

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120115\
 Data File : BF083388.D
 Acq On : 1 Dec 2015 22:58
 Operator : UM/IZ
 Sample : G4593-08
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S24-15.5

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.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	3.904	191	194	204	rBV	61732	120789	2.18%	0.221%
2	4.464	241	243	248	rBV	34891	63547	1.15%	0.117%
3	4.659	257	260	275	rVB	868490	1544062	27.92%	2.831%
4	5.139	300	302	317	rBV	4528842	5107130	92.35%	9.364%
5	6.225	394	397	399	rBV	3841104	5464769	98.82%	10.019%
6	6.327	403	406	408	rBV	3909644	5104946	92.31%	9.360%
7	6.545	423	425	428	rBV	995614	969249	17.53%	1.777%
8	6.705	437	439	441	rBV	3894023	3483982	63.00%	6.388%
9	7.116	473	475	478	rBV	2495588	3232583	58.45%	5.927%
10	7.276	486	489	495	rBV	187960	285831	5.17%	0.524%
11	7.836	535	538	540	rBV	1389596	1272907	23.02%	2.334%
12	8.922	630	633	635	rBV	4446222	5530215	100.00%	10.139%
13	9.322	666	668	670	rBV	1152263	1071054	19.37%	1.964%
14	9.448	675	679	681	rBV2	26670	59290	1.07%	0.109%
15	9.539	685	687	689	rBV	134383	132557	2.40%	0.243%
16	9.585	689	691	693	rBV	1320537	1383126	25.01%	2.536%
17	9.791	705	709	711	rBV3	42238	75851	1.37%	0.139%
18	10.042	723	731	732	rBV	224142	324860	5.87%	0.596%
19	10.122	737	738	740	rVB	50995	56916	1.03%	0.104%
20	10.259	747	750	751	rBV	86063	122611	2.22%	0.225%
21	10.282	751	752	756	rVB3	38337	68472	1.24%	0.126%
22	10.374	757	760	763	rVB	3264975	3374993	61.03%	6.188%
23	10.522	771	773	777	rBV	202097	339699	6.14%	0.623%
24	10.968	809	812	814	rBV3	33898	59830	1.08%	0.110%
25	11.014	814	816	818	rBV2	150873	170386	3.08%	0.312%
26	11.116	823	825	830	rBV2	986586	2035850	36.81%	3.733%
27	11.208	830	833	836	rVV	135478	180406	3.26%	0.331%
28	11.528	858	861	864	rVB	66071	103366	1.87%	0.190%
29	11.734	876	879	880	rBV	110232	175665	3.18%	0.322%
30	11.768	880	882	885	rVV	88121	129480	2.34%	0.237%
31	11.837	885	888	889	rVV	461205	728074	13.17%	1.335%
32	11.859	889	890	892	rVV	143363	185236	3.35%	0.340%
33	11.894	892	893	899	rVB2	65071	90624	1.64%	0.166%
34	12.111	910	912	914	rBV	49175	65413	1.18%	0.120%

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35	12.145	914	915	920	rVB3	32356	81966	1.48%	0.150%
36	12.385	932	936	938	rVB2	43570	69122	1.25%	0.127%
37	12.465	939	943	946	rVV2	78845	172358	3.12%	0.316%
38	12.648	957	959	962	rBV	456927	673024	12.17%	1.234%
39	12.785	966	971	974	rBV4	45281	95061	1.72%	0.174%
40	12.899	979	981	984	rBV	130699	234308	4.24%	0.430%
41	12.957	984	986	989	rVV	462886	775217	14.02%	1.421%
42	13.220	1005	1009	1012	rVV	3551463	4696483	84.92%	8.611%
43	13.300	1012	1016	1018	rVB2	30764	63022	1.14%	0.116%
44	13.448	1027	1029	1032	rVB	57015	86833	1.57%	0.159%
45	13.597	1040	1042	1044	rBV2	55907	80072	1.45%	0.147%
46	14.362	1105	1109	1111	rBV	35289	55943	1.01%	0.103%
47	14.408	1111	1113	1117	rVV2	30986	59931	1.08%	0.110%
48	14.671	1133	1136	1139	rBV	79611	192162	3.47%	0.352%
49	14.740	1139	1142	1145	rVV2	775748	1400348	25.32%	2.567%
50	14.785	1145	1146	1148	rVB	168729	139995	2.53%	0.257%
51	15.357	1193	1196	1199	rBV	32839	64942	1.17%	0.119%
52	16.077	1256	1259	1263	rBV4	22093	58522	1.06%	0.107%
53	16.237	1270	1273	1276	rVB	187559	253763	4.59%	0.465%
54	16.294	1276	1278	1281	rVV	122262	261472	4.73%	0.479%
55	16.340	1281	1282	1286	rVB	58371	57000	1.03%	0.105%
56	16.443	1286	1291	1294	rBV2	26593	55344	1.00%	0.101%
57	16.671	1308	1311	1314	rBV	68309	103818	1.88%	0.190%
58	16.740	1314	1317	1320	rVB	116876	164540	2.98%	0.302%
59	16.808	1320	1323	1329	rBV	607468	941984	17.03%	1.727%
60	17.060	1339	1345	1351	rBV7	15817	64481	1.17%	0.118%
61	18.191	1440	1444	1449	rVB2	63464	130683	2.36%	0.240%
62	18.569	1473	1477	1483	rBV	53629	115959	2.10%	0.213%
63	18.900	1500	1506	1515	rVB3	18556	97300	1.76%	0.178%
64	20.077	1603	1609	1619	rBV7	18415	103195	1.87%	0.189%
65	20.466	1639	1643	1651	rVB3	19084	79573	1.44%	0.146%

