

Data Path : Z:\HPCHEM1\BNA F\DATA\BF010716\
 Data File : BF084318.D
 Acq On : 8 Jan 2016 00:38
 Operator : SJ/UM
 Sample : H1028-08
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SW005

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	5.058	257	260	267	rVB	1523500	2365406	29.09%	3.250%
2	5.481	294	297	299	rBV	7206640	7781336	95.71%	10.693%
3	6.510	384	387	390	rVB	5748531	7799131	95.93%	10.717%
4	6.636	395	398	401	rBV	6232527	7615342	93.67%	10.465%
5	6.864	416	418	421	rVB	1921108	1565152	19.25%	2.151%
6	7.024	429	432	434	rBV	6777850	5930257	72.94%	8.149%
7	7.436	465	468	470	rBV	5133135	5501699	67.67%	7.560%
8	8.156	528	531	533	rBV	2145216	2125213	26.14%	2.920%
9	9.230	622	625	628	rBV	8407741	8130234	100.00%	11.172%
10	9.619	657	659	662	rBV	719345	996888	12.26%	1.370%
11	9.905	682	684	687	rVB	2502886	2241242	27.57%	3.080%
12	10.339	718	722	724	rBV2	141070	151851	1.87%	0.209%
13	10.556	738	741	745	rBV	51679	100396	1.23%	0.138%
14	10.693	750	753	756	rBV2	4464129	5493453	67.57%	7.549%
15	10.808	762	763	770	rVB	156929	297524	3.66%	0.409%
16	11.265	801	803	804	rBV	135467	142646	1.75%	0.196%
17	11.390	812	814	816	rBV	2563993	2244977	27.61%	3.085%
18	11.448	816	819	823	rVB	40792	85201	1.05%	0.117%
19	12.979	950	953	956	rBV	7267529	7480527	92.01%	10.280%
20	13.882	1029	1032	1042	rBV	267641	492491	6.06%	0.677%
21	14.019	1042	1044	1047	rBV2	1620408	1847216	22.72%	2.538%
22	15.448	1167	1169	1172	rBV	916076	1392662	17.13%	1.914%
23	16.134	1224	1229	1240	rVB	121823	495827	6.10%	0.681%
24	16.968	1298	1302	1309	rBV	44277	96503	1.19%	0.133%
25	17.643	1357	1361	1365	rBV	46390	116682	1.44%	0.160%
26	18.431	1425	1430	1437	rBV	47663	147972	1.82%	0.203%
27	19.380	1508	1513	1519	rBV2	38103	132795	1.63%	0.182%

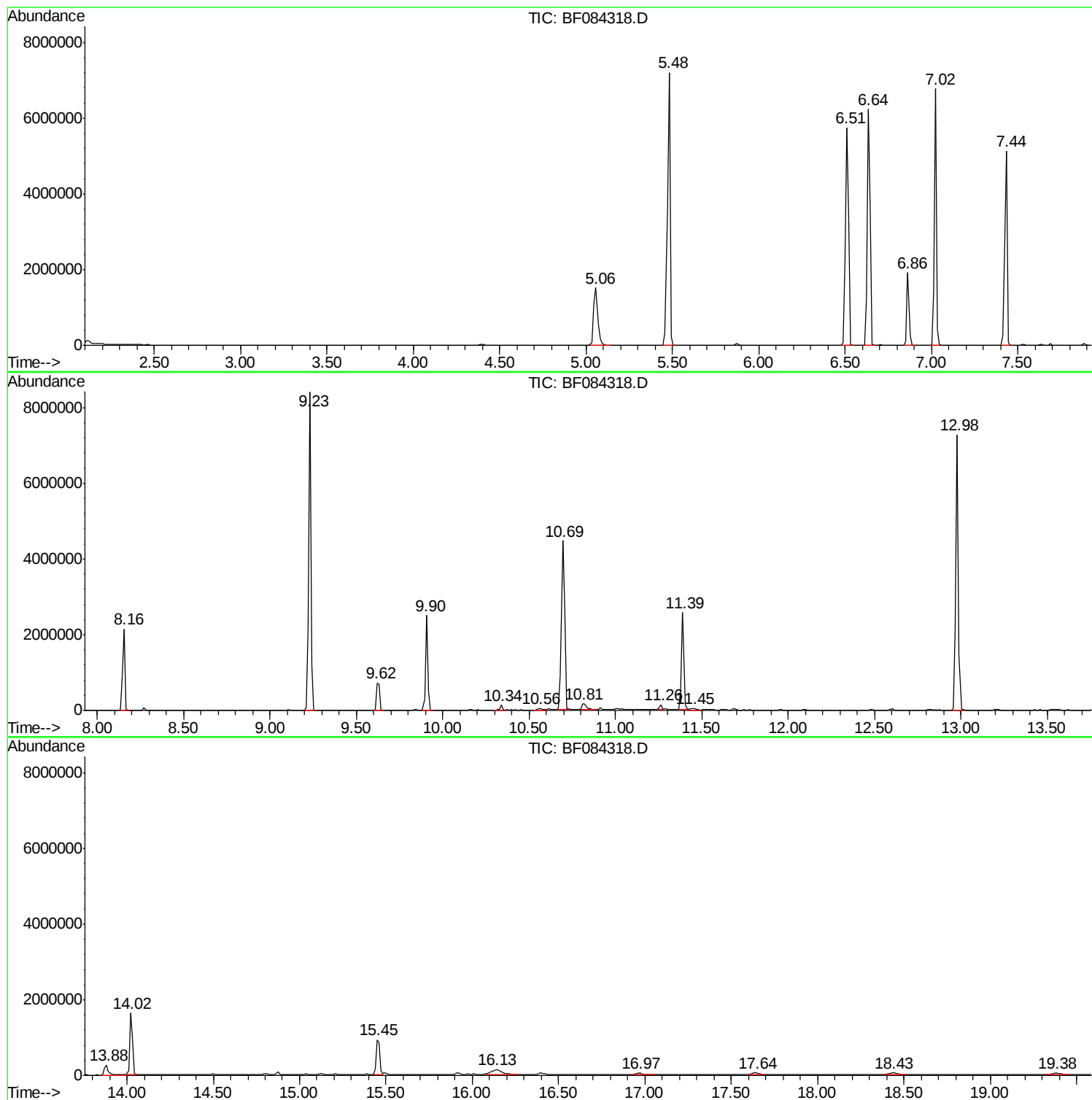
