

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041719\
 Data File : BM019951.D
 Acq On : 17 Apr 2019 22:55
 Operator : JU/SJ
 Sample : K2417-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 FK-PS-01-045

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:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM041519.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	5.140	360	366	391	rBV	3572099	5130405	92.07%	9.914%
2	6.187	538	544	552	rBV	205368	293449	5.27%	0.567%
3	6.722	628	635	661	rBV	3193145	5227228	93.81%	10.102%
4	6.940	666	672	677	rVB	66770	105201	1.89%	0.203%
5	7.063	686	693	706	rBV	3616056	5572240	100.00%	10.768%
6	7.528	765	772	781	rBV	627757	995964	17.87%	1.925%
7	7.839	818	825	841	rVB	2486766	3922860	70.40%	7.581%
8	8.698	964	971	997	rBV	1608945	3060136	54.92%	5.914%
9	10.304	1237	1244	1253	rBV	611165	1160796	20.83%	2.243%
10	10.369	1253	1255	1265	rVB5	36321	77035	1.38%	0.149%
11	12.798	1661	1668	1682	rBV	3252135	4977441	89.33%	9.619%
12	13.668	1810	1816	1829	rBV	346781	581411	10.43%	1.124%
13	14.192	1899	1905	1919	rBV2	817619	1333663	23.93%	2.577%
14	14.421	1936	1944	1950	rBV2	43727	62883	1.13%	0.122%
15	15.704	2155	2162	2179	rBV	2454948	3720491	66.77%	7.190%
16	16.956	2370	2375	2386	rBV2	831662	1319516	23.68%	2.550%
17	17.851	2522	2527	2537	rBV	278794	474264	8.51%	0.917%
18	18.562	2644	2648	2653	rVB	139692	182992	3.28%	0.354%
19	19.068	2724	2734	2740	rVB	396055	521840	9.36%	1.008%
20	19.145	2742	2747	2764	rBV	109707	256402	4.60%	0.495%
21	19.262	2764	2767	2775	rBV4	33920	56463	1.01%	0.109%
22	19.603	2819	2825	2830	rBV	4370782	5087841	91.31%	9.832%
23	20.062	2899	2903	2912	rVB6	81923	140277	2.52%	0.271%
24	20.239	2930	2933	2939	rVB4	120164	153745	2.76%	0.297%
25	20.339	2946	2950	2953	rBV	76660	94828	1.70%	0.183%
