

Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
 Data File : BG022251.D
 Acq On : 9 Jun 2016 2:46
 Operator : UM/SJ
 Sample : H3381-10
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 W-53

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_G\METHODS\8270-BG060616.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	5.188	393	400	410	rVB	14281	28176	1.09%	0.213%
2	5.663	474	481	492	rBV	210364	398147	15.45%	3.006%
3	6.586	631	638	648	rVB2	16561	33623	1.30%	0.254%
4	7.279	745	756	769	rBV	143257	293084	11.37%	2.213%
5	7.649	811	819	835	rBV	451196	910445	35.33%	6.875%
6	8.008	870	880	888	rVB2	18963	43988	1.71%	0.332%
7	8.119	888	899	908	rBV	91707	182202	7.07%	1.376%
8	8.443	942	954	970	rBV	440644	927198	35.98%	7.001%
9	8.754	1000	1007	1011	rBV3	16125	31784	1.23%	0.240%
10	8.807	1011	1016	1024	rVB2	24017	48527	1.88%	0.366%
11	9.224	1079	1087	1091	rBV2	57100	111603	4.33%	0.843%
12	9.289	1091	1098	1110	rVB	337858	718881	27.90%	5.428%
13	9.571	1141	1146	1156	rVB	14083	29988	1.16%	0.226%
14	9.747	1170	1176	1183	rVB	60504	111991	4.35%	0.846%
15	10.111	1232	1238	1244	rBV3	21893	51411	1.99%	0.388%
16	10.323	1262	1274	1282	rBV2	151703	303169	11.76%	2.289%
17	10.934	1366	1378	1391	rBV	150663	422546	16.40%	3.191%
18	11.039	1391	1396	1403	rVB	35457	69138	2.68%	0.522%
19	11.398	1451	1457	1465	rVB4	20722	42482	1.65%	0.321%
20	11.509	1469	1476	1481	rBV3	20185	44303	1.72%	0.335%
21	11.833	1524	1531	1540	rVB3	25212	51974	2.02%	0.392%
22	12.032	1553	1565	1573	rBV	19656	41366	1.61%	0.312%
23	12.115	1573	1579	1585	rBV4	20787	41784	1.62%	0.316%
24	12.467	1630	1639	1646	rVV3	53874	105805	4.11%	0.799%
25	12.802	1685	1696	1704	rBV2	59594	149441	5.80%	1.128%
26	12.867	1704	1707	1719	rVB9	19345	49134	1.91%	0.371%
27	13.366	1784	1792	1799	rBV	1179015	2028487	78.71%	15.317%
28	14.071	1901	1912	1917	rBV3	28380	69350	2.69%	0.524%
29	14.747	2017	2027	2035	rBV2	262137	491344	19.07%	3.710%
30	16.233	2273	2280	2291	rBV	802486	1322739	51.33%	9.988%
31	17.491	2487	2494	2506	rVB2	292505	510929	19.83%	3.858%
32	20.082	2928	2935	2948	rBV	1765865	2577040	100.00%	19.459%
33	21.762	3213	3221	3230	rBV	264114	518060	20.10%	3.912%
34	25.058	3771	3782	3797	rVB2	145822	483553	18.76%	3.651%

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ClientSampleId :
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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

Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

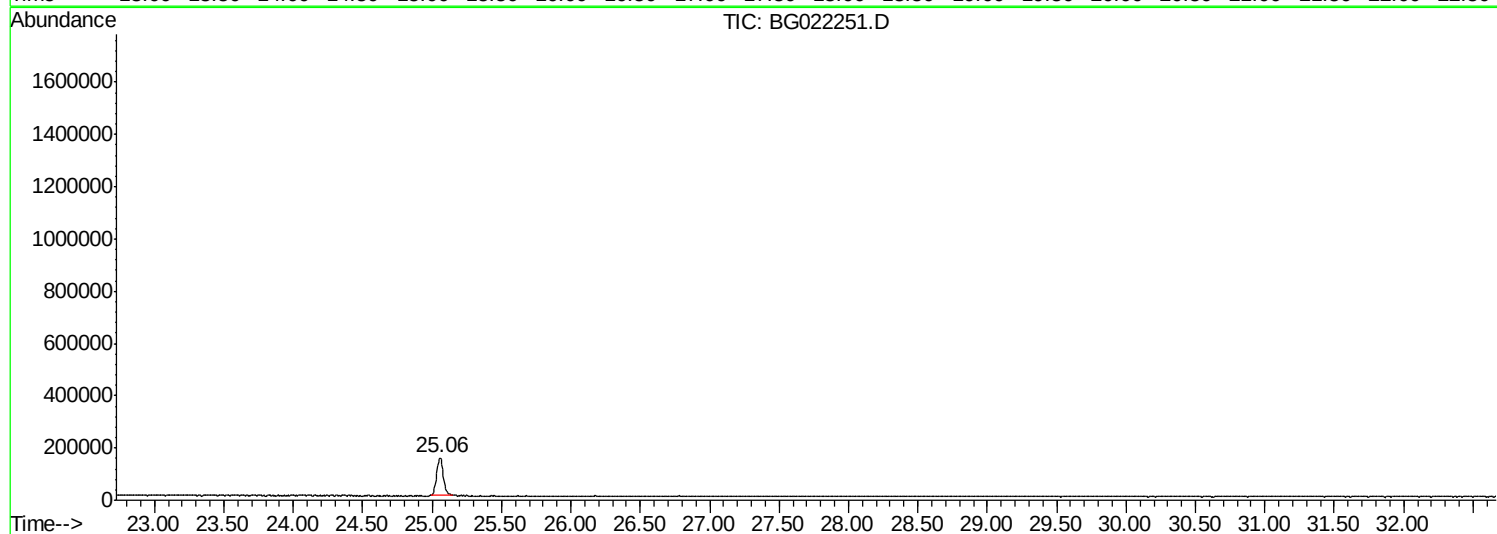
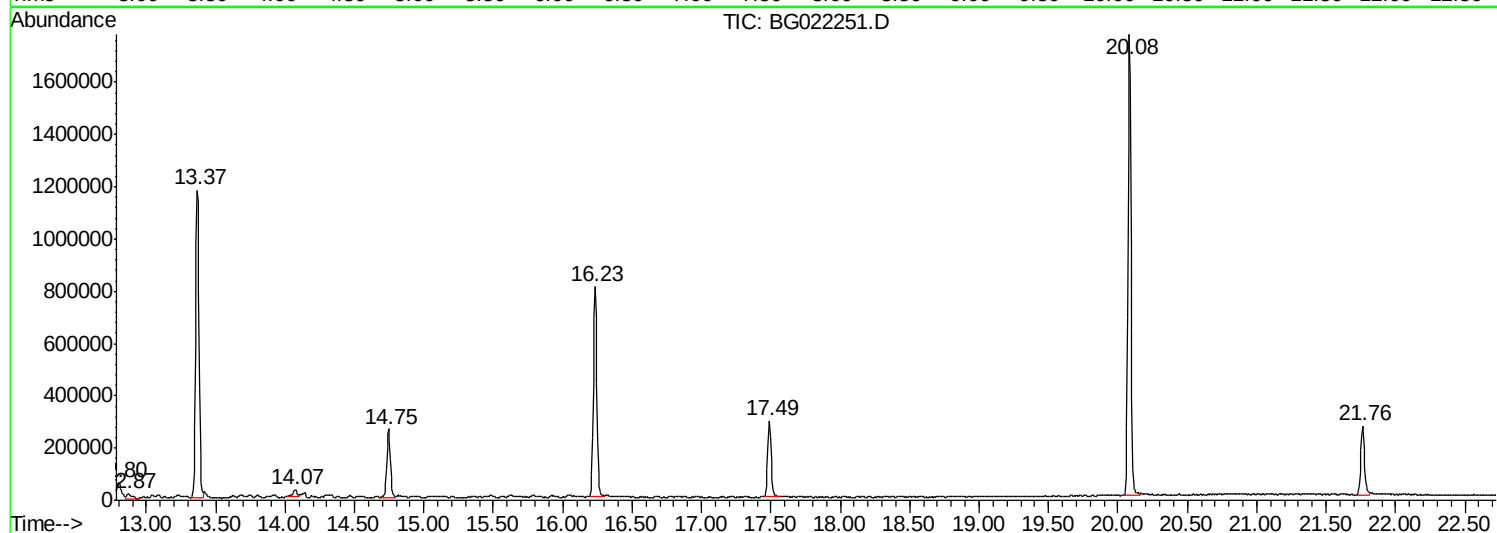
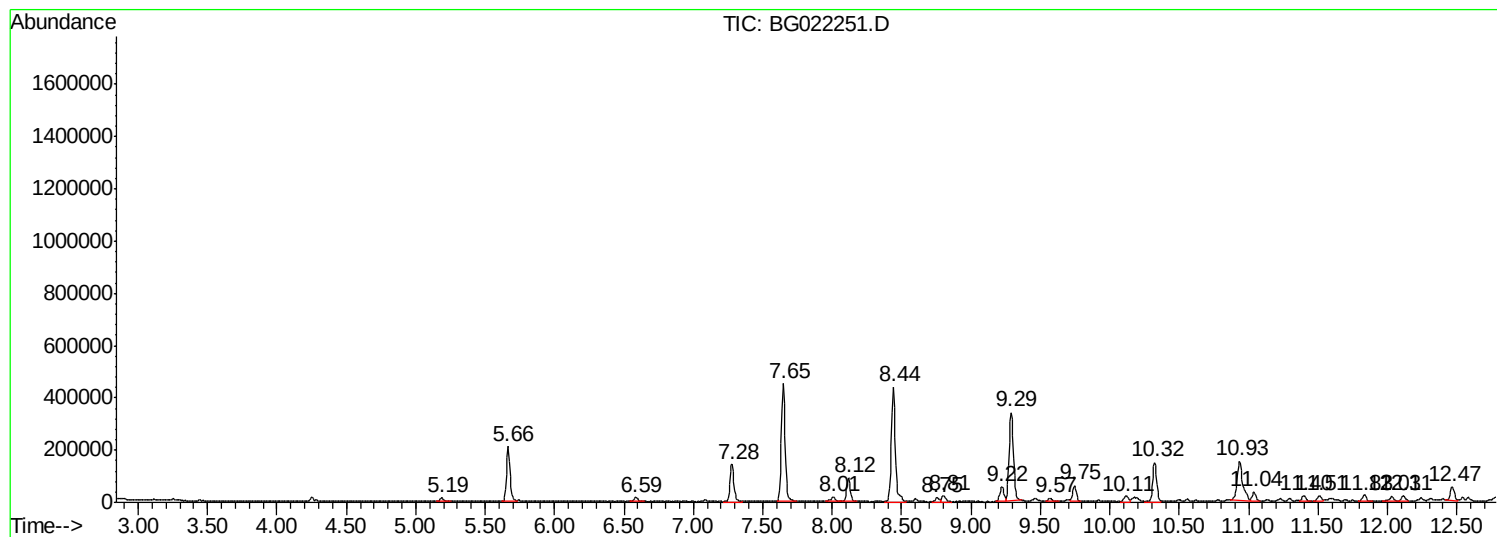
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Quant Title :

