

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
 Data File : BF085560.D
 Acq On : 19 Mar 2016 7:59
 Operator : UM/SJ
 Sample : H1804-03
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TRACK-D-GRID-7B

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-BF031816.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	3.819	131	134	138	rBV	269939	410037	9.93%	1.048%
2	5.064	241	243	248	rBV	47857	75680	1.83%	0.193%
3	5.316	263	265	267	rBV	69670	77642	1.88%	0.198%
4	5.419	272	274	275	rBV	135749	150920	3.65%	0.386%
5	5.453	275	277	280	rVB	2516342	3340077	80.88%	8.539%
6	5.682	294	297	300	rVB	110225	144930	3.51%	0.371%
7	6.493	365	368	370	rBV	2566830	3702517	89.66%	9.466%
8	6.619	376	379	382	rBV	3004637	3311078	80.18%	8.465%
9	6.676	382	384	387	rVB2	67195	82753	2.00%	0.212%
10	6.847	397	399	402	rBV	667470	754226	18.26%	1.928%
11	7.007	410	413	416	rBV	2839737	2685876	65.04%	6.867%
12	7.110	421	422	426	rVB	162223	144332	3.50%	0.369%
13	7.419	446	449	451	rBV	2241247	2343041	56.74%	5.990%
14	7.499	454	456	464	rVB	36317	65810	1.59%	0.168%
15	7.887	486	490	492	rBV2	41358	76239	1.85%	0.195%
16	8.139	509	512	516	rBV2	1099708	1332274	32.26%	3.406%
17	8.848	572	574	578	rBB	60810	54538	1.32%	0.139%
18	9.213	603	606	609	rBV	3656695	3914251	94.79%	10.007%
19	9.602	638	640	642	rVB	406583	527832	12.78%	1.349%
20	9.819	657	659	663	rBV	31637	45676	1.11%	0.117%
21	9.888	663	665	668	rVB	1159874	1082383	26.21%	2.767%
22	10.185	689	691	697	rBV	27825	49934	1.21%	0.128%
23	10.322	699	703	705	rBV2	131597	180387	4.37%	0.461%
24	10.539	719	722	726	rBV	43893	78556	1.90%	0.201%
25	10.676	731	734	737	rBV2	1985034	2589781	62.71%	6.621%
26	10.791	743	744	749	rBV	106151	189390	4.59%	0.484%
27	11.248	782	784	785	rBV	64580	73708	1.78%	0.188%
28	11.374	793	795	796	rBV	1349688	1182159	28.63%	3.022%
29	11.396	796	797	798	rVV	115200	86862	2.10%	0.222%
30	11.442	798	801	804	rVB3	37834	63796	1.54%	0.163%
31	11.899	840	841	844	rBV	168927	230864	5.59%	0.590%
32	11.979	847	848	852	rVB2	30066	48385	1.17%	0.124%
33	12.585	898	901	902	rBV	84755	82632	2.00%	0.211%
34	12.688	907	910	916	rBV4	126754	187267	4.53%	0.479%

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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

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35	12.779	916	918	919	rVV	301149	250043	6.06%	0.639%
36	12.814	919	921	922	rVB	75398	86073	2.08%	0.220%
37	12.962	931	934	936	rBV	4028590	4129467	100.00%	10.557%
38	13.442	971	976	978	rBV	1307394	2266073	54.88%	5.793%
39	13.511	981	982	991	rVB	233857	315957	7.65%	0.808%
40	13.865	1011	1013	1020	rBV	86105	196236	4.75%	0.502%
41	14.002	1023	1025	1028	rBV	1010383	1106296	26.79%	2.828%
42	14.117	1033	1035	1039	rVV	187510	319596	7.74%	0.817%
43	14.174	1039	1040	1047	rVB2	28514	51850	1.26%	0.133%
44	15.431	1147	1150	1157	rBV	705894	859246	20.81%	2.197%
45	15.877	1186	1189	1194	rVB	32560	49705	1.20%	0.127%
46	16.357	1227	1231	1237	rVB	35878	67343	1.63%	0.172%
47	16.917	1276	1280	1286	rBV	22861	51464	1.25%	0.132%