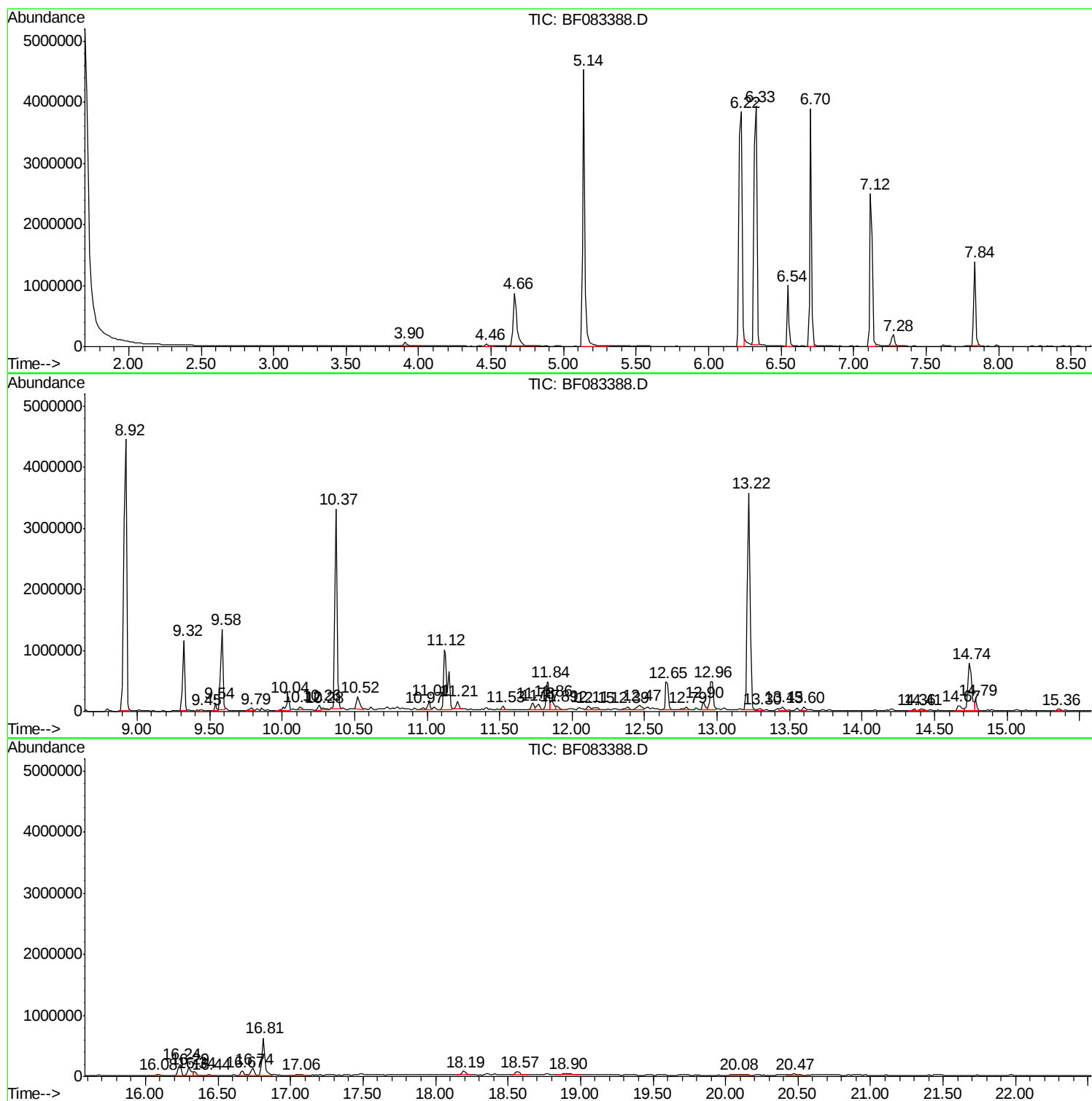
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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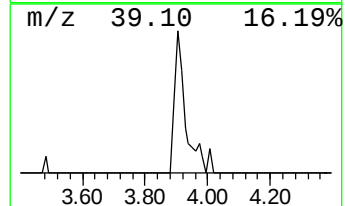
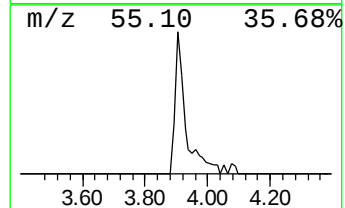
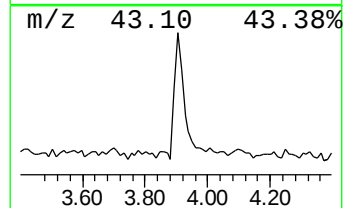
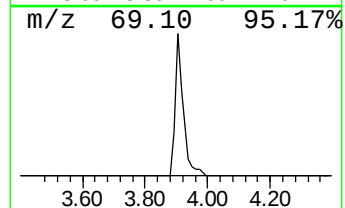
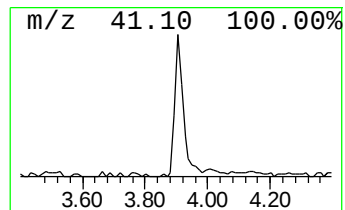
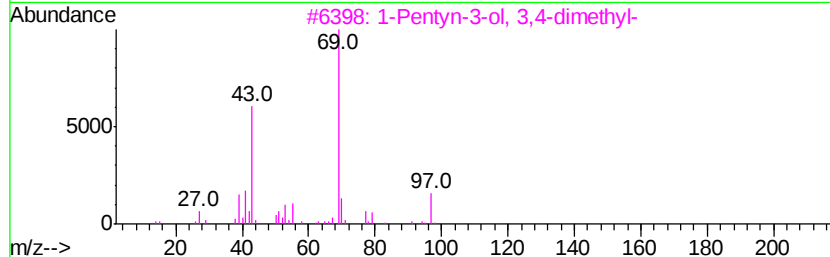
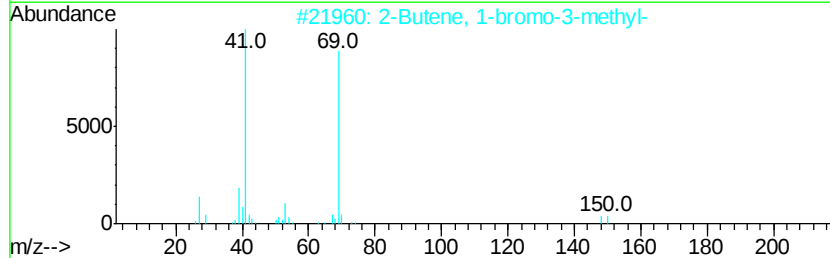
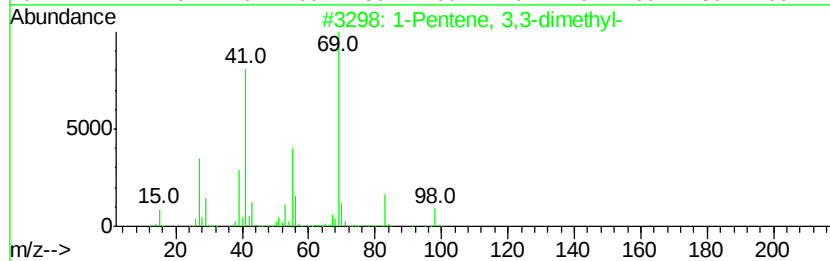
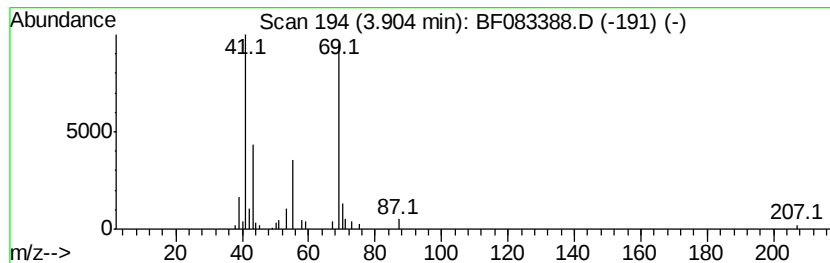
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 1-Pentene, 3,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.90	2.49 ng	120789	1,4-Dichlorobenzene-d4	6.54

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Pentene, 3,3-dimethyl-	98	C7H14	003404-73-7	50
2			2-Butene, 1-bromo-3-methyl-	148	C5H9Br	000870-63-3	45
3			1-Pentyn-3-ol, 3,4-dimethyl-	112	C7H12O	001482-15-1	42
4			1-Butene, 3,3-dimethyl-	84	C6H12	000558-37-2	39
5			2-Butenoic acid, 4-nitrophenyl e...	207	C10H9NO4	014617-88-0	38



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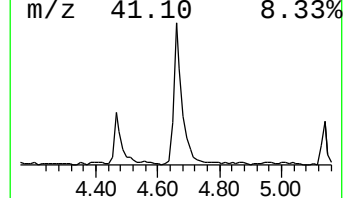
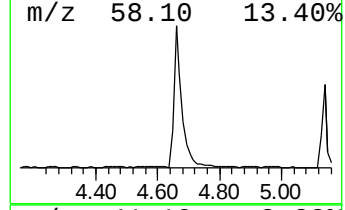
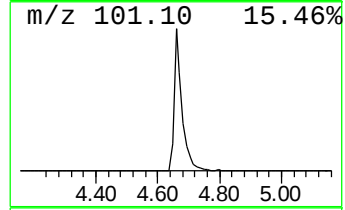
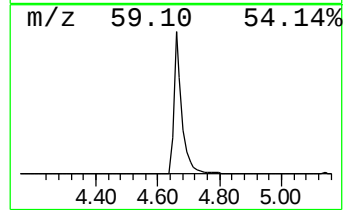
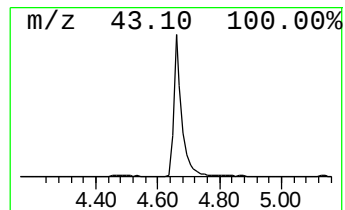
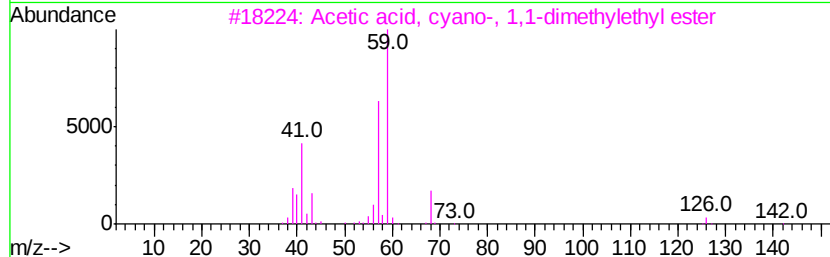
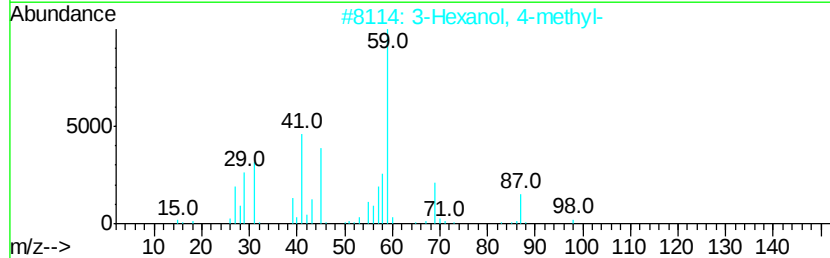
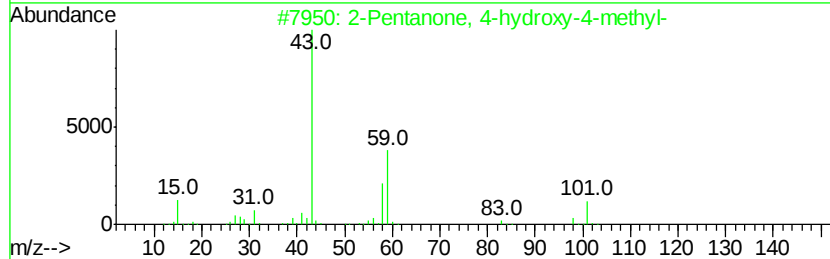
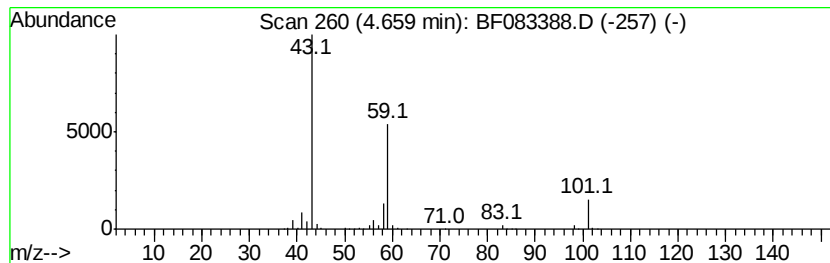
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.66	31.86 ng	1544060	1,4-Dichlorobenzene-d4	6.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
4		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
5		3-Pentanol	88	C5H12O	000584-02-1	9