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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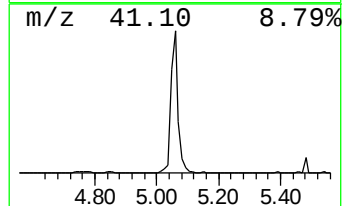
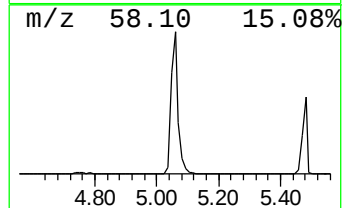
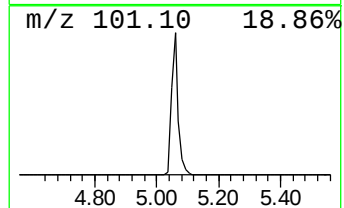
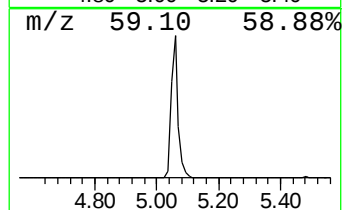
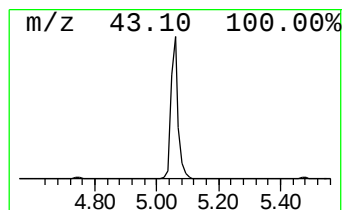
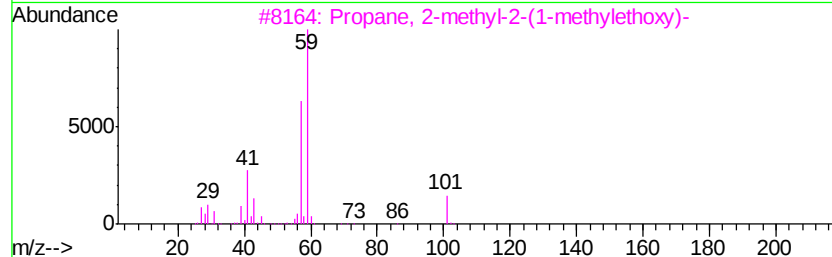
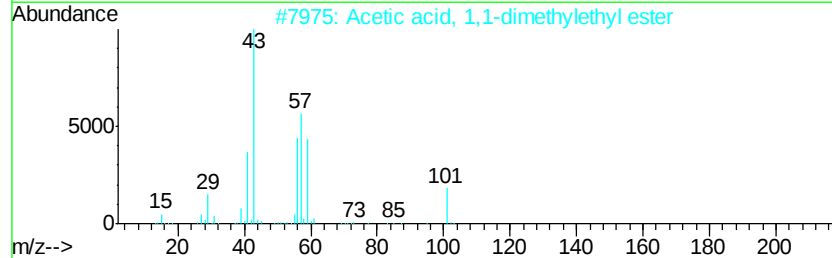
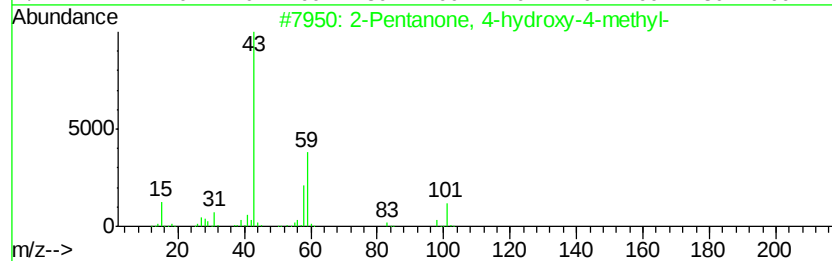
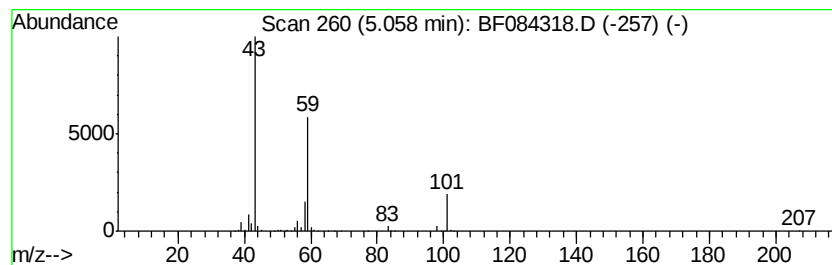
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.06	30.23 ng	2365410	1,4-Dichlorobenzene-d4	6.86

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53	
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39	
3	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33	
4	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33	
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25	



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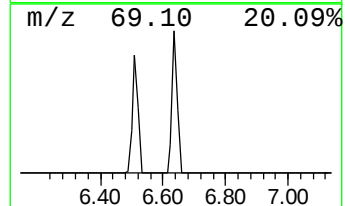