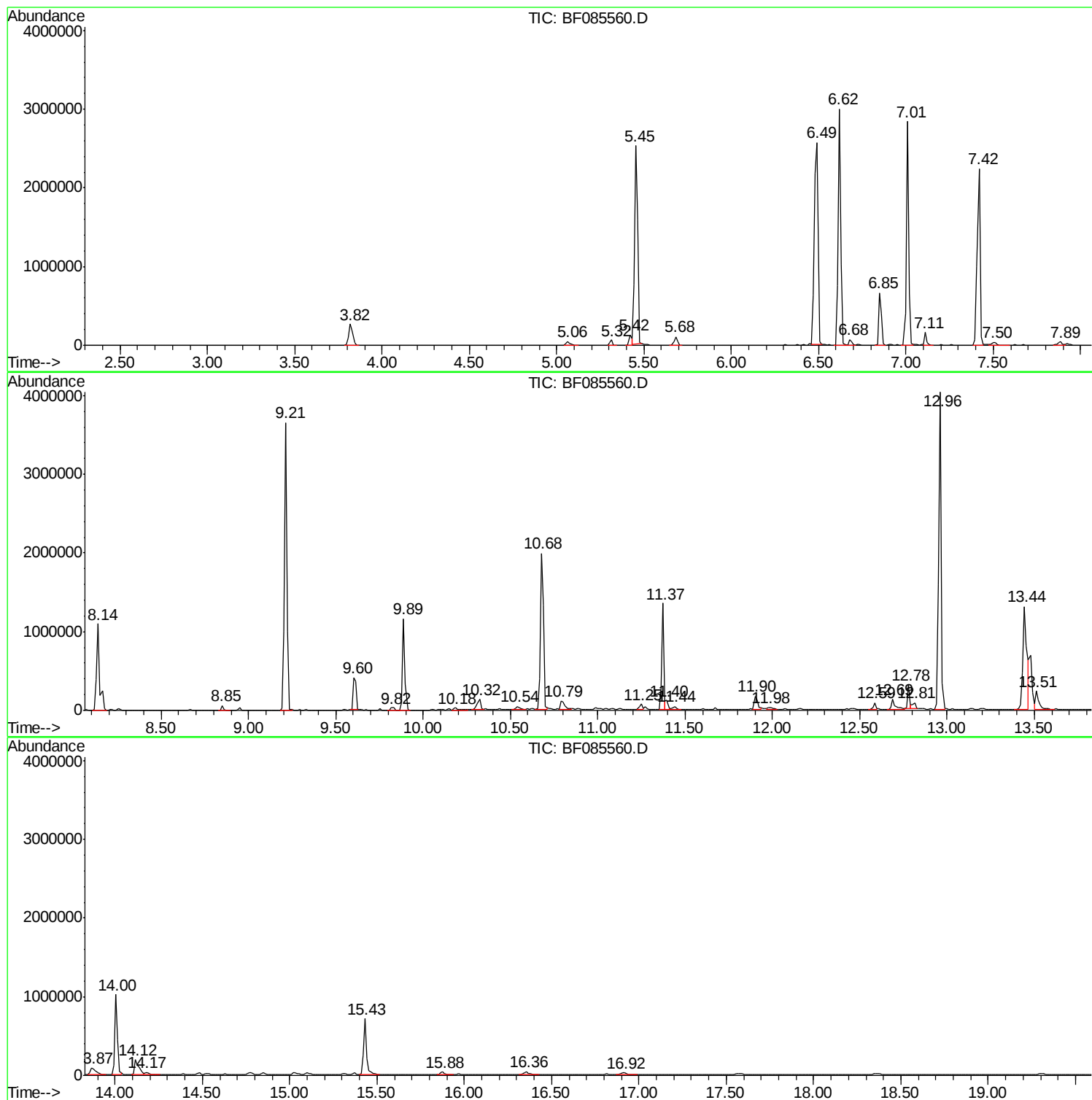
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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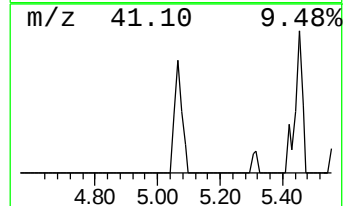
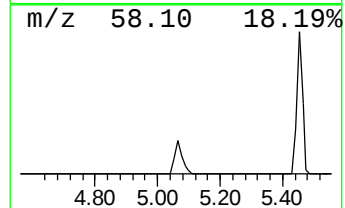
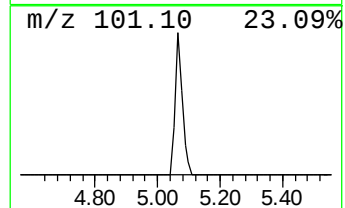
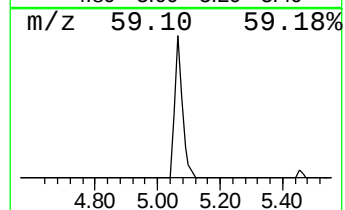
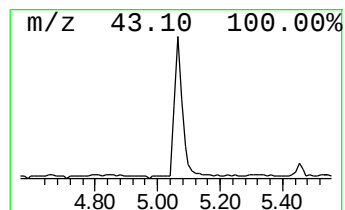
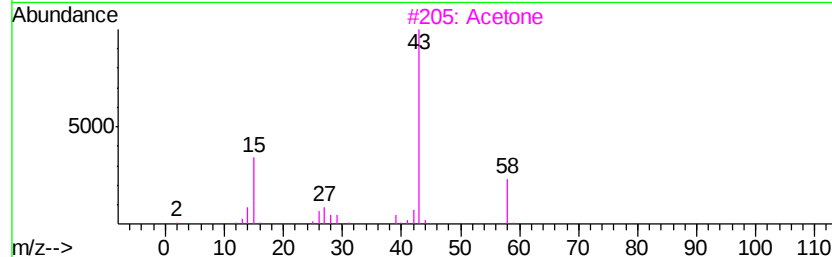
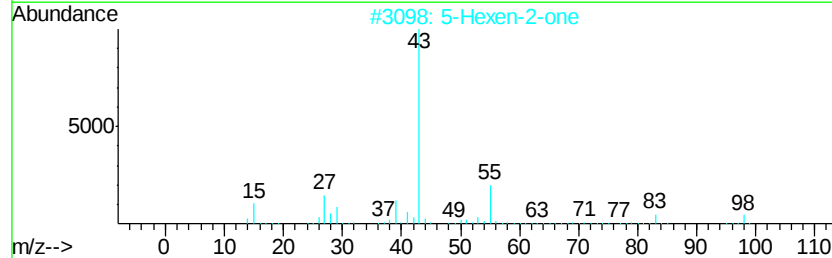
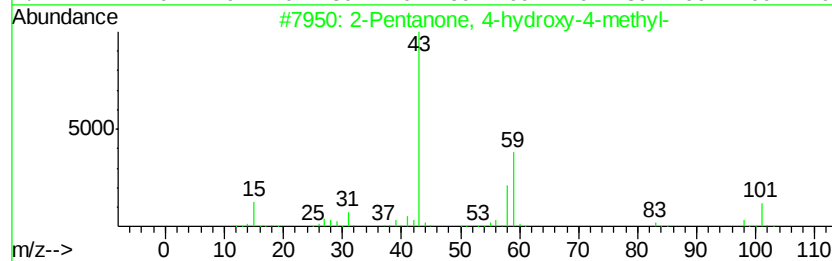
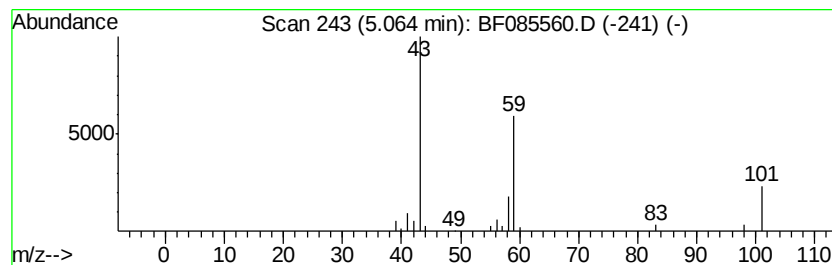
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.06	2.01 ng	75680	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	56
2			5-Hexen-2-one	98	C6H10O	000109-49-9	10
3			Acetone	58	C3H6O	000067-64-1	9
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9
5			1-Butanol, 3-methoxy-	104	C5H12O2	002517-43-3	9



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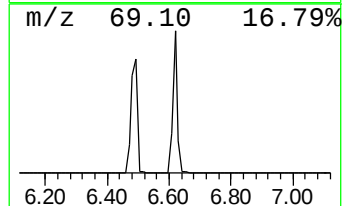
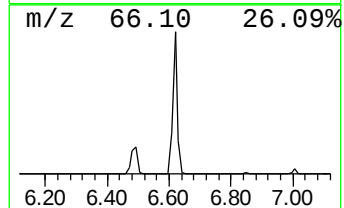
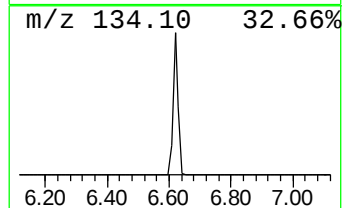
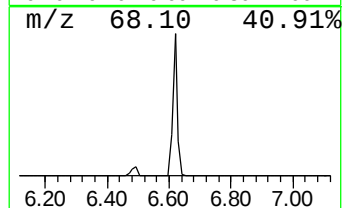
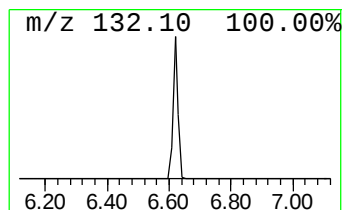
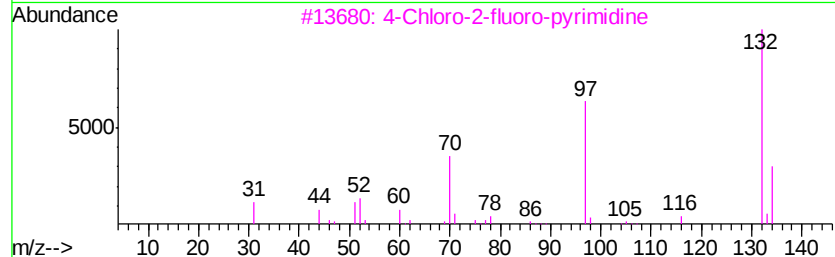
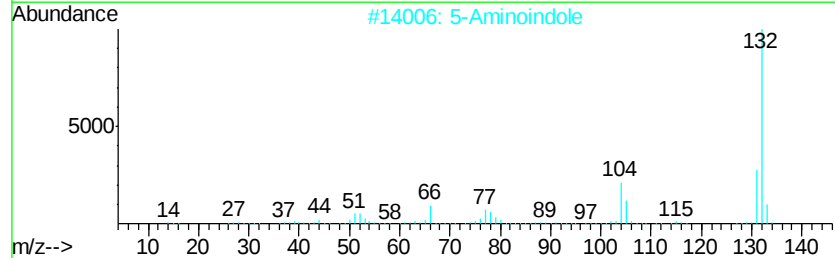
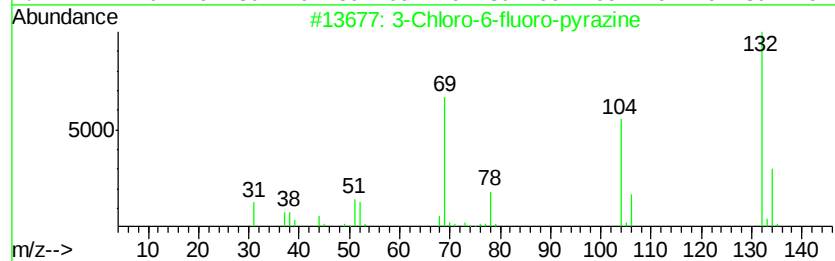
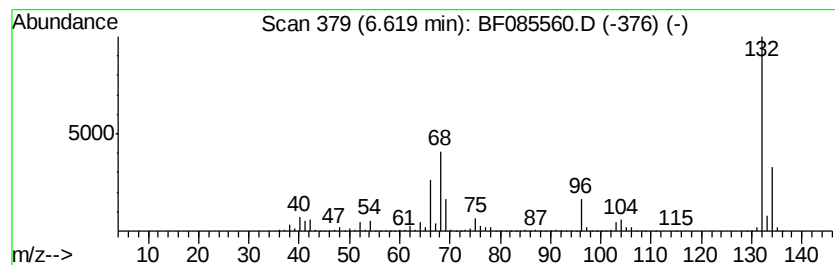
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	87.80 ng	3311080	1,4-Dichlorobenzene-d4	6.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	27
2	5-Aminoindole			132	C8H8N2	005192-03-0	12
3	4-Chloro-2-fluoro-pyrimidine			132	C4H2ClFN2	051422-00-5	9
4	2H-Cyclopenta[d]pyridazine, 2-me...			132	C8H8N2	022291-85-6	9
5	1H-Benzimidazole, 2-methyl-			132	C8H8N2	000615-15-6	9



