

Data Path : Z:\HPCHEM1\BNA F\DATA\BF072417\
 Data File : BF096999.D
 Acq On : 24 Jul 2017 21:47
 Operator : SJ/JU
 Sample : I4361-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RAMP-D-4-(0-2)

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-BF071317.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.346	73	78	86	rVB	108526	163875	4.41%	0.534%
2	4.675	470	474	491	rVB	274501	465938	12.54%	1.519%
3	5.128	546	551	554	rBV	3117637	3430198	92.32%	11.180%
4	6.181	725	730	744	rBV	2907512	3715468	100.00%	12.110%
5	6.293	744	749	752	rBV	3148720	3317560	89.29%	10.813%
6	6.510	783	786	793	rBV	827362	762943	20.53%	2.487%
7	6.669	809	813	817	rBV	2525701	2478993	66.72%	8.080%
8	7.081	878	883	886	rBV	2289501	2480714	66.77%	8.086%
9	7.792	1000	1004	1008	rBV	1076890	967664	26.04%	3.154%
10	8.875	1182	1188	1191	rBV	3127490	3485402	93.81%	11.360%
11	9.269	1251	1255	1259	rBV	339768	286141	7.70%	0.933%
12	9.539	1296	1301	1305	rBV	1113608	1014789	27.31%	3.308%
13	9.969	1367	1374	1376	rBV3	25648	37812	1.02%	0.123%
14	9.998	1376	1379	1382	rVB	51279	45830	1.23%	0.149%
15	10.328	1430	1435	1439	rBV	1798110	2031413	54.67%	6.621%
16	10.463	1455	1458	1465	rBV	85754	104329	2.81%	0.340%
17	10.910	1529	1534	1537	rBV2	41699	40051	1.08%	0.131%
18	11.016	1548	1552	1555	rBV	1171020	1039206	27.97%	3.387%
19	11.569	1642	1646	1655	rBV2	125403	135548	3.65%	0.442%
20	12.604	1816	1822	1825	rBV2	2687735	3100510	83.45%	10.106%
21	13.510	1972	1976	1985	rVB	126373	170725	4.59%	0.556%
22	13.639	1994	1998	2004	rVB	883409	822055	22.13%	2.679%
23	14.992	2223	2228	2234	rVB	496288	583577	15.71%	1.902%

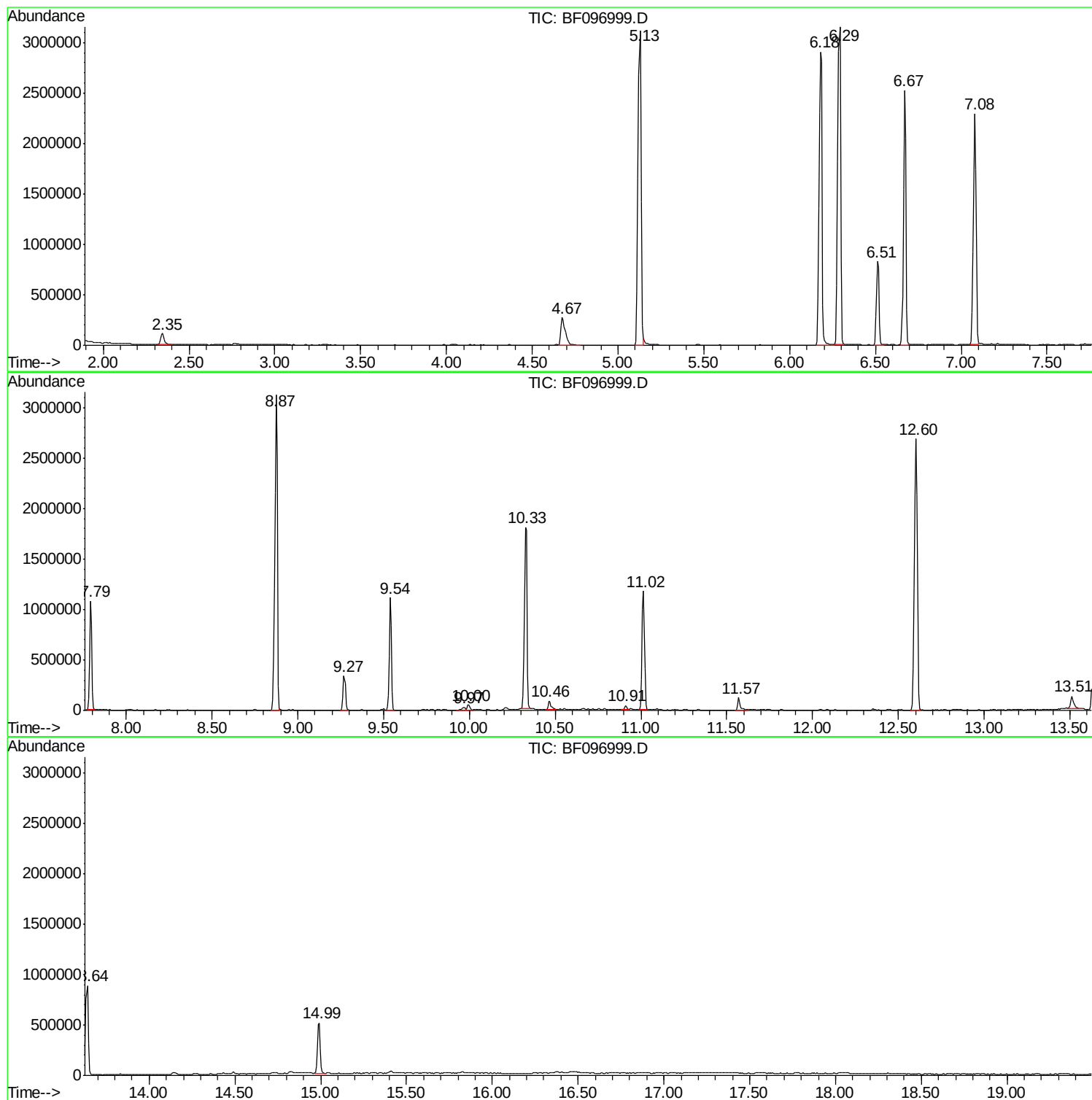
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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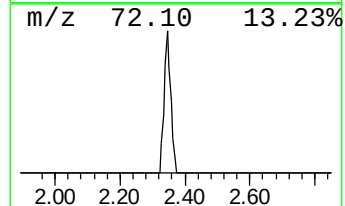
