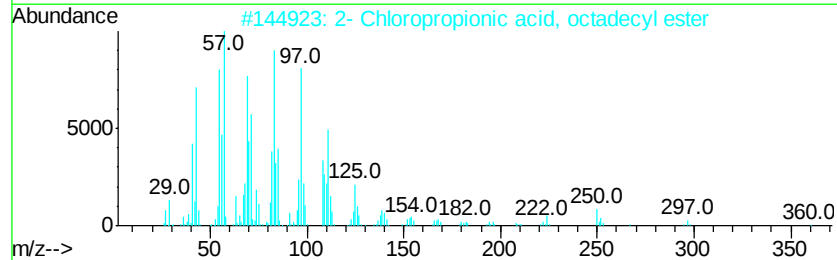
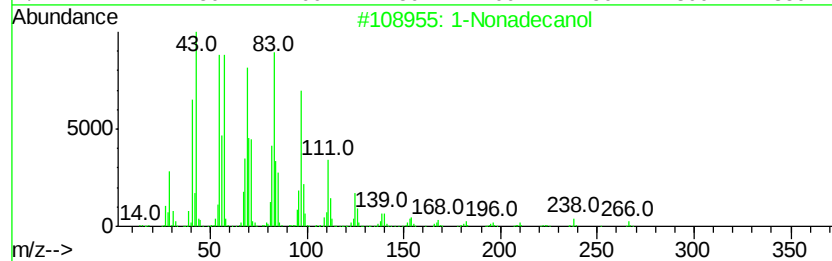
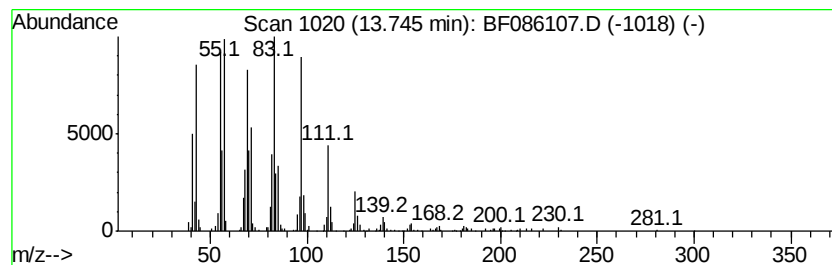
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1-Nonadecanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.75	7.16 ng	400750	Chrysene-d12	13.89

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Nonadecanol	284	C19H40O	001454-84-8	93
2			2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	93
3			Cyclopentadecane	210	C15H30	000295-48-7	93
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 : BF086107.D
Acq On : 6 Apr 2016 19:35
Operator : UM/SJ
Sample : H2254-04
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-17