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



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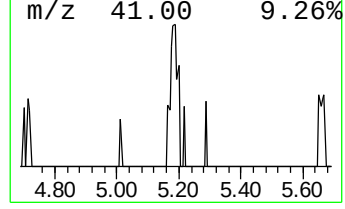
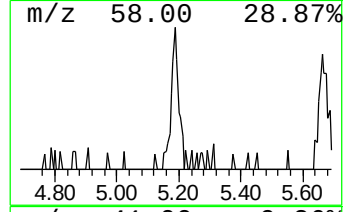
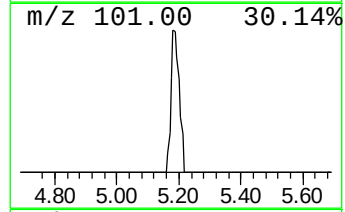
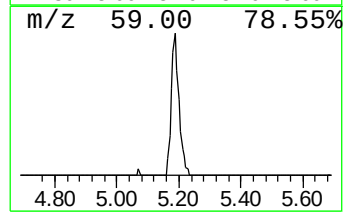
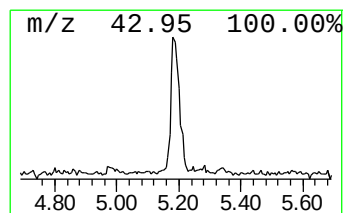
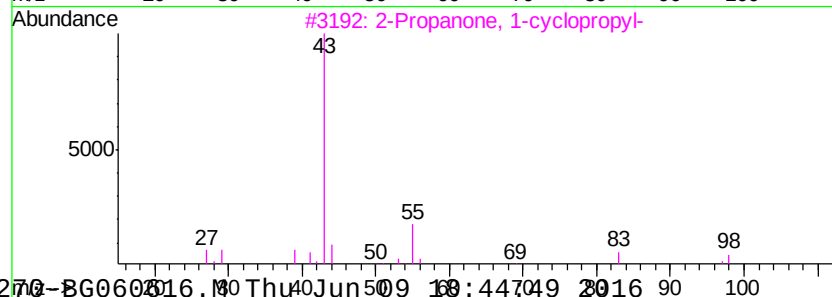
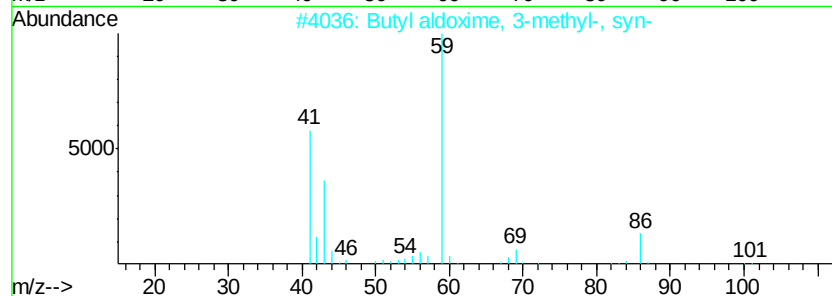
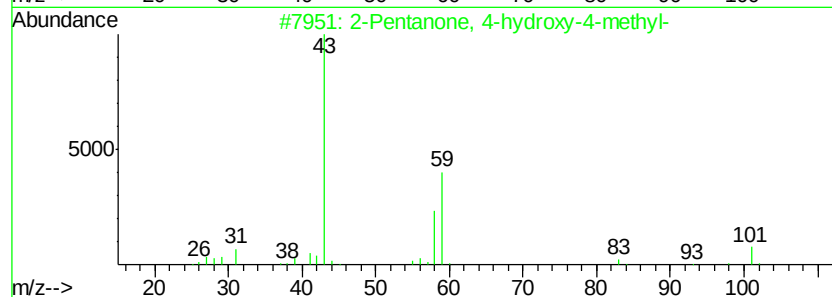
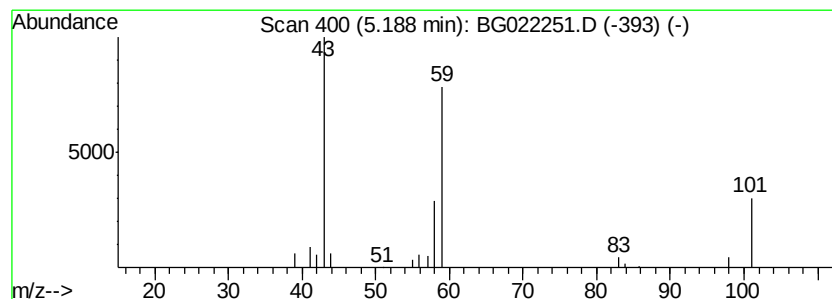
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Quant Title :

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.19	3.09 ng	28176	1,4-Dichlorobenzene-d4	8.12	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25
3	2-Propanone, 1-cyclopropyl-	98	C6H10O	004160-75-2	9
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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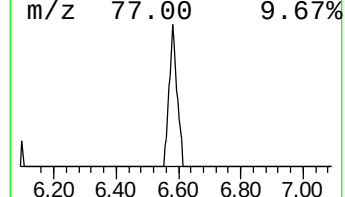
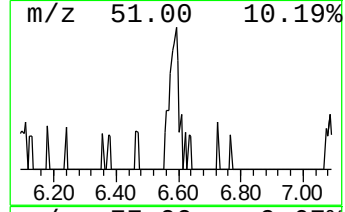
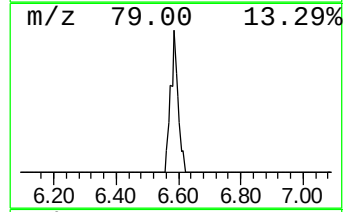
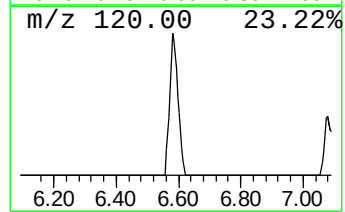
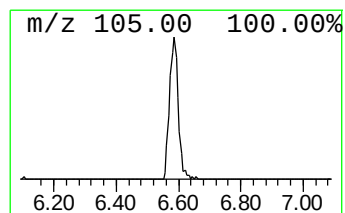
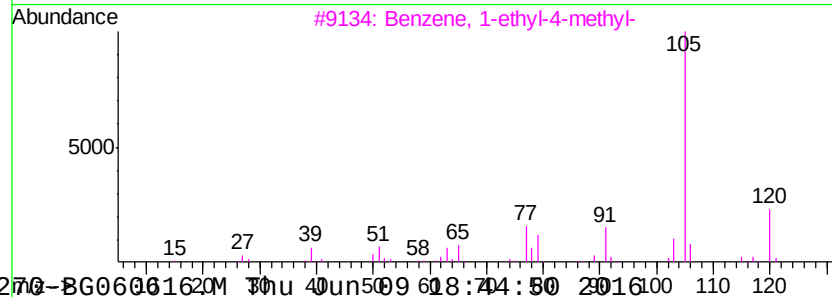
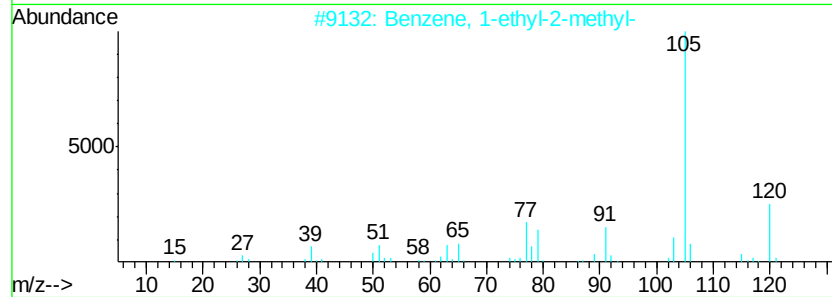
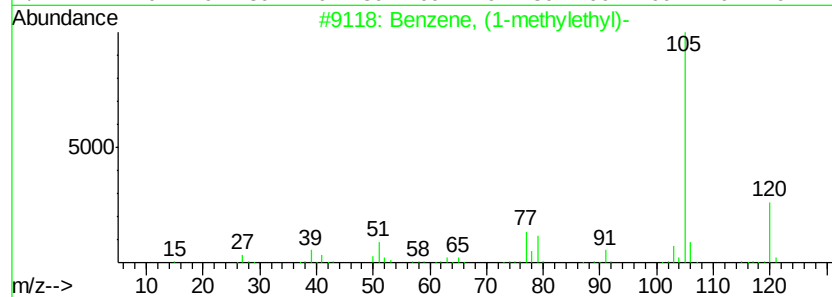
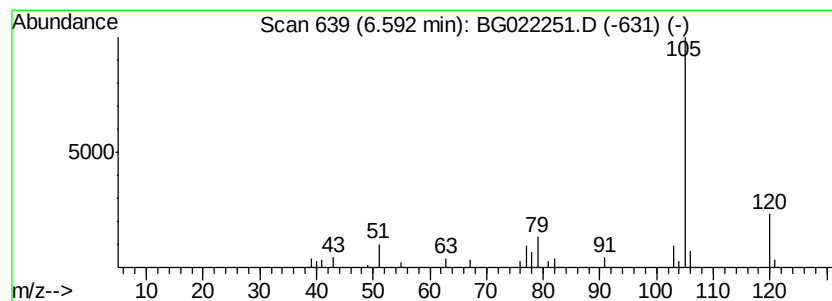
Quant Title :