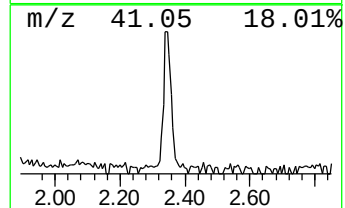
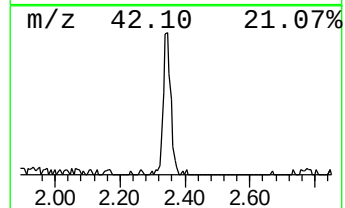
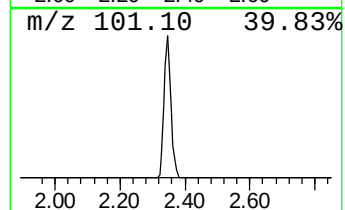
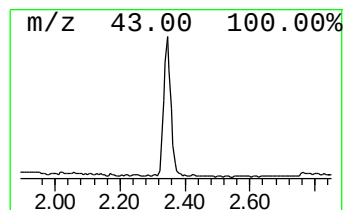
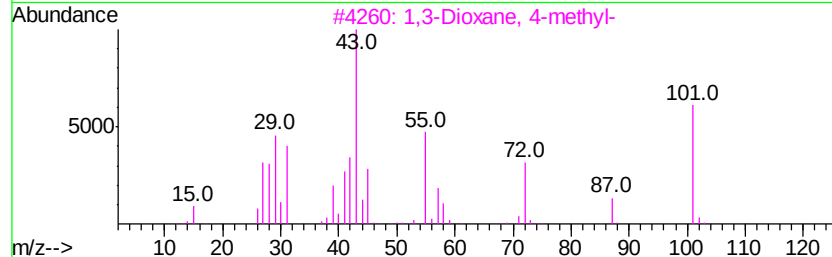
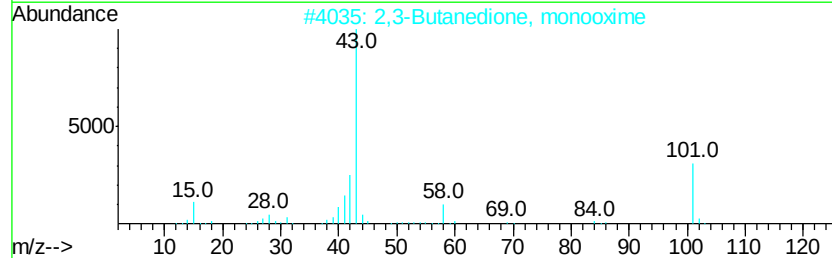
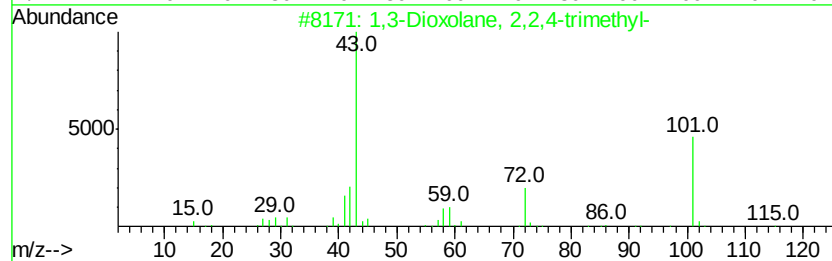
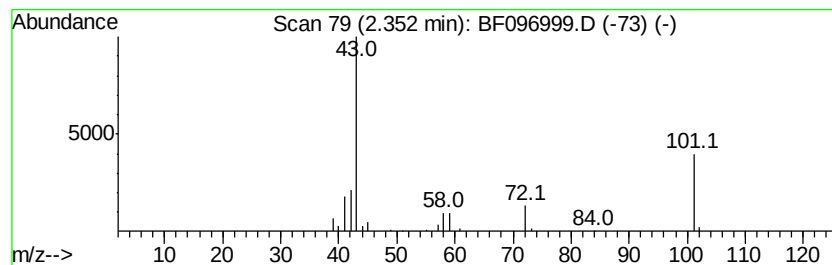
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF071317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 1,3-Dioxolane, 2,2,4-trimet... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.35	4.30 ng	163875	1,4-Dichlorobenzene-d4	6.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Dioxolane, 2,2,4-trimethyl-	116	C6H12O2	001193-11-9	83
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	38
3		1,3-Dioxane, 4-methyl-	102	C5H10O2	001120-97-4	35
4		1,3-Dioxolane-2-methanol, 2,4-di...	132	C6H12O3	053951-43-2	28