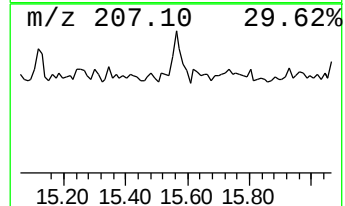
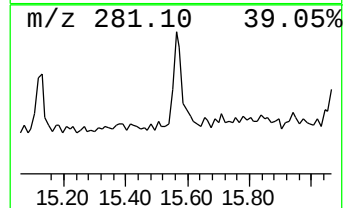
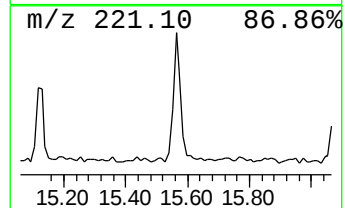
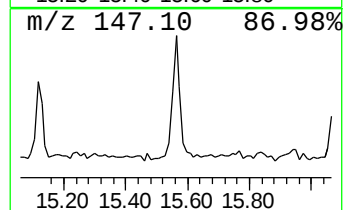
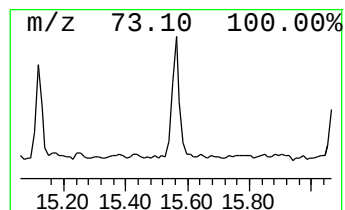
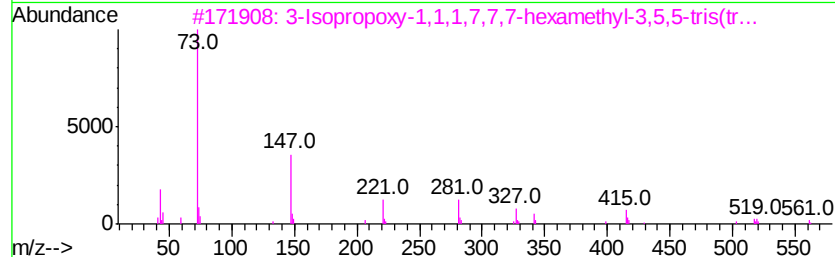
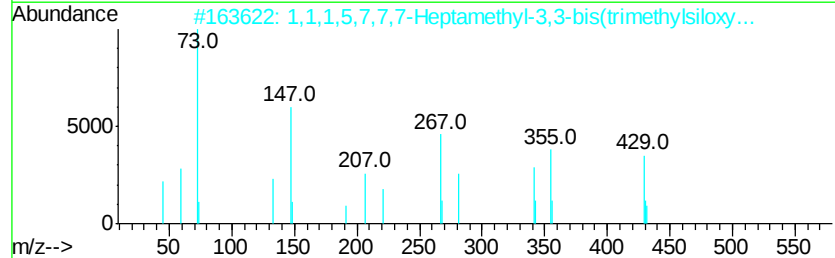
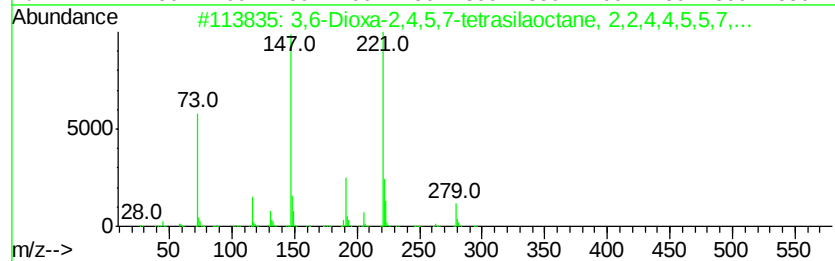
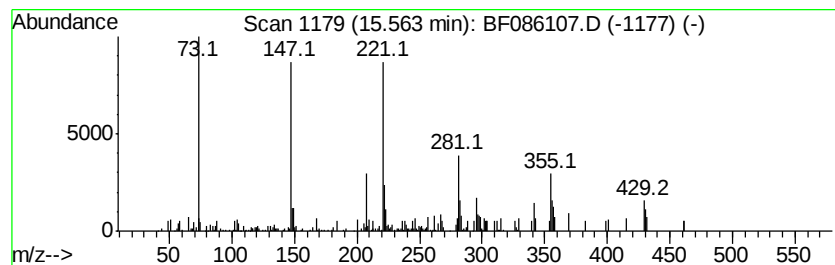
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 3,6-Dioxa-2,4,5,7-tetrasila... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.56	2.59 ng	119860	Perylene-d12	15.30

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H3002Si4	004342-25-0	50
2			1,1,1,5,7,7,7-Heptamethyl-3,3-bi...	444	C13H4005Si6	038147-00-1	38
3			3-Isopropoxv-1,1,1,7,7,7-hexamet...	576	C18H5207Si7	071579-69-6	32
4			N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	30
5			Pentasiloxane, dodecamethyl-	384	C12H3604Si5	000141-63-9	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
Data File : BF086107.D
Acq On : 6 Apr 2016 19:35
Operator : UM/SJ
Sample : H2254-04
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-17