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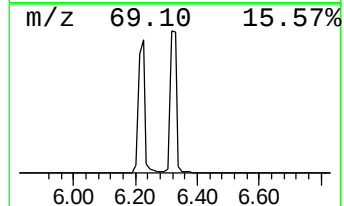
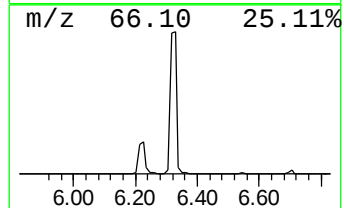
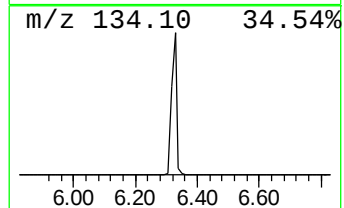
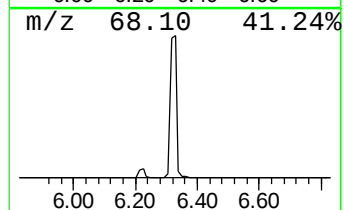
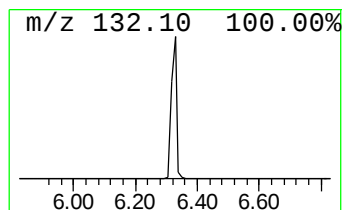
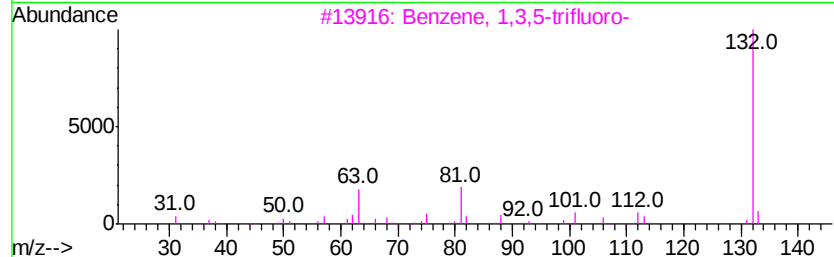
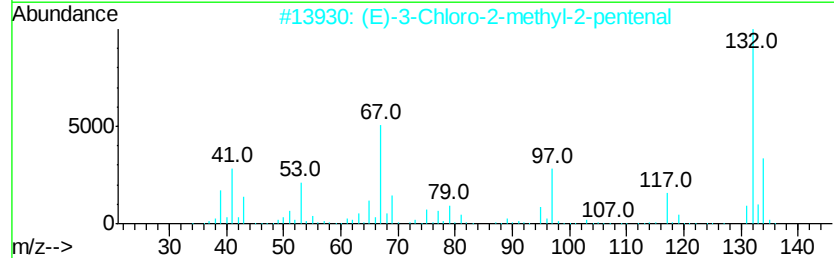
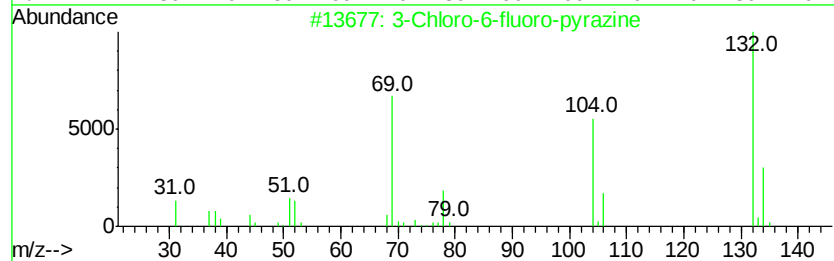
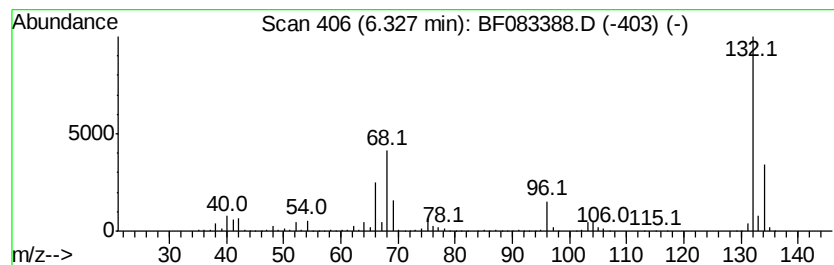
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.33 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.33	105.34 ng	5104950	1,4-Dichlorobenzene-d4	6.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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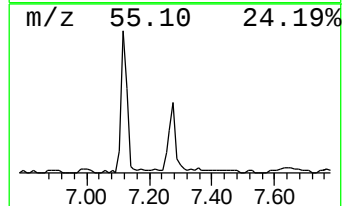
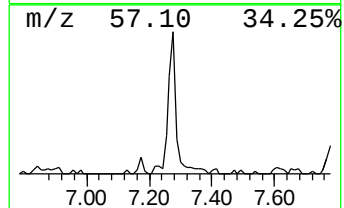
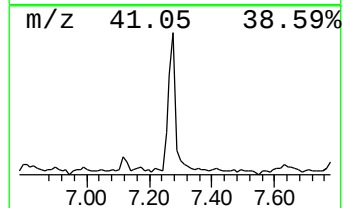
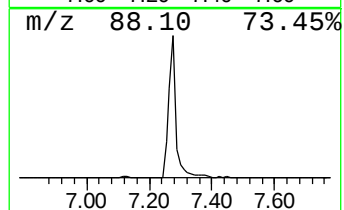
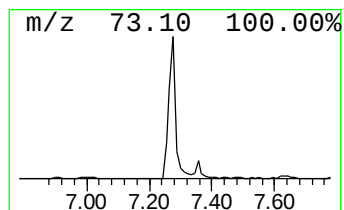
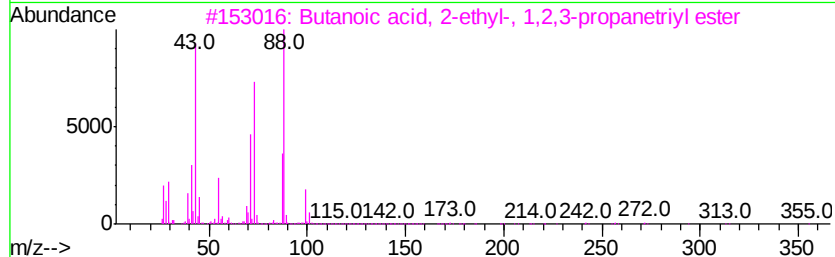
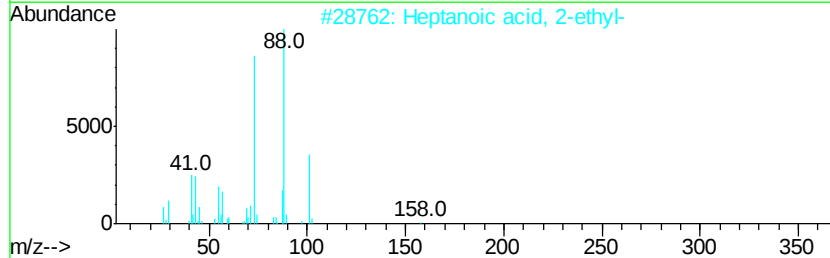
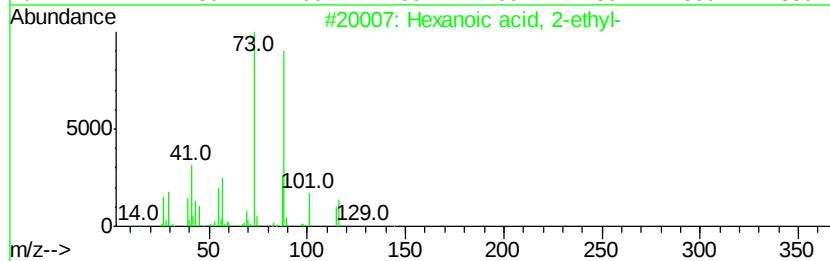
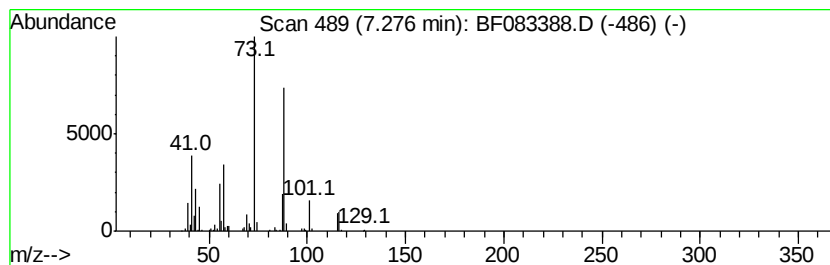
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 Hexanoic acid, 2-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.28	4.49 ng	285831	Naphthalene-d8	7.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
2		Heptanoic acid, 2-ethyl-	158	C9H18O2	003274-29-1	59
3		Butanoic acid, 2-ethyl-, 1,2,3-p...	386	C21H38O6	056554-54-2	53
4		Acetic acid, diethyl-	116	C6H12O2	000088-09-5	40
5		Octanoic acid, 2-methyl-, methyl...	172	C10H20O2	002177-86-8	40