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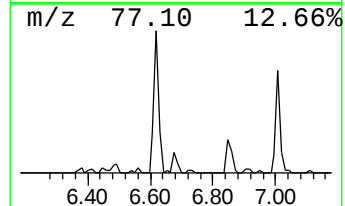
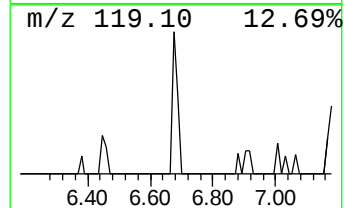
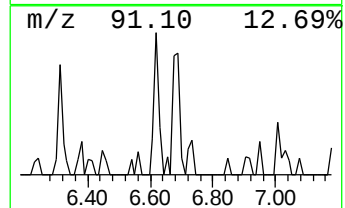
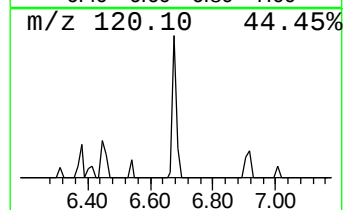
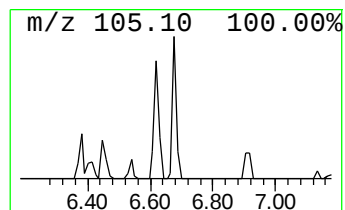
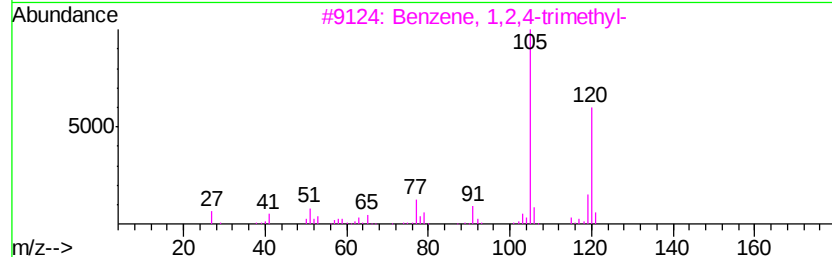
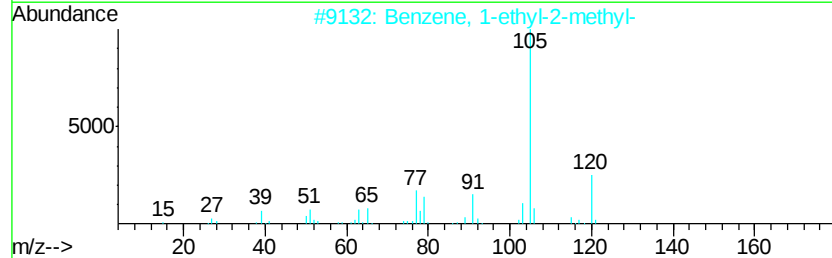
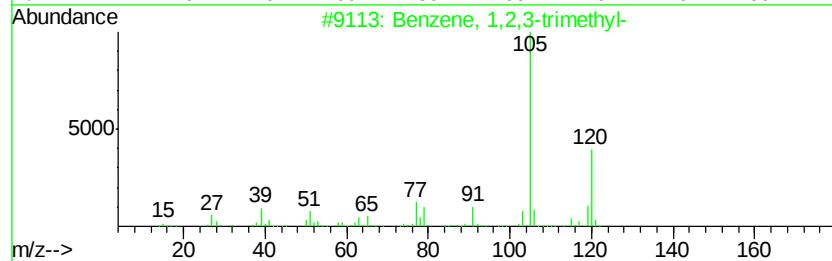
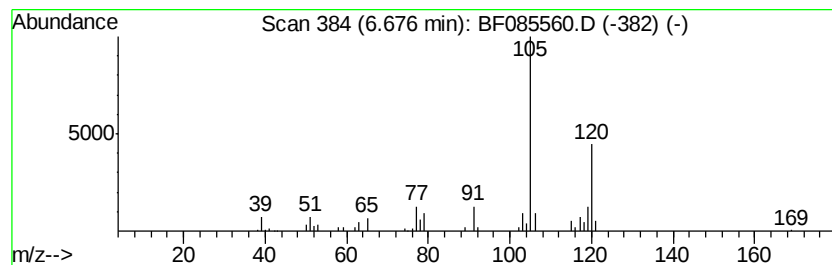
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.68	2.19 ng	82753	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	90
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



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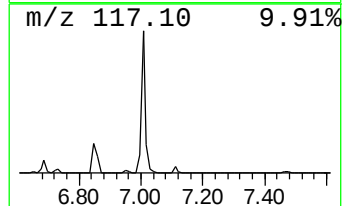
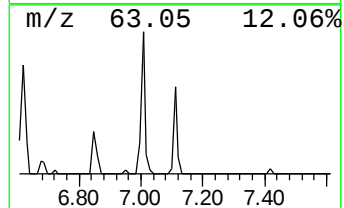
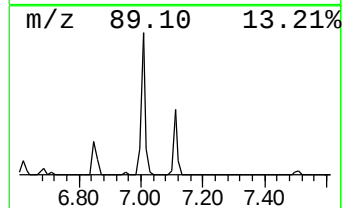
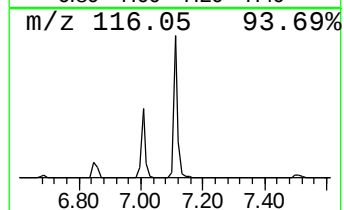
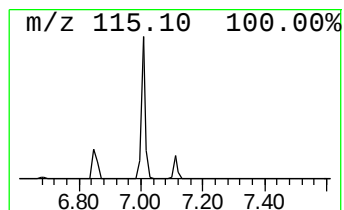
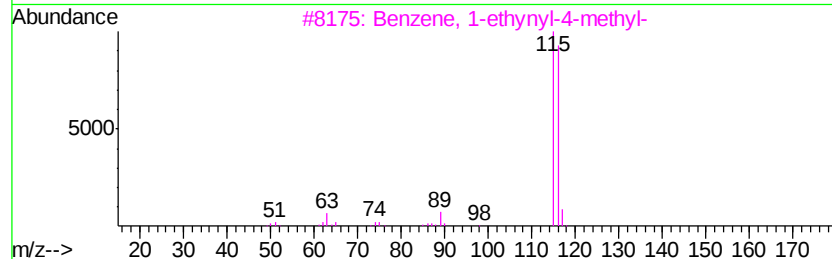
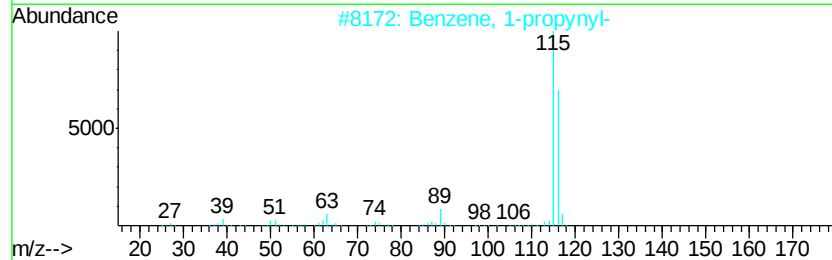
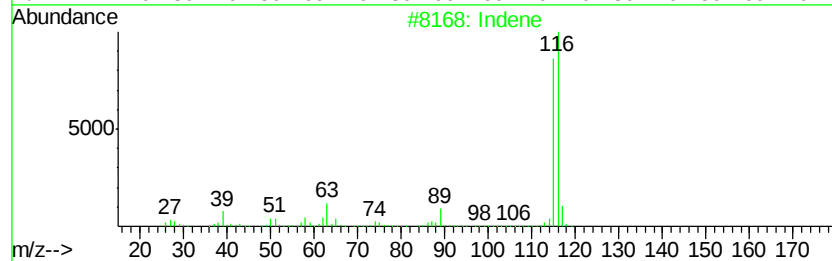
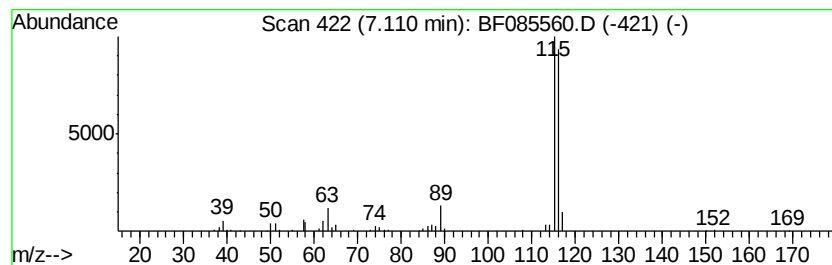
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 8 Indene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.11	3.83 ng	144332	1,4-Dichlorobenzene-d4	6.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indene	116	C9H8	000095-13-6	97
2			Benzene, 1-propynyl-	116	C9H8	000673-32-5	95
3			Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	91
4			Benzene, 1-ethynyl-2-methyl-	116	C9H8	000766-47-2	91
5			Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	91



