

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011016\
 Data File : BF084358.D
 Acq On : 10 Jan 2016 12:06
 Operator : SJ/UM
 Sample : H1043-05
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-3S(0-10)

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-BF123115.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.047	256	259	267	rVB	1151997	1791833	25.77%	2.418%
2	5.470	294	296	299	rBV	5848339	5882839	84.62%	7.940%
3	6.510	383	387	389	rBV	4845704	6952256	100.00%	9.383%
4	6.625	395	397	400	rBV	5522108	6007867	86.42%	8.108%
5	6.853	415	417	419	rBV	2055719	1663167	23.92%	2.245%
6	7.013	428	431	433	rBV	5684634	5181939	74.54%	6.994%
7	7.425	464	467	469	rBV	4347994	4788492	68.88%	6.463%
8	7.527	475	476	483	rVB	68612	104405	1.50%	0.141%
9	8.145	527	530	532	rBV	1915227	2195112	31.57%	2.963%
10	9.219	621	624	626	rBV	7409701	6587374	94.75%	8.890%
11	9.619	656	659	660	rBV	977731	1253377	18.03%	1.692%
12	9.825	675	677	681	rBV	154773	218072	3.14%	0.294%
13	9.893	681	683	685	rVV	2667871	2405510	34.60%	3.246%
14	9.928	685	686	688	rVB	119493	91350	1.31%	0.123%
15	10.065	694	698	699	rBV	45615	72949	1.05%	0.098%
16	10.088	699	700	704	rVB3	44365	90202	1.30%	0.122%
17	10.328	720	721	723	rVB	404695	321176	4.62%	0.433%
18	10.442	729	731	734	rVB	82968	99656	1.43%	0.134%
19	10.545	737	740	744	rBV	113249	210768	3.03%	0.284%
20	10.682	749	752	755	rBV	4097850	4397818	63.26%	5.935%
21	10.796	759	762	769	rVB2	327887	624892	8.99%	0.843%
22	10.991	777	779	784	rVB	65644	124115	1.79%	0.168%
23	11.254	798	802	803	rBV	229255	257959	3.71%	0.348%
24	11.276	803	804	807	rVB	93363	92106	1.32%	0.124%
25	11.379	811	813	814	rBV	2701708	2434999	35.02%	3.286%
26	11.448	817	819	821	rVB	220841	233265	3.36%	0.315%
27	11.608	830	833	837	rBV2	393800	526219	7.57%	0.710%
28	11.676	837	839	841	rVB	122700	100671	1.45%	0.136%
29	11.882	854	857	858	rBV	76584	91978	1.32%	0.124%
30	11.905	858	859	861	rVB	151751	122263	1.76%	0.165%
31	11.985	864	866	871	rVV2	146732	233861	3.36%	0.316%
32	12.088	871	875	876	rVV	43887	77672	1.12%	0.105%
33	12.168	880	882	887	rVB2	54003	130088	1.87%	0.176%
34	12.431	903	905	911	rBV	904768	1201103	17.28%	1.621%

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35	12.511	911	912	916	rVB3	55874	74487	1.07%	0.101%
36	12.591	916	919	921	rBV	1050619	1067271	15.35%	1.440%
37	12.819	935	939	940	rBV	736713	1041555	14.98%	1.406%
38	12.854	940	942	946	rVB	492099	636652	9.16%	0.859%
39	12.968	949	952	954	rVB	6393855	6078216	87.43%	8.203%
40	13.139	964	967	969	rVB	106178	191699	2.76%	0.259%
41	13.208	969	973	975	rBV2	118931	225974	3.25%	0.305%
42	13.242	975	976	980	rVB3	72002	125964	1.81%	0.170%
43	13.414	989	991	993	rBV	186202	191153	2.75%	0.258%
44	13.540	997	1002	1006	rVB6	52459	126113	1.81%	0.170%
45	13.654	1006	1012	1014	rVB3	43164	106391	1.53%	0.144%
46	13.757	1018	1021	1023	rBV	57549	97267	1.40%	0.131%
47	13.871	1028	1031	1039	rVB2	478900	869782	12.51%	1.174%
48	14.008	1040	1043	1051	rVB2	1625798	2824812	40.63%	3.812%
49	14.122	1051	1053	1054	rBV	65237	73431	1.06%	0.099%
50	14.397	1073	1077	1079	rBV	82144	137774	1.98%	0.186%
51	14.488	1083	1085	1087	rBV3	53597	90043	1.30%	0.122%
52	14.774	1108	1110	1115	rBV6	29833	77302	1.11%	0.104%
53	14.865	1115	1118	1123	rVB2	102090	124318	1.79%	0.168%
54	15.025	1130	1132	1138	rBV	259977	580415	8.35%	0.783%
55	15.128	1138	1141	1143	rVV3	40704	94794	1.36%	0.128%
56	15.323	1155	1158	1161	rVB	139336	183518	2.64%	0.248%
57	15.380	1161	1163	1166	rVB	207226	246815	3.55%	0.333%
58	15.437	1166	1168	1171	rBV	963127	1419692	20.42%	1.916%
59	15.974	1211	1215	1224	rVB5	41718	176806	2.54%	0.239%
60	16.831	1287	1290	1295	rBV2	78026	167285	2.41%	0.226%
61	17.243	1323	1326	1334	rVB	74796	165985	2.39%	0.224%
62	17.608	1354	1358	1362	rVB	33004	80609	1.16%	0.109%
63	18.409	1423	1428	1432	rBV2	34218	105527	1.52%	0.142%
64	19.346	1506	1510	1519	rVB3	34816	146754	2.11%	0.198%