26	20.397	2957	2960	2965	rVB6	62423	94831	1.70%	0.183%
27	20.721	3011	3015	3021	rBV7	76732	117401	2.11%	0.227%
28	20.844	3032	3036	3037	rBV	312435	318401	5.71%	0.615%
29	20.868	3037	3040	3049	rVB	700312	937369	16.82%	1.811%
30	21.174	3086	3092	3098	rBV	1255037	1768116	31.73%	3.417%
31	21.392	3126	3129	3137	rBV2	104152	153052	2.75%	0.296%
32	21.721	3182	3185	3187	rBV	95807	107865	1.94%	0.208%
33	21.750	3187	3190	3202	rVB2	199580	322439	5.79%	0.623%
34	22.650	3339	3343	3349	rBV	264507	388730	6.98%	0.751%

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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

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35	23.362	3458	3464	3477	rVB	1086642	2008359	36.04%	3.881%
36	23.915	3552	3558	3568	rVB2	441550	1020679	18.32%	1.972%

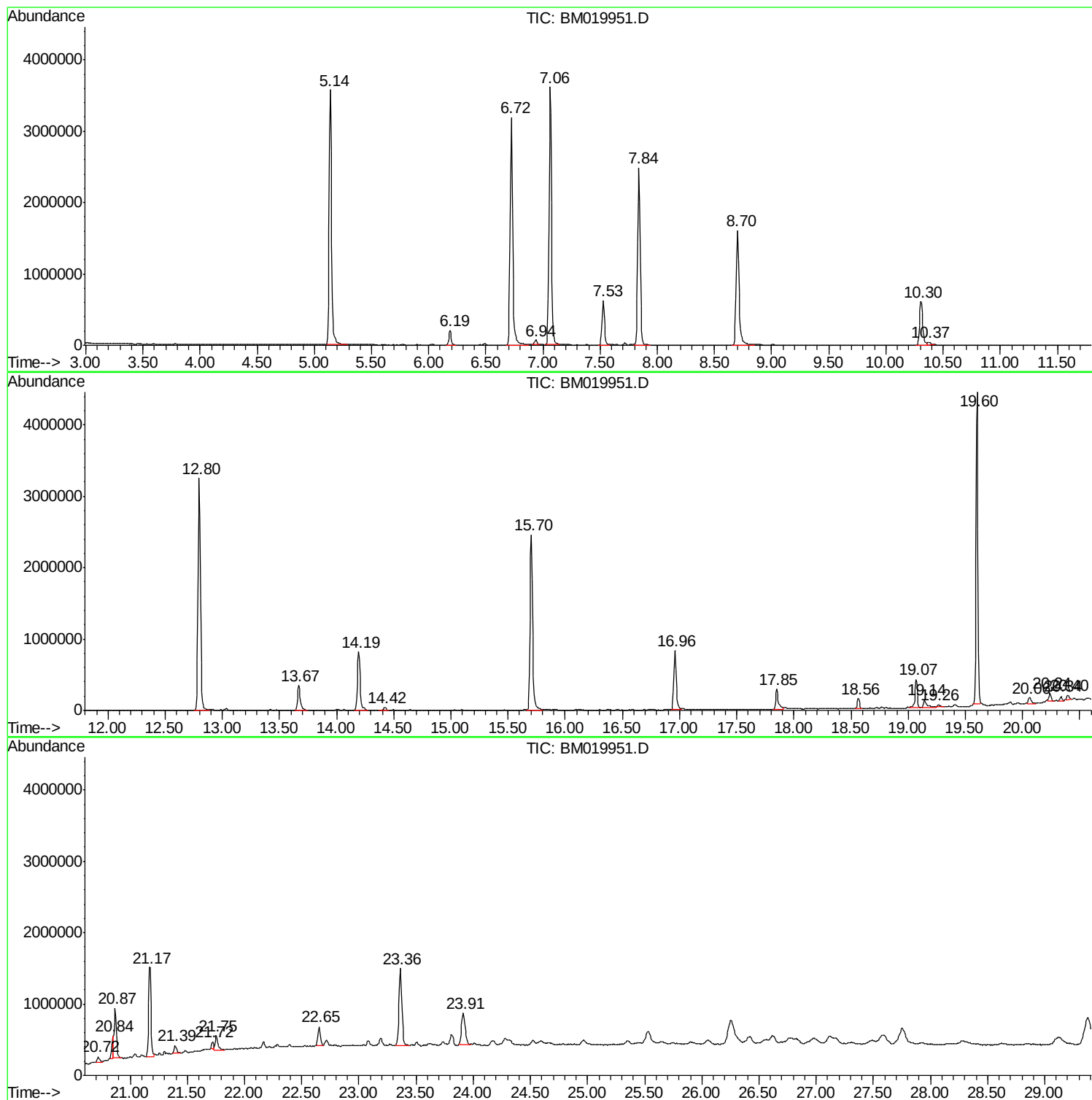
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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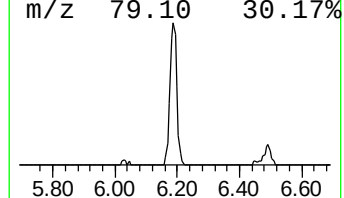
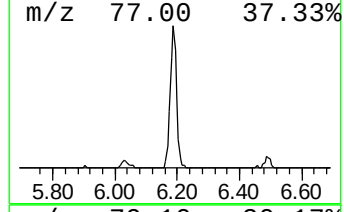
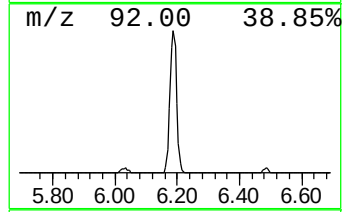
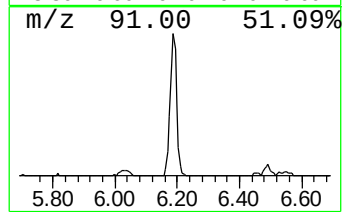
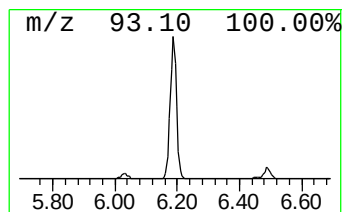
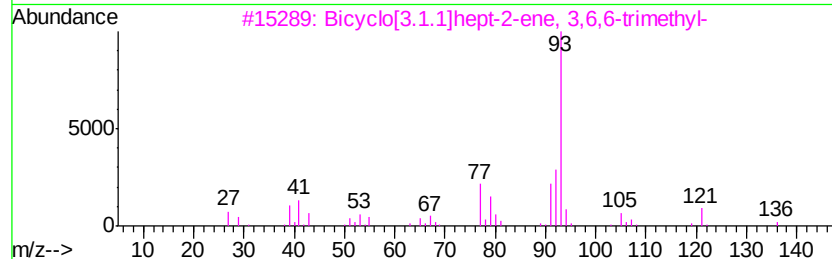
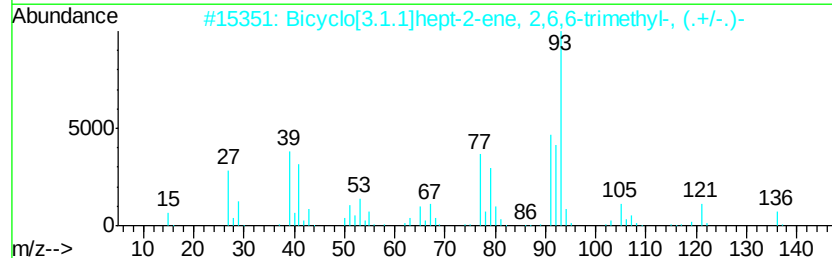
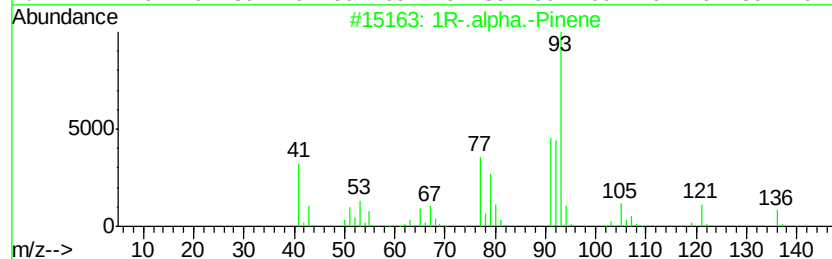
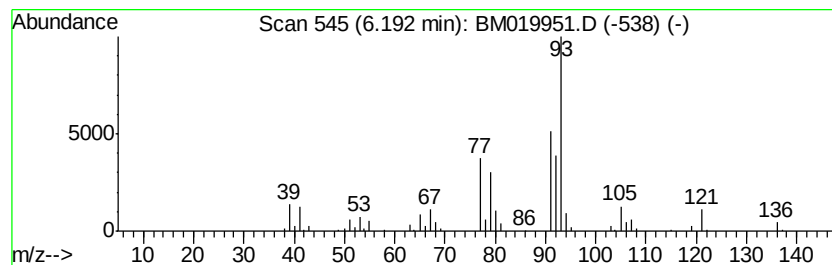
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 1R-.alpha.-Pinene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.19	5.89 ng	293449	1,4-Dichlorobenzene-d4	7.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1R-.alpha.-Pinene	136	C10H16	007785-70-8	95
2		Bicyclo[3.1.1]hept-2-ene, 2,6,6-...	136	C10H16	002437-95-8	95
3		Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	94
4		Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000508-32-7	91