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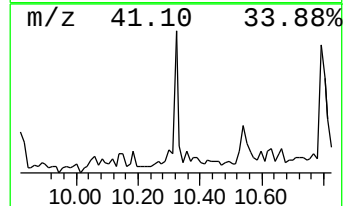
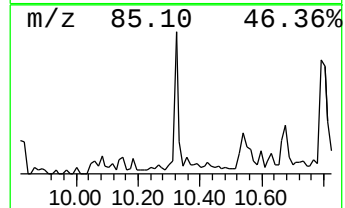
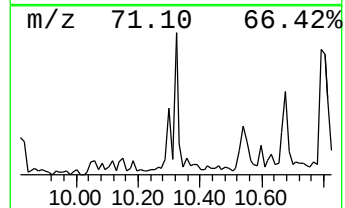
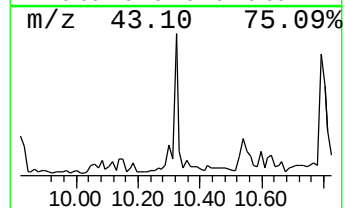
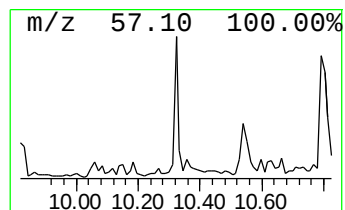
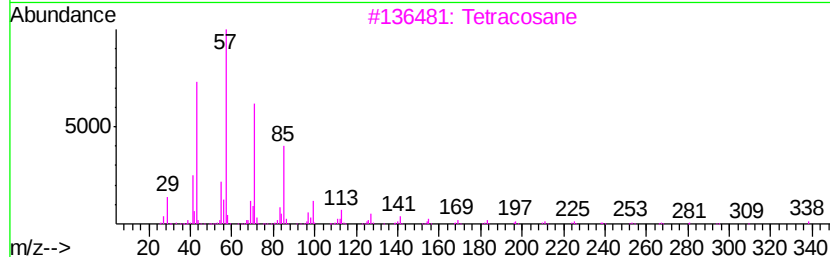
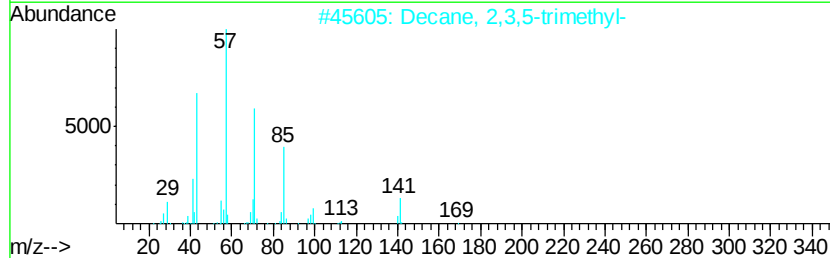
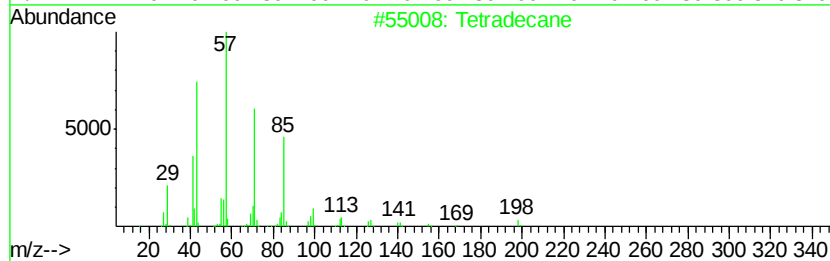
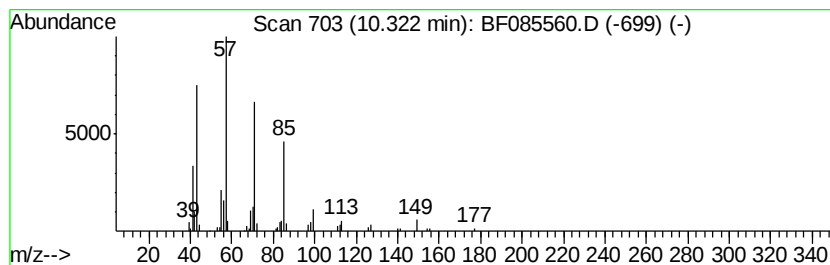
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TRACK-D-GRID-7B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetradecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.32	3.33 ng	180387	Acenaphthene-d10	9.89
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tetradecane	198 C14H30	000629-59-4 86
2		Decane, 2,3,5-trimethyl-	184 C13H28	062238-11-3 83
3		Tetracosane	338 C24H50	000646-31-1 78
4		Pentadecane	212 C15H32	000629-62-9 78
5		Eicosane	282 C20H42	000112-95-8 72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
Data File : BF085560.D
Acq On : 19 Mar 2016 7:59
Operator : UM/SJ
Sample : H1804-03
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TRACK-D-GRID-7B