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

Peak Number 2 Benzene, (1-methylethyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.59	3.69 ng	33623	1,4-Dichlorobenzene-d4	8.12

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	91
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	86
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	86
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	86
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	83



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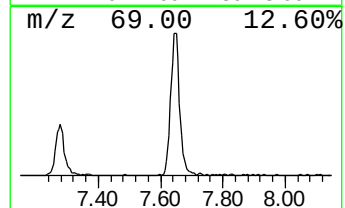
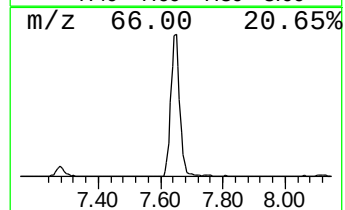
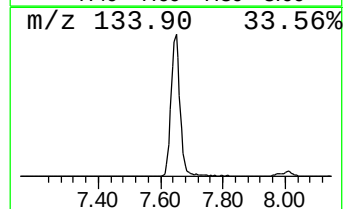
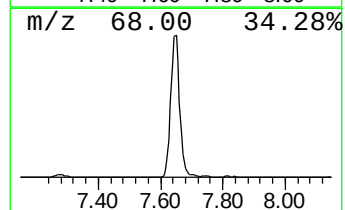
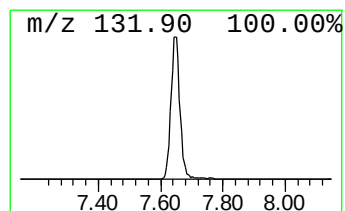
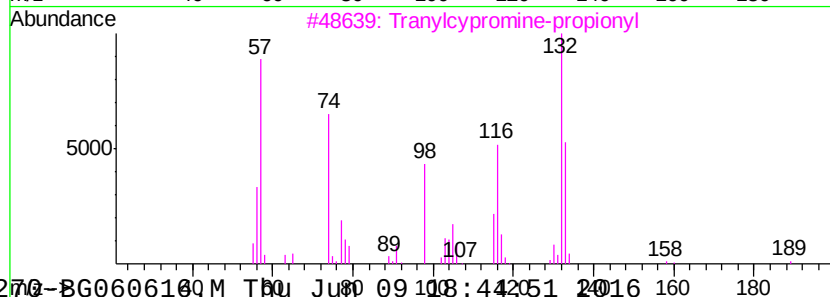
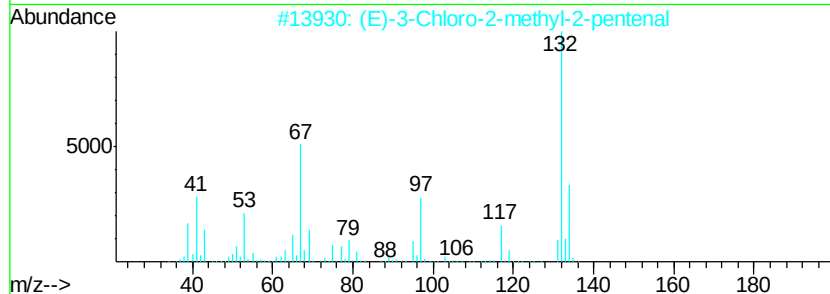
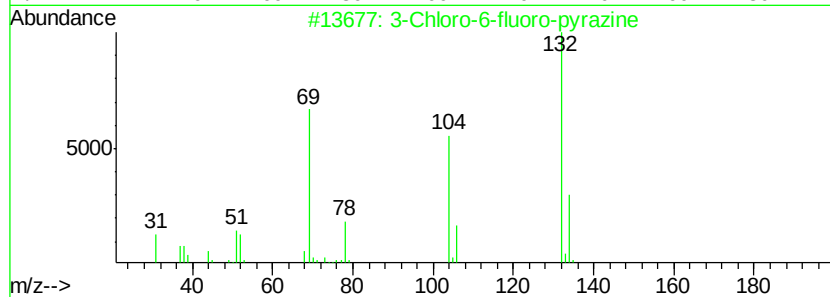
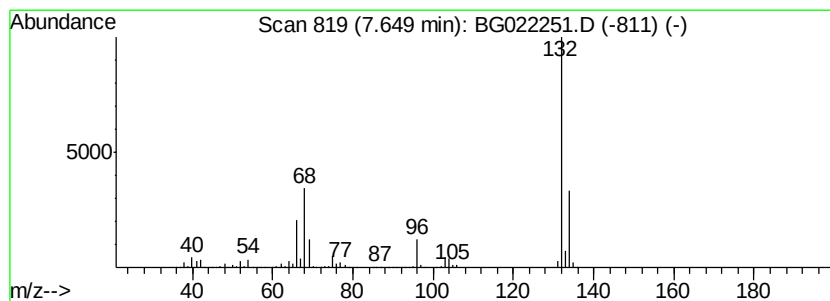
Quant Title :

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

Peak Number 3 unknown7.65 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.65	99.94 ng	910445	1,4-Dichlorobenzene-d4	8.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
3	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4	Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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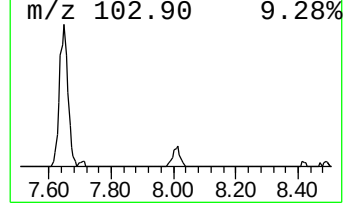
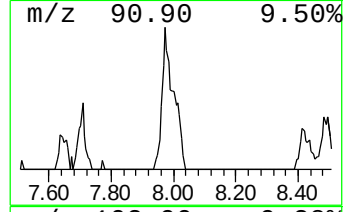
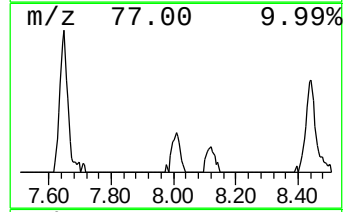
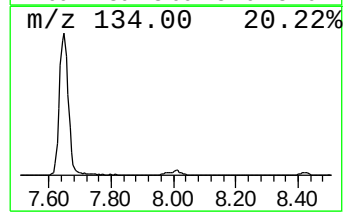
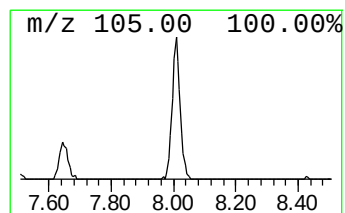
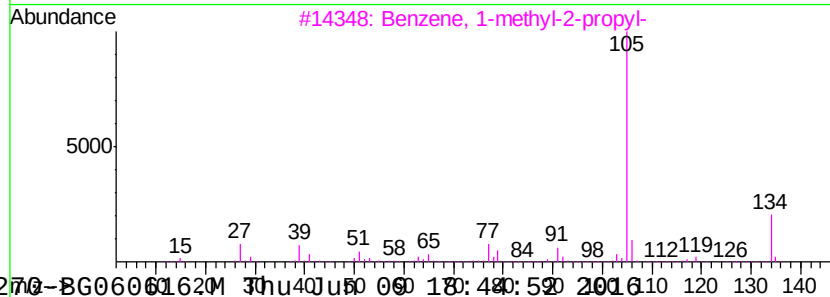
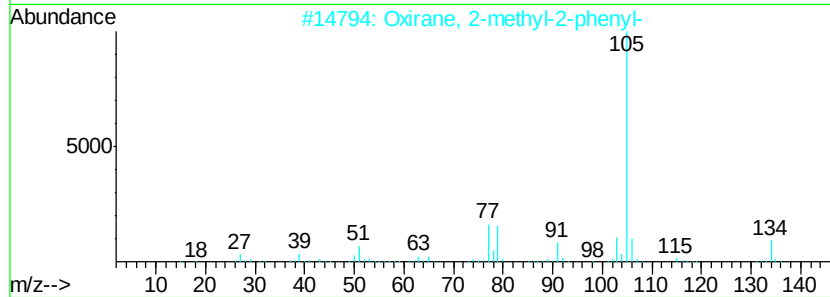
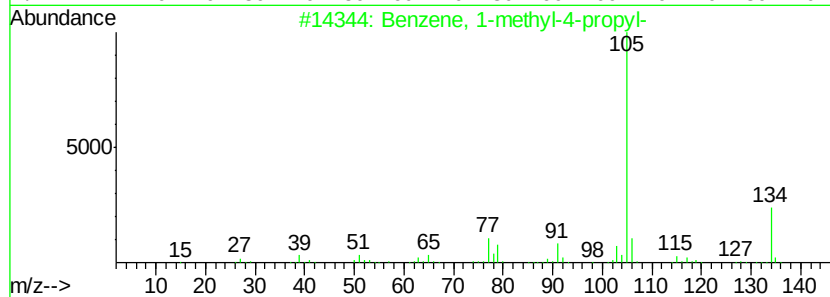
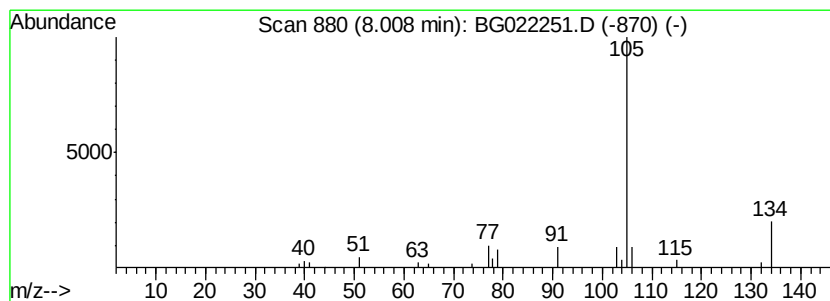
Quant Title :

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

Peak Number 4 Benzene, 1-methyl-4-propyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.01	4.83 ng	43988	1,4-Dichlorobenzene-d4	8.12

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	86
2			Oxirane, 2-methyl-2-phenyl-	134	C9H10O	002085-88-3	86
3			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	80
4			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	80
5			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	80