5		.alpha.-Pinene	136	C10H16	000080-56-8	91



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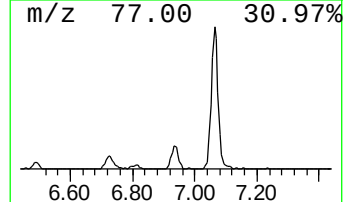
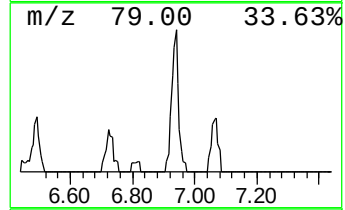
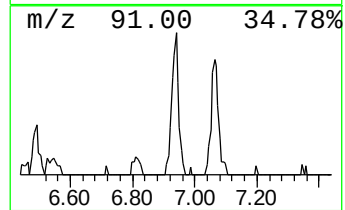
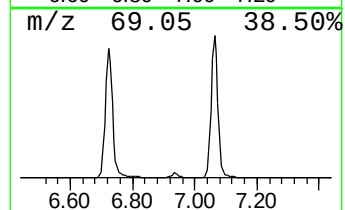
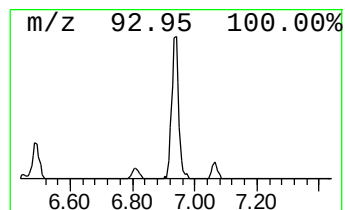
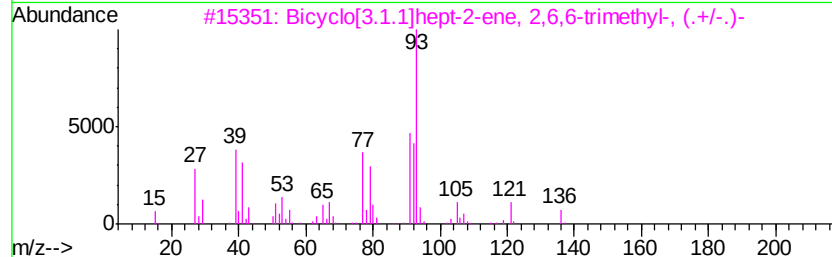
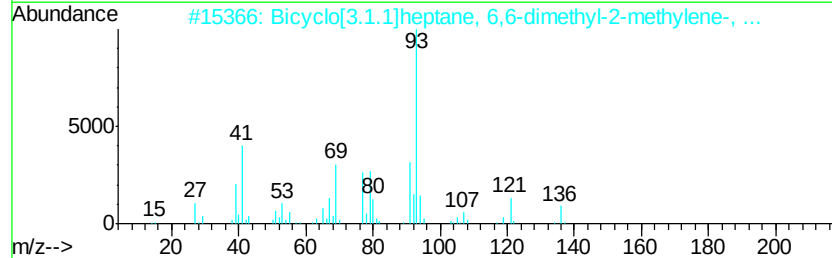
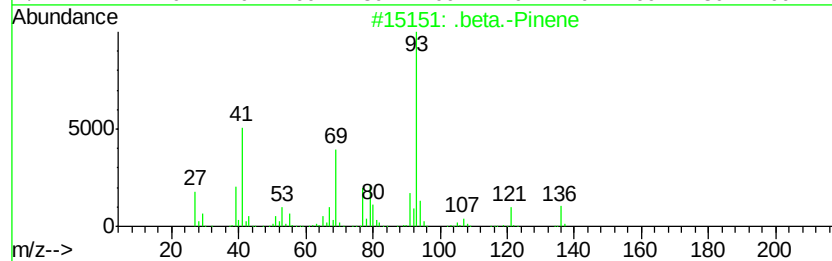
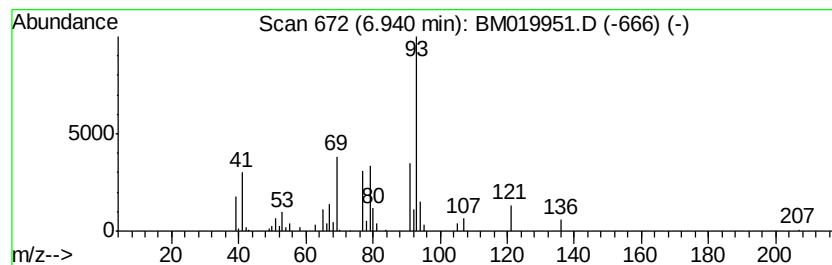
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 .beta.-Pinene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.94	2.11 ng	105201	1,4-Dichlorobenzene-d4	7.53

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.beta.-Pinene	136	C10H16	000127-91-3	90
2			Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	83
3			Bicyclo[3.1.1]hept-2-ene, 2,6,6-...	136	C10H16	002437-95-8	83
4			Cyclopentene, 3-isopropenyl-5,5-...	136	C10H16	1000162-25-4	83
5			Bicyclo[3.1.0]hexane, 4-methylen...	136	C10H16	003387-41-5	78



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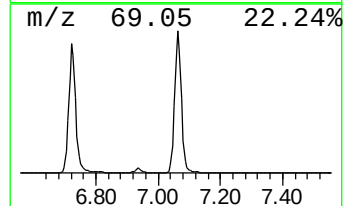
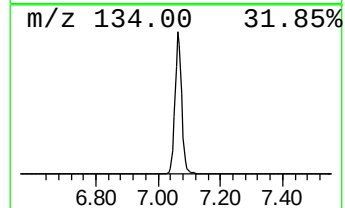
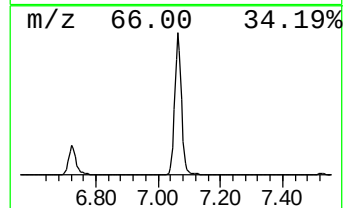
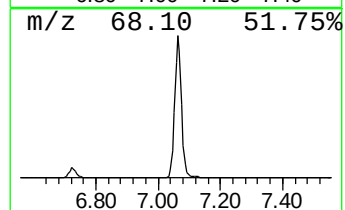
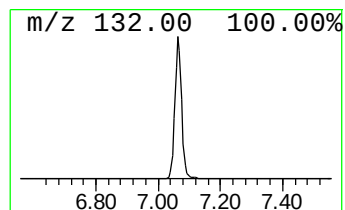
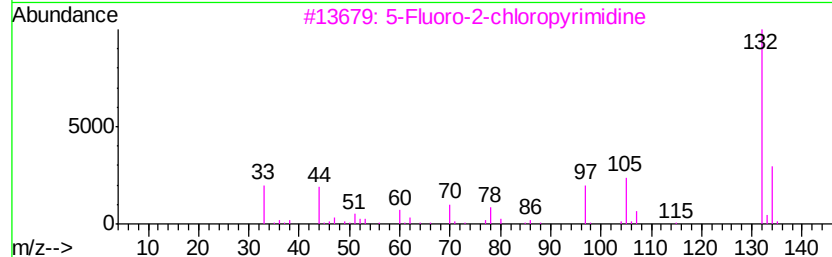
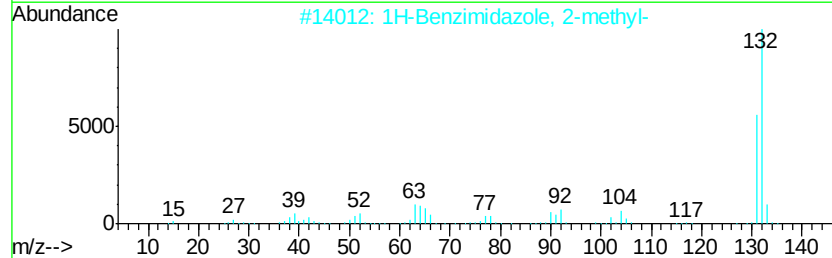
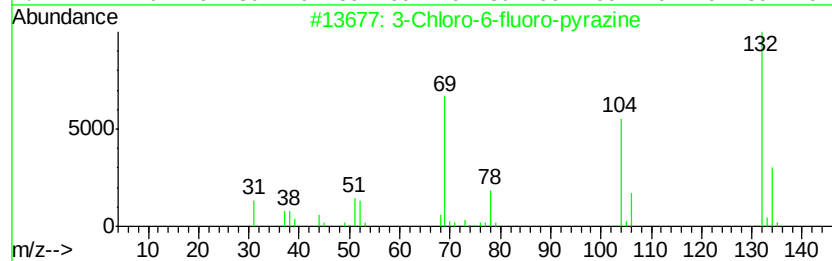
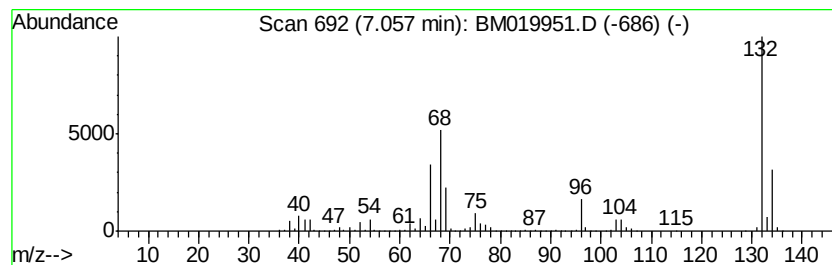
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown7.06 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.06	111.90 ng	5572240	1,4-Dichlorobenzene-d4	7.53

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	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	Xenon	132	Xe	007440-63-3	9