5		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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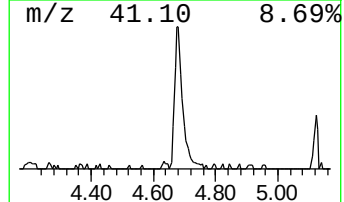
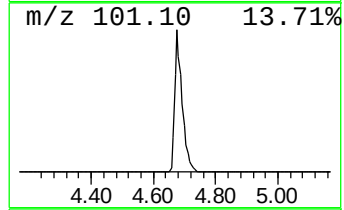
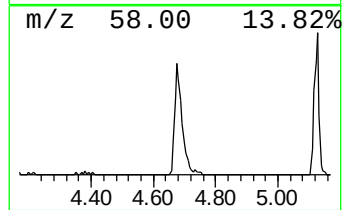
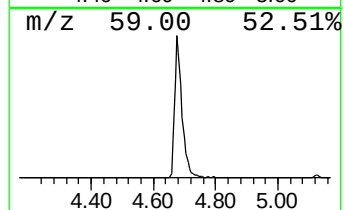
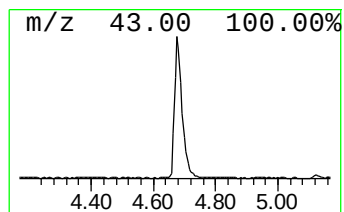
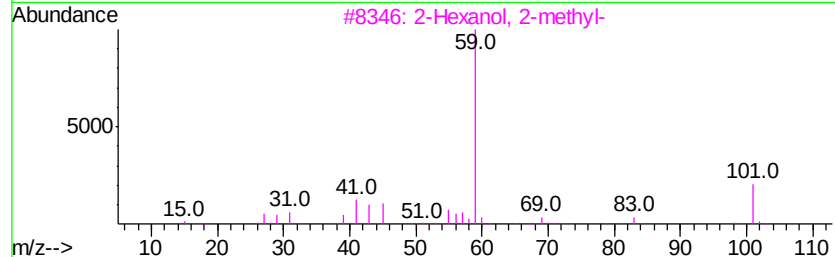
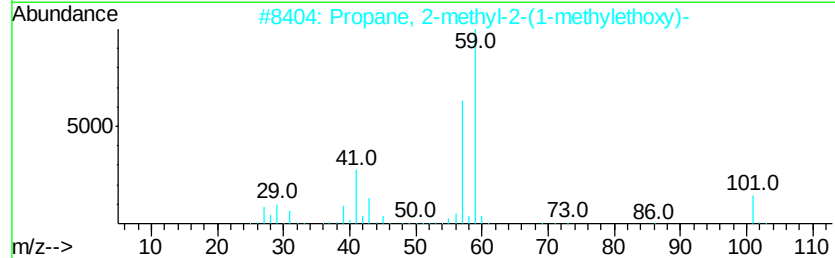
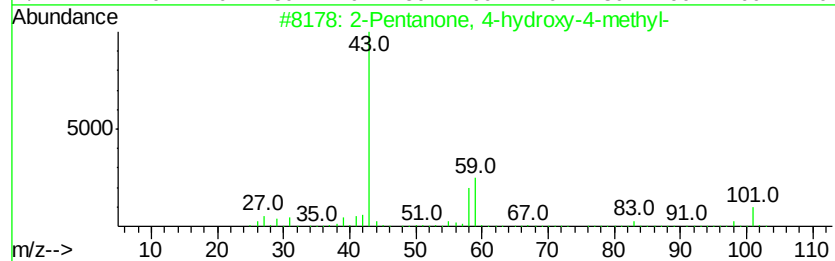
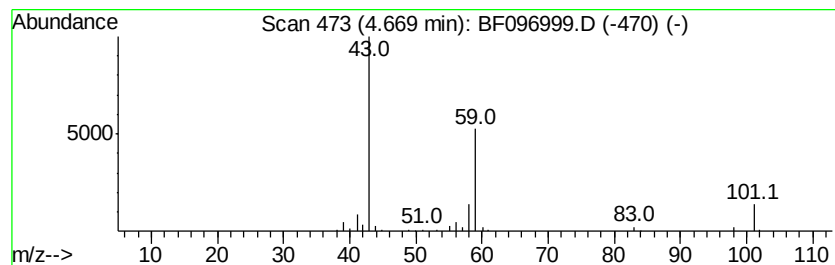
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.67	12.21 ng	465938	1,4-Dichlorobenzene-d4	6.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	53
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	25
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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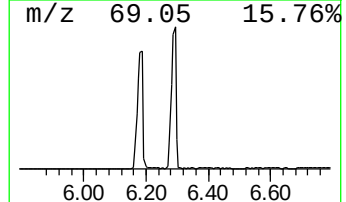
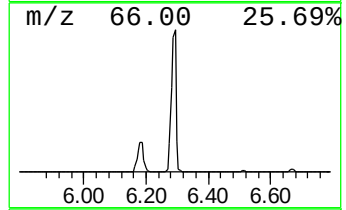
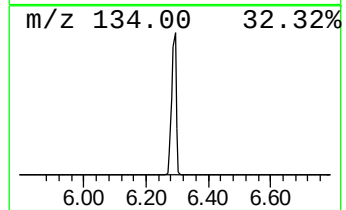
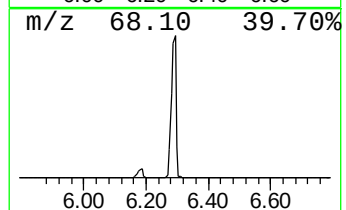