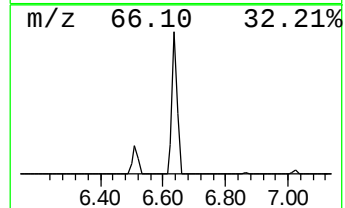
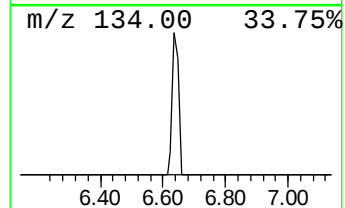
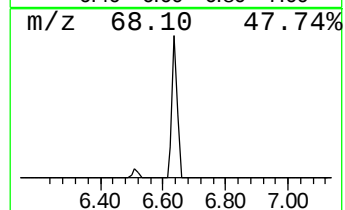
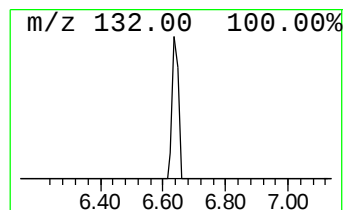
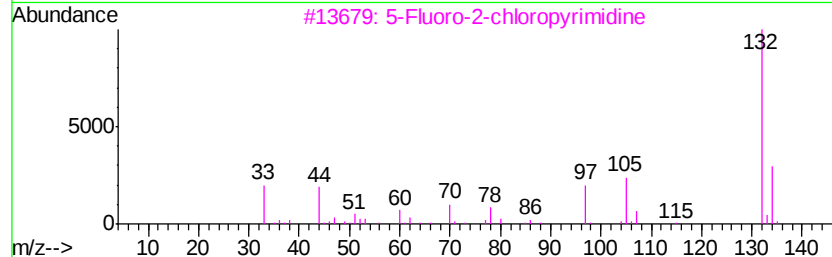
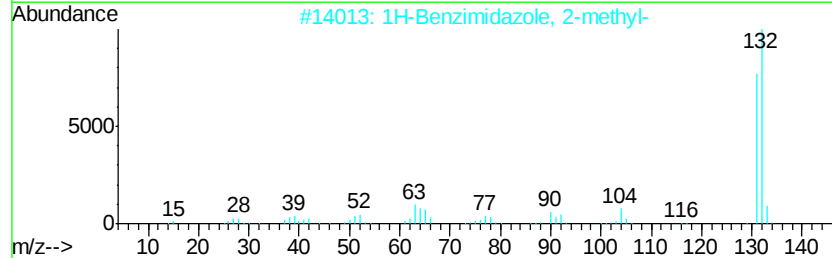
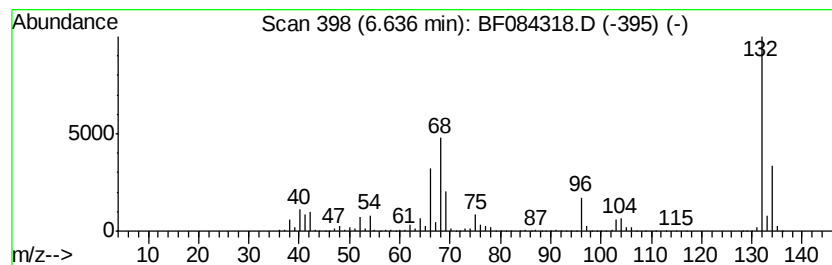
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.64 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.64	97.31 ng	7615340	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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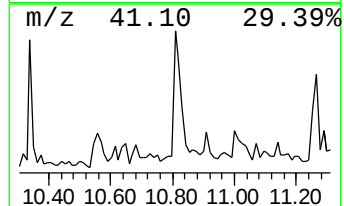
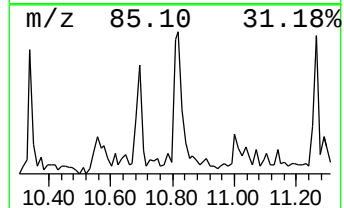
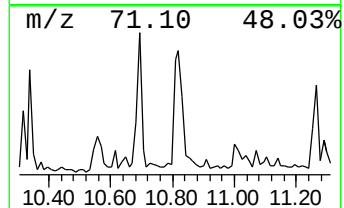
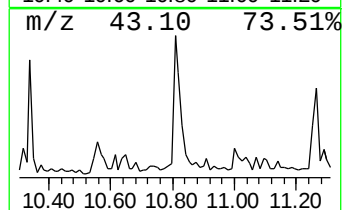
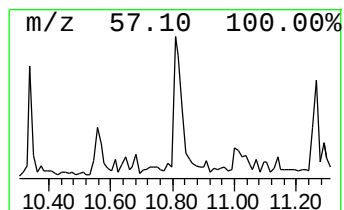
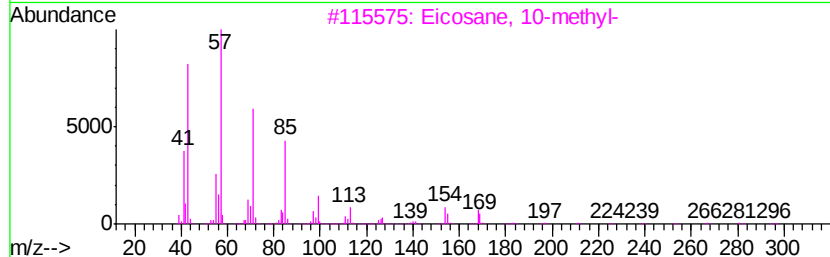
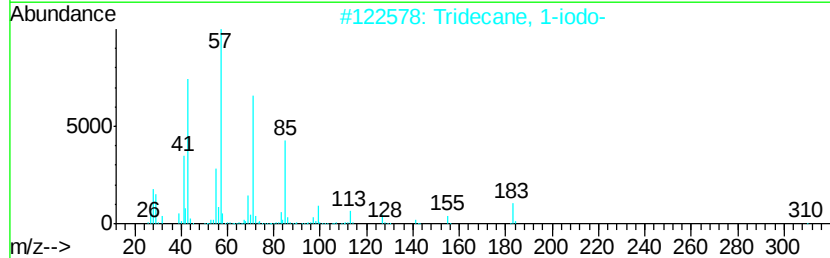
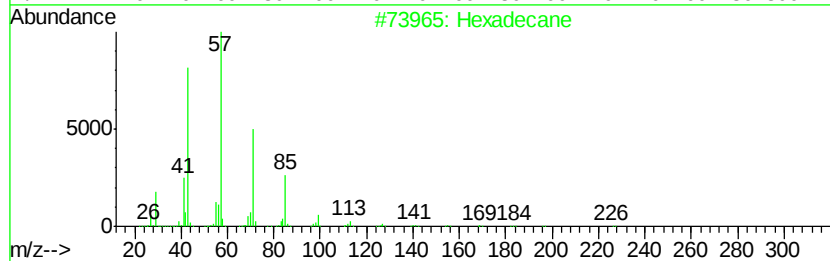
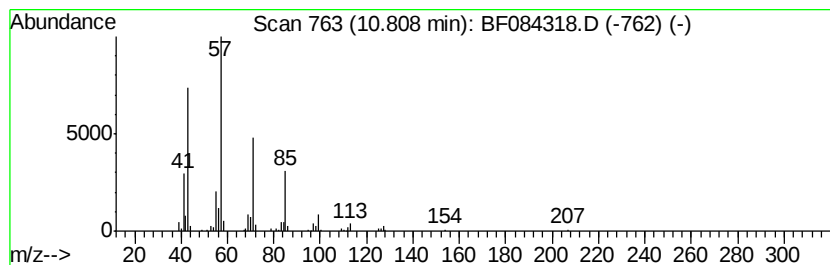