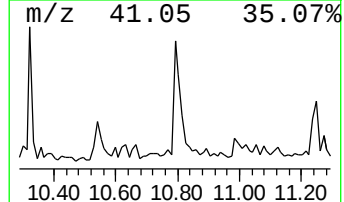
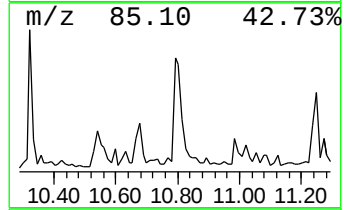
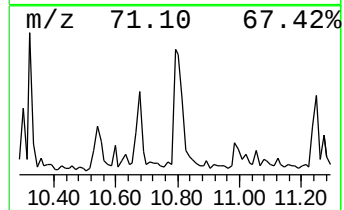
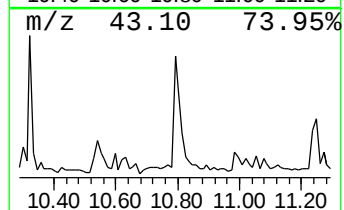
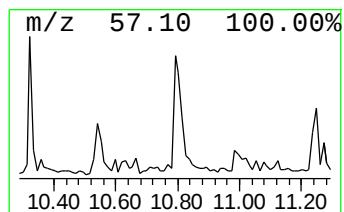
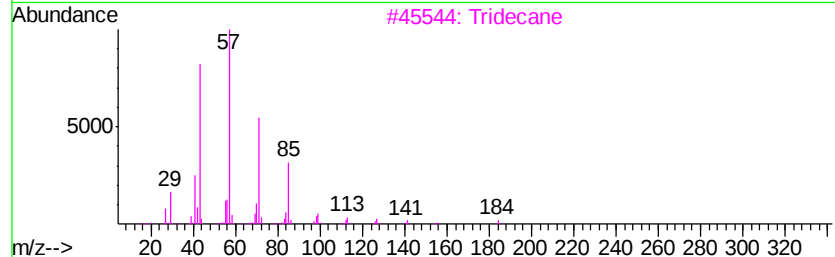
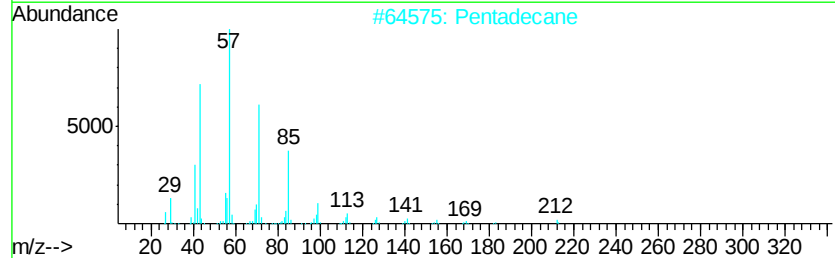
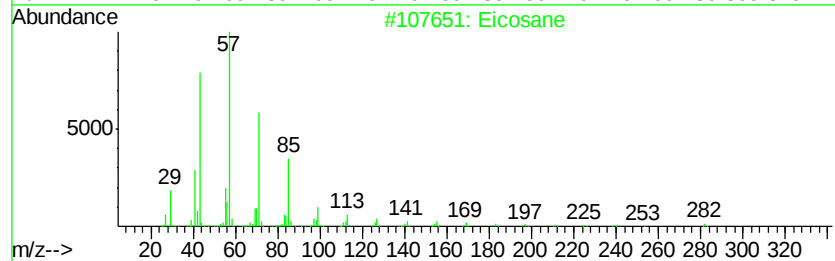
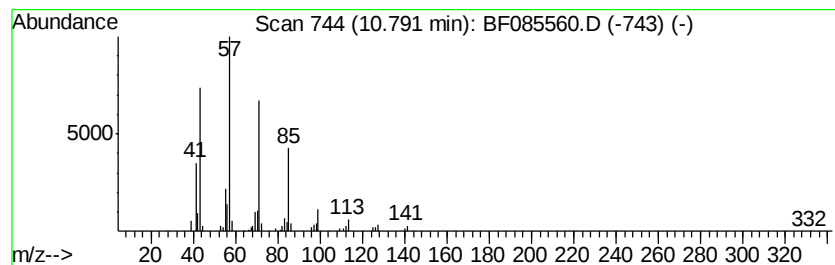
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Eicosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.79	3.20 ng	189390	Phenanthrene-d10	11.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	86
2		Pentadecane	212	C15H32	000629-62-9	80
3		Tridecane	184	C13H28	000629-50-5	80
4		Dodecane, 2-methyl-	184	C13H28	001560-97-0	78
5		Nonane, 5-butyl-	184	C13H28	017312-63-9	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
Data File : BF085560.D
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ALS Vial : 24 Sample Multiplier: 1

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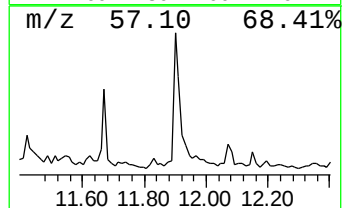
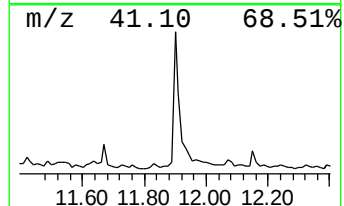
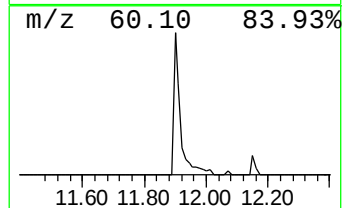
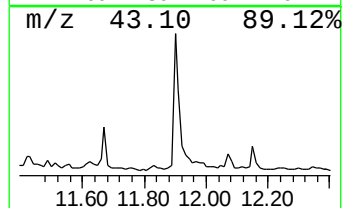
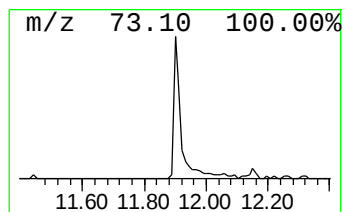
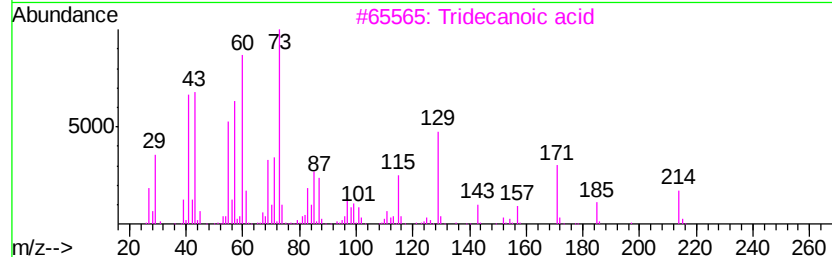
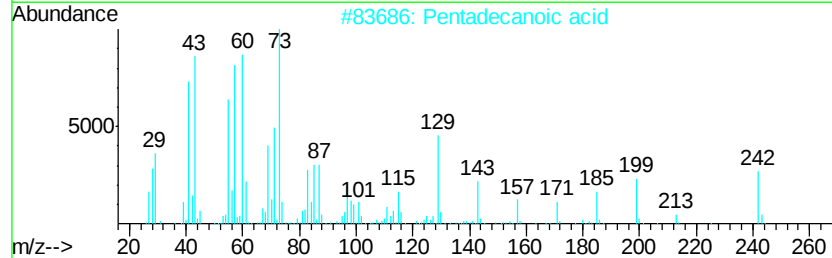
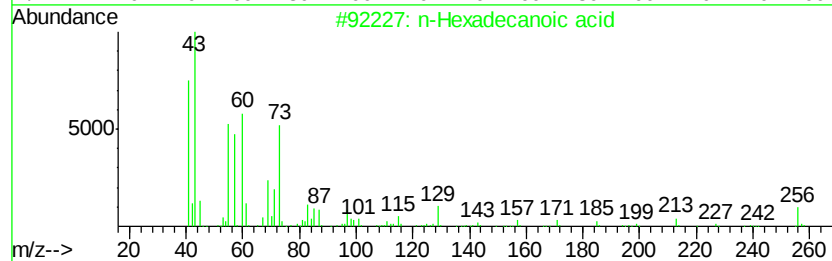
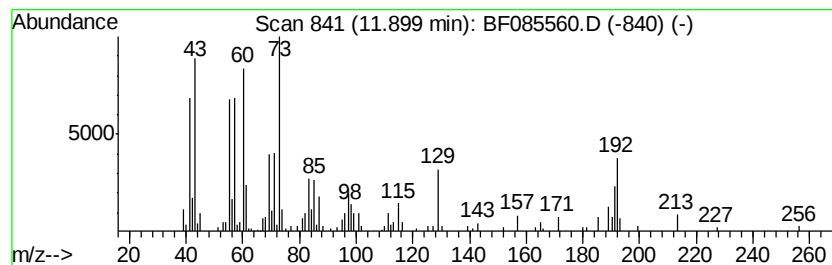
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.90	3.91 ng	230864	Phenanthrene-d10	11.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	64
3		Tridecanoic acid	214	C13H26O2	000638-53-9	52
4		Dodecane, 1,2-dibromo-	326	C12H24Br2	055334-42-4	15
5		Hexadecane	226	C16H34	000544-76-3	11