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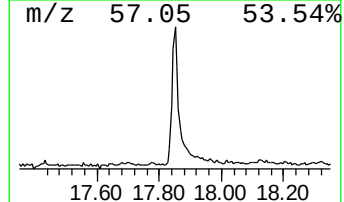
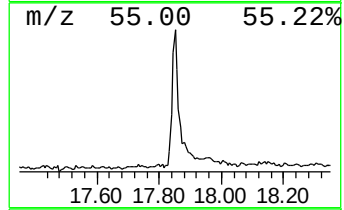
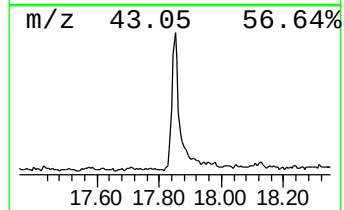
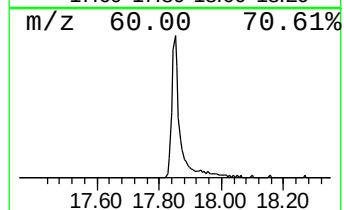
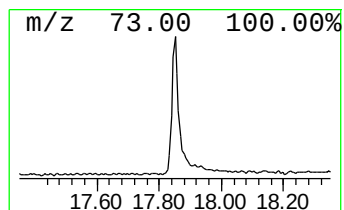
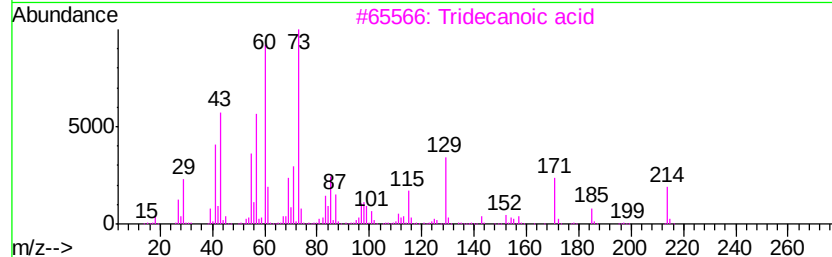
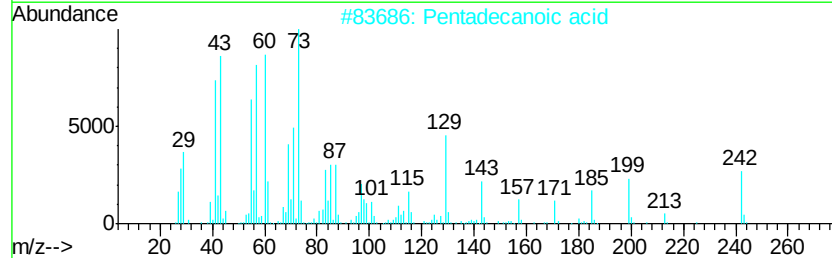
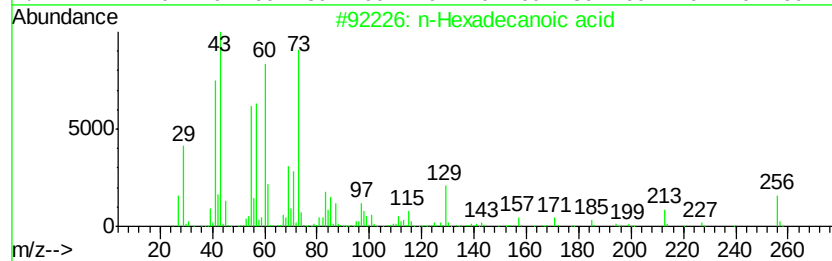
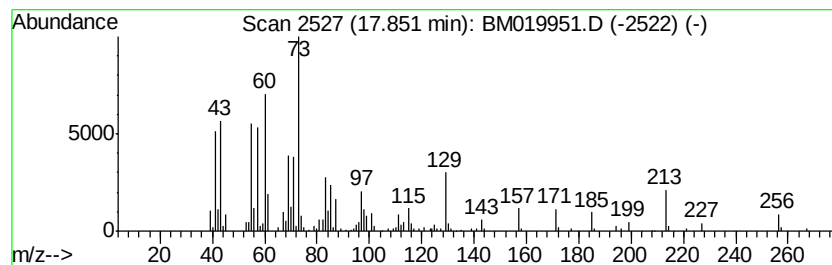
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.
17.85	7.19 ng	474264	Phenanthrene-d10		16.96
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	64
3	Tridecanoic acid	214	C13H26O2	000638-53-9	52
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	52
5	Dodecanoic acid	200	C12H24O2	000143-07-7	50



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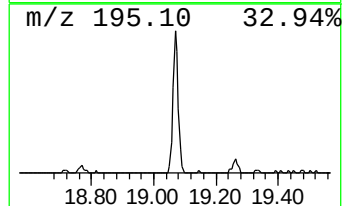
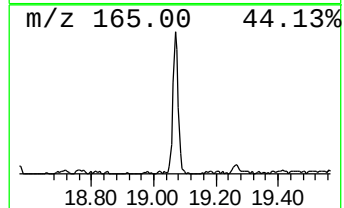
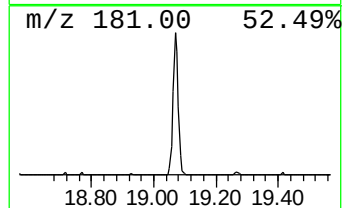
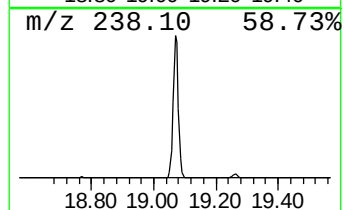
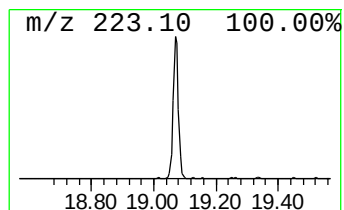
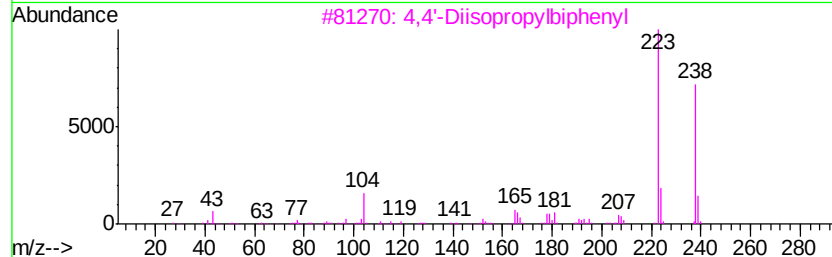
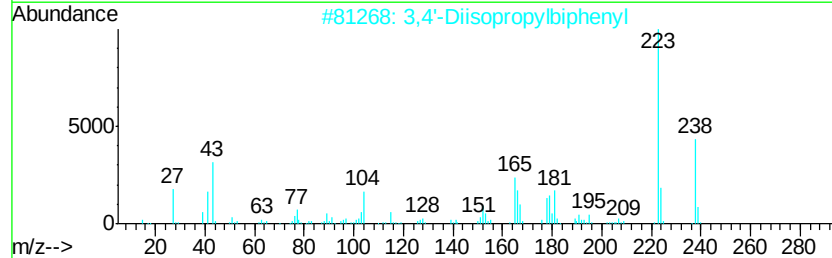
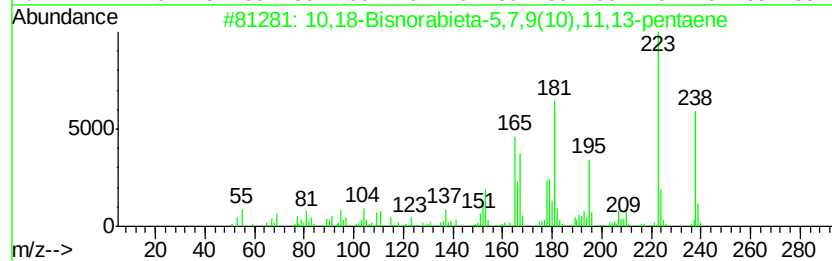
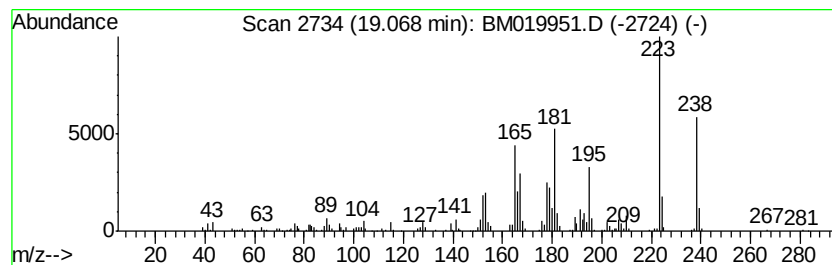
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.07	5.90 ng	521840	Chrysene-d12	21.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99
2		3,4'-Diisopropylbiphenyl	238	C18H22	061434-46-6	58
3		4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	50
4		Anthracene, 9,10-dimethoxy-	238	C16H14O2	002395-97-3	46
5		9-Acridinecarboxylic acid	223	C14H9NO2	005336-90-3	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041719\
Data File : BM019951.D
Acq On : 17 Apr 2019 22:55
Operator : JU/SJ
Sample : K2417-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