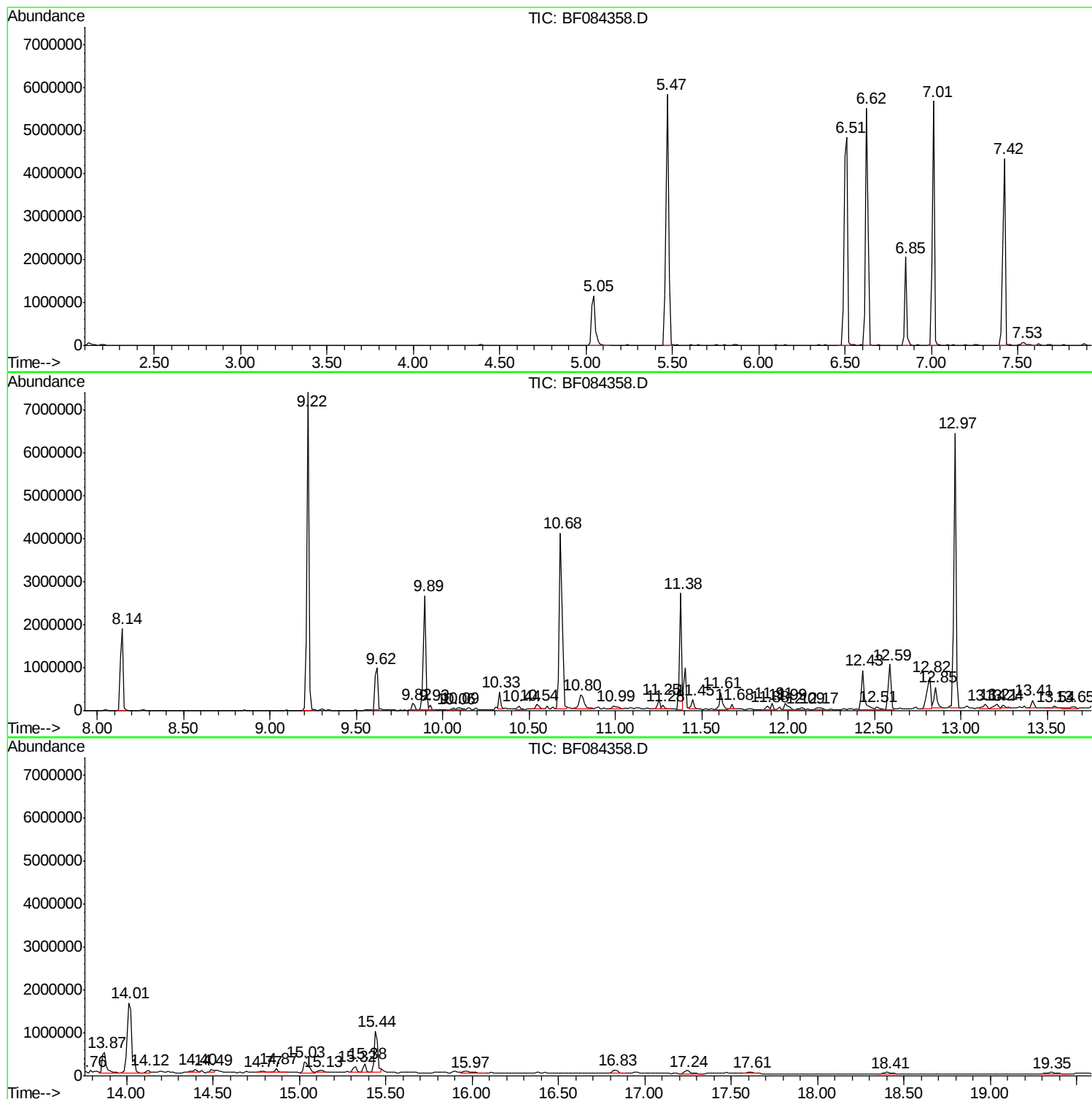
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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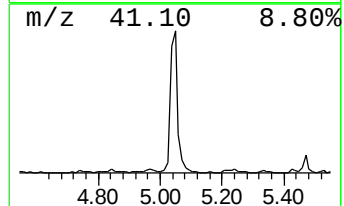
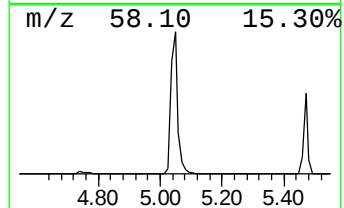
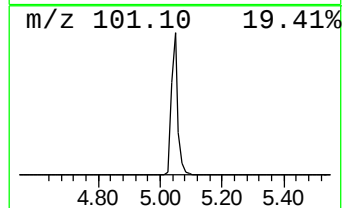
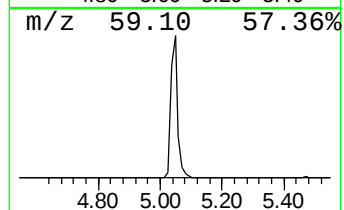
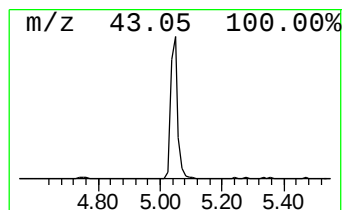
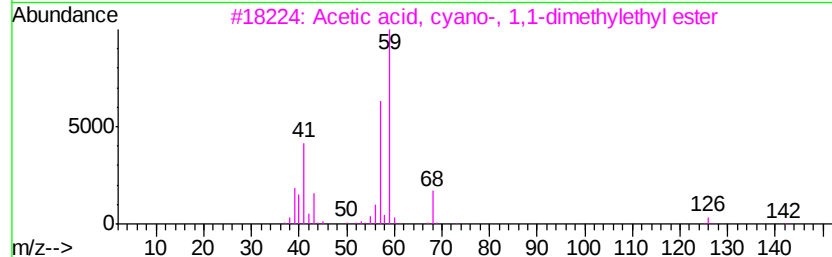
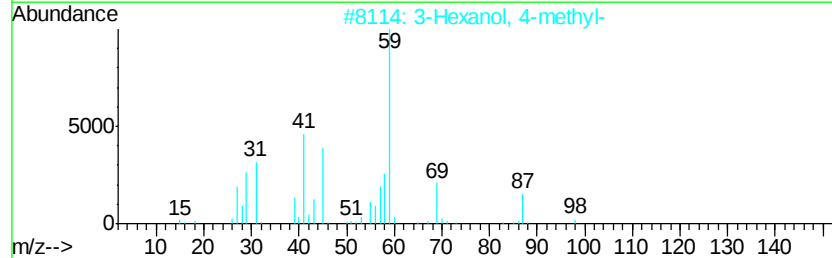
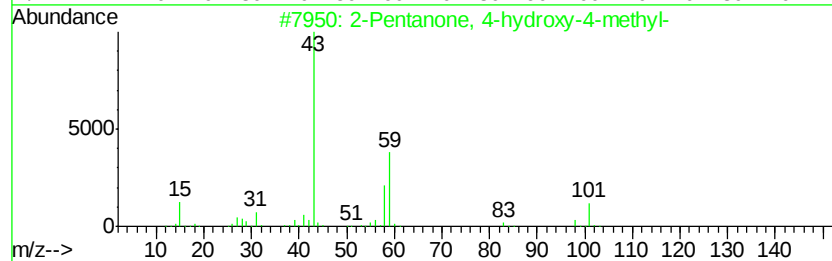
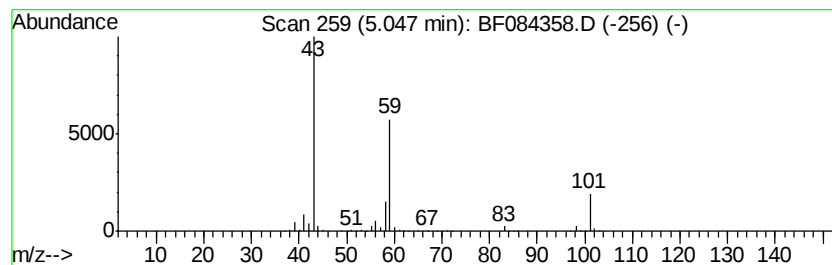
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown5.05 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	21.55 ng	1791830	1,4-Dichlorobenzene-d4	6.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	47
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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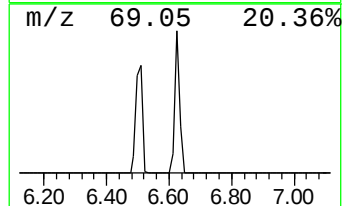
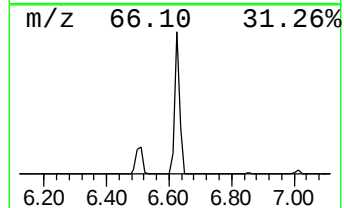
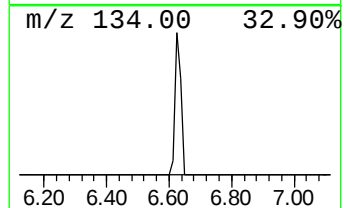
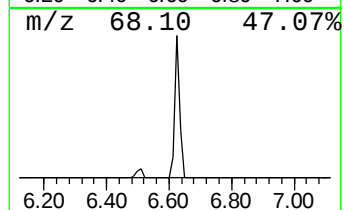
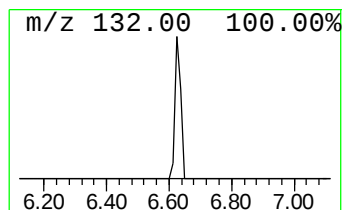
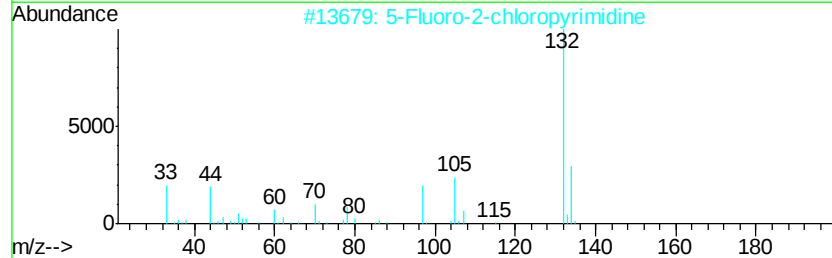
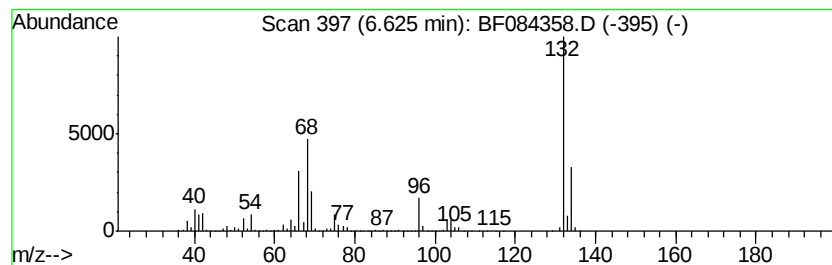
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	72.25 ng	6007870	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5		1,2,3-Trifluorobenzene	132	C6H3F3	001489-53-8	9



