

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121915\
 Data File : BF083941.D
 Acq On : 19 Dec 2015 4:43
 Operator : UM/SJ
 Sample : G4840-06
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W-48

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-BF121815.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	2.373	18	25	27	rBV	22954	50436	2.35%	0.377%
2	3.893	155	158	161	rBV	27444	38586	1.80%	0.288%
3	4.579	215	218	221	rBB	16457	25179	1.17%	0.188%
4	4.830	238	240	241	rBV	34650	36072	1.68%	0.270%
5	4.853	241	242	245	rVB	18654	23704	1.10%	0.177%
6	5.047	257	259	263	rBV	72086	103581	4.83%	0.774%
7	5.482	294	297	299	rBV	678588	684886	31.92%	5.120%
8	6.510	385	387	389	rBV	358459	430222	20.05%	3.216%
9	6.636	396	398	400	rBV	1365804	1219723	56.85%	9.118%
10	6.819	412	414	416	rBB	40214	40220	1.87%	0.301%
11	6.865	416	418	420	rBV	429422	347880	16.21%	2.601%
12	7.025	429	432	434	rBV	1497412	1338610	62.39%	10.007%
13	7.185	445	446	448	rBV	43575	41656	1.94%	0.311%
14	7.425	464	467	469	rBV	725730	948691	44.22%	7.092%
15	7.516	473	475	476	rBV	32335	31446	1.47%	0.235%
16	7.562	476	479	481	rVB	36742	35903	1.67%	0.268%
17	7.722	491	493	494	rVB	32677	36841	1.72%	0.275%
18	7.825	500	502	505	rVB2	108091	135516	6.32%	1.013%
19	7.893	505	508	510	rBB	21724	26212	1.22%	0.196%
20	8.145	528	530	532	rBV	387411	440165	20.52%	3.291%
21	8.202	532	535	537	rVB2	131559	159507	7.43%	1.192%
22	8.533	562	564	565	rVB2	31765	25020	1.17%	0.187%
23	8.785	582	586	587	rBV2	18304	30065	1.40%	0.225%
24	8.990	602	604	607	rVB2	24839	33750	1.57%	0.252%
25	9.070	607	611	615	rBV3	22599	46366	2.16%	0.347%
26	9.230	622	625	627	rVB	2314060	2145427	100.00%	16.039%
27	9.619	657	659	662	rVB	32520	29859	1.39%	0.223%
28	9.905	681	684	686	rBV	580281	508381	23.70%	3.801%
29	10.339	719	722	724	rBV	21918	23165	1.08%	0.173%
30	10.693	750	753	755	rBV	1086413	1005797	46.88%	7.519%
31	11.391	811	814	816	rBV	405019	432644	20.17%	3.234%
32	11.596	829	832	834	rBV	39961	36122	1.68%	0.270%
33	12.751	931	933	935	rBV	60591	72608	3.38%	0.543%
34	12.968	950	952	955	rBV	1331312	1701161	79.29%	12.717%

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35	13.745	1018	1020	1025	rBV2	37704	58203	2.71%	0.435%
36	13.871	1029	1031	1034	rBV2	31472	50281	2.34%	0.376%
37	14.020	1042	1044	1047	rVB	439521	473859	22.09%	3.542%
38	14.637	1095	1098	1102	rVB3	18817	38448	1.79%	0.287%
39	15.460	1168	1170	1175	rVB	325960	470453	21.93%	3.517%