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Misc :
ALS Vial : 13 Sample Multiplier: 1

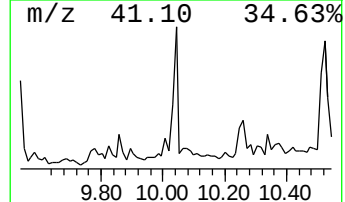
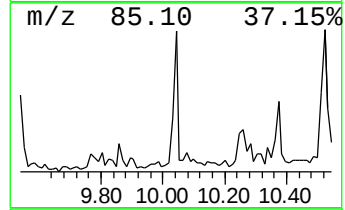
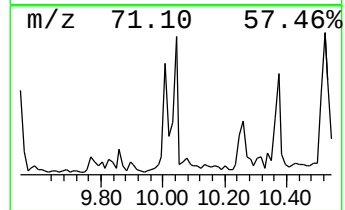
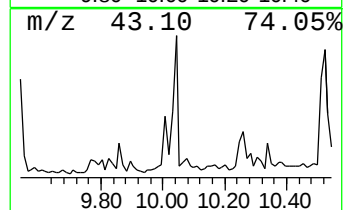
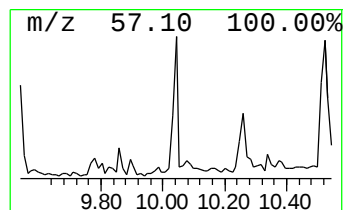
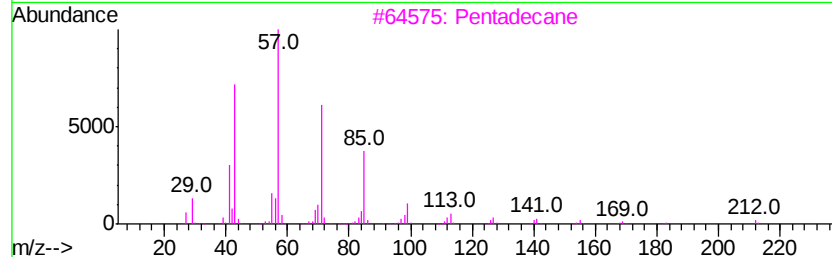
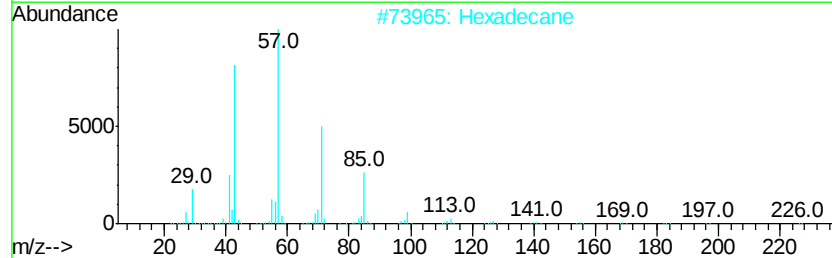
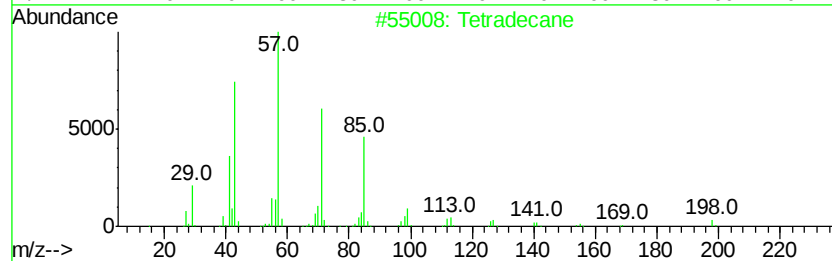
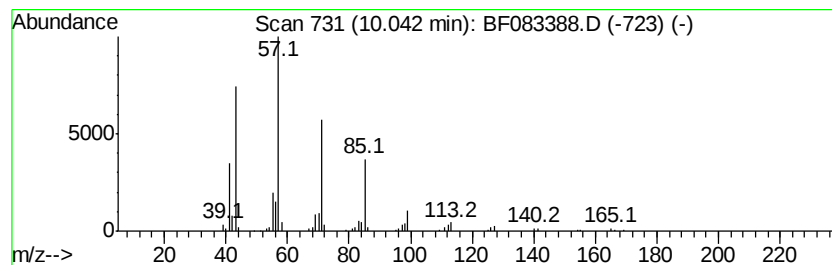
Instrument :
BNA_F
ClientSampleId :
S24-15.5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Tetradecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.04	4.70 ng	324860	Acenaphthene-d10		9.58
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	90
2	Hexadecane	226	C16H34	000544-76-3	90
3	Pentadecane	212	C15H32	000629-62-9	90
4	Tridecane	184	C13H28	000629-50-5	86
5	Eicosane	282	C20H42	000112-95-8	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120115\
Data File : BF083388.D
Acq On : 1 Dec 2015 22:58
Operator : UM/IZ
Sample : G4593-08
Misc :
ALS Vial : 13 Sample Multiplier: 1