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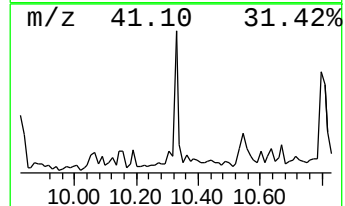
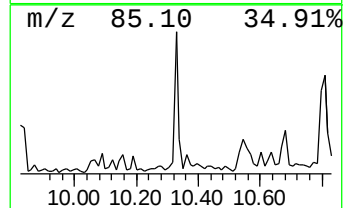
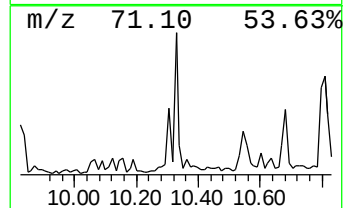
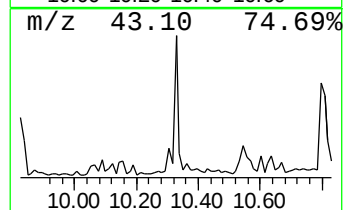
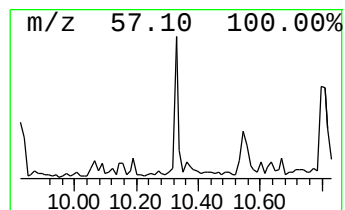
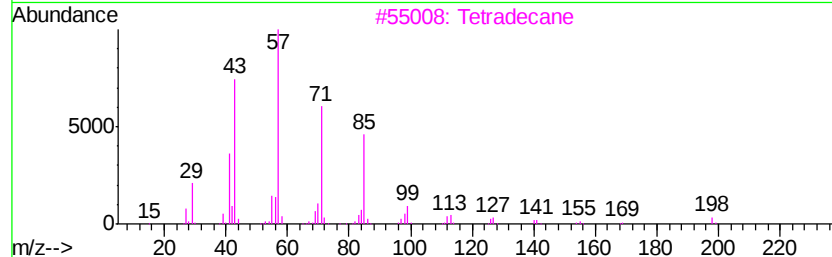
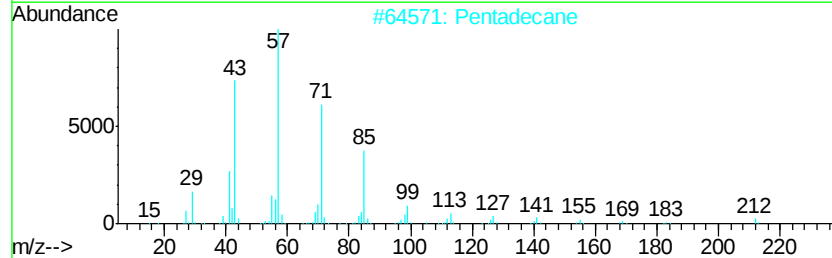
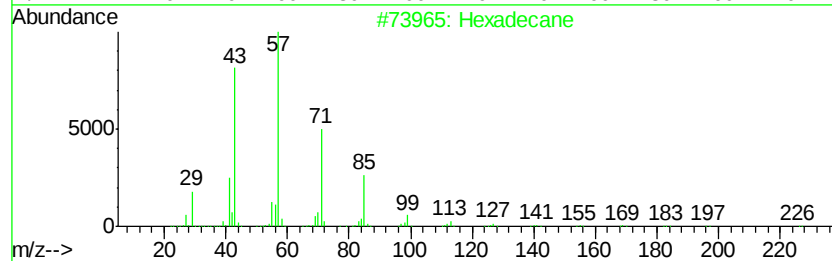
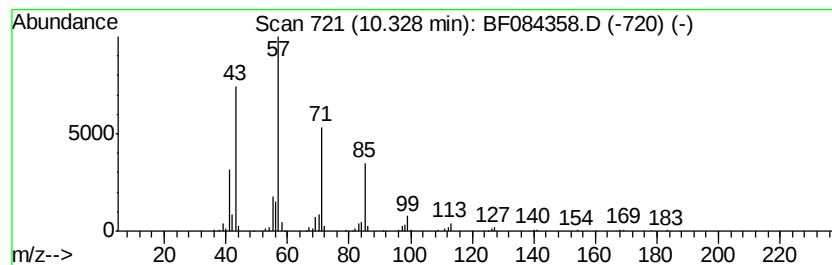
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 Hexadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.33	2.67 ng	321176	Acenaphthene-d10	9.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	91
2		Pentadecane	212	C15H32	000629-62-9	90
3		Tetradecane	198	C14H30	000629-59-4	90
4		Tridecane	184	C13H28	000629-50-5	83
5		Eicosane	282	C20H42	000112-95-8	72