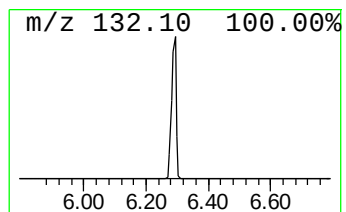
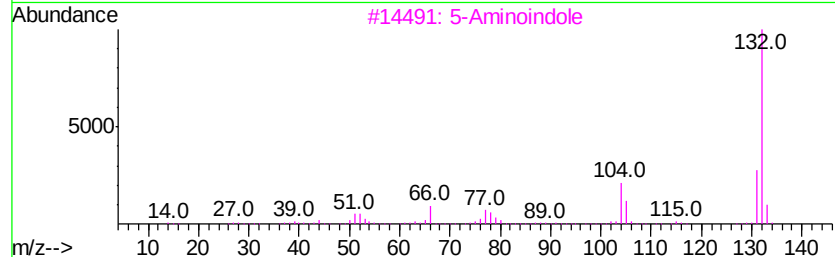
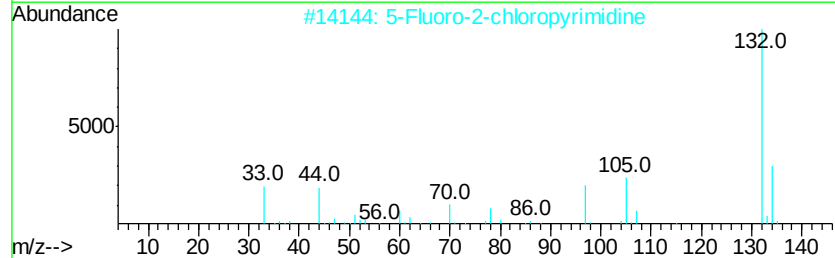
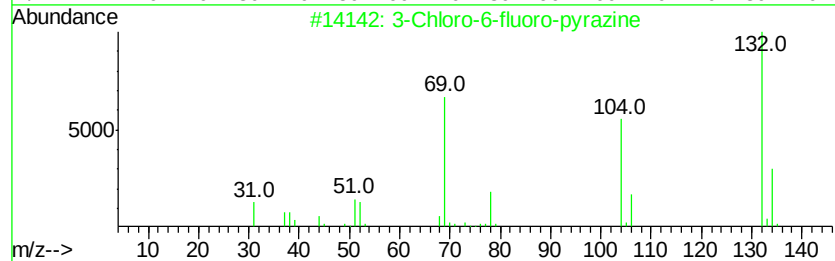
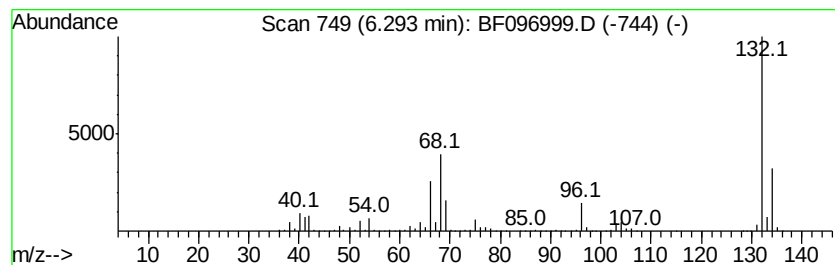
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Peak Number 3 unknown6.29 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.29	86.97 ng	3317560	1,4-Dichlorobenzene-d4	6.51

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	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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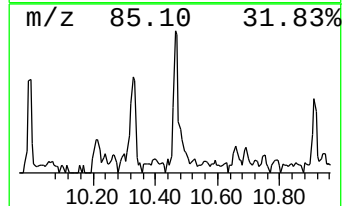
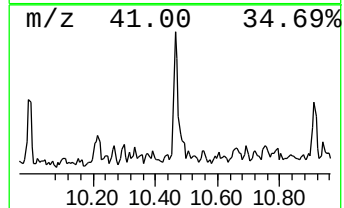
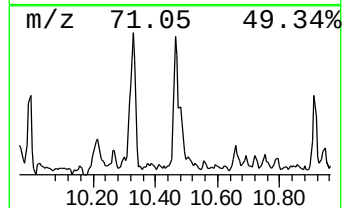
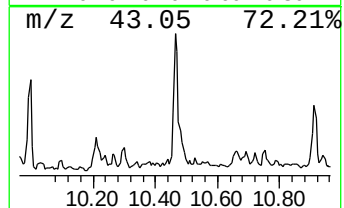
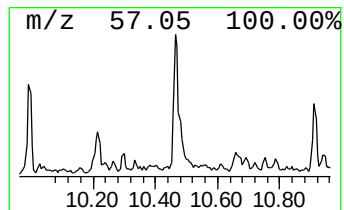
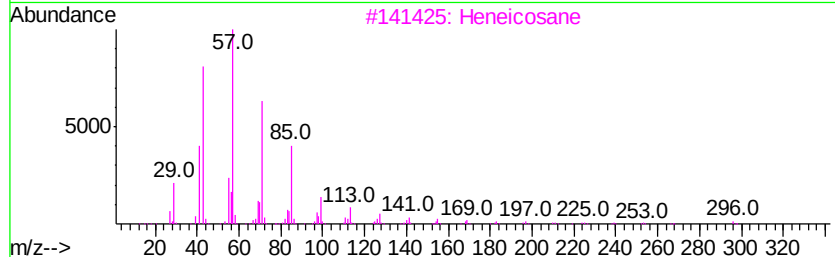
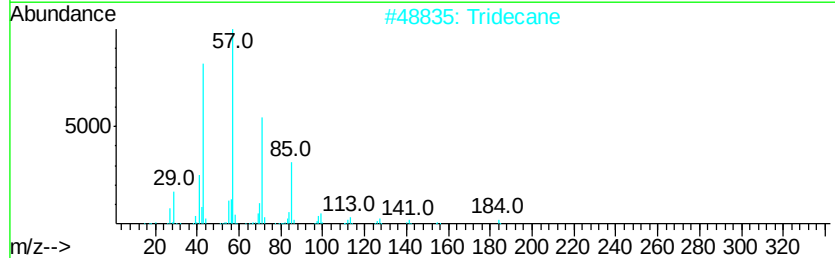
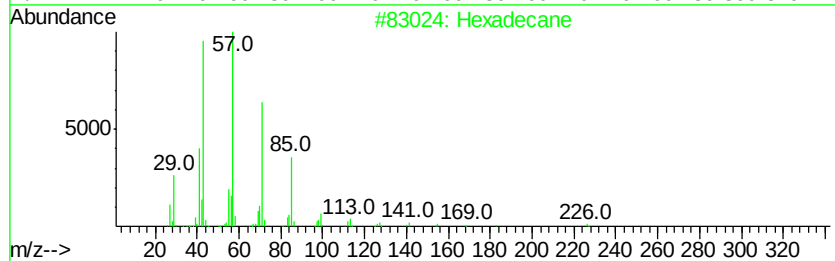
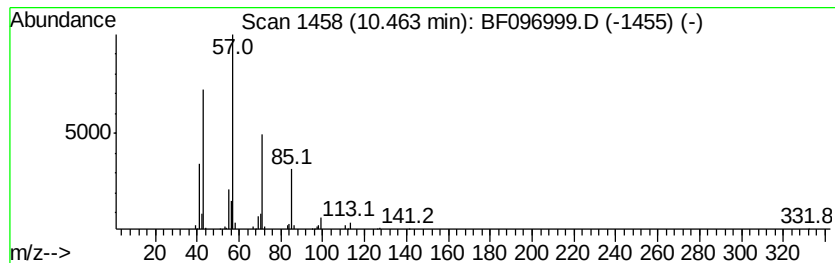