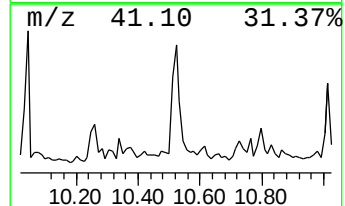
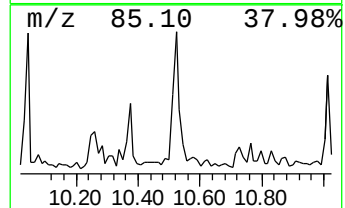
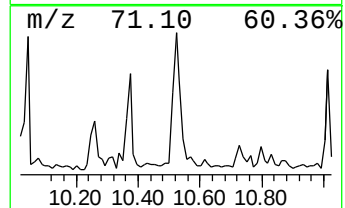
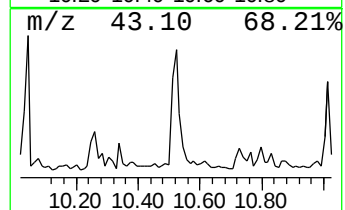
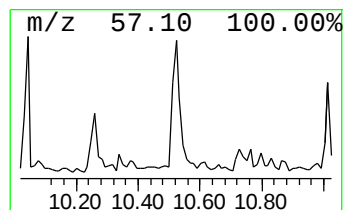
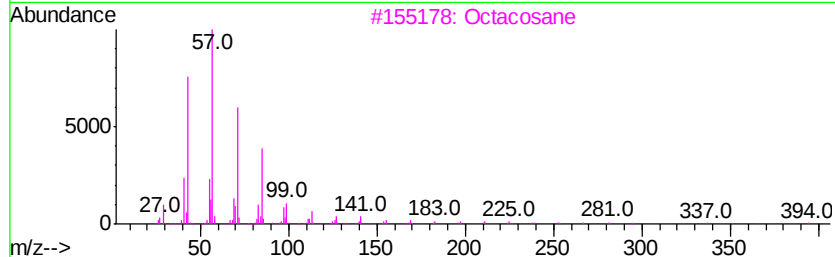
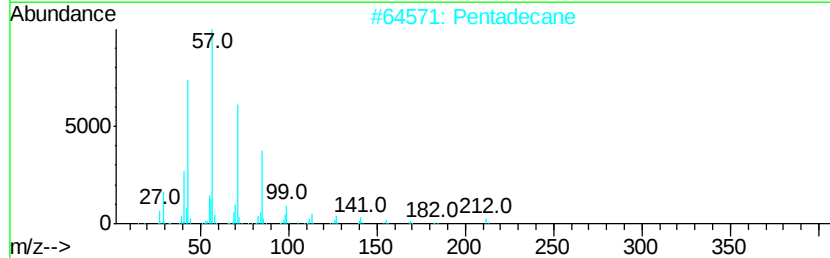
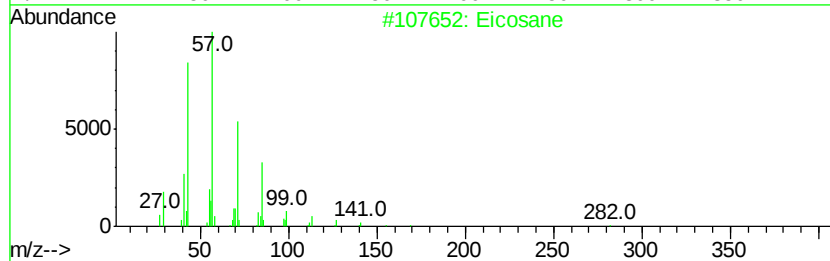
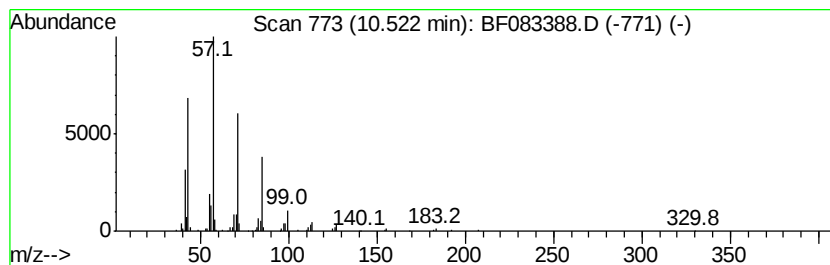
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Eicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.		
10.52	3.34 ng	339699	Phenanthrene-d10		11.13		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	91
2			Pentadecane	212	C15H32	000629-62-9	90
3			Octacosane	394	C28H58	000630-02-4	90
4			Heptadecane	240	C17H36	000629-78-7	90
5			Hexadecane	226	C16H34	000544-76-3	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120115\
Data File : BF083388.D
Acq On : 1 Dec 2015 22:58
Operator : UM/IZ
Sample : G4593-08
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S24-15.5

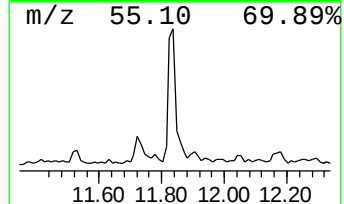
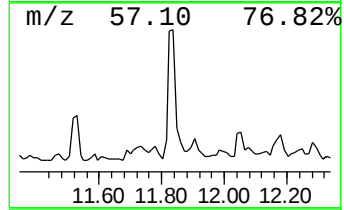
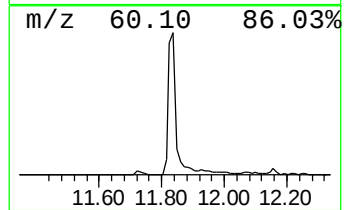
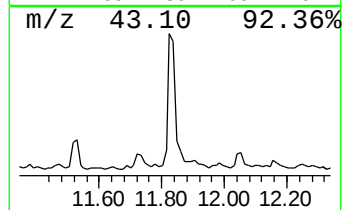
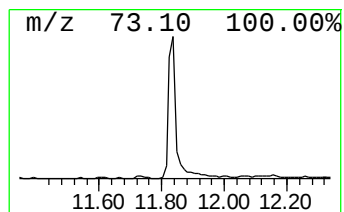
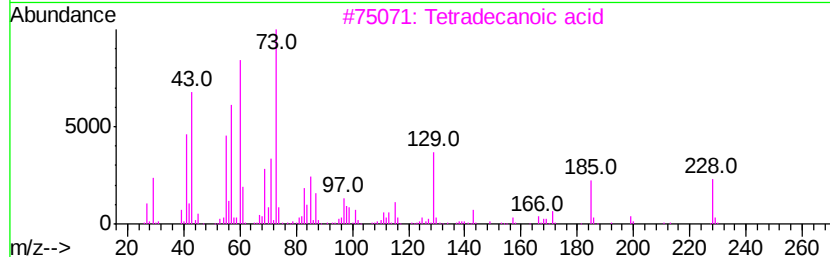
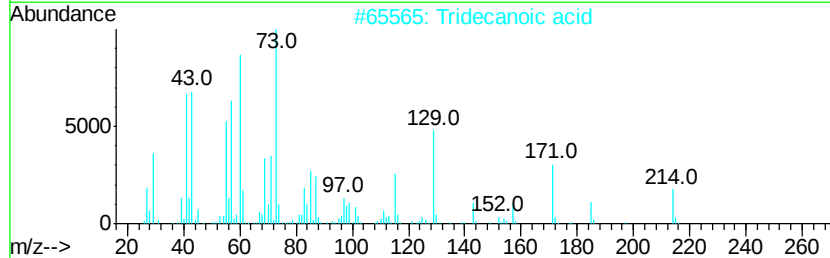
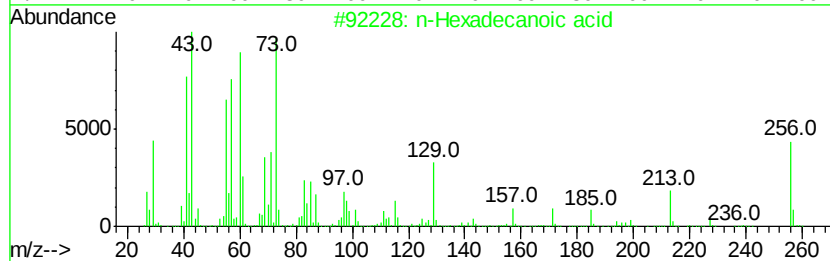
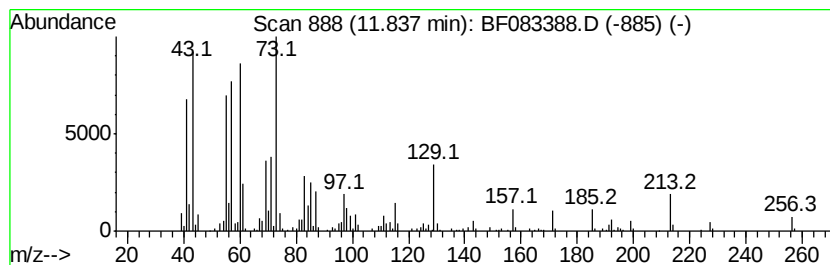
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.84	7.15 ng	728074	Phenanthrene-d10	11.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tridecanoic acid	214	C13H26O2	000638-53-9	95
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	90
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	86
5		Undecane	156	C11H24	001120-21-4	11



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120115\
Data File : BF083388.D
Acq On : 1 Dec 2015 22:58
Operator : UM/IZ
Sample : G4593-08
Misc :
ALS Vial : 13 Sample Multiplier: 1