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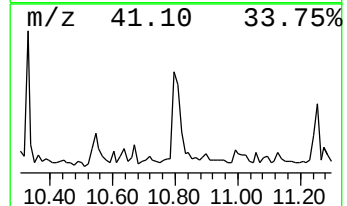
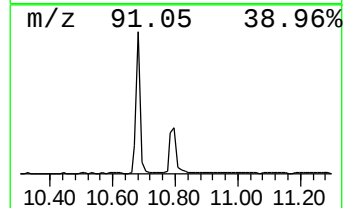
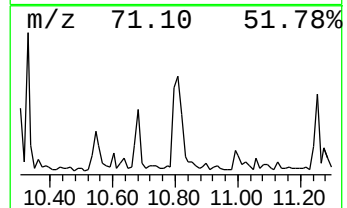
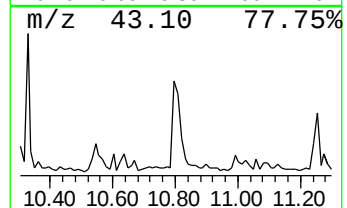
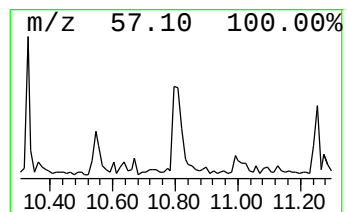
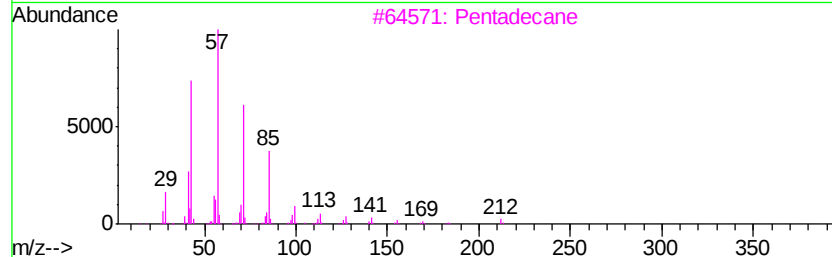
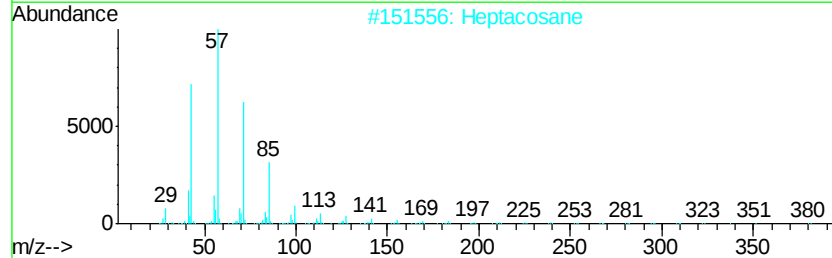
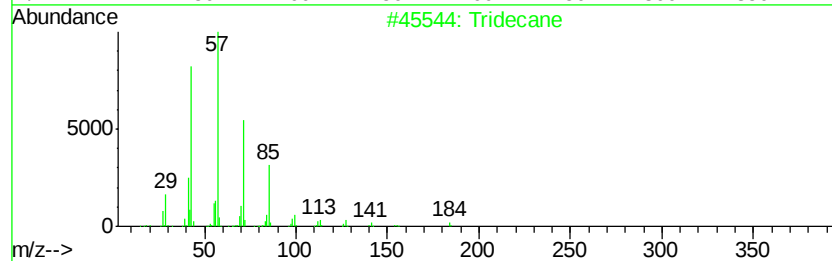
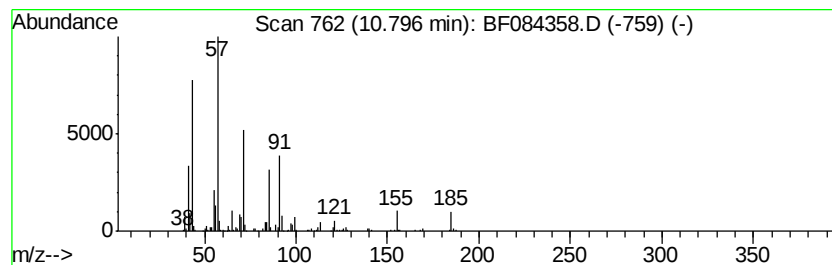
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Peak Number 5 Tridecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.80	5.13 ng	624892	Phenanthrene-d10	11.38	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	76
2	Heptacosane	380	C27H56	000593-49-7	72
3	Pentadecane	212	C15H32	000629-62-9	72
4	Hexadecane	226	C16H34	000544-76-3	68
5	Eicosane	282	C20H42	000112-95-8	64