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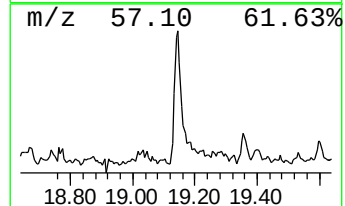
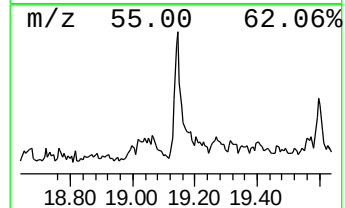
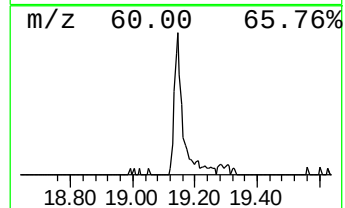
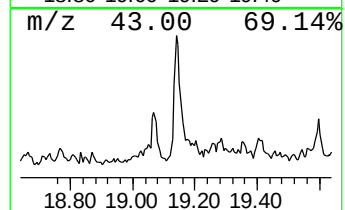
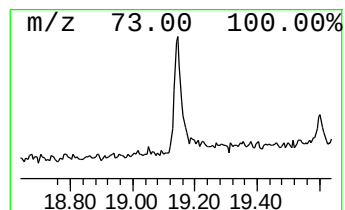
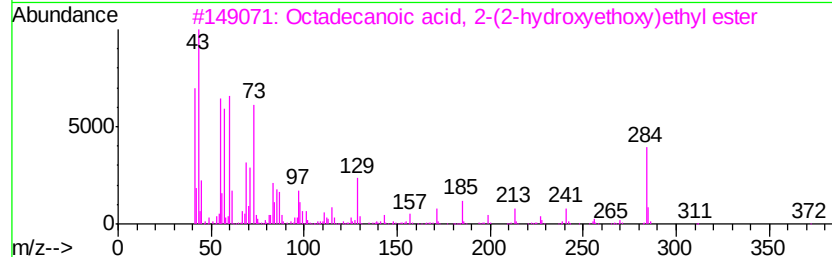
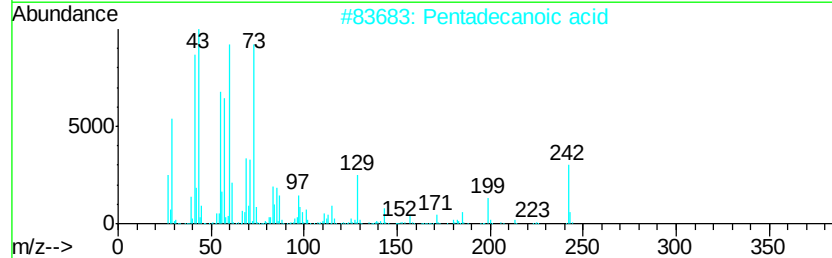
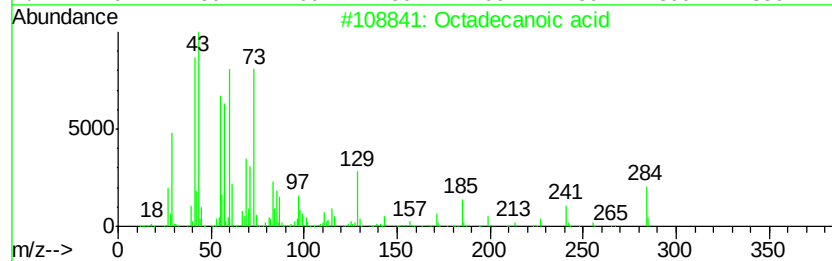
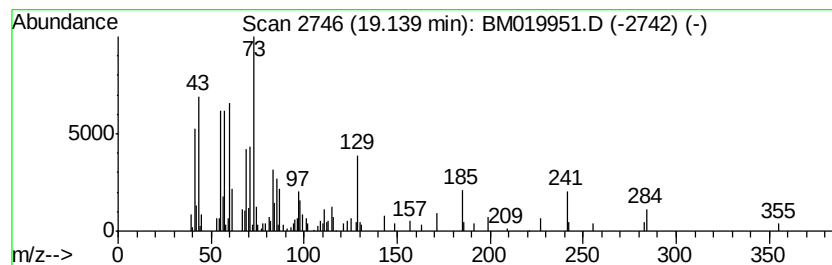
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.14	2.90 ng	256402	Chrysene-d12	21.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	95
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3		Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	87
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	87
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	56



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041719\
Data File : BM019951.D
Acq On : 17 Apr 2019 22:55
Operator : JU/SJ
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ClientSampleId :
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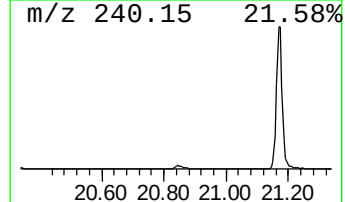
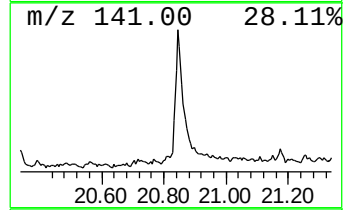
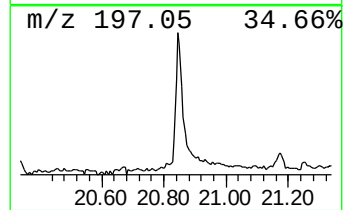
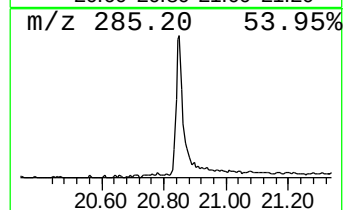
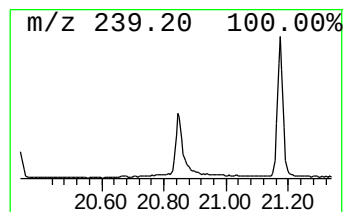
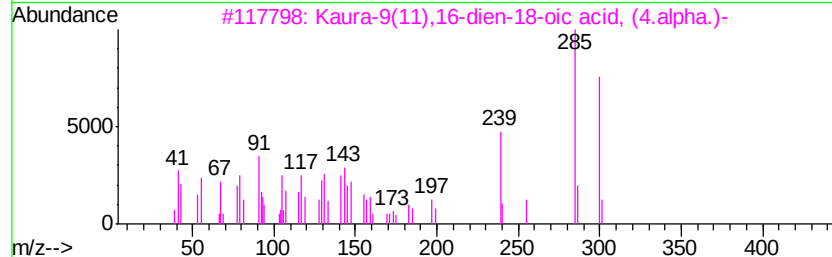
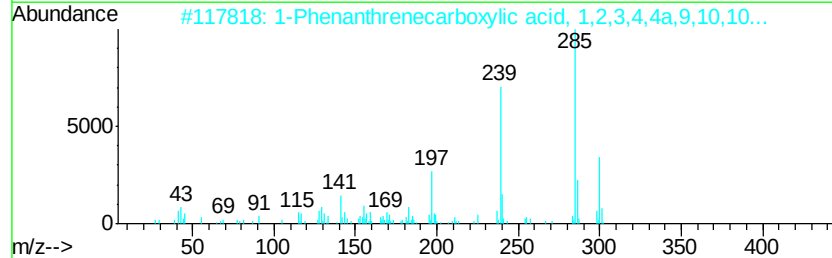
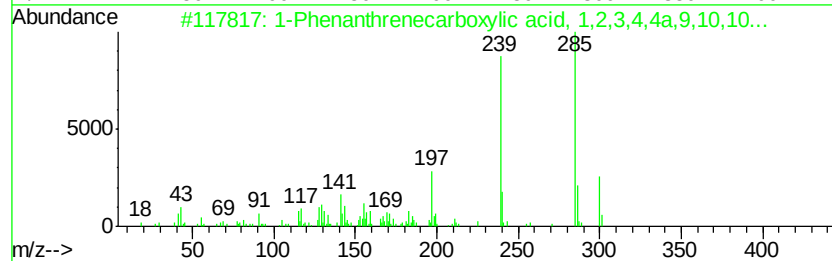
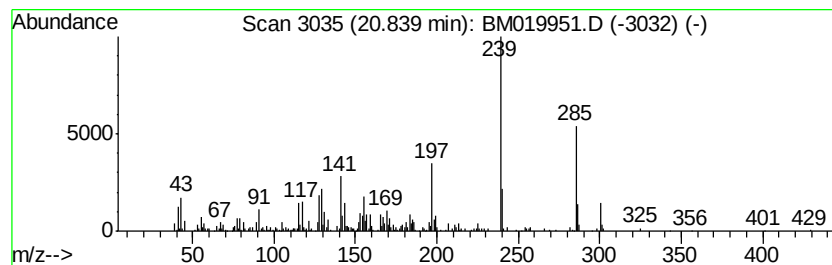
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Phenanthrenecarboxylic ac... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.84	3.60 ng	318401	Chrysene-d12	21.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenanthrenecarboxylic acid, 1...	300	C20H28O2	001740-19-8	93
2		1-Phenanthrenecarboxylic acid, 1...	300	C20H28O2	005155-70-4	93