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
Data File : BF085560.D
Acq On : 19 Mar 2016 7:59
Operator : UM/SJ
Sample : H1804-03
Misc :
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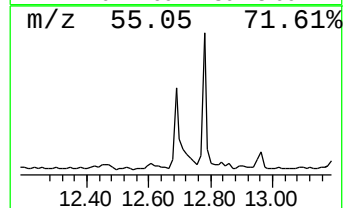
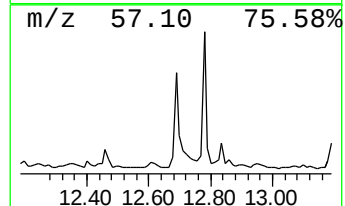
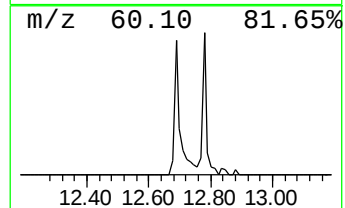
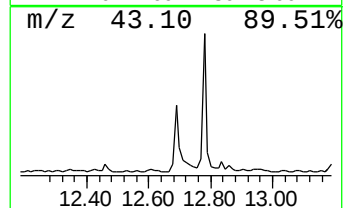
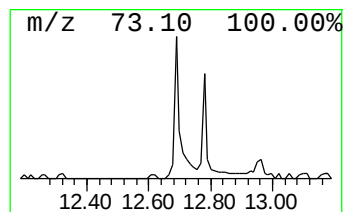
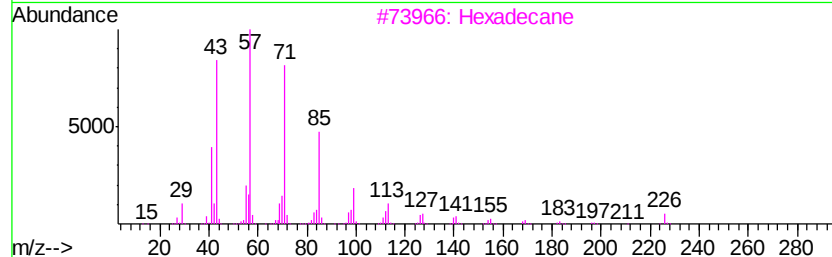
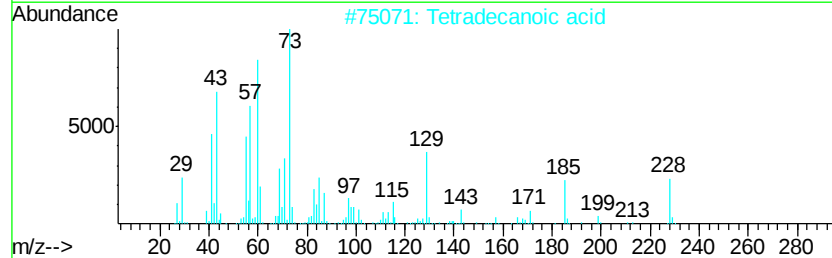
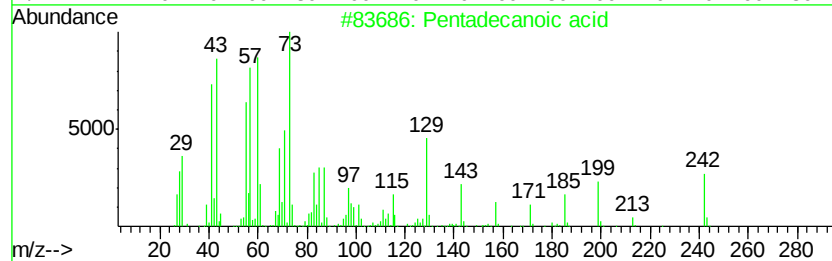
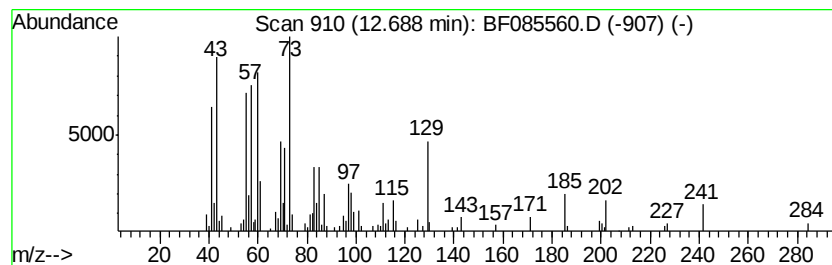
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Pentadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.69	3.39 ng	187267	Chrysene-d12	14.00		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecanoic acid	242	C15H30O2	001002-84-2	86
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
3		Hexadecane	226	C16H34	000544-76-3	25
4		Tetracontane, 3,5,24-trimethyl-	605	C43H88	055162-61-3	18
5		tert-Hexadecanethiol	258	C16H34S	025360-09-2	15



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
Data File : BF085560.D
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Operator : UM/SJ
Sample : H1804-03
Misc :
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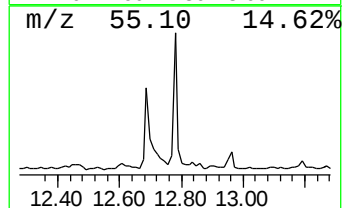
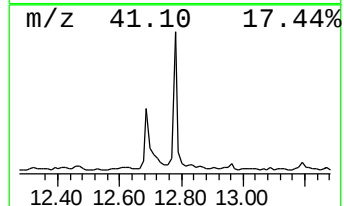
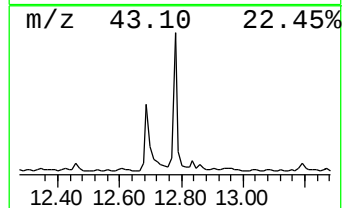
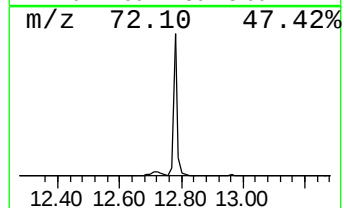
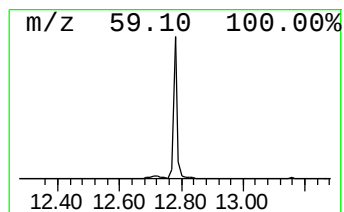
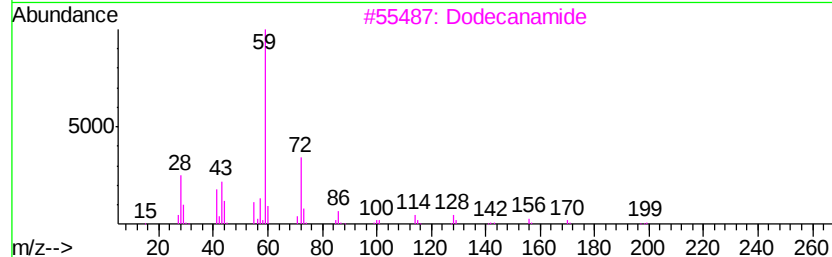
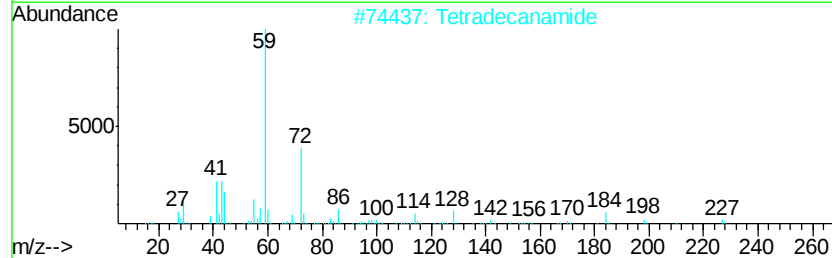
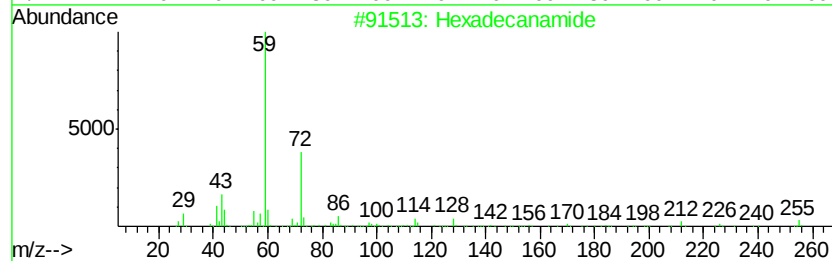
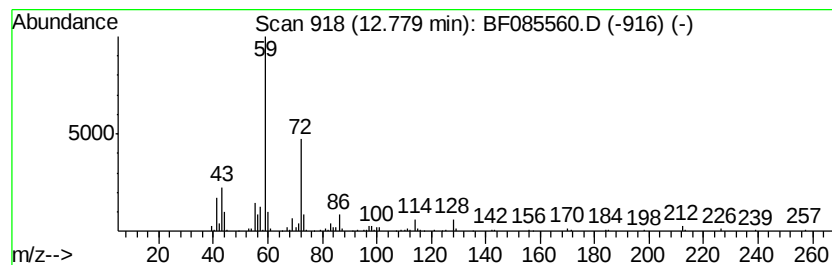
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hexadecanamide Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.78	4.52 ng	250043	Chrysene-d12		14.00
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanamide	255	C16H33NO	000629-54-9	91
2	Tetradecanamide	227	C14H29NO	000638-58-4	90
3	Dodecanamide	199	C12H25NO	001120-16-7	72
4	Pentadecanamide, 15-bromo-	319	C15H30BrNO	1000163-86-1	64
5	Octanamide	143	C8H17NO	000629-01-6	56