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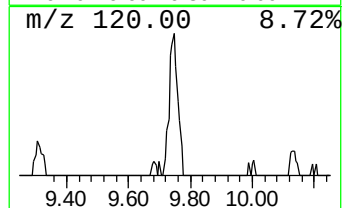
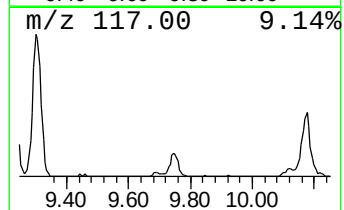
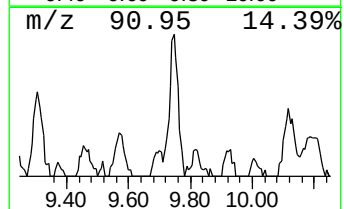
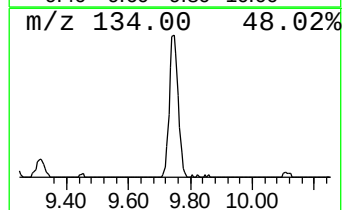
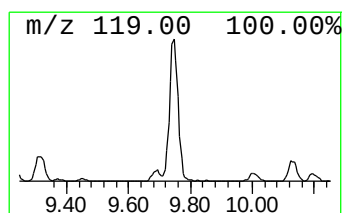
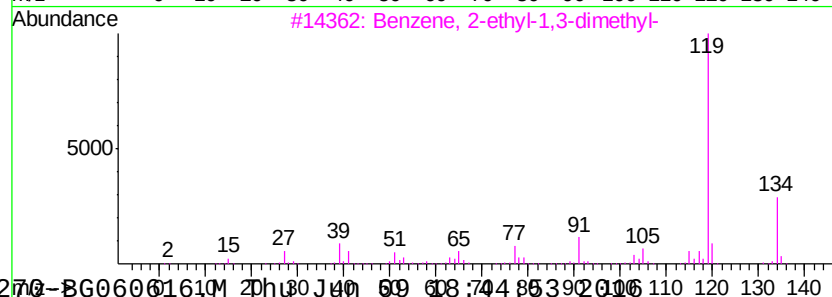
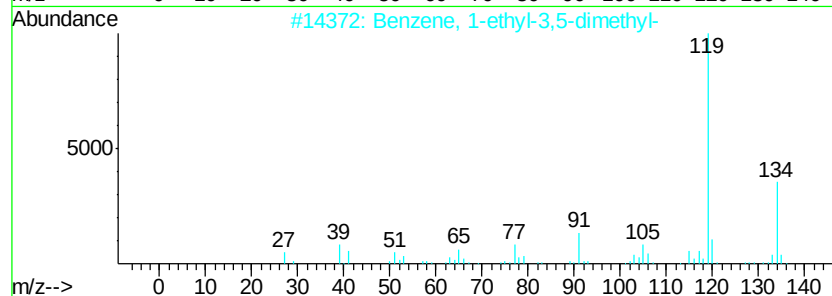
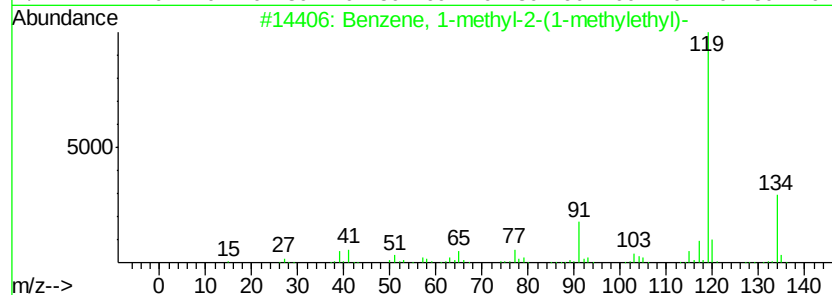
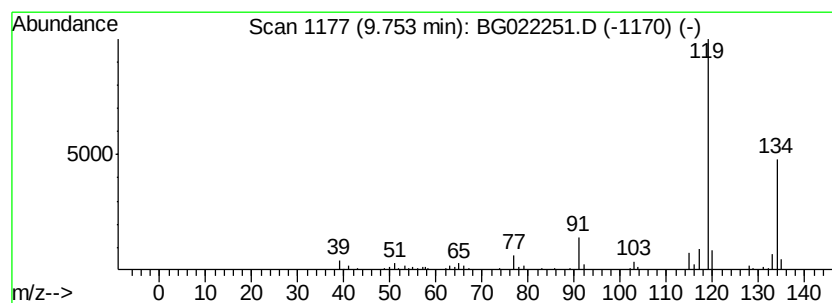
Quant Title :

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

Peak Number 7 Benzene, 1-methyl-2-(1-meth... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.75	5.30 ng	111991	Napthalene-d8	10.93

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene,	1-methyl-2-(1-methyleth...	134	C10H14		000527-84-4	91
2	Benzene,	1-ethyl-3,5-dimethyl-	134	C10H14		000934-74-7	91
3	Benzene,	2-ethyl-1,3-dimethyl-	134	C10H14		002870-04-4	91
4	Benzene,	1-methyl-3-(1-methyleth...	134	C10H14		000535-77-3	91
5	Benzene,	1,2,3,5-tetramethyl-	134	C10H14		000527-53-7	91



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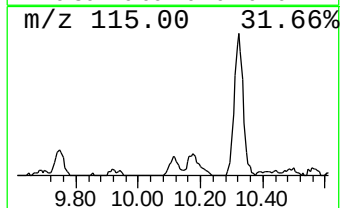
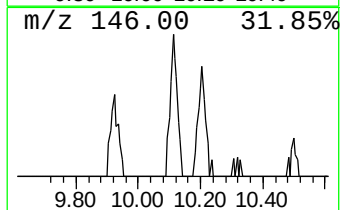
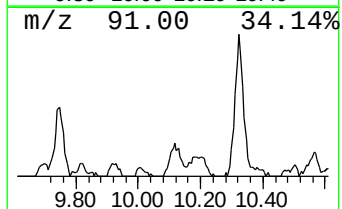
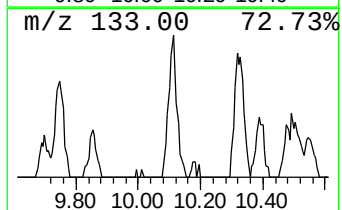
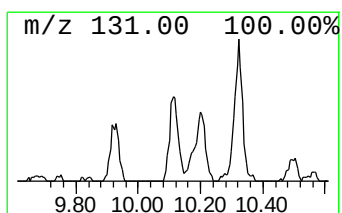
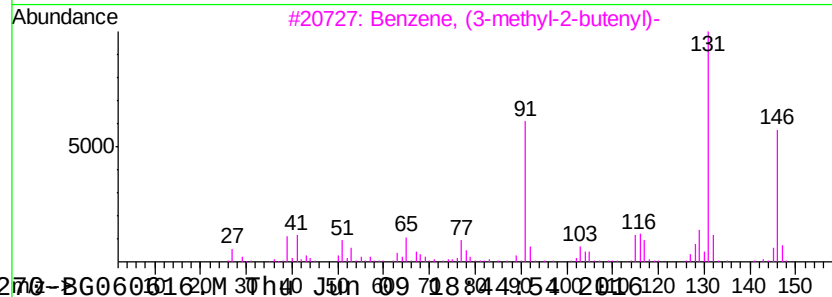
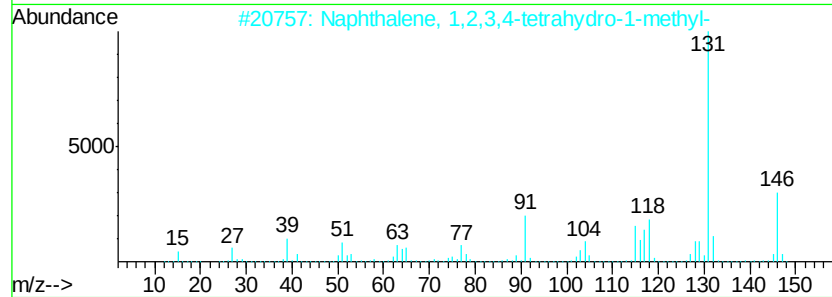
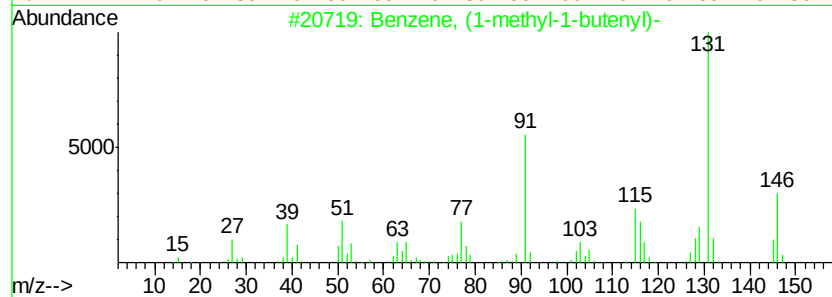
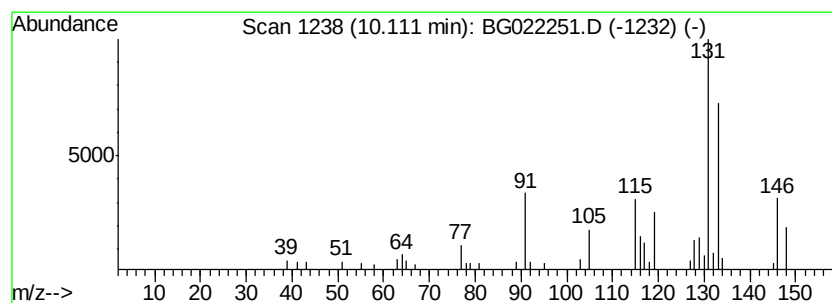
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ClientSampleId :
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Quant Title :

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

Peak Number 8 Benzene, (1-methyl-1-butenyl)- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.11	2.43 ng	51411	Naphthalene-d8	10.93	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	60
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	60
3	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	60
4	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	58
5	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	53



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Operator : UM/SJ
Sample : H3381-10
Misc :
ALS Vial : 27 Sample Multiplier: 1