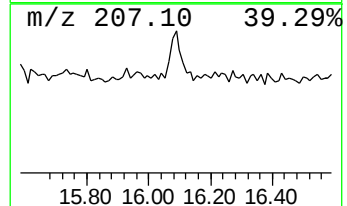
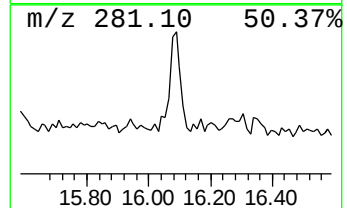
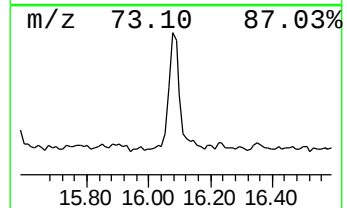
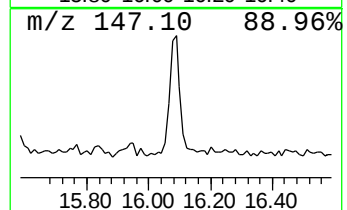
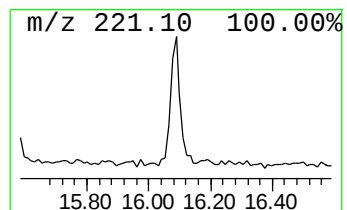
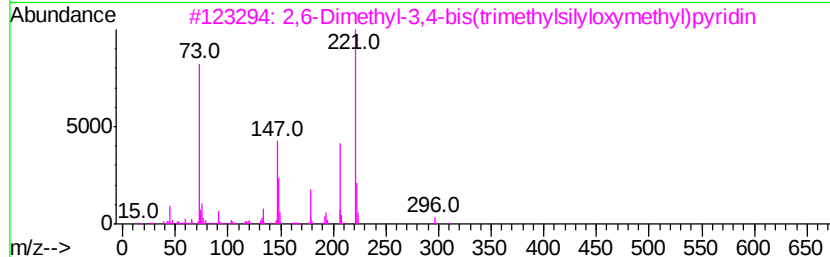
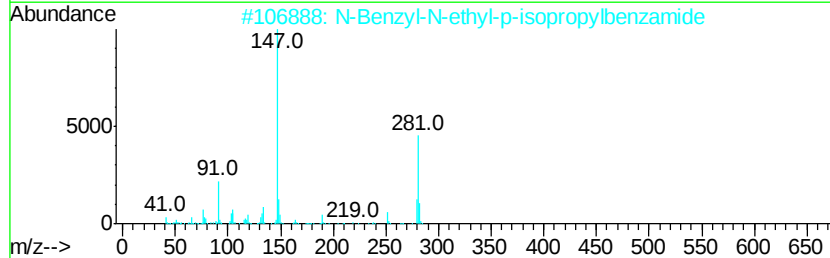
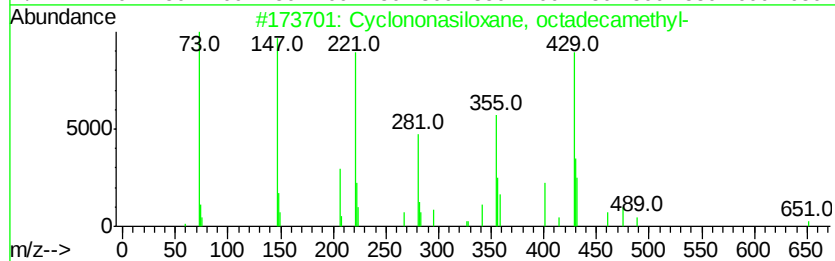
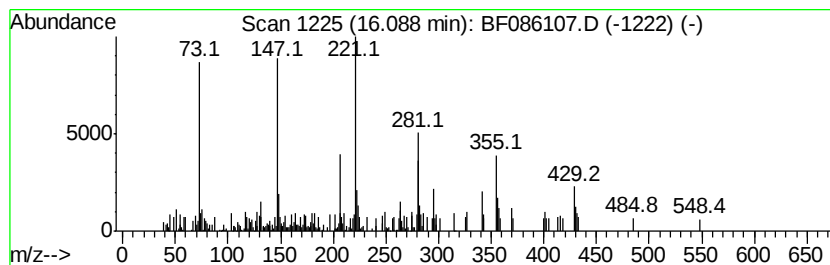
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Cyclononasiloxane, octadeca... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.09	2.76 ng	127843	Perylene-d12	15.30

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclononasiloxane, octadecamethyl-	666	C18H5409Si9	000556-71-8	90
2			N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	25
3			2,6-Dimethyl-3,4-bis(trimethylsilo...	311	C15H29N02Si2	1000079-52-1	12
4			Cyclohexasiloxane, dodecamethyl-	444	C12H3606Si6	000540-97-6	11
5			Pentasiloxane, dodecamethyl-	384	C12H3604Si5	000141-63-9	10



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040616\
Data File : BF086107.D
Acq On : 6 Apr 2016 19:35
Operator : UM/SJ
Sample : H2254-04
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SP-17