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
Data File : BF085560.D
Acq On : 19 Mar 2016 7:59
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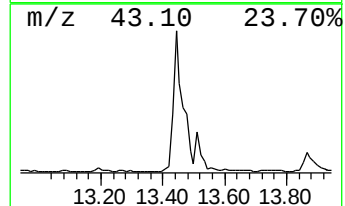
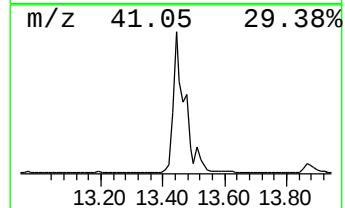
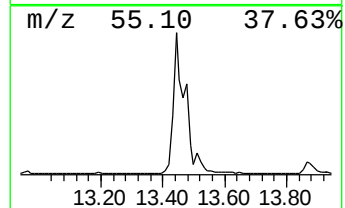
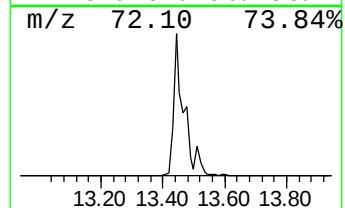
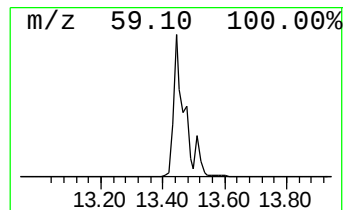
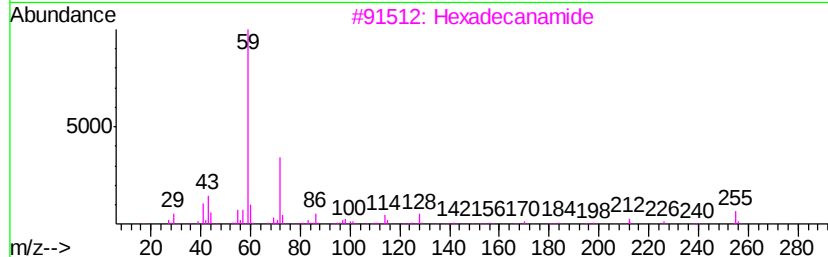
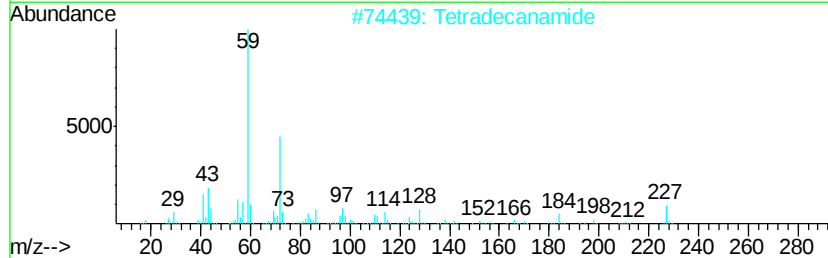
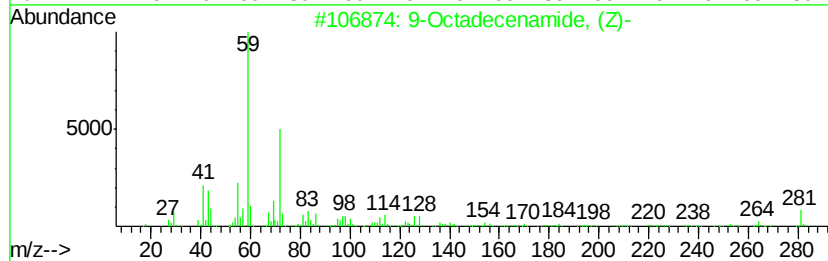
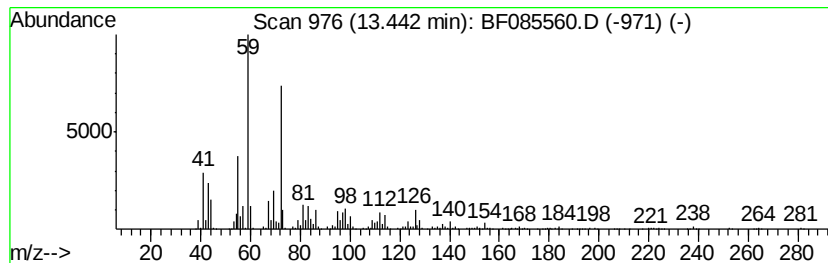
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 9-Octadecenamide, (Z)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.44	40.97 ng	2266070	Chrysene-d12	14.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	72
2		Tetradecanamide	227	C14H29NO	000638-58-4	50
3		Hexadecanamide	255	C16H33NO	000629-54-9	50
4		Nonanamide	157	C9H19NO	001120-07-6	43
5		Undecanamide, 11-bromo-	263	C11H22BrNO	005875-26-3	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
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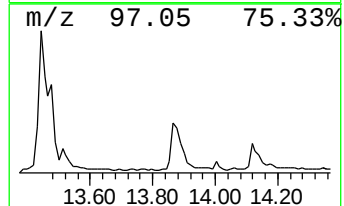
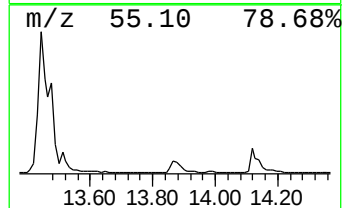
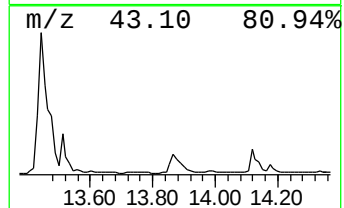
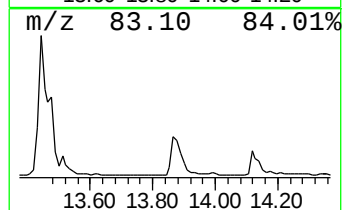
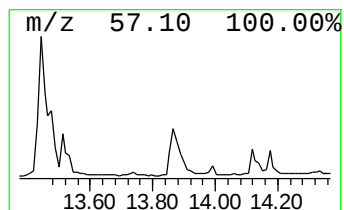
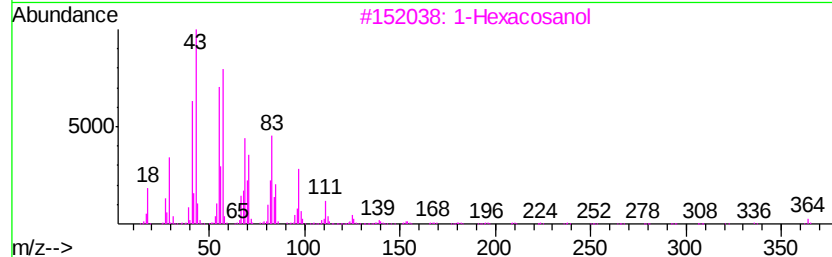
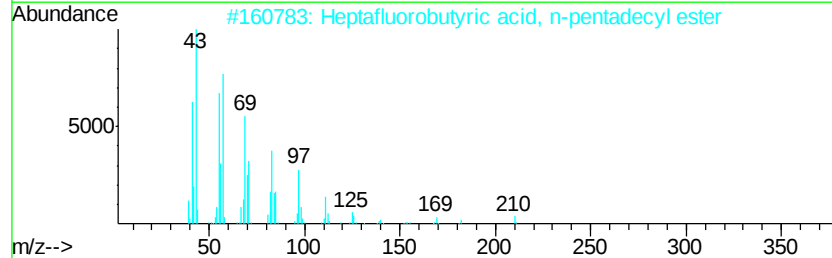
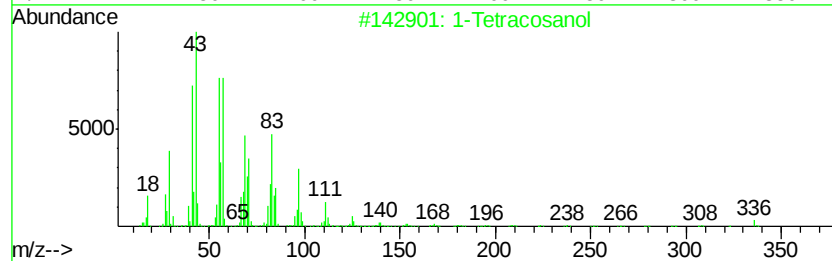
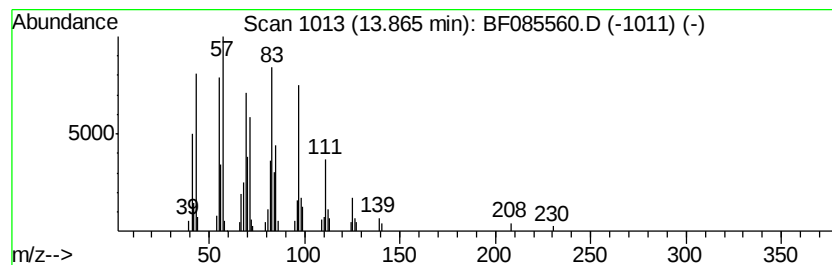
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 1-Tetracosanol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.		
13.87	3.55 ng	196236	Chrysene-d12		14.00		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Tetracosanol	354	C24H50O	000506-51-4	91
2			Heptafluorobutyric acid, n-penta...	424	C19H31F7O2	1000216-79-3	91
3			1-Hexacosanol	382	C26H54O	000506-52-5	91
4			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	90
5			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
Data File : BF085560.D
Acq On : 19 Mar 2016 7:59
Operator : UM/SJ
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Misc :
ALS Vial : 24 Sample Multiplier: 1