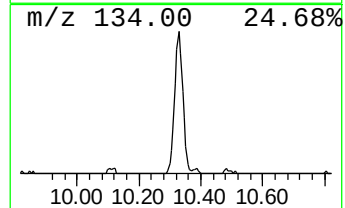
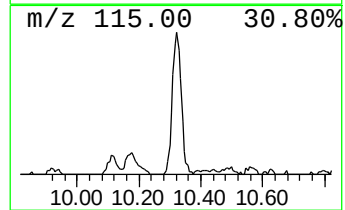
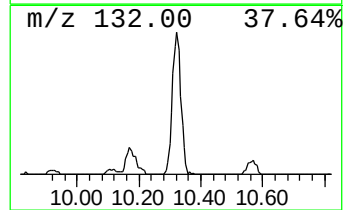
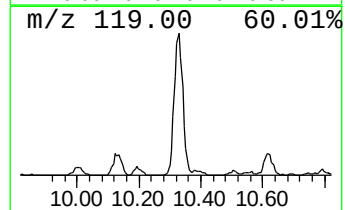
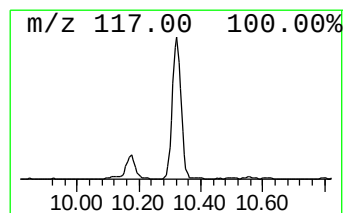
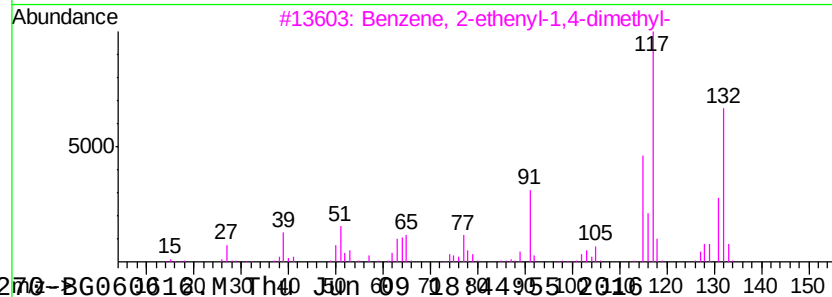
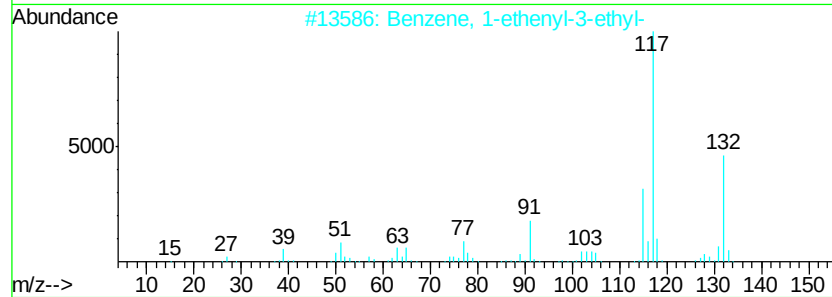
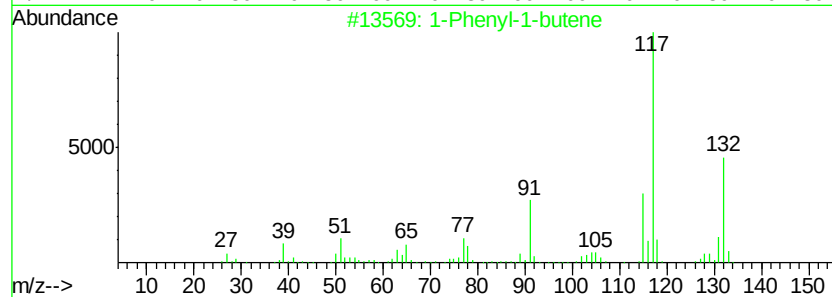
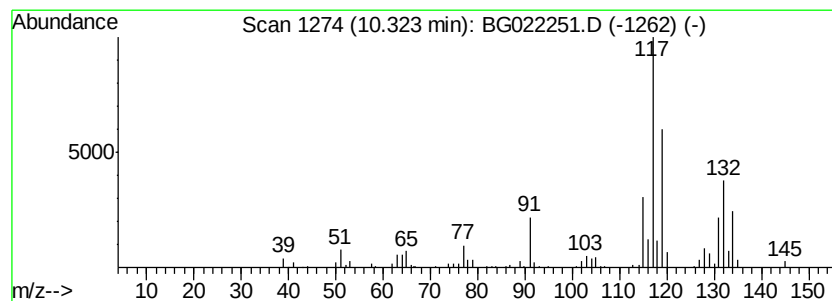
Instrument :
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ClientSampleId :
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Quant Title :

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

Peak Number 9 1-Phenyl-1-butene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.32	14.35 ng	303169	Napthalene-d8	10.93
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8 92
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4 91
3	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6 90
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1 86
5	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7 64



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Data File : BG022251.D
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Sample : H3381-10
Misc :
ALS Vial : 27 Sample Multiplier: 1

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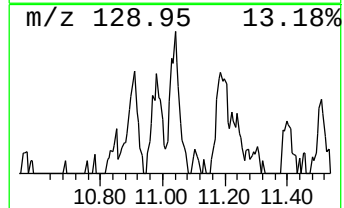
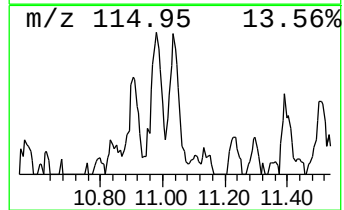
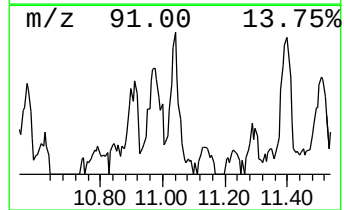
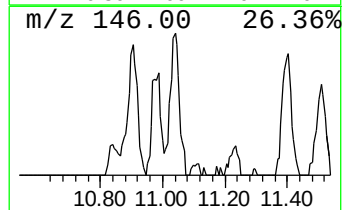
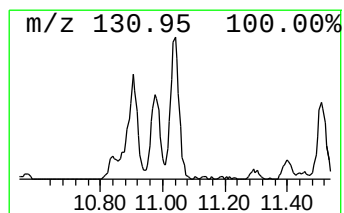
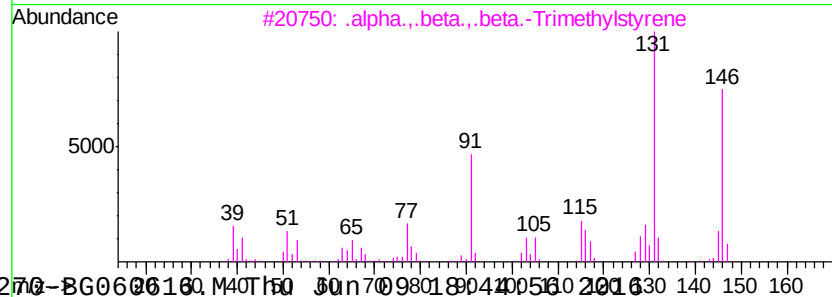
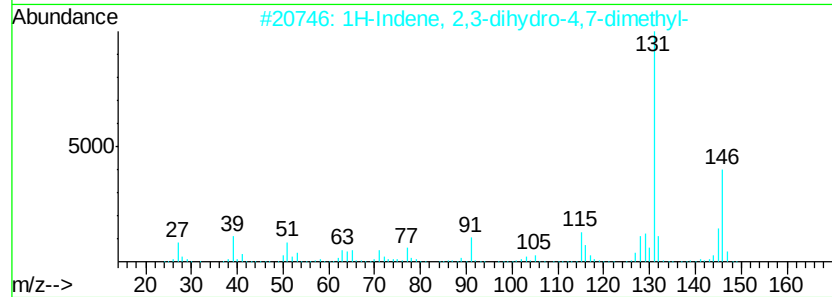
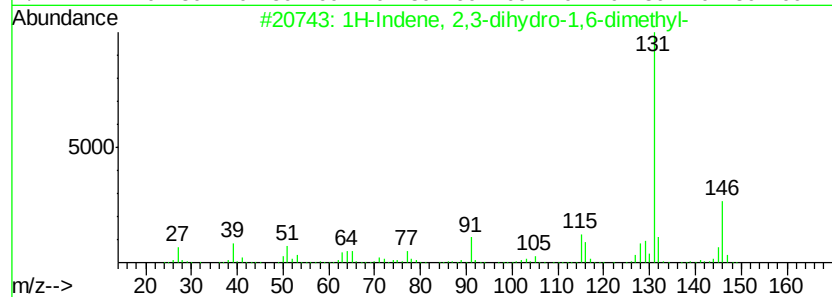
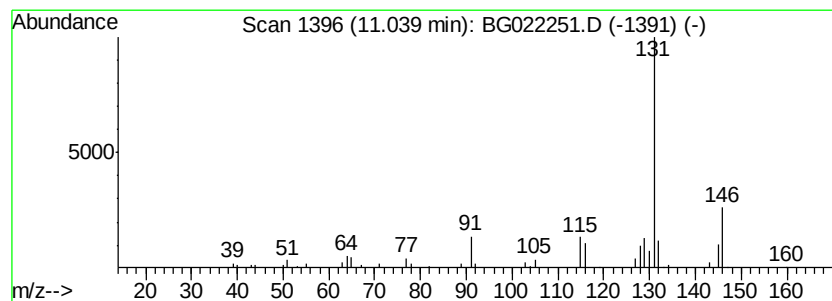
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.04	3.27 ng	69138	Napthalene-d8	10.93

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	95
2			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
3			.alpha.,.beta.,.beta.-Trimethyls...	146	C11H14	000769-57-3	91
4			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91
5			1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	91



Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
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Sample : H3381-10
Misc :
ALS Vial : 27 Sample Multiplier: 1