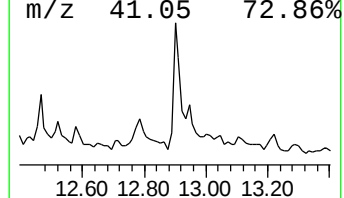
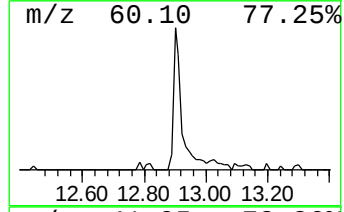
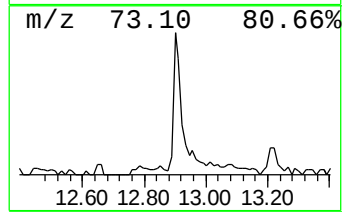
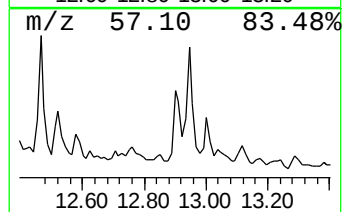
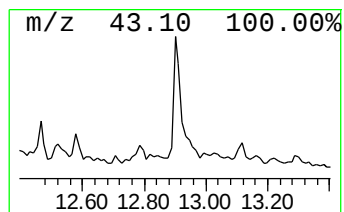
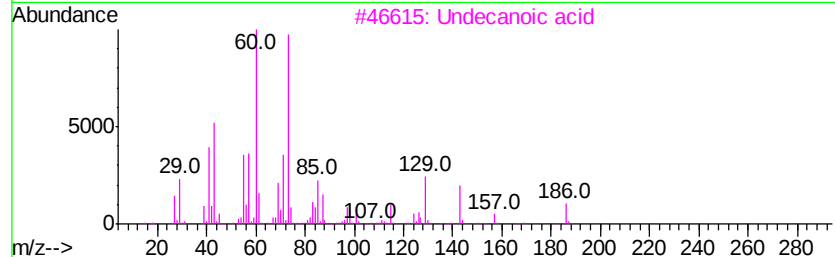
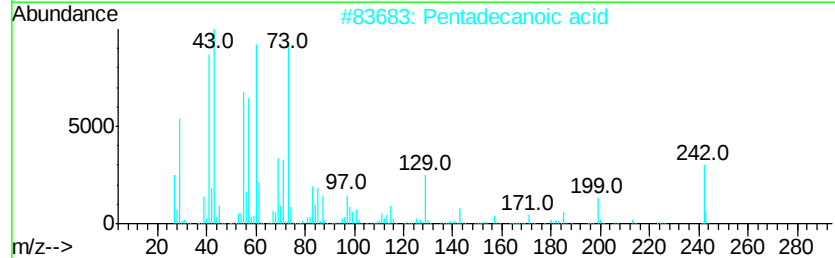
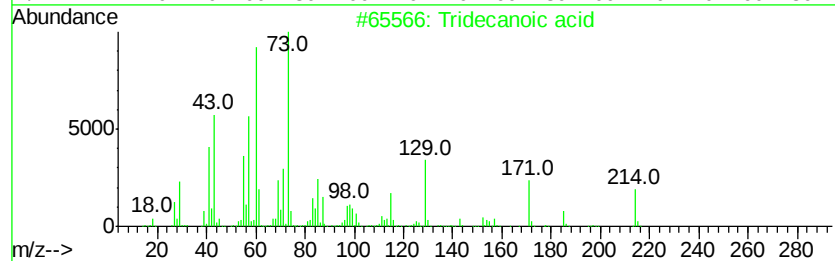
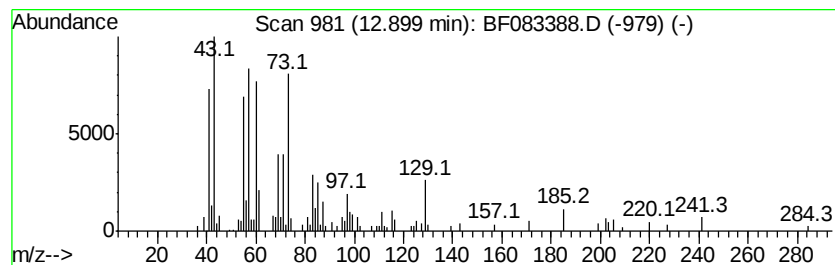
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tridecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.90	2.30 ng	234308	Phenanthrene-d10		11.13
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecanoic acid	214	C13H26O2	000638-53-9	72
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	64
3	Undecanoic acid	186	C11H22O2	000112-37-8	64
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120115\
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Acq On : 1 Dec 2015 22:58
Operator : UM/IZ
Sample : G4593-08
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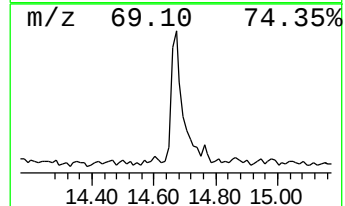
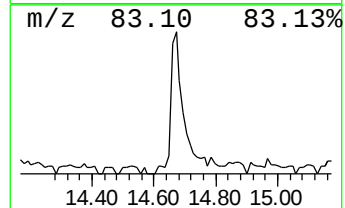
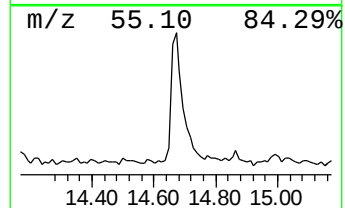
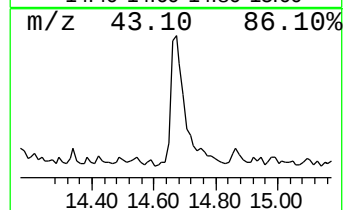
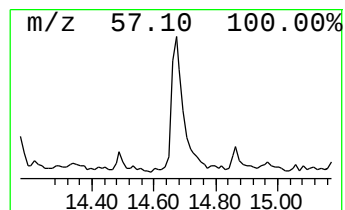
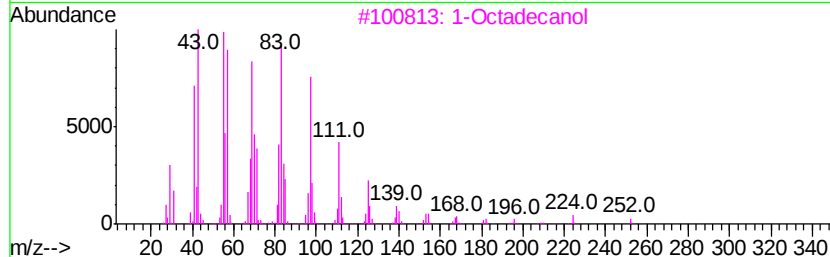
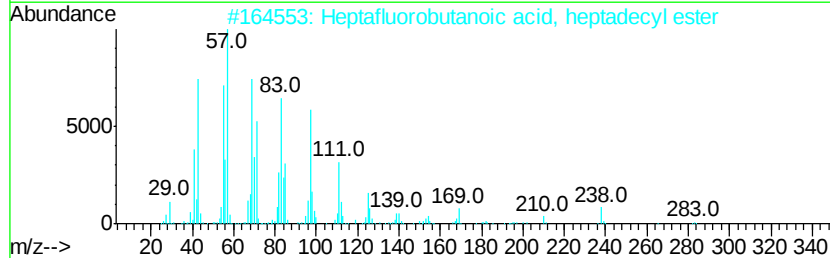
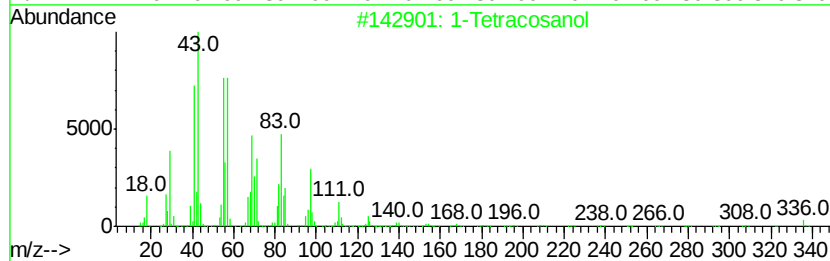
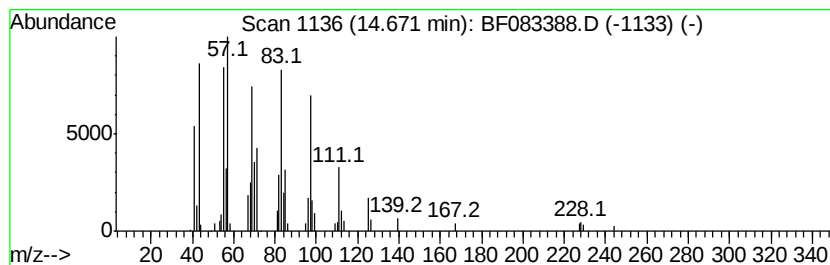
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1-Tetracosanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.67	2.74 ng	192162	Chrysene-d12	14.74
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Tetracosanol		354 C24H50O	000506-51-4 91
2	Heptafluorobutanoic acid, heptad...		452 C21H35F7O2	1000282-97-3 87
3	1-Octadecanol		270 C18H38O	000112-92-5 87
4	11-Tricosene		322 C23H46	052078-56-5 86
5	Bromoacetic acid, hexadecyl ester		362 C18H35BrO2	005454-48-8 83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120115\
Data File : BF083388.D
Acq On : 1 Dec 2015 22:58
Operator : UM/IZ
Sample : G4593-08
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S24-15.5