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Hexadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.46	2.01 ng	104329	Phenanthrene-d10	11.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	91
2		Tridecane	184	C13H28	000629-50-5	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Pentadecane	212	C15H32	000629-62-9	83
5		Tetradecane	198	C14H30	000629-59-4	83



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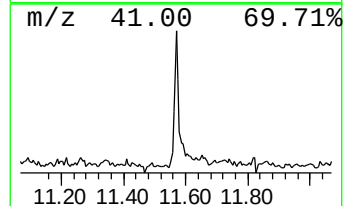
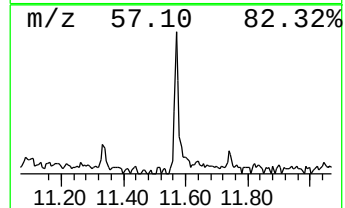
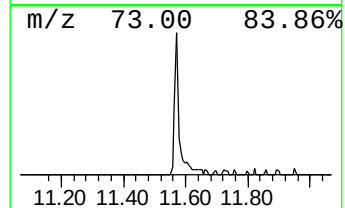
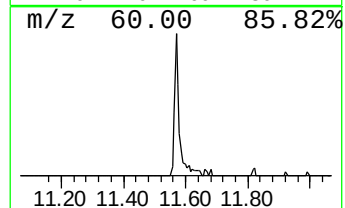
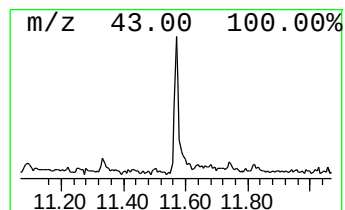
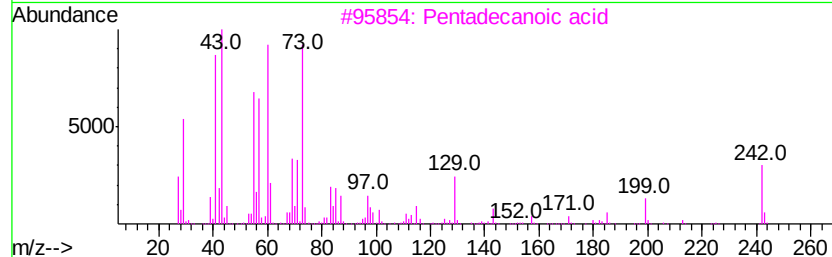
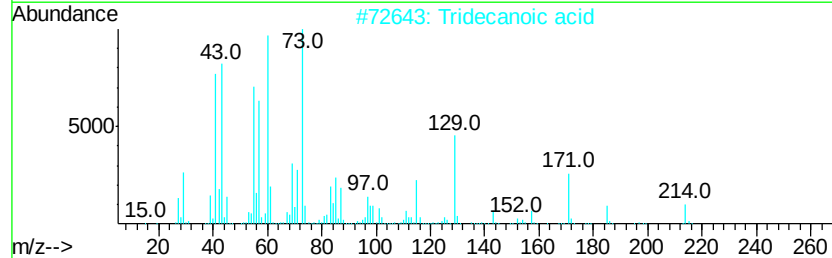
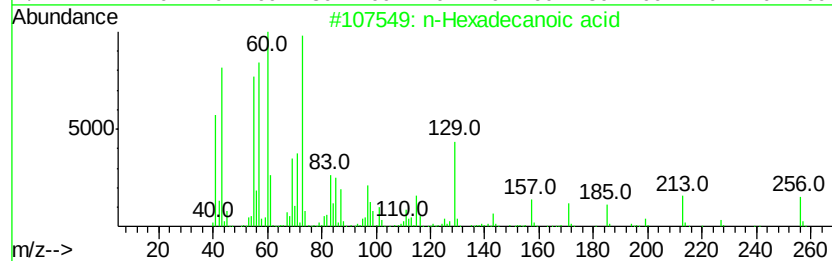
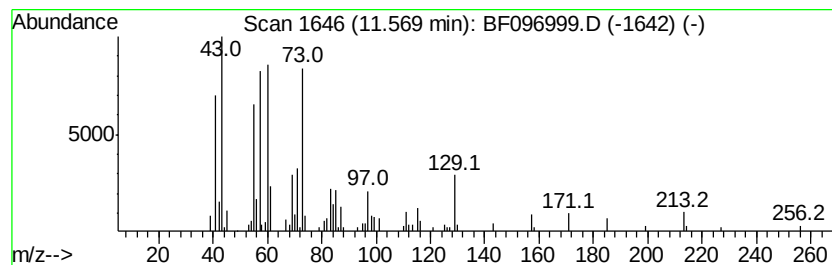
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.57	2.61 ng	135548	Phenanthrene-d10		11.02	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	90