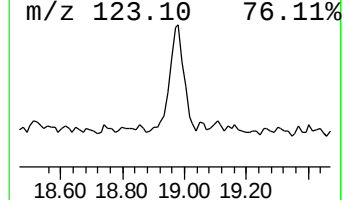
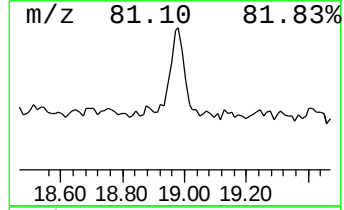
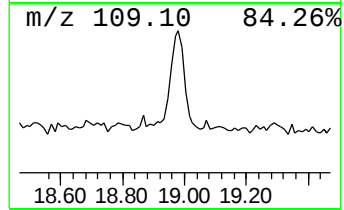
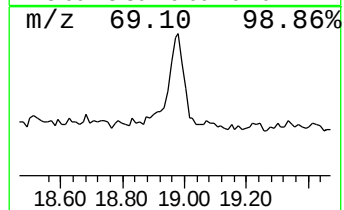
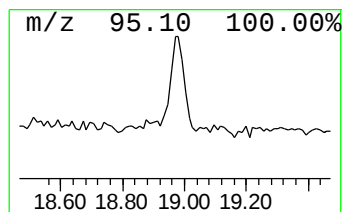
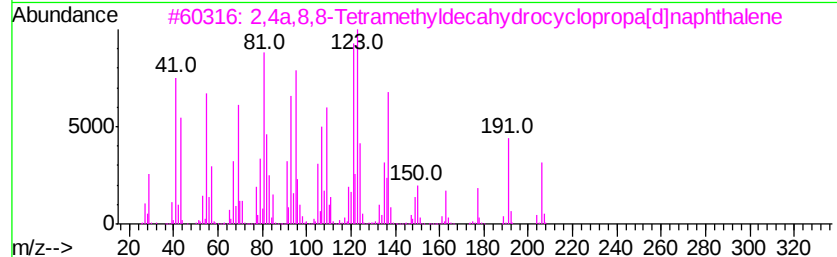
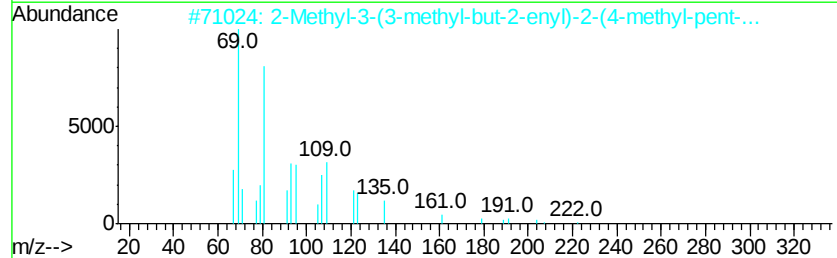
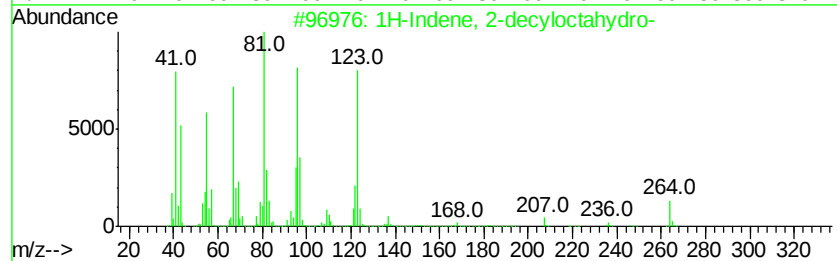
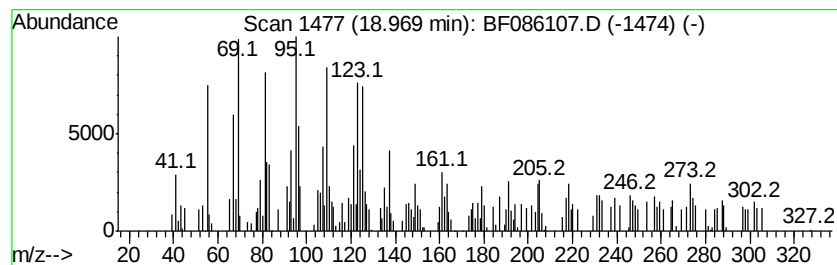
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 1H-Indene, 2-decyloctahydro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.97	5.32 ng	246503	Perylene-d12	15.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2-decvloctahvdro-	264	C19H36	055044-34-3	51
2		2-Methyl-3-(3-methyl-but-2-enyl)...	222	C15H26O	1000144-10-2	46
3		2,4a,8,8-Tetramethyldecahvdrocyc...	206	C15H26	074022-04-1	42
4		1-Formyl-2,2-dimethyl-3-trans-(3...	220	C15H24O	1000144-09-7	38
5		2,2,6-Trimethyl-1-(2-methyl-cycl...	218	C15H22O	1000188-72-8	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
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Sample : H2254-04
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-17

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
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2-Pentanone, 4-hy...	4.91	6.6	ng	303588	1	6.74	914865	20.0
unknown6.51	6.51	56.0	ng	2562750	1	6.74	914865	20.0
n-Hexadecanoic acid	11.80	3.1	ng	162534	4	11.27	1058940	20.0
1-Nonadecanol	13.75	7.2	ng	400750	5	13.89	1119330	20.0
3,6-Dioxa-2,4,5,7...	15.56	2.6	ng	119860	6	15.30	926344	20.0
Cyclononasiloxane...	16.09	2.8	ng	127843	6	15.30	926344	20.0
1H-Indene, 2-decy...	18.97	5.3	ng	246503	6	15.30	926344	20.0