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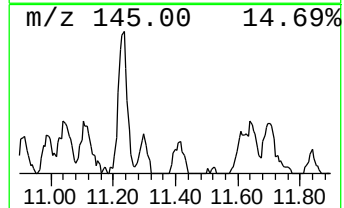
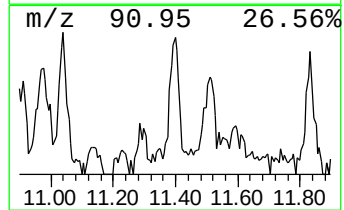
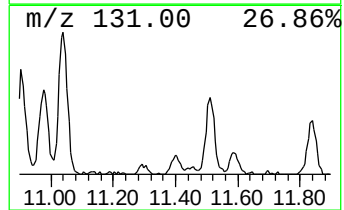
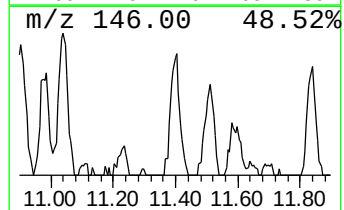
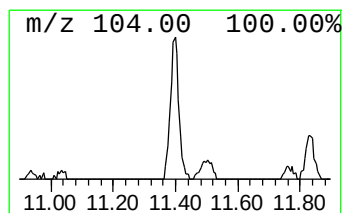
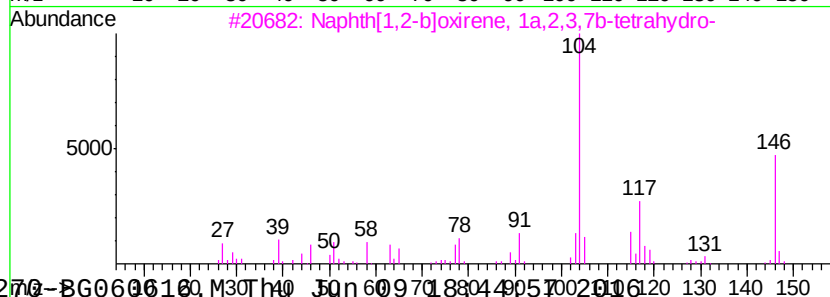
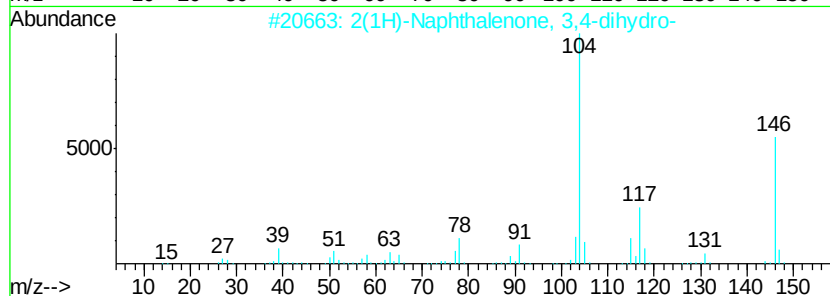
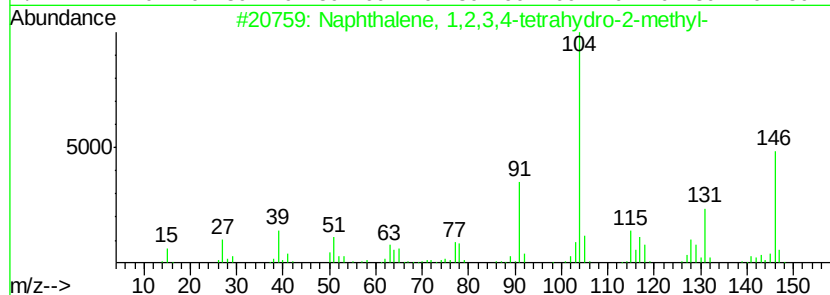
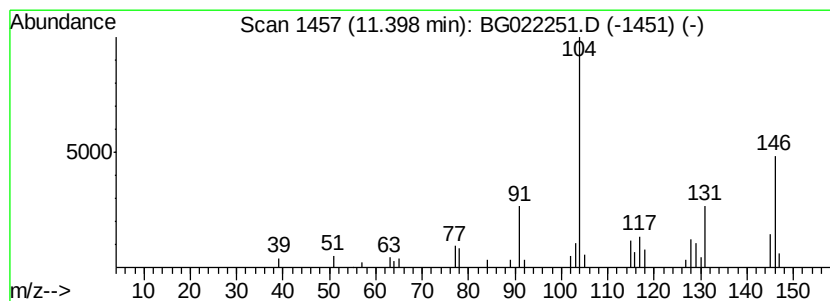
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.40	2.01 ng	42482	Naphthalene-d8	10.93

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	93
2		2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	53
3		Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	50
4		2-Propenoic acid, 2-phenylethyl ...	176	C11H12O2	003530-36-7	35
5		2-Bromopropionic acid, 2-phenyle...	256	C11H13BrO2	043216-34-8	25



Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
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Sample : H3381-10
Misc :
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Instrument :
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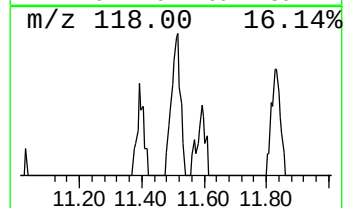
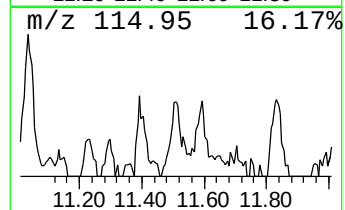
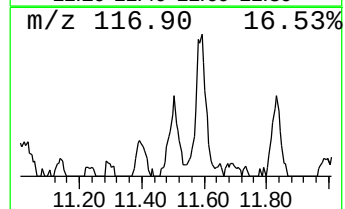
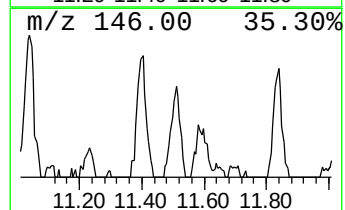
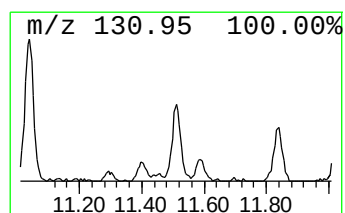
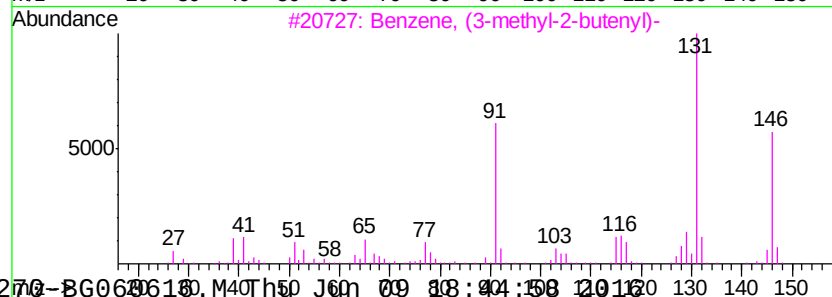
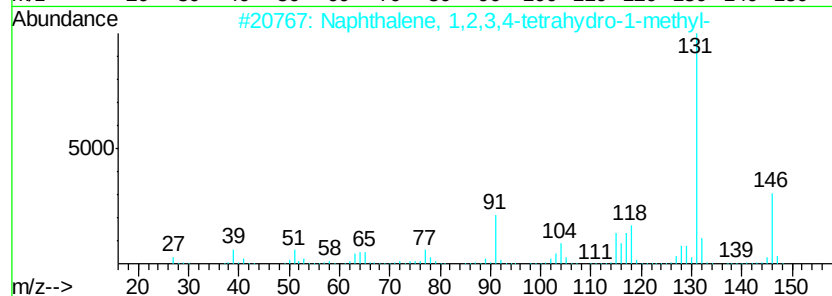
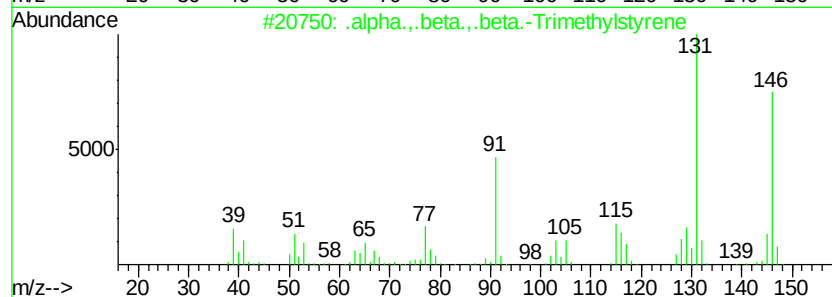
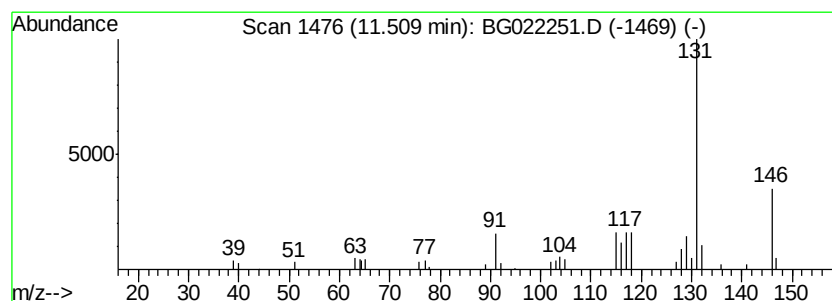
Quant Title :

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

Peak Number 12 .alpha.,.beta.,.beta.-Trime... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.51	2.10 ng	44303	Naphthalene-d8	10.93

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.,.beta.,.beta.-Trimethyls...	146	C11H14	000769-57-3	94
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	94
3	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	93
4	1H-Indene, 2,3-dihydro-2,2-dimethyl-	146	C11H14	020836-11-7	93
5	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90



Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
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Sample : H3381-10
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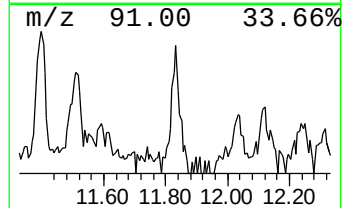
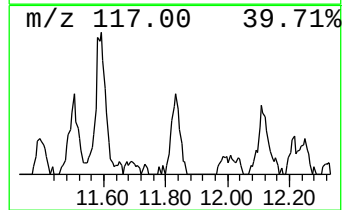
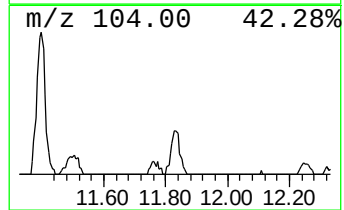
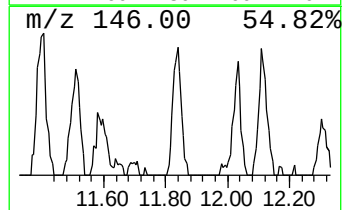
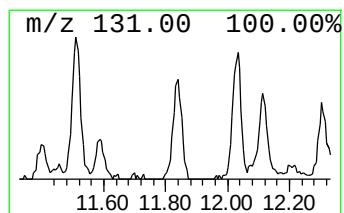
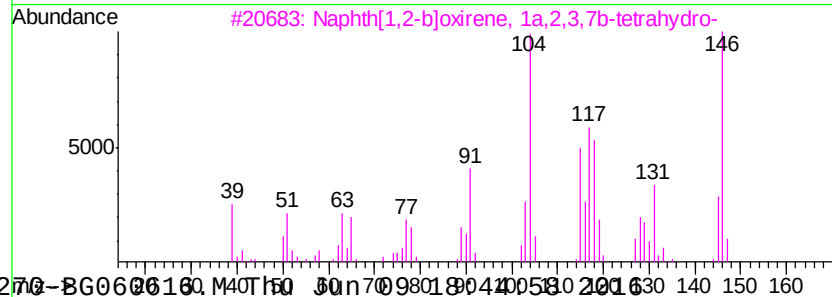
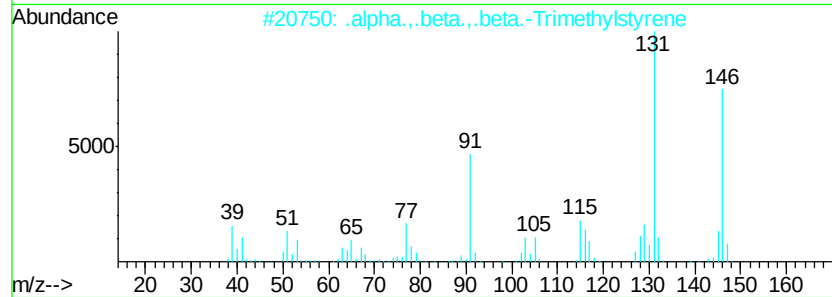
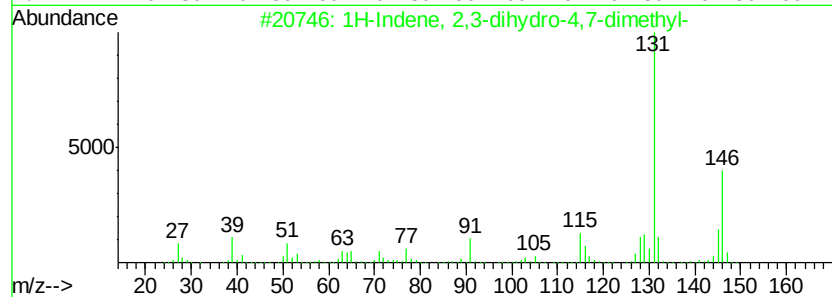
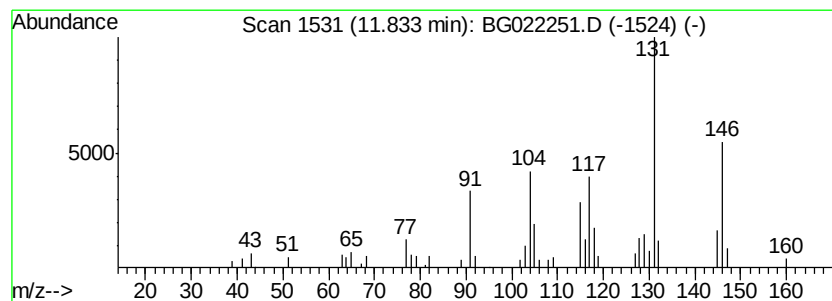
Instrument :
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ClientSampleId :
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Quant Title :

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

Peak Number 13 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.83	2.46 ng	51974	Napthalene-d8	10.93	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
2	.alpha.,.beta.,.beta.-Trimethyls...	146	C11H14	000769-57-3	91
3	Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	83
4	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	76
5	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70



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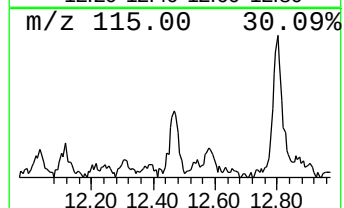
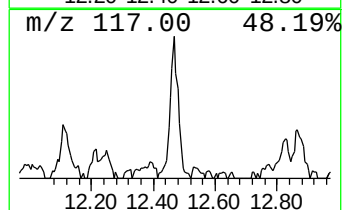
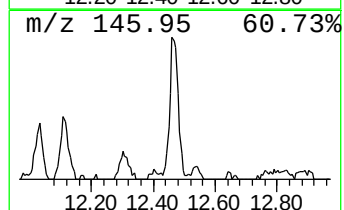
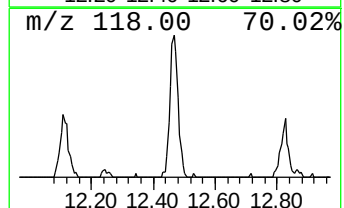
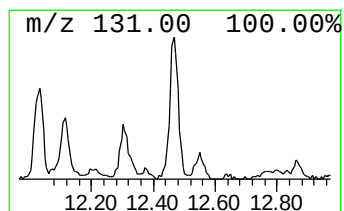
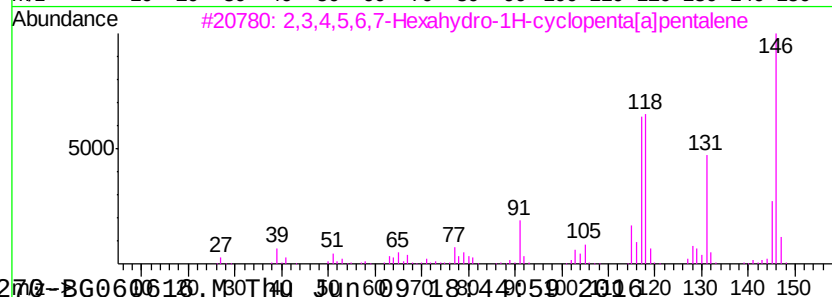
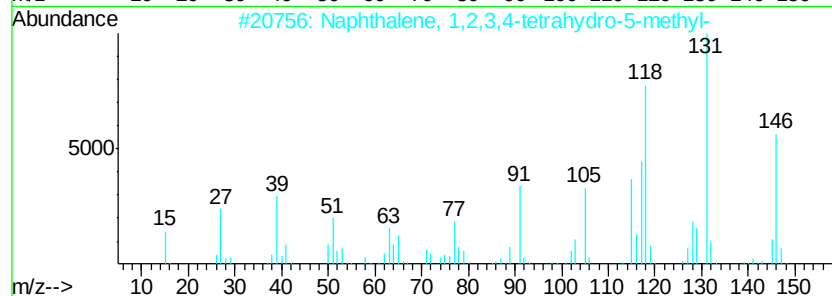
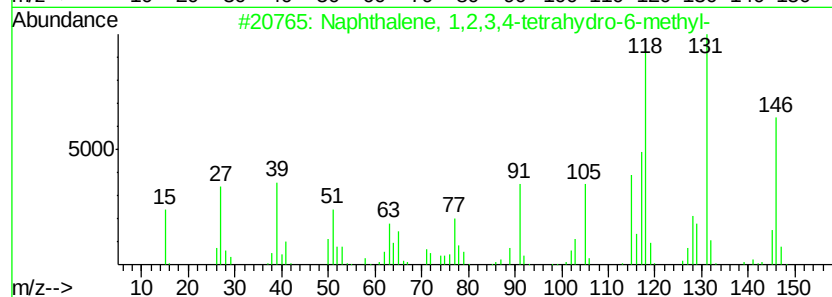
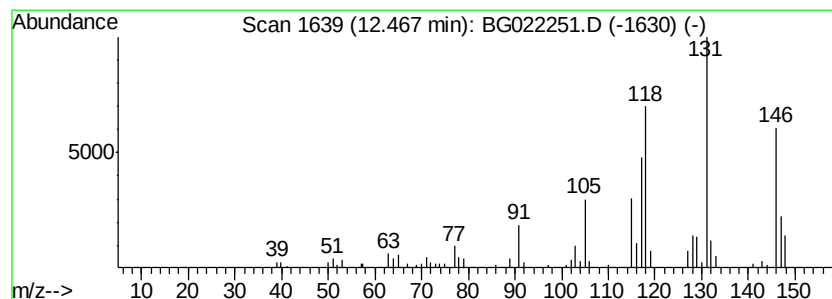
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Quant Title :

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

Peak Number 14 Naphthalene, 1,2,3,4-tetra... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.47	5.01 ng	105805	Naphthalene-d8	10.93	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	95
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	93
3	2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1000189-31-0	68
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	60
5	.alpha.,.beta.,.beta.-Trimethyls...	146	C11H14	000769-57-3	60



Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
Data File : BG022251.D
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Sample : H3381-10
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
W-53