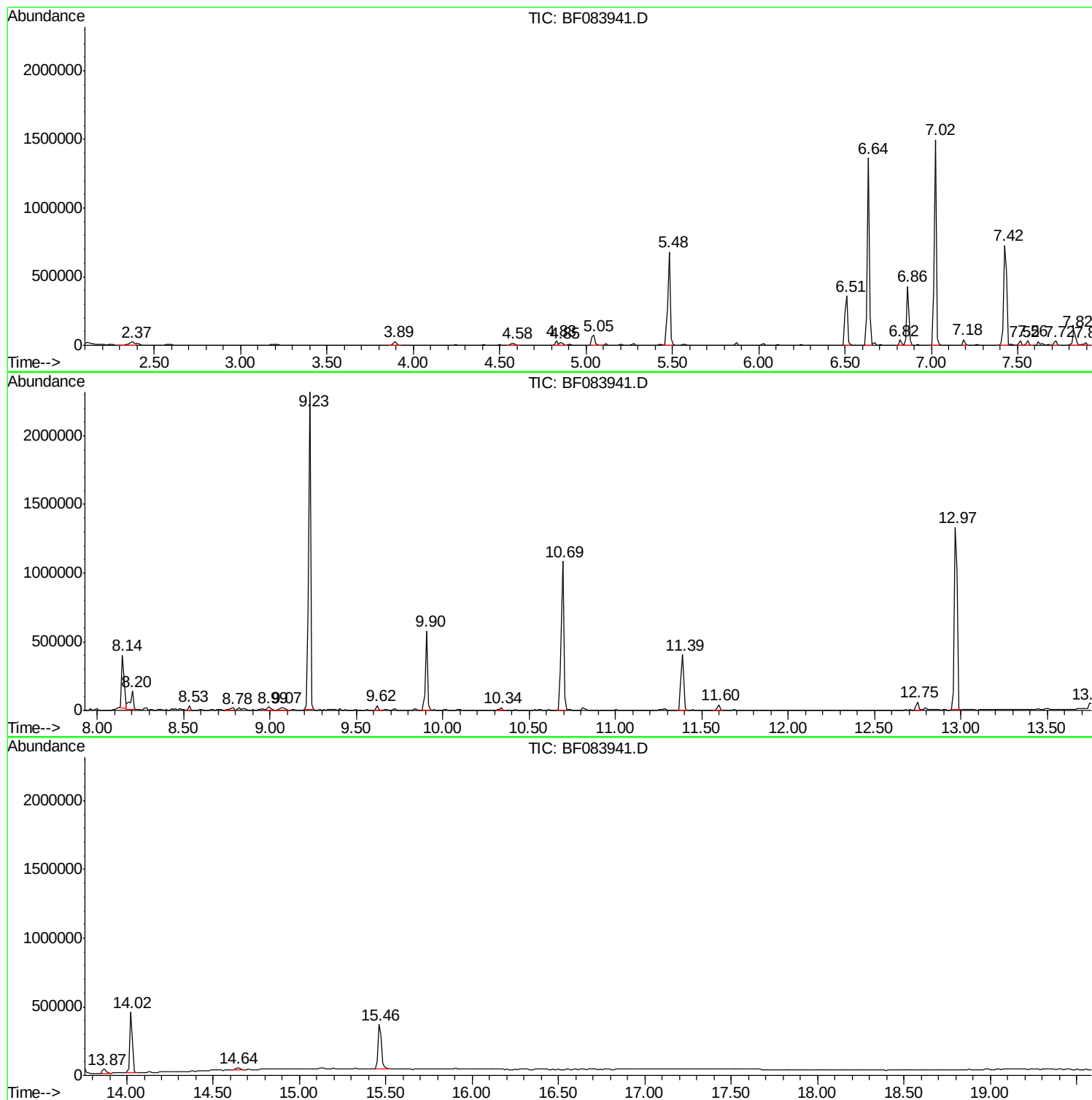
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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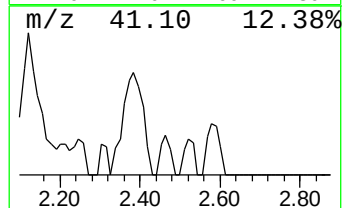
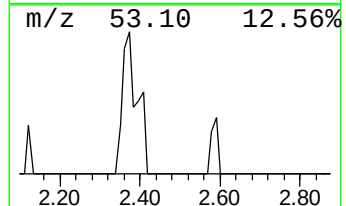
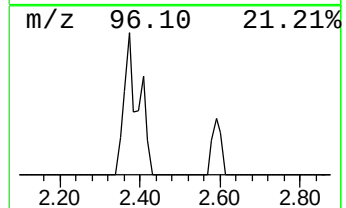
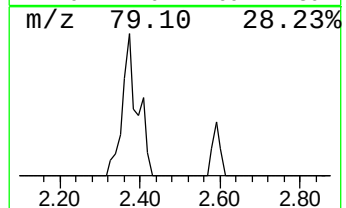
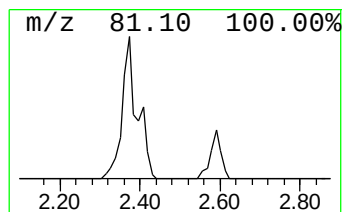
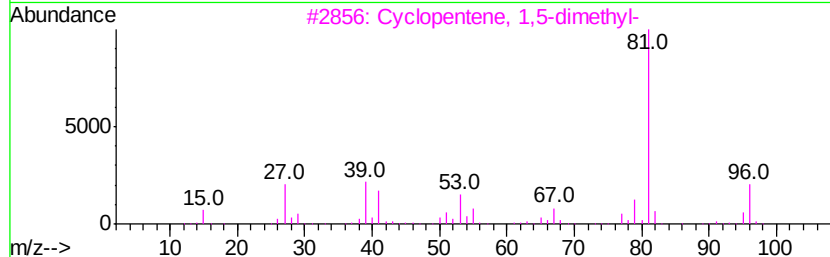
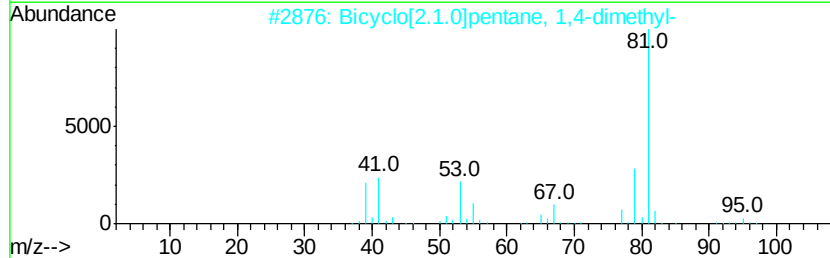
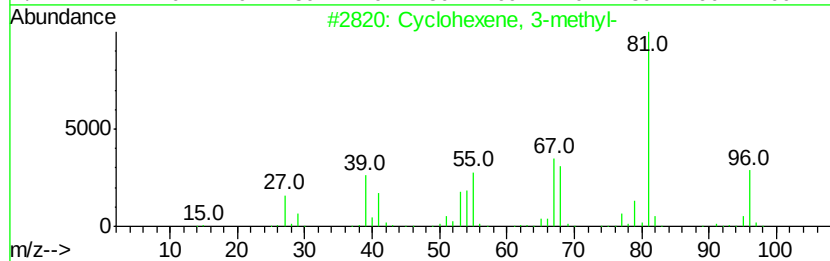
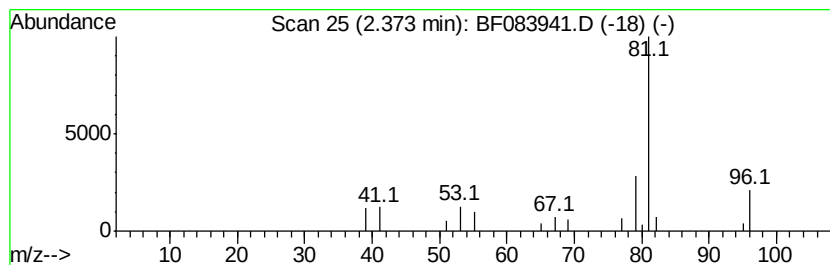