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Peak Number 4 Hexadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.81	2.65 ng	297524	Phenanthrene-d10	11.39	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	91
2	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90
3	Eicosane, 10-methyl-	296	C21H44	054833-23-7	90
4	Eicosane	282	C20H42	000112-95-8	90
5	Pentadecane	212	C15H32	000629-62-9	83



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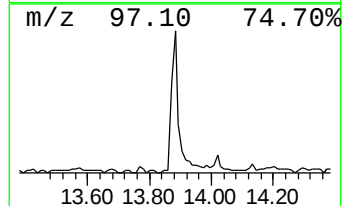
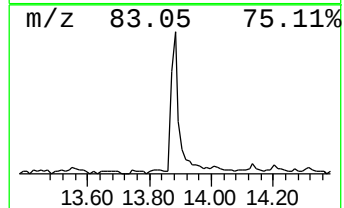
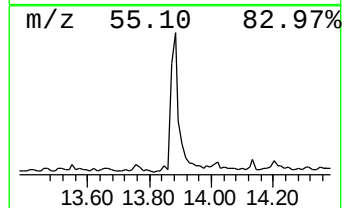
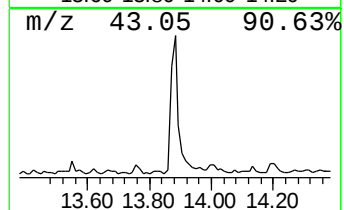
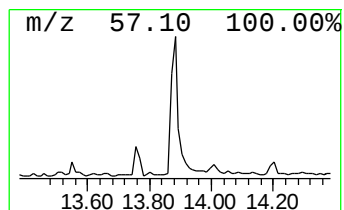
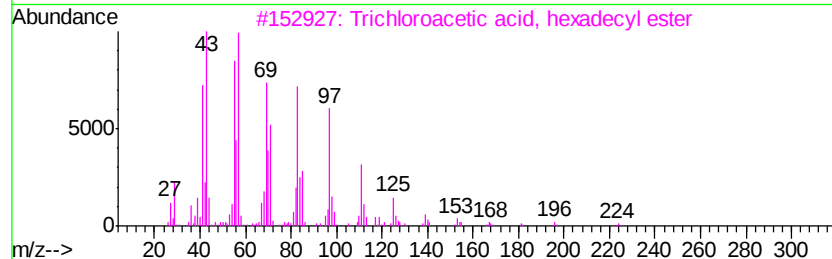
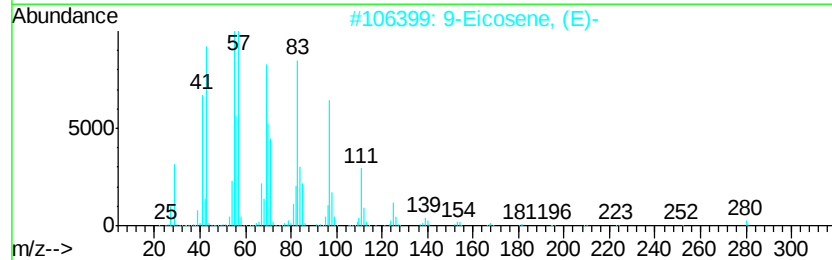
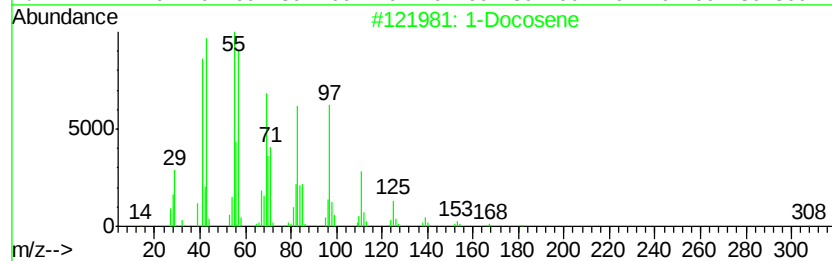
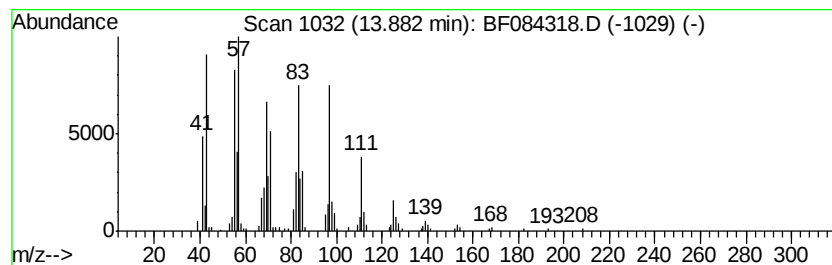
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 1-Docosene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.88	5.33 ng	492491	Chrysene-d12	14.02