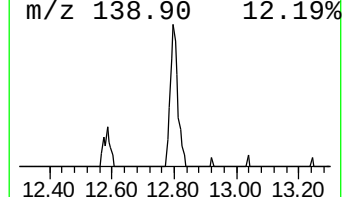
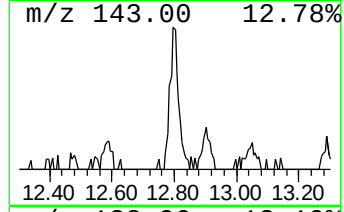
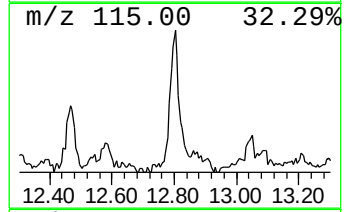
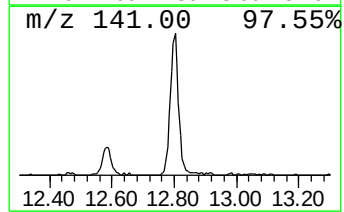
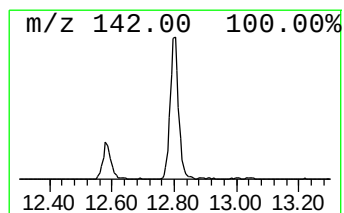
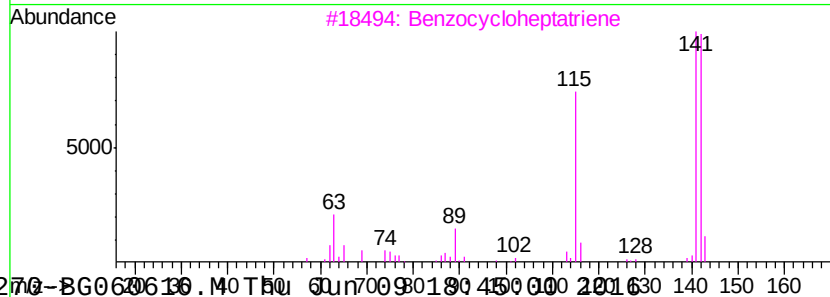
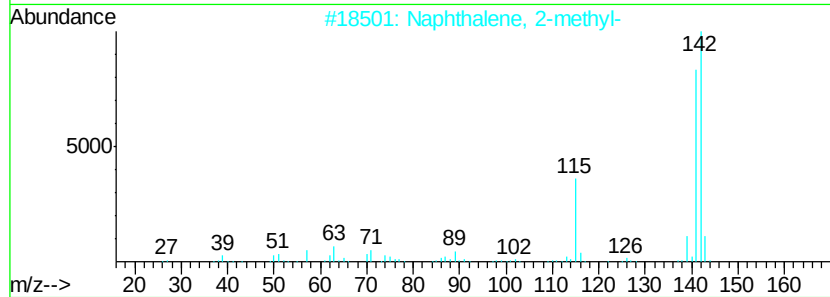
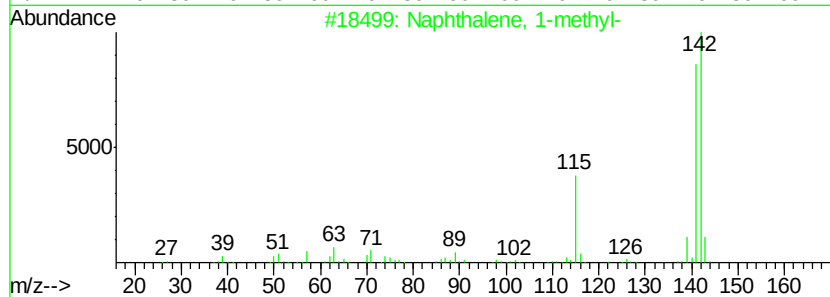
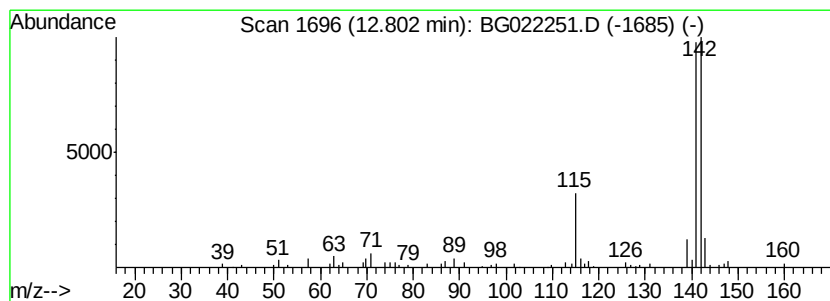
Quant Title :

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

Peak Number 15 Naphthalene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.80	7.07 ng	149441	Naphthalene-d8	10.93

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	97
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
3		Benzocycloheptatriene	142	C11H10	000264-09-5	91
4		1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	91
5		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	90



Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
Data File : BG022251.D
Acq On : 9 Jun 2016 2:46
Operator : UM/SJ
Sample : H3381-10
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
W-53

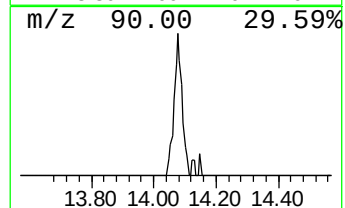
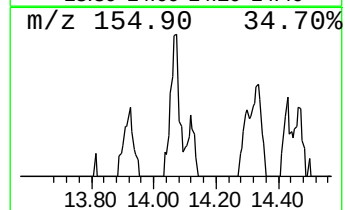
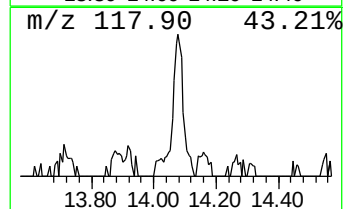
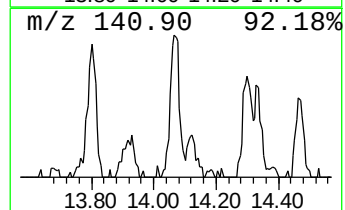
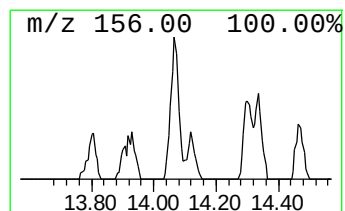
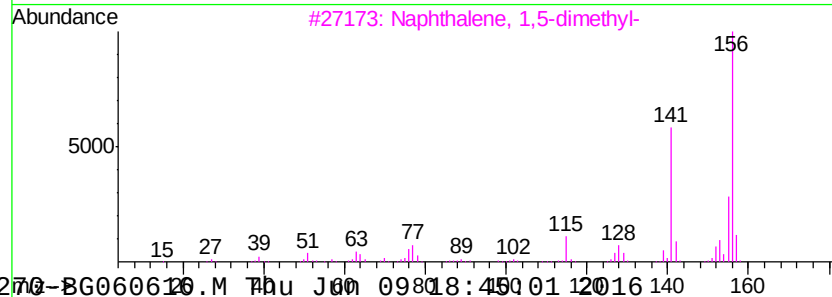
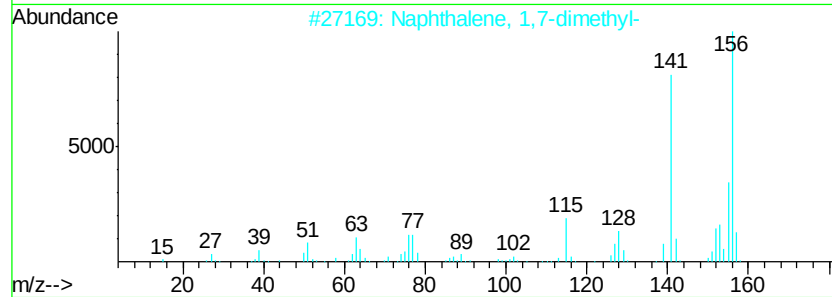
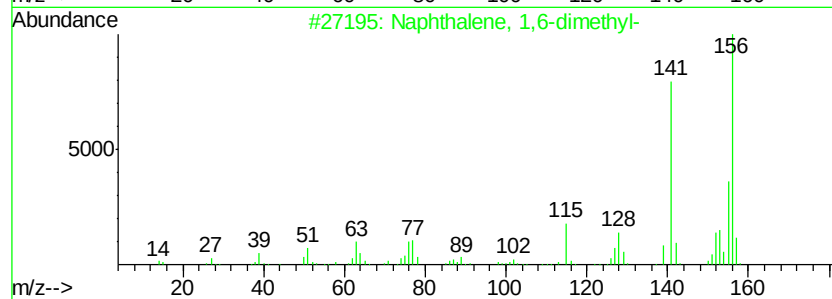
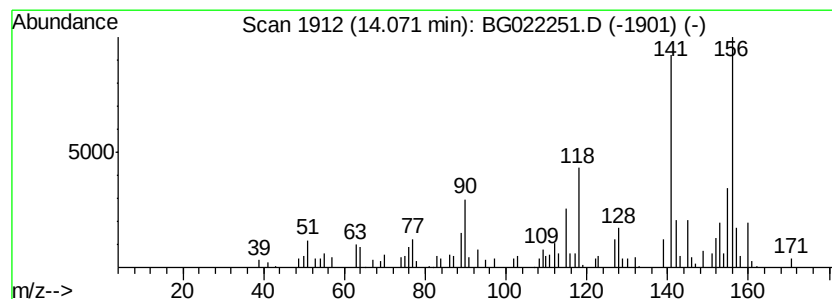
Quant Title :

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

Peak Number 16 Naphthalene, 1,6-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.07	2.82 ng	69350	Acenaphthene-d10	14.75

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	78
2			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	78
3			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	64
4			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	60
5			Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	60



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060816\
Data File : BG022251.D
Acq On : 9 Jun 2016 2:46
Operator : UM/SJ
Sample : H3381-10
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
W-53

Quant Title :

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...	5.19	3.1	ng	28176	1	8.12	182202	20.0
Benzene, (1-methy...	6.59	3.7	ng	33623	1	8.12	182202	20.0
unknown7.65	7.65	99.9	ng	910445	1	8.12	182202	20.0
Benzene, 1-methyl...	8.01	4.8	ng	43988	1	8.12	182202	20.0
Benzene, 1-methyl...	9.75	5.3	ng	111991	2	10.93	422546	20.0
Benzene, (1-methy...	10.11	2.4	ng	51411	2	10.93	422546	20.0
1-Phenyl-1-butene	10.32	14.4	ng	303169	2	10.93	422546	20.0
1H-Indene, 2,3-di...	11.04	3.3	ng	69138	2	10.93	422546	20.0
Naphthalene, 1,2,...	11.40	2.0	ng	42482	2	10.93	422546	20.0
.alpha.,.beta.,.b...	11.51	2.1	ng	44303	2	10.93	422546	20.0
1H-Indene, 2,3-di...	11.83	2.5	ng	51974	2	10.93	422546	20.0
Naphthalene, 1,2,...	12.47	5.0	ng	105805	2	10.93	422546	20.0
Naphthalene, 1-me...	12.80	7.1	ng	149441	2	10.93	422546	20.0
Naphthalene, 1,6-...	14.07	2.8	ng	69350	3	14.75	491344	20.0