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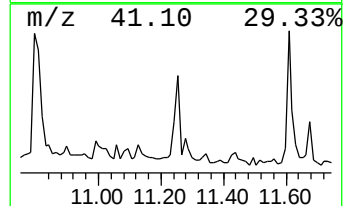
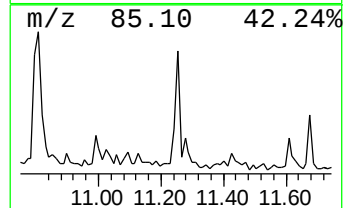
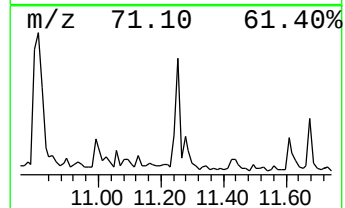
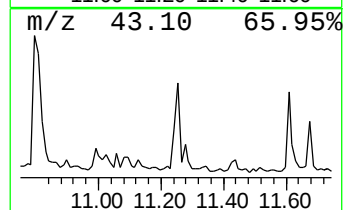
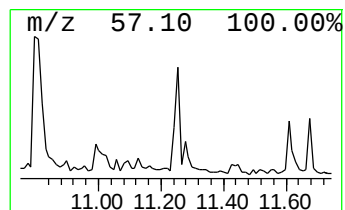
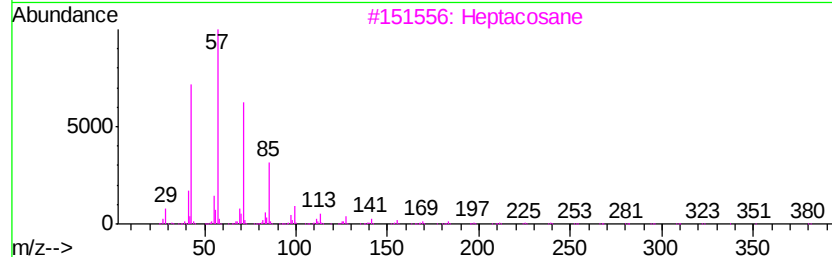
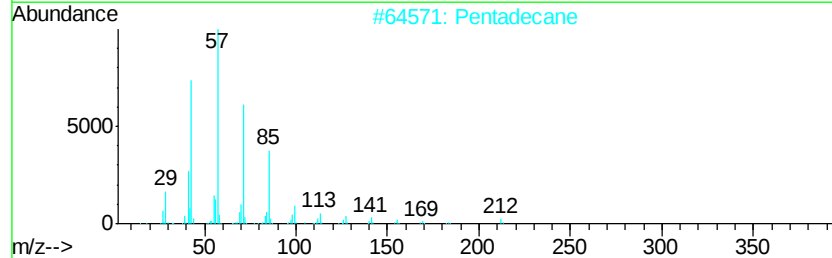
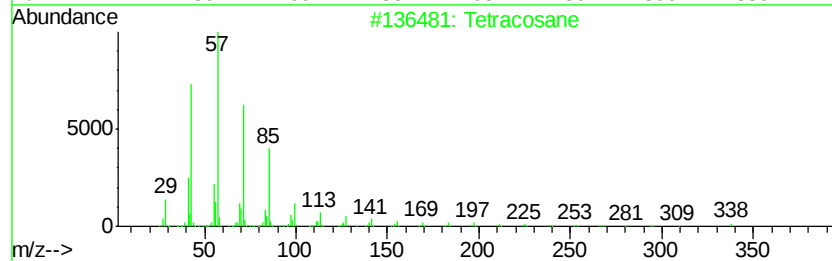
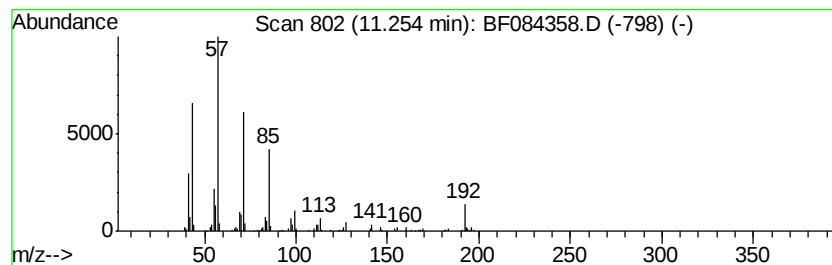
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ClientSampleId :
SB-3S(0-10)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Tetracosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.25	2.12 ng	257959	Phenanthrene-d10		11.38	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	87
2		Pentadecane	212	C15H32	000629-62-9	87
3		Heptacosane	380	C27H56	000593-49-7	87
4		Octacosane	394	C28H58	000630-02-4	87
5		Heneicosane	296	C21H44	000629-94-7	86



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011016\
Data File : BF084358.D
Acq On : 10 Jan 2016 12:06
Operator : SJ/UM
Sample : H1043-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-3S(0-10)

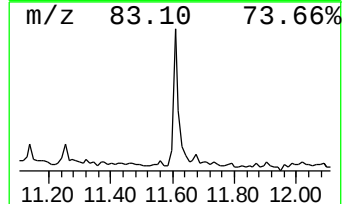
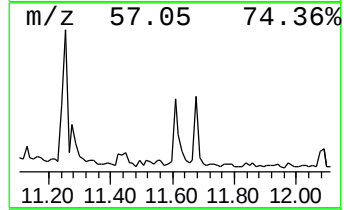
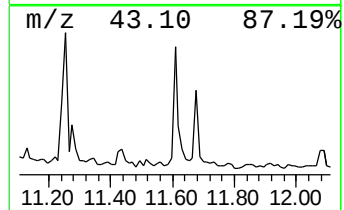
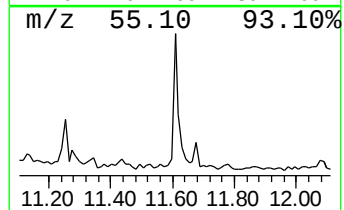
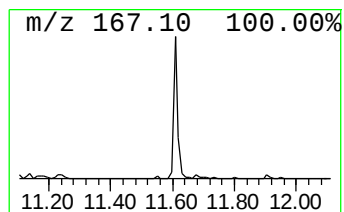
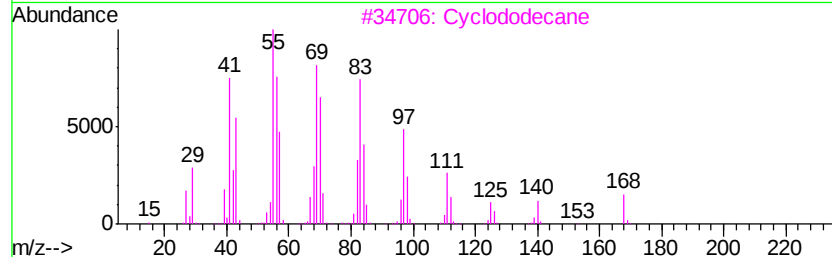
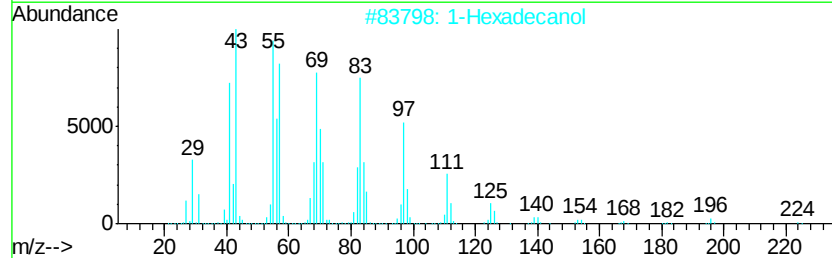
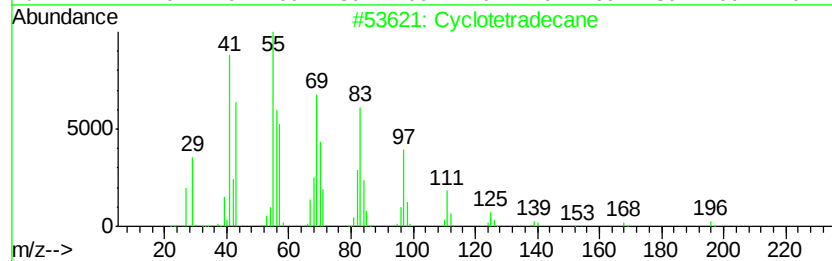
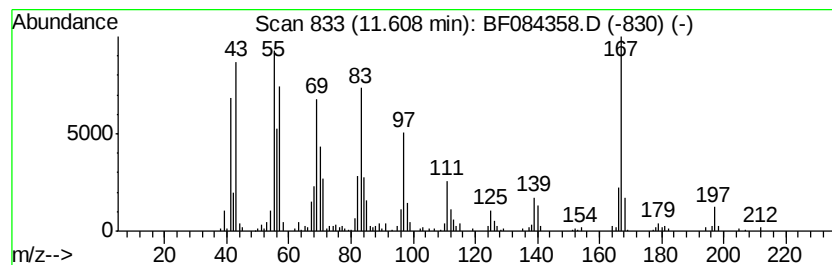
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Cyclotetradecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.61	4.32 ng	526219	Phenanthrene-d10	11.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetradecane	196	C14H28	000295-17-0	95
2	1-Hexadecanol	242	C16H34O	036653-82-4	91
3	Cyclododecane	168	C12H24	000294-62-2	90
4	1-Pentadecanol	228	C15H32O	000629-76-5	83
5	Hexadecen-1-ol, trans-9-	240	C16H32O	064437-47-4	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011016\
Data File : BF084358.D
Acq On : 10 Jan 2016 12:06
Operator : SJ/UM
Sample : H1043-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-3S(0-10)