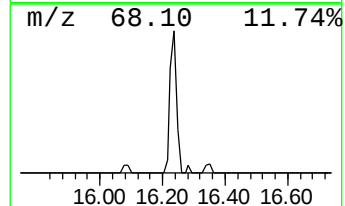
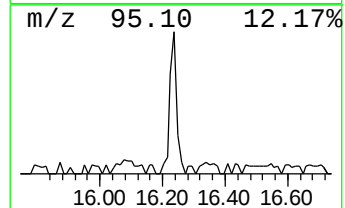
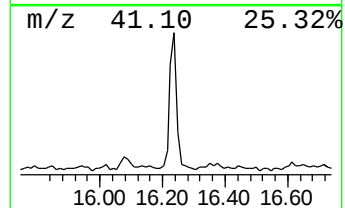
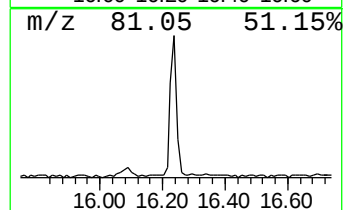
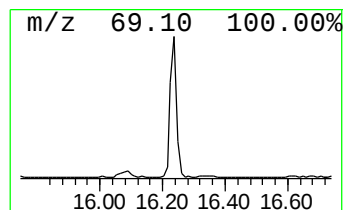
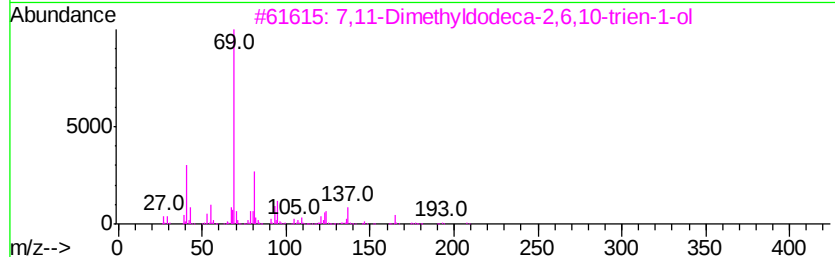
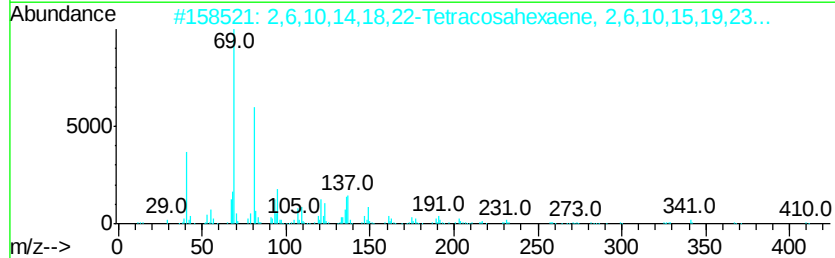
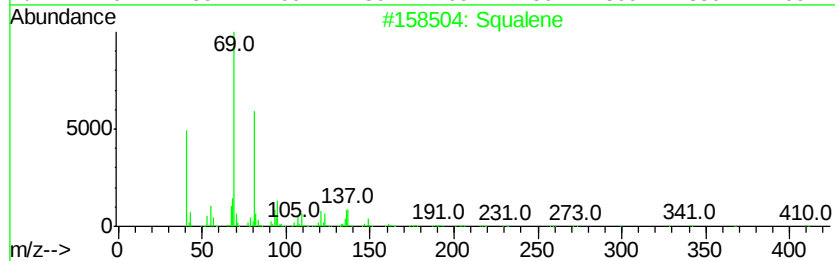
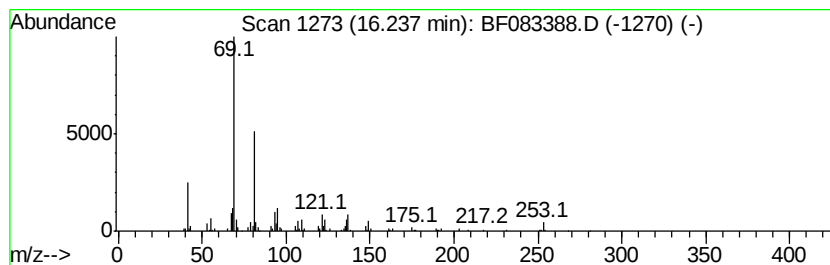
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Squalene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.24	5.39 ng	253763	Perylene-d12	16.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	007683-64-9	91
2		2,6,10,14,18,22-Tetracosahexaene...	410	C30H50	000111-02-4	90
3		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	83
4		4,9,13,17-Tetramethyl-4,8,12,16-...	316	C22H36O	056882-09-8	83
5		3,7,11-Tridecatrienitrile, 4,8...	231	C16H25N	006006-01-5	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120115\
Data File : BF083388.D
Acq On : 1 Dec 2015 22:58
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Sample : G4593-08
Misc :
ALS Vial : 13 Sample Multiplier: 1

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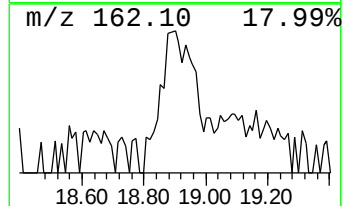
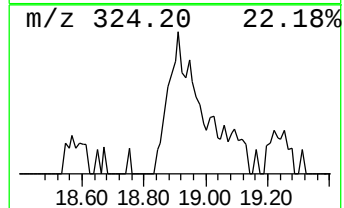
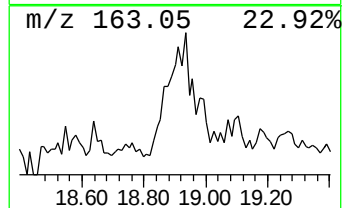
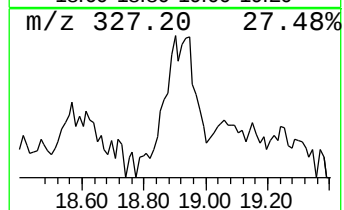
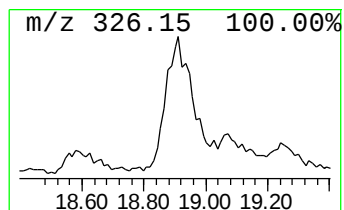
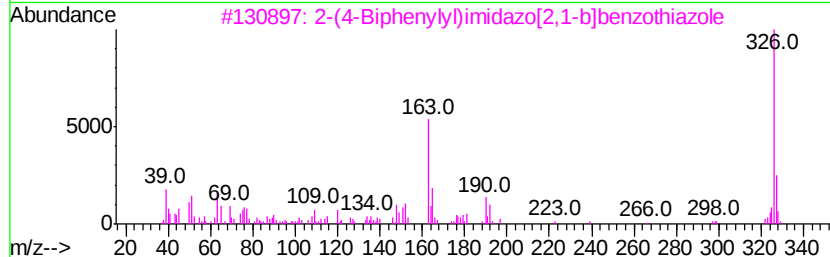
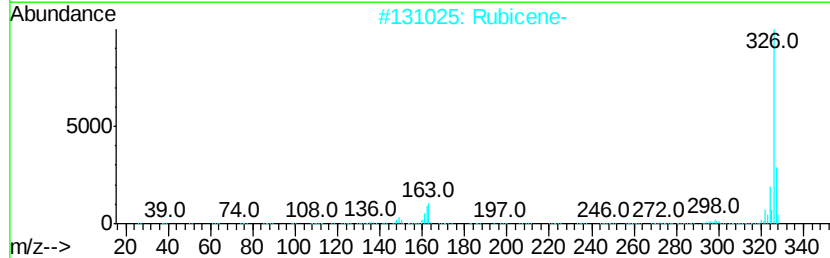
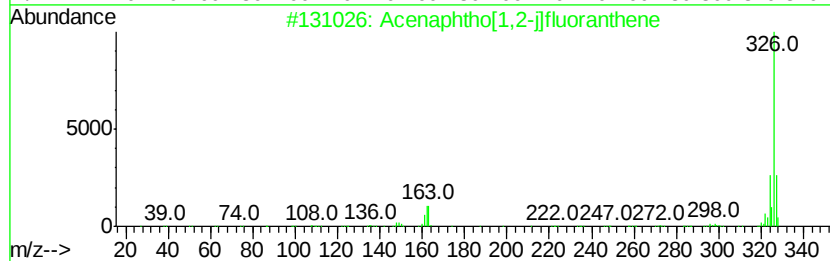
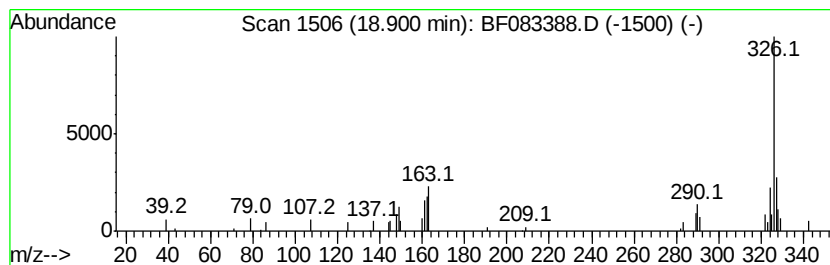
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Acenaphtho[1,2-j]fluoranthene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.90	2.07 ng	97300	Perylene-d12	16.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Acenaphtho[1,2-il]fluoranthene	326	C26H14	000193-21-5	70
2		Rubicene-	326	C26H14	000197-61-5	70
3		2-(4-Biphenylvl)imidazo[2,1-b]be...	326	C21H14N2S	038956-29-5	59
4		Silane, (estra-1,3,5(10),16-tetr...	326	C21H30OSi	069688-27-3	50
5		Benzothiophene-3-carbonitrile, 4...	326	C18H18N2O2S	329722-10-3	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120115\
Data File : BF083388.D
Acq On : 1 Dec 2015 22:58
Operator : UM/IZ
Sample : G4593-08
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S24-15.5