3		Kaura-9(11),16-dien-18-oic acid,...	300	C20H28O2	022338-67-6	72
4		1-Phenanthrenecarboxylic acid, 1...	314	C21H30O2	001235-74-1	49
5		Indole, 3-imidazolo[2,1-b]thiazo...	239	C13H9N3S	292855-05-1	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041719\
Data File : BM019951.D
Acq On : 17 Apr 2019 22:55
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Sample : K2417-01
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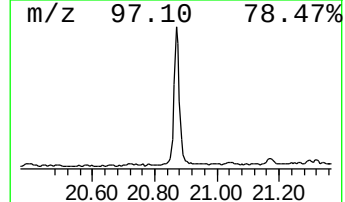
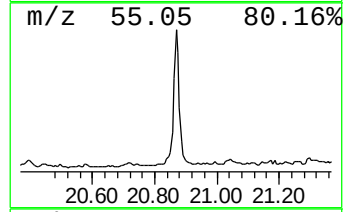
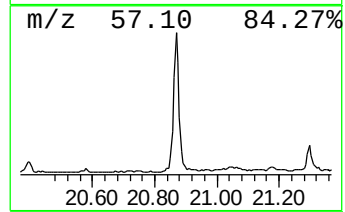
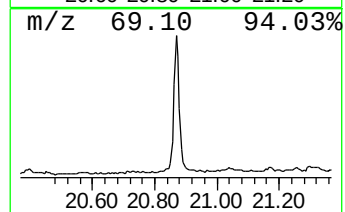
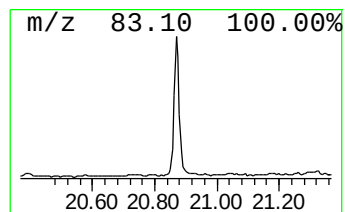
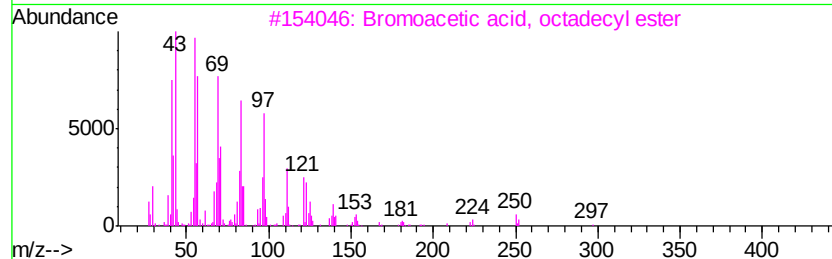
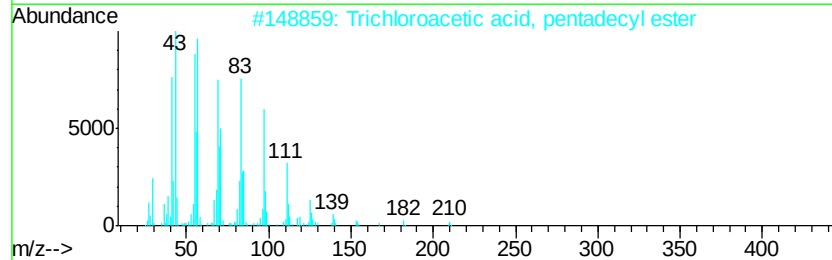
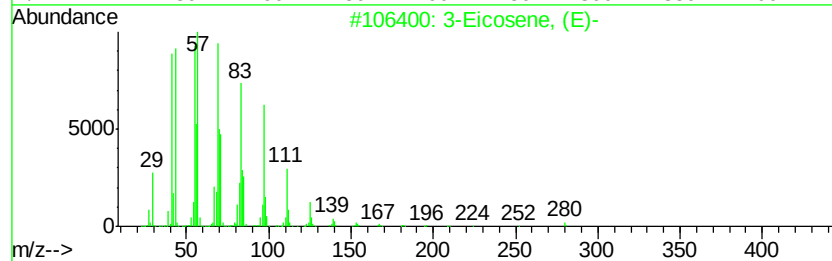
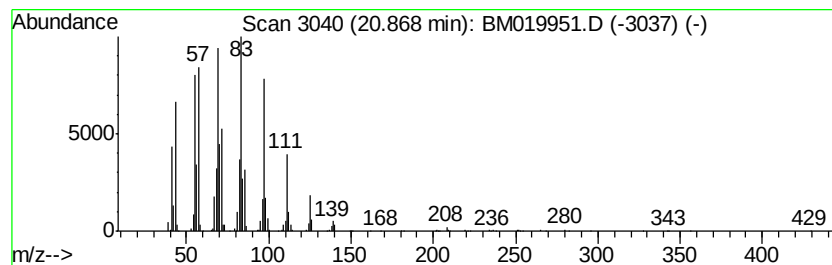
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 3-Eicosene, (E)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.87	10.60 ng	937369	Chrysene-d12	21.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Eicosene, (E)-	280	C20H40	074685-33-9	91
2		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91
3		Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	91
4		1-Heptadecene	238	C17H34	006765-39-5	90
5		1-Hexadecanol	242	C16H34O	036653-82-4	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041719\
Data File : BM019951.D
Acq On : 17 Apr 2019 22:55
Operator : JU/SJ
Sample : K2417-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

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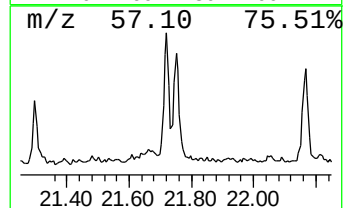
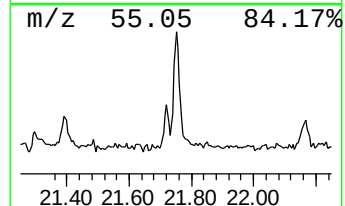
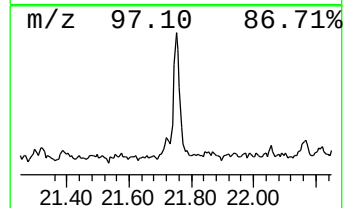
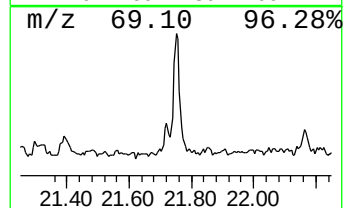
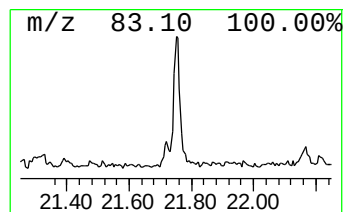
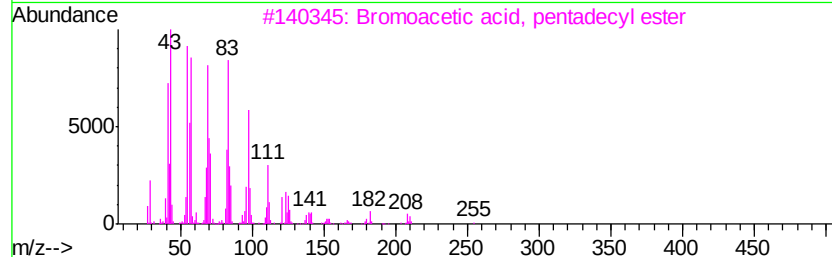
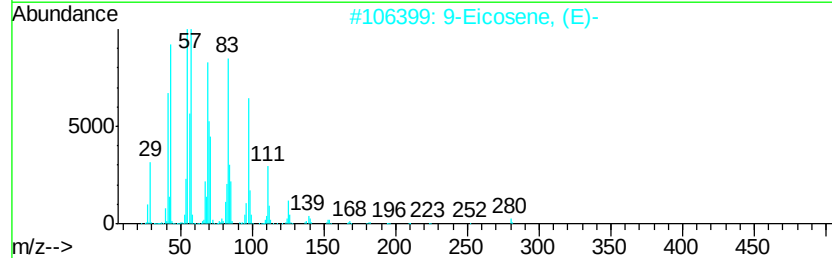