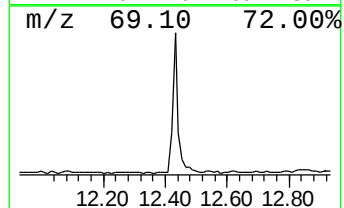
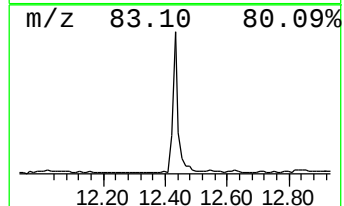
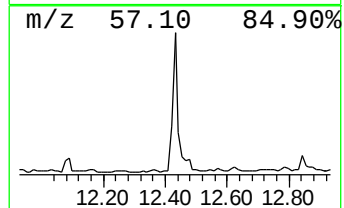
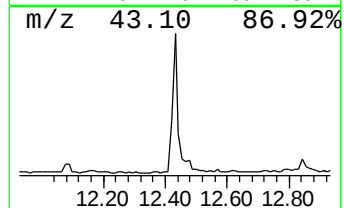
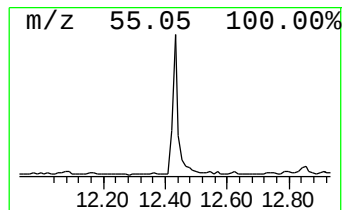
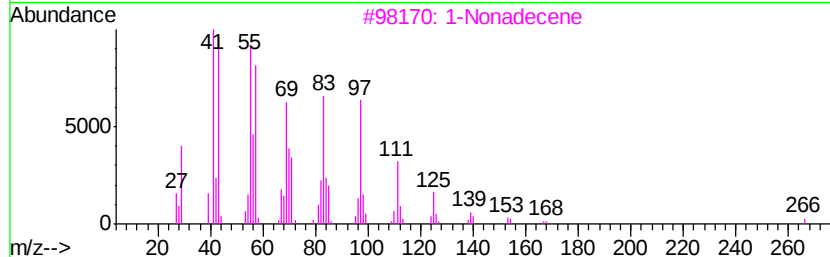
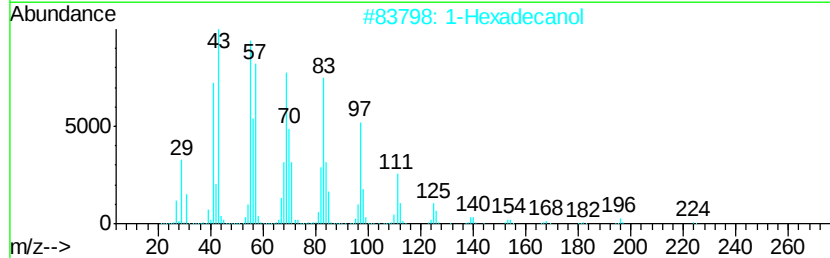
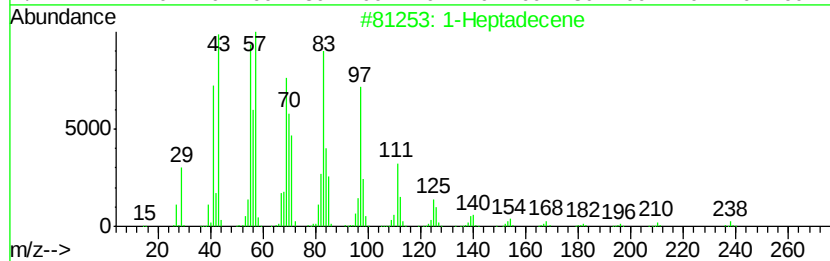
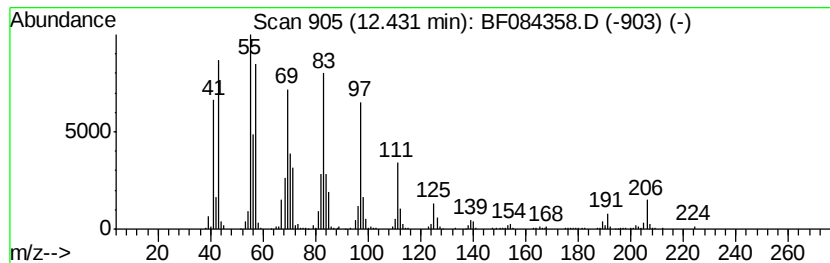
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1-Heptadecene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.43	9.87 ng	1201100	Phenanthrene-d10	11.38

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptadecene		238	C17H34	006765-39-5	91
2	1-Hexadecanol		242	C16H34O	036653-82-4	91
3	1-Nonadecene		266	C19H38	018435-45-5	91
4	5-Eicosene, (E)-		280	C20H40	074685-30-6	91
5	9-Eicosene, (E)-		280	C20H40	074685-29-3	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011016\
Data File : BF084358.D
Acq On : 10 Jan 2016 12:06
Operator : SJ/UM
Sample : H1043-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