2	Tridecanoic acid		214	C13H26O2	000638-53-9	89
3	Pentadecanoic acid		242	C15H30O2	001002-84-2	72
4	Tetradecanoic acid		228	C14H28O2	000544-63-8	58
5	Undecanoic acid		186	C11H22O2	000112-37-8	53



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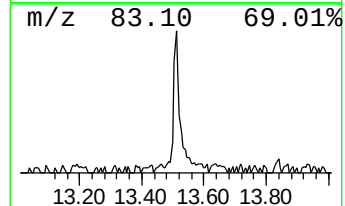
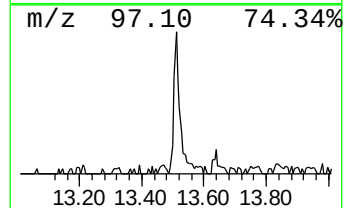
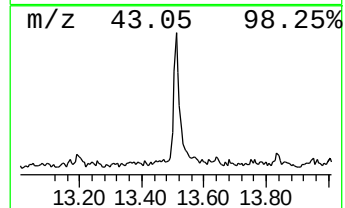
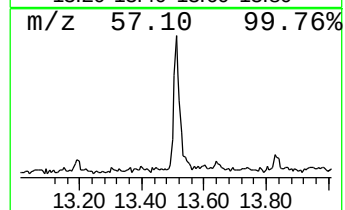
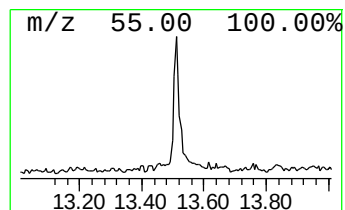
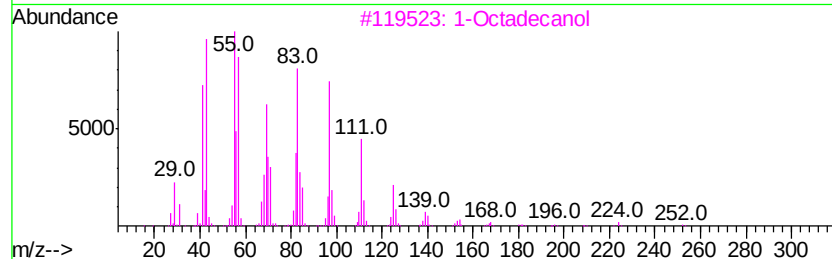
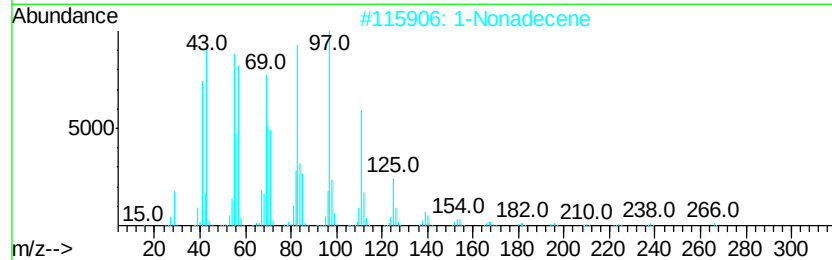
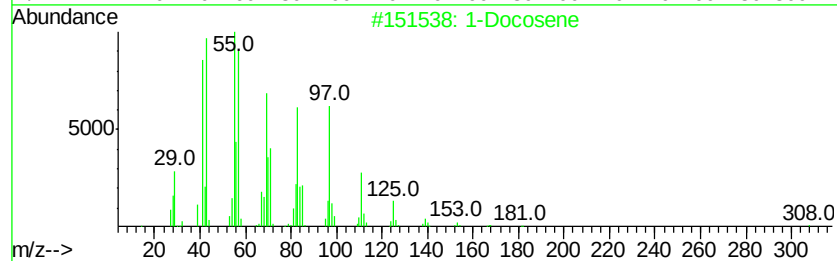
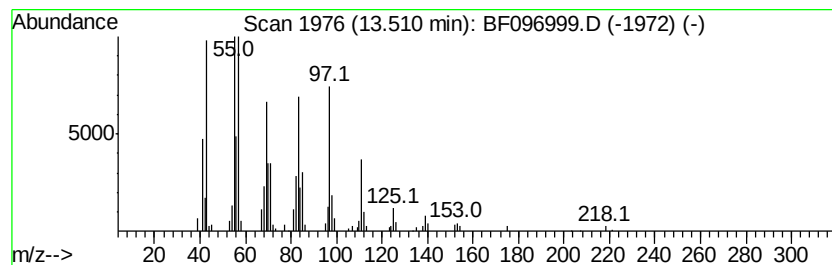
Instrument :
BNA_F
ClientSampleId :
RAMP-D-4-(0-2)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF071317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 1-Docosene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.51	4.15 ng	170725	Chrysene-d12		13.64
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	91
2	1-Nonadecene	266	C19H38	018435-45-5	91
3	1-Octadecanol	270	C18H38O	000112-92-5	87
4	1-Heneicosanol	312	C21H44O	015594-90-8	87
5	Trichloroacetic acid, tetradecyl...	358	C16H29Cl3O2	074339-52-9	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072417\
Data File : BF096999.D
Acq On : 24 Jul 2017 21:47
Operator : SJ/JU
Sample : I4361-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RAMP-D-4-(0-2)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1,3-Dioxolane, 2,...	2.35	4.3	ng	163875	1	6.51	762943	20.0
2-Pentanone, 4-hy...	4.67	12.2	ng	465938	1	6.51	762943	20.0
unknown6.29	6.29	87.0	ng	3317560	1	6.51	762943	20.0
Hexadecane	10.46	2.0	ng	104329	4	11.02	1039210	20.0
n-Hexadecanoic acid	11.57	2.6	ng	135548	4	11.02	1039210	20.0
1-Docosene	13.51	4.2	ng	170725	5	13.64	822055	20.0