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Docosene		308 C22H44	001599-67-3 91
2	9-Eicosene, (E)-		280 C20H40	074685-29-3 90
3	Trichloroacetic acid, hexadecyl ...		386 C18H33Cl3O2	074339-54-1 90
4	5-Eicosene, (E)-		280 C20H40	074685-30-6 87
5	17-Pentatriacontene		491 C35H70	006971-40-0 86



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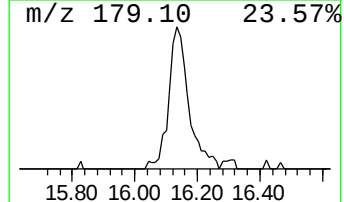
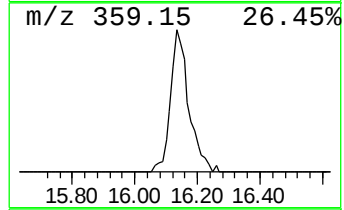
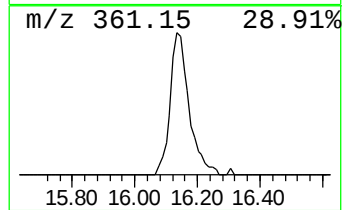
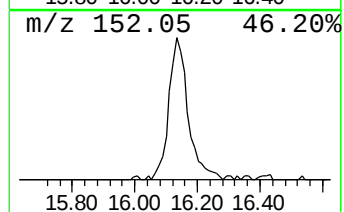
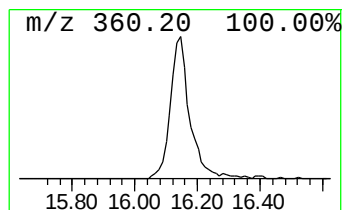
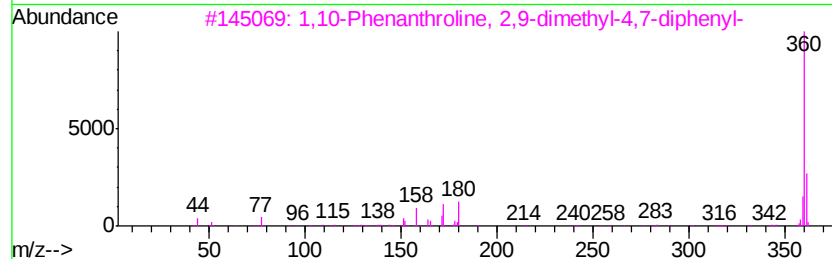
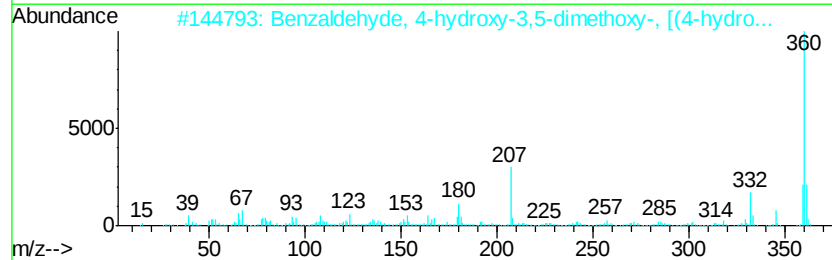
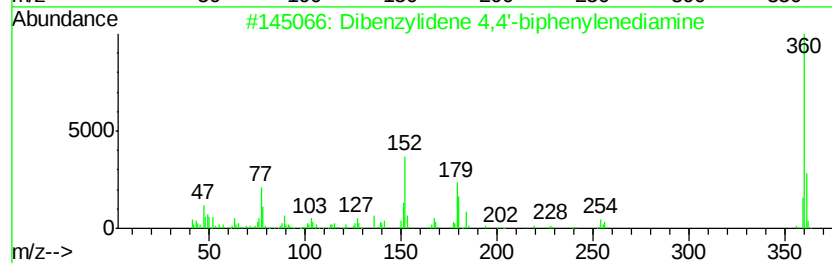
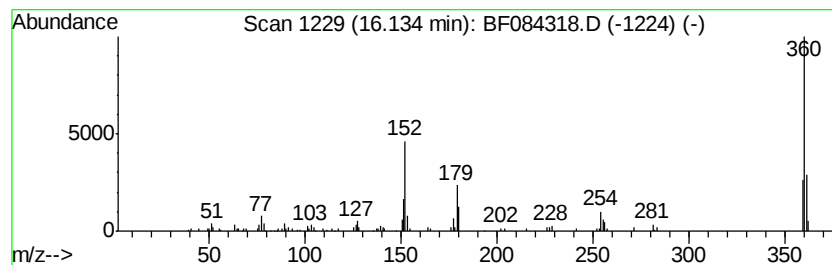