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ClientSampleId :
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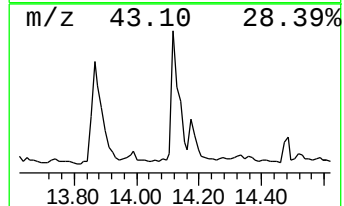
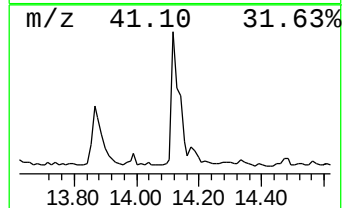
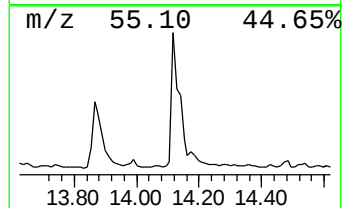
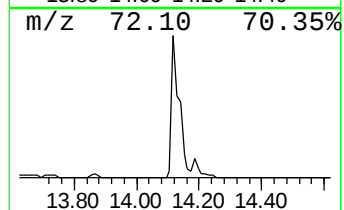
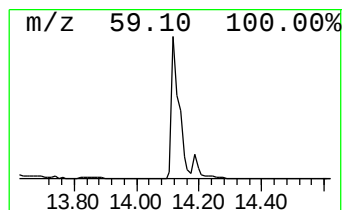
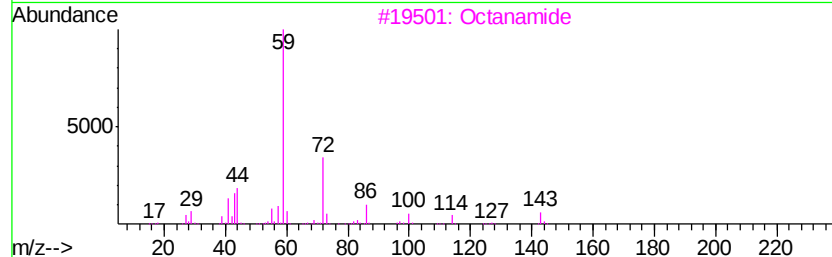
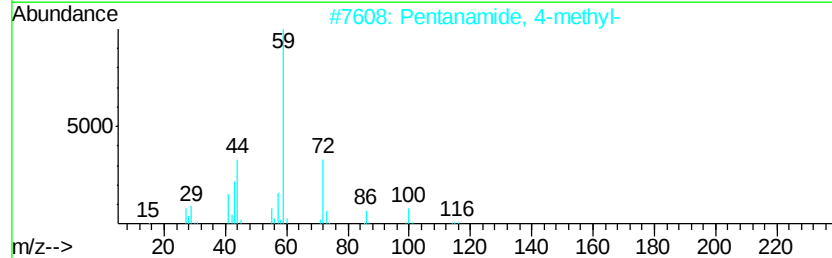
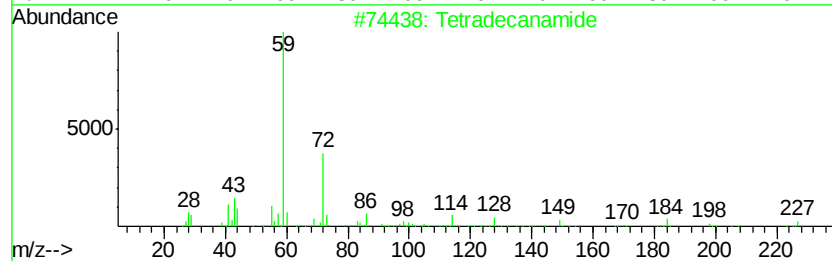
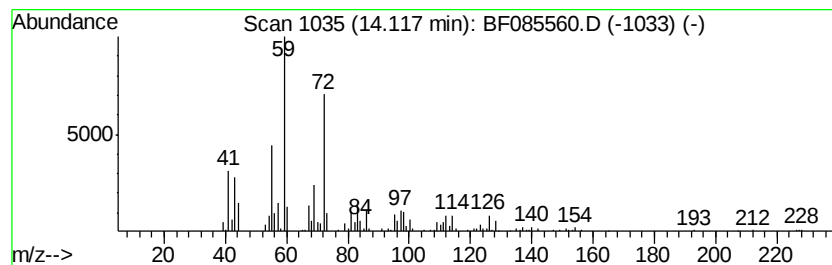
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Tetradecanamide Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.12	5.78 ng	319596	Chrysene-d12	14.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecanamide	227	C14H29NO	000638-58-4	59
2		Pentanamide, 4-methyl-	115	C6H13NO	001119-29-5	59
3		Octanamide	143	C8H17NO	000629-01-6	53
4		Dodecanamide	199	C12H25NO	001120-16-7	53
5		Heptanamide, 4-ethyl-5-methyl-	171	C10H21NO	054789-40-1	45



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031816\
Data File : BF085560.D
Acq On : 19 Mar 2016 7:59
Operator : UM/SJ
Sample : H1804-03
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TRACK-D-GRID-7B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.06	2.0	ng	75680	1	6.85	754226	20.0
unknown6.62	6.62	87.8	ng	3311080	1	6.85	754226	20.0
Benzene, 1,2,3-tr...	6.68	2.2	ng	82753	1	6.85	754226	20.0
Indene	7.11	3.8	ng	144332	1	6.85	754226	20.0
Tetradecane	10.32	3.3	ng	180387	3	9.89	1082380	20.0
Eicosane	10.79	3.2	ng	189390	4	11.37	1182160	20.0
n-Hexadecanoic acid	11.90	3.9	ng	230864	4	11.37	1182160	20.0
Pentadecanoic acid	12.69	3.4	ng	187267	5	14.00	1106300	20.0
Hexadecanamide	12.78	4.5	ng	250043	5	14.00	1106300	20.0
9-Octadecenamide,...	13.44	41.0	ng	2266070	5	14.00	1106300	20.0
1-Tetracosanol	13.87	3.5	ng	196236	5	14.00	1106300	20.0
Tetradecanamide	14.12	5.8	ng	319596	5	14.00	1106300	20.0