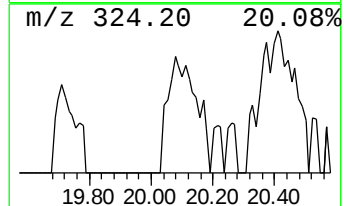
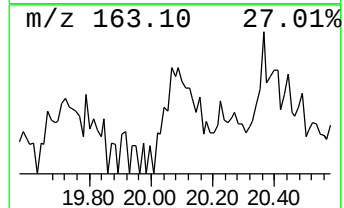
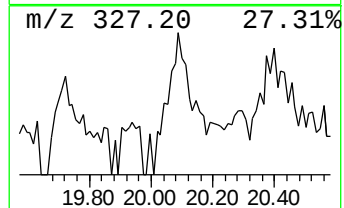
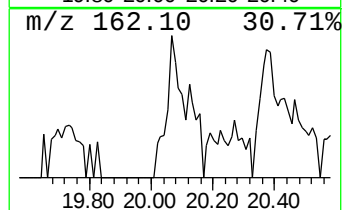
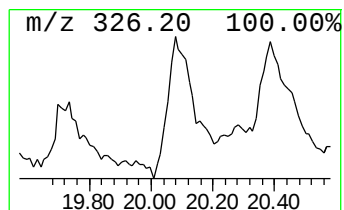
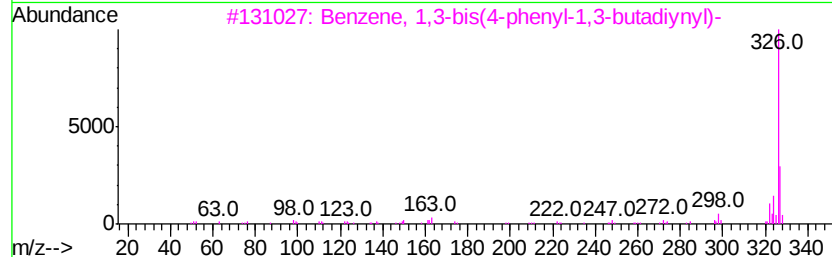
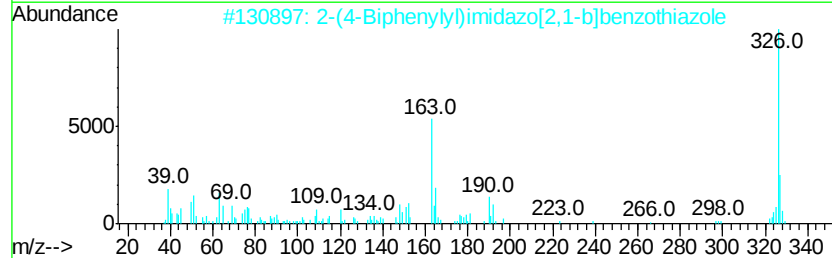
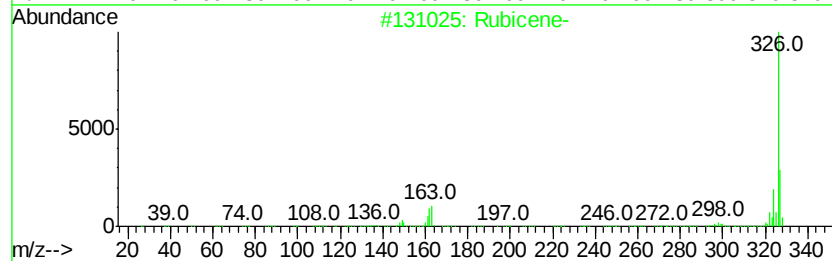
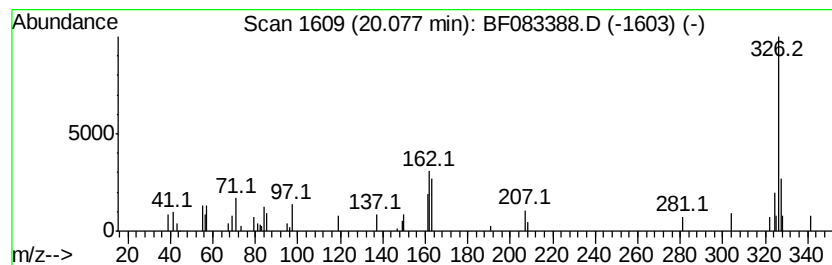
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Rubicene- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.08	2.19 ng	103195	Perylene-d12	16.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Rubicene-	326	C26H14	000197-61-5	72
2		2-(4-Biphenylvl)imidazo[2,1-b]be...	326	C21H14N2S	038956-29-5	50
3		Benzene, 1,3-bis(4-phenyl-1,3-bu...	326	C26H14	037902-13-9	43
4		Silane, (estra-1,3,5(10),16-tetr...	326	C21H30OSi	069688-27-3	43
5		9-Anthracenamine, N-[(4-nitrope...	326	C21H14N2O2	127387-74-0	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120115\
Data File : BF083388.D
Acq On : 1 Dec 2015 22:58
Operator : UM/IZ
Sample : G4593-08
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S24-15.5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.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
1-Pentene, 3,3-di...	3.90	2.5	ng	120789	1	6.54	969249	20.0
2-Pentanone, 4-hy...	4.66	31.9	ng	1544060	1	6.54	969249	20.0
unknown6.33	6.33	105.3	ng	5104950	1	6.54	969249	20.0
Hexanoic acid, 2-...	7.28	4.5	ng	285831	2	7.84	1272910	20.0
Tetradecane	10.04	4.7	ng	324860	3	9.58	1383130	20.0
Eicosane	10.52	3.3	ng	339699	4	11.13	2035850	20.0
n-Hexadecanoic acid	11.84	7.2	ng	728074	4	11.13	2035850	20.0
Tridecanoic acid	12.90	2.3	ng	234308	4	11.13	2035850	20.0
1-Tetracosanol	14.67	2.7	ng	192162	5	14.74	1400350	20.0
Squalene	16.24	5.4	ng	253763	6	16.81	941984	20.0
Acenaphtho[1,2-j]...	18.90	2.1	ng	97300	6	16.81	941984	20.0
Rubcene-	20.08	2.2	ng	103195	6	16.81	941984	20.0