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Cyclohexene, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.37	2.90 ng	50436	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexene, 3-methyl-	96	C7H12	000591-48-0	72
2		Bicyclo[2.1.0]pentane, 1,4-dimet...	96	C7H12	017065-18-8	72
3		Cyclopentene, 1,5-dimethyl-	96	C7H12	016491-15-9	72
4		Cyclopentane, 1-methyl-2-methylene-	96	C7H12	041158-41-2	72
5		1,4-Pentadiene, 3,3-dimethyl-	96	C7H12	001112-35-2	72



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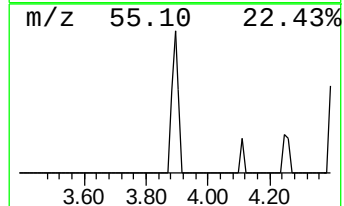
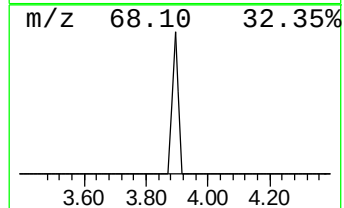
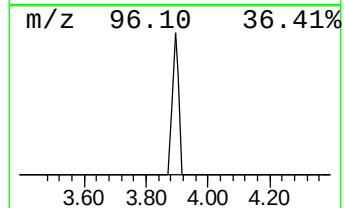
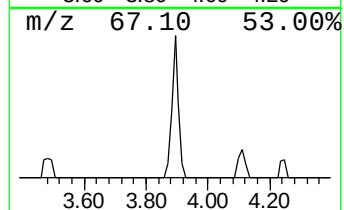