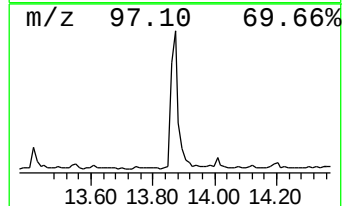
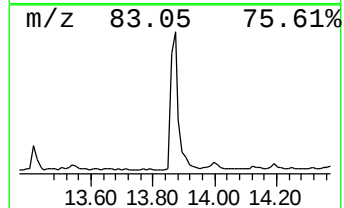
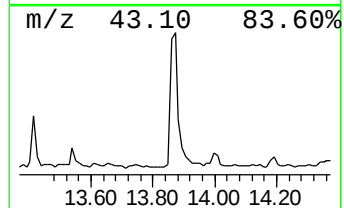
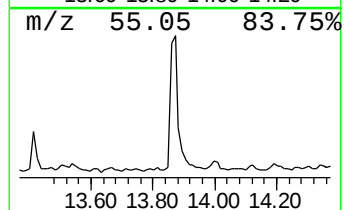
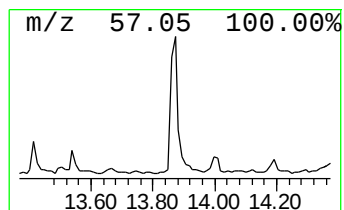
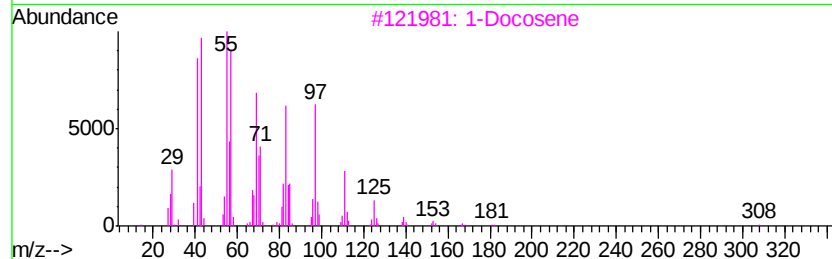
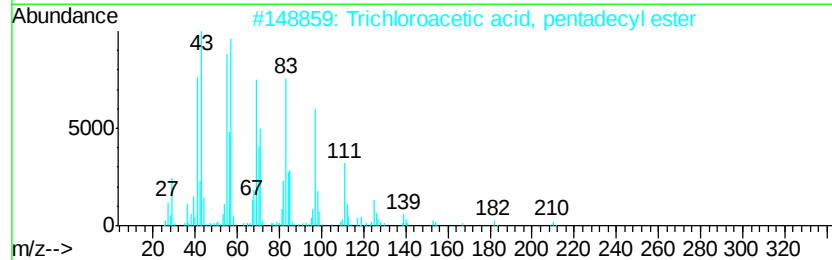
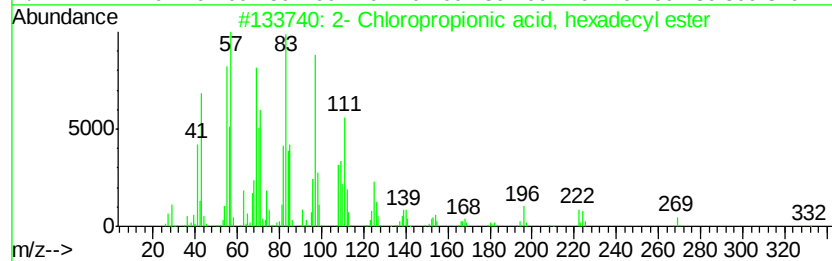
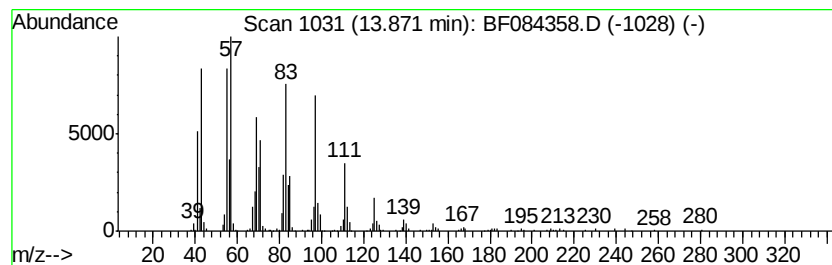
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ClientSampleId :
SB-3S(0-10)

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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 2- Chloropropionic acid, he... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.87	6.16 ng	869782	Chrysene-d12	14.02		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-	Chloropropionic acid, hexadec...	332	C19H37ClO2	086711-81-1	91
2		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91
3		1-Docosene	308	C22H44	001599-67-3	91
4		Trifluoroacetic acid, n-octadecv...	366	C20H37F3O2	079392-43-1	91
5		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011016\
Data File : BF084358.D
Acq On : 10 Jan 2016 12:06
Operator : SJ/UM
Sample : H1043-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-3S(0-10)

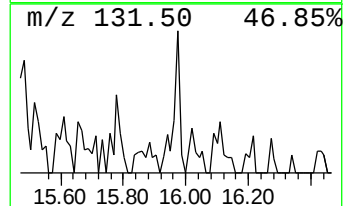
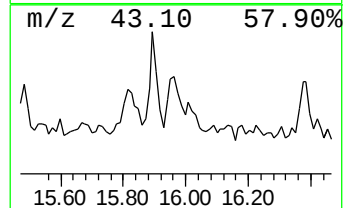
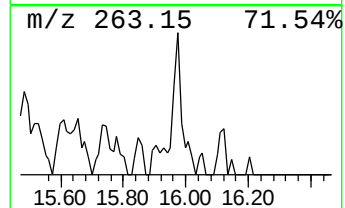
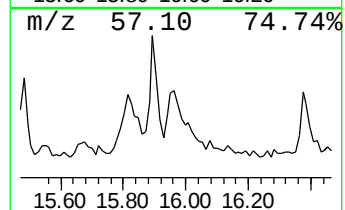
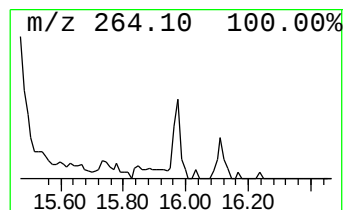
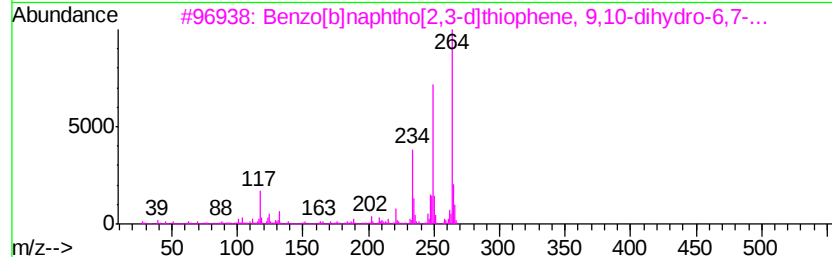
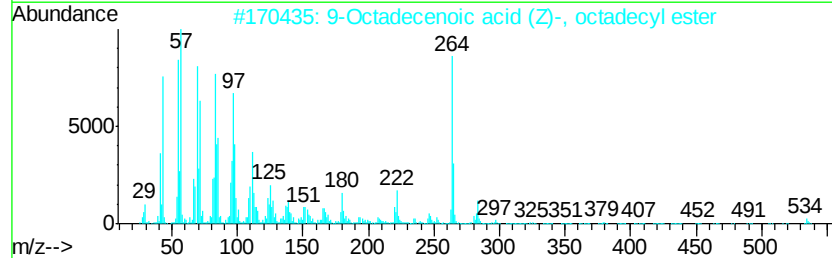
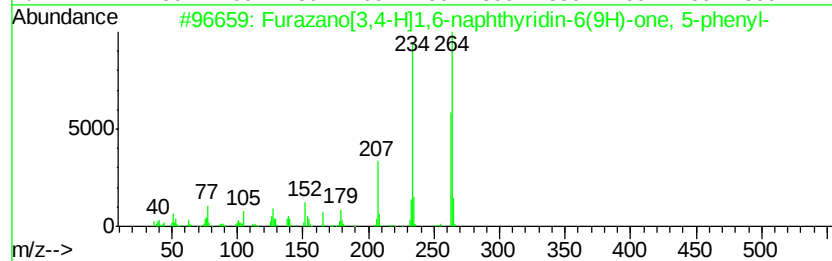
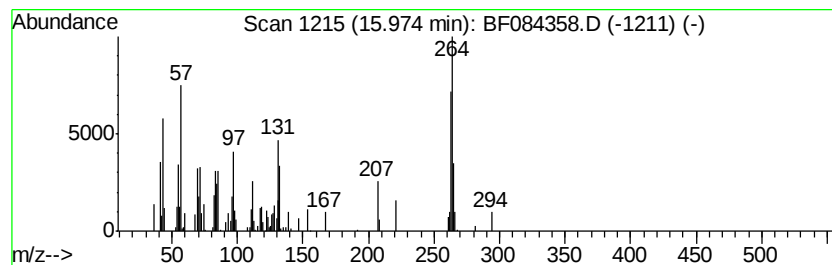
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown15.97 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.97	2.49 ng	176806	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Furazano[3,4-H]1,6-naphthvridin-...	264	C14H8N4O2	296245-19-7	32
2		9-Octadecenoic acid (Z)-, octade...	535	C36H70O2	017673-49-3	32
3		Benzo[b]naphtho[2,3-d]thiophene,...	264	C18H16S	024964-02-1	30
4		3-Hydroxycinnamic acid, ethyl es...	264	C14H20O3Si	1000071-70-0	22
5		9H-Purin-6-amine, N,9-bis(trimet...	279	C11H21N5Si2	017995-04-9	22