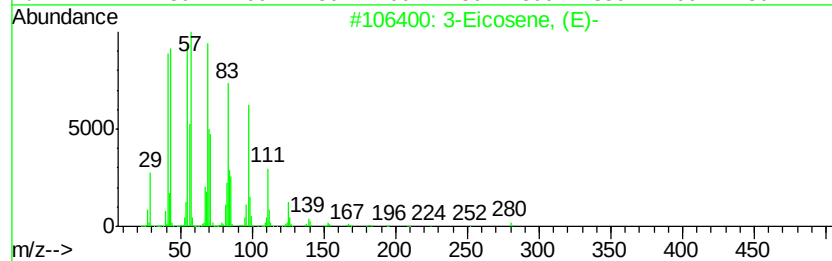
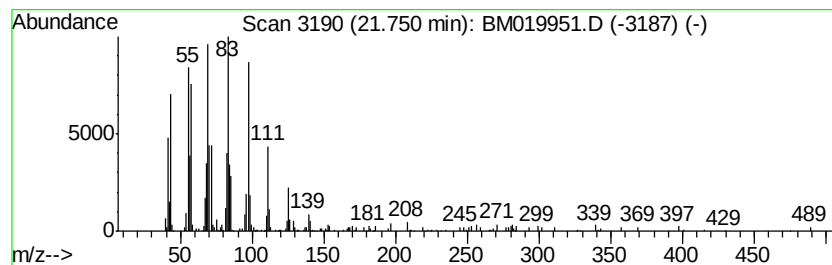
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 9-Eicosene, (E)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.75	3.65 ng	322439	Chrysene-d12	21.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Eicosene, (E)-	280	C20H40	074685-33-9	96
2		9-Eicosene, (E)-	280	C20H40	074685-29-3	94
3		Bromoacetic acid, pentadecyl ester	348	C17H33BrO2	131143-01-6	91
4		3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	91
5		1-Eicosanol	298	C20H42O	000629-96-9	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041719\
Data File : BM019951.D
Acq On : 17 Apr 2019 22:55
Operator : JU/SJ
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Misc :
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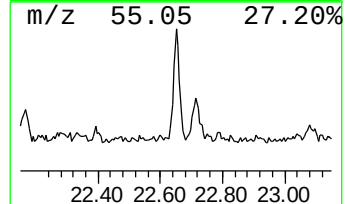
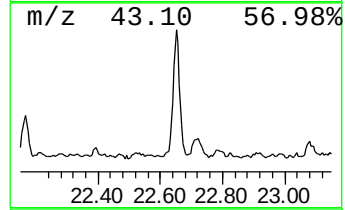
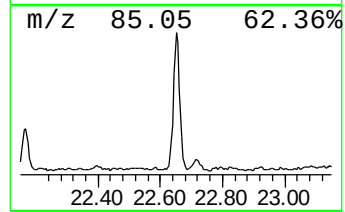
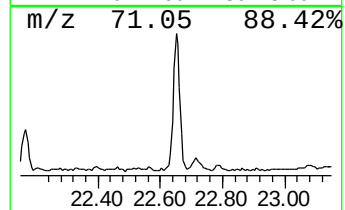
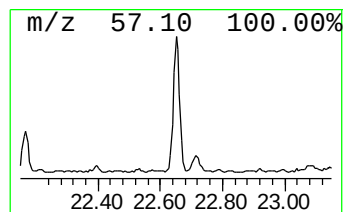
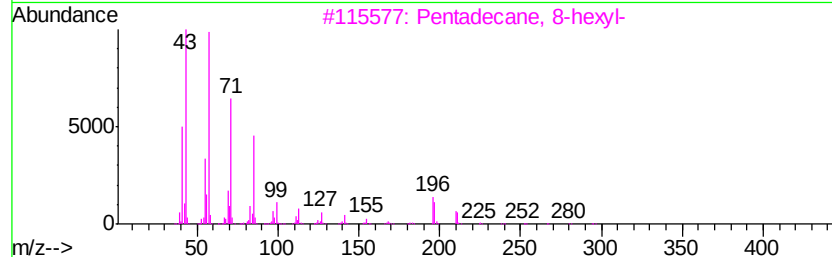
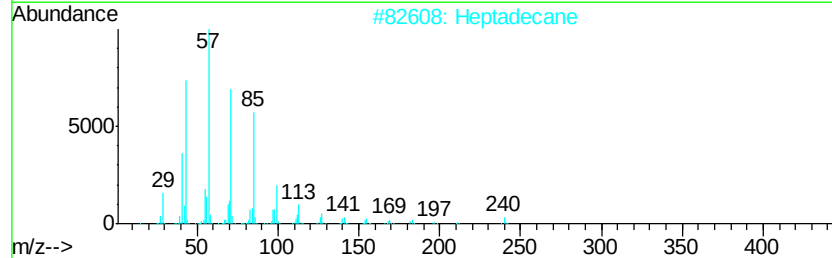
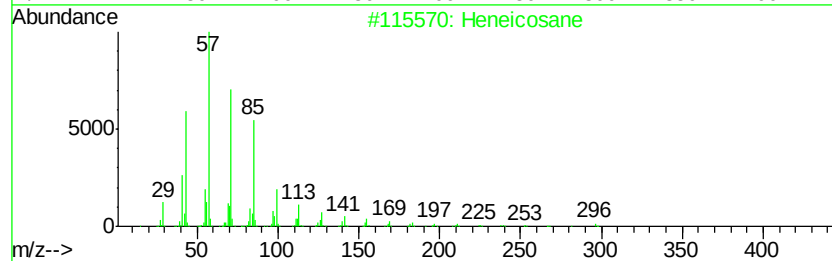
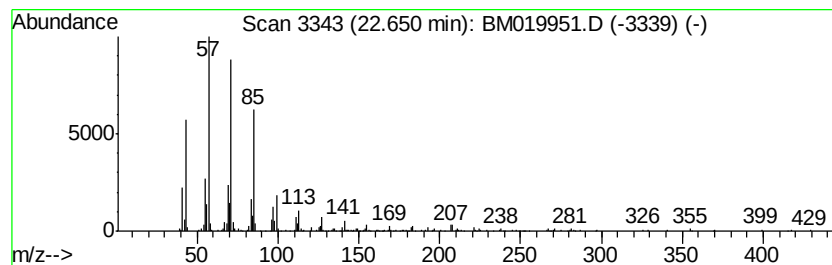
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.65	3.87 ng	388730	Perylene-d12	23.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Heptadecane	240	C17H36	000629-78-7	91
3		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	90
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
5		Tridecanol, 2-ethyl-2-methyl-	242	C16H34O	1000115-66-1	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041719\
Data File : BM019951.D
Acq On : 17 Apr 2019 22:55
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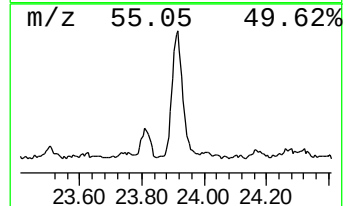
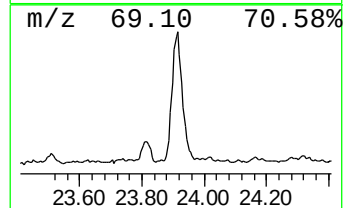