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 6 Dibenzylidene 4,4'-biphenyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.13	7.12 ng	495827	Perylene-d12	15.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dibenzylidene 4,4'-biphenylenedi...	360	C26H20N2	006311-48-4	91
2		Benzaldehyde, 4-hydroxy-3,5-dime...	360	C18H20N2O6	014414-32-5	49
3		1,10-Phenanthroline, 2,9-dimethv...	360	C26H20N2	004733-39-5	47
4		N,N'-di-2-Naphthyl-p-phenylenedi...	360	C26H20N2	000093-46-9	46
5		1,3,2-Dioxaborolane, bis(spiro[4...	360	C20H13B06	1000150-23-7	43



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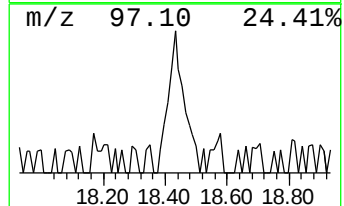
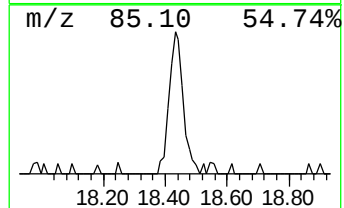
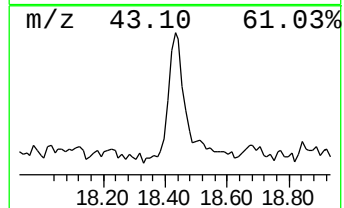
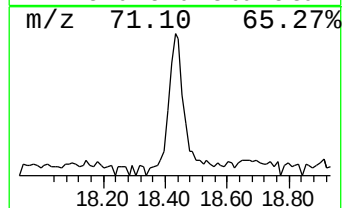
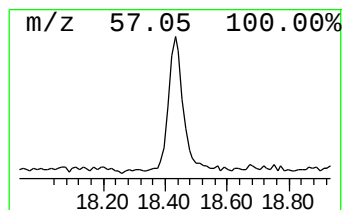
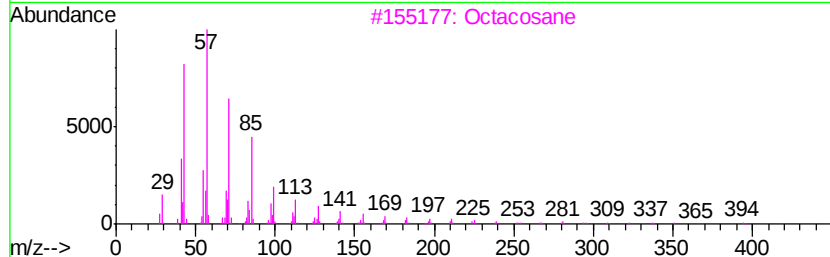
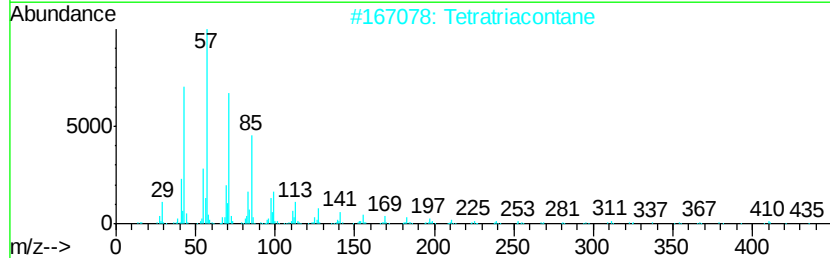
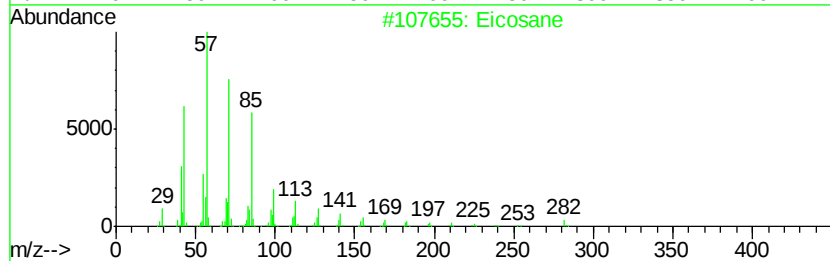
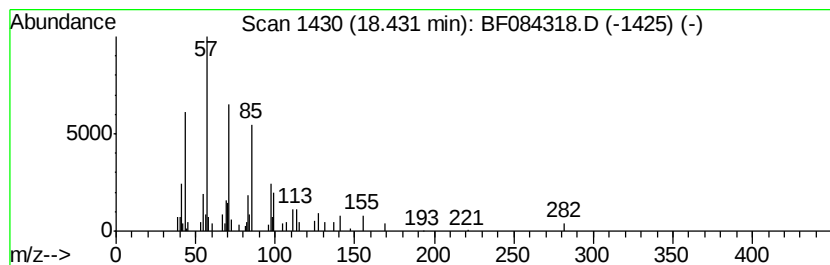
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Eicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.43	2.13 ng	147972	Perylene-d12	15.46

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	94
2	Tetratriacontane		479	C34H70	014167-59-0	90
3	Octacosane		394	C28H58	000630-02-4	87
4	Pentacosane		352	C25H52	000629-99-2	83
5	Tetratetracontane		619	C44H90	007098-22-8	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010716\
Data File : BF084318.D
Acq On : 8 Jan 2016 00:38
Operator : SJ/UM
Sample : H1028-08
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SW005

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
2-Pentanone, 4-hy...	5.06	30.2	ng	2365410	1	6.86	1565150	20.0
unknown6.64	6.64	97.3	ng	7615340	1	6.86	1565150	20.0
Hexadecane	10.81	2.6	ng	297524	4	11.39	2244980	20.0
1-Docosene	13.88	5.3	ng	492491	5	14.02	1847220	20.0
Dibenzylidene 4,4...	16.13	7.1	ng	495827	6	15.46	1392660	20.0
Eicosane	18.43	2.1	ng	147972	6	15.46	1392660	20.0