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011016\
Data File : BF084358.D
Acq On : 10 Jan 2016 12:06
Operator : SJ/UM
Sample : H1043-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-3S(0-10)

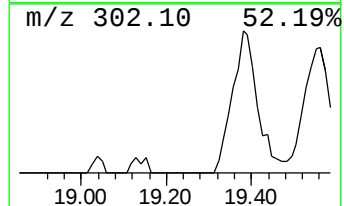
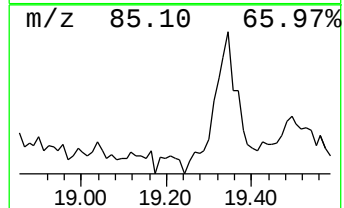
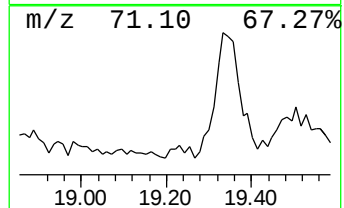
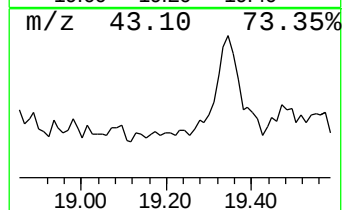
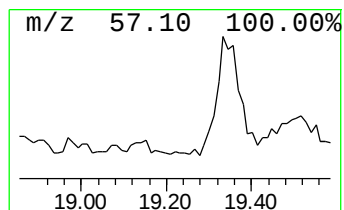
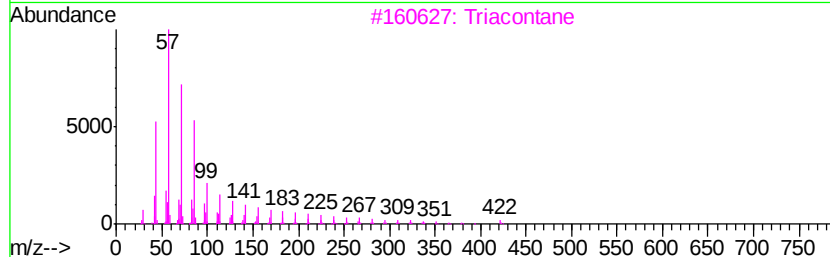
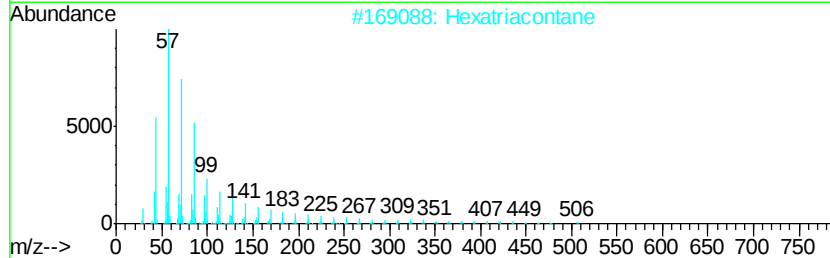
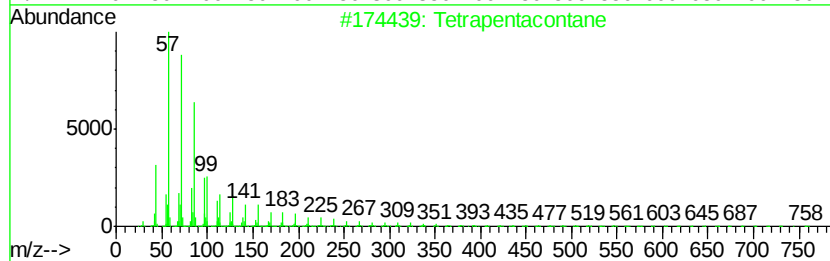
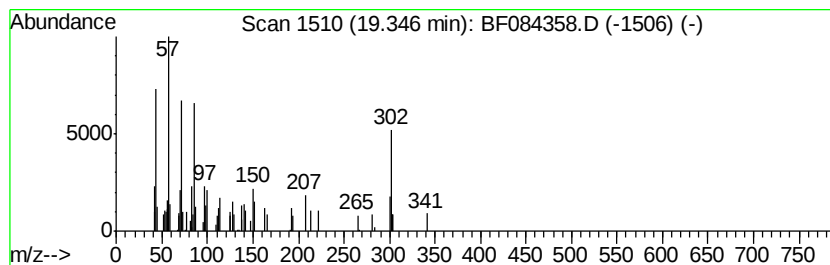
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown19.35 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.35	2.07 ng	146754	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetrapentacontane	759	C54H110	005856-66-6	43
2		Hexatriacontane	507	C36H74	000630-06-8	38
3		Triacontane	422	C30H62	000638-68-6	38
4		2-Bromo dodecane	248	C12H25Br	013187-99-0	38
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011016\
Data File : BF084358.D
Acq On : 10 Jan 2016 12:06
Operator : SJ/UM
Sample : H1043-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-3S(0-10)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.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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unknown6.62	6.62	72.3	ng	6007870	1	6.85	1663170	20.0
Hexadecane	10.33	2.7	ng	321176	3	9.89	2405510	20.0
Tridecane	10.80	5.1	ng	624892	4	11.38	2435000	20.0
Tetracosane	11.25	2.1	ng	257959	4	11.38	2435000	20.0
Cyclotetradecane	11.61	4.3	ng	526219	4	11.38	2435000	20.0
1-Heptadecene	12.43	9.9	ng	1201100	4	11.38	2435000	20.0
2-Chloropropioni...	13.87	6.2	ng	869782	5	14.02	2824810	20.0
unknown15.97	15.97	2.5	ng	176806	6	15.44	1419690	20.0
unknown19.35	19.35	2.1	ng	146754	6	15.44	1419690	20.0