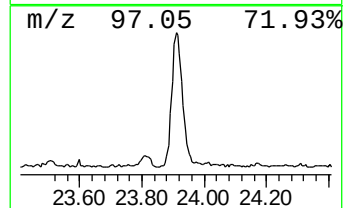
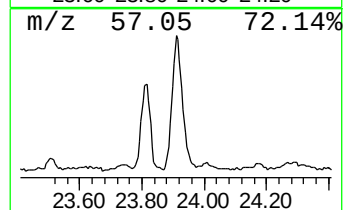
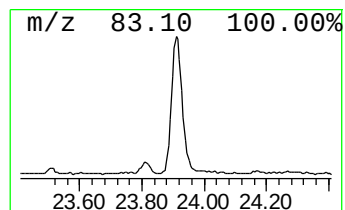
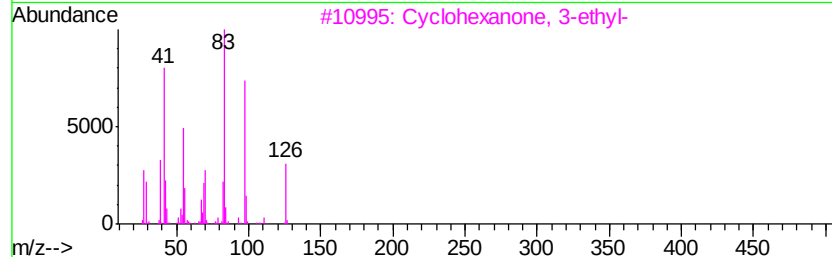
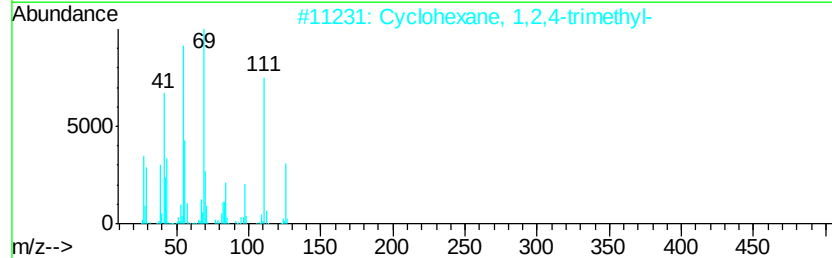
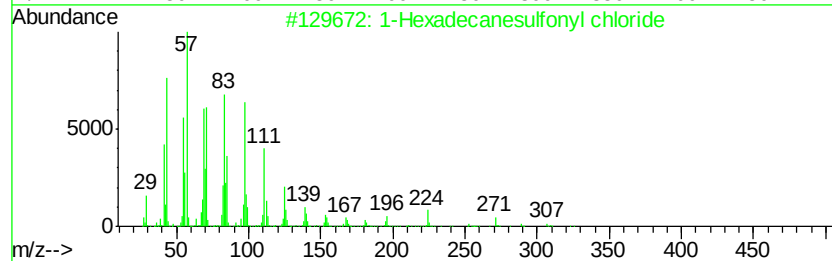
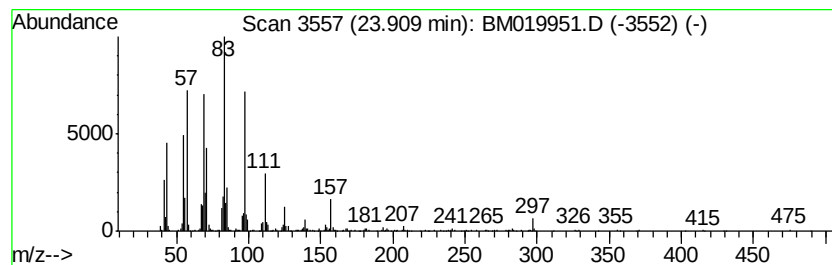
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown23.91 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.91	10.16 ng	1020680	Perylene-d12	23.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	38
2		Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	30
3		Cyclohexanone, 3-ethyl-	126	C8H14O	022461-89-8	30
4		Cyclohexane, 1-ethyl-2-propyl-	154	C11H22	062238-33-9	25
5		Octadecane, 1-(ethenyloxy)-	296	C20H40O	000930-02-9	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041719\
 Data File : BM019951.D
 Acq On : 17 Apr 2019 22:55
 Operator : JU/SJ
 Sample : K2417-01
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 ALS Vial : 19 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM041519.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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.beta.-Pinene	6.94	2.1	ng	105201	1	7.53	995964	20.0
unknown7.06	7.06	111.9	ng	5572240	1	7.53	995964	20.0
n-Hexadecanoic acid	17.85	7.2	ng	474264	4	16.96	1319520	20.0
4b,8-Dimethyl-2-i...	18.56	2.8	ng	182992	4	16.96	1319520	20.0
10,18-Bisnorabiet...	19.07	5.9	ng	521840	5	21.17	1768120	20.0
Octadecanoic acid	19.14	2.9	ng	256402	5	21.17	1768120	20.0
1-Phenanthrenecar...	20.84	3.6	ng	318401	5	21.17	1768120	20.0
3-Eicosene, (E)-	20.87	10.6	ng	937369	5	21.17	1768120	20.0
9-Eicosene, (E)-	21.75	3.6	ng	322439	5	21.17	1768120	20.0
Heneicosane	22.65	3.9	ng	388730	6	23.36	2008360	20.0
unknown23.91	23.91	10.2	ng	1020680	6	23.36	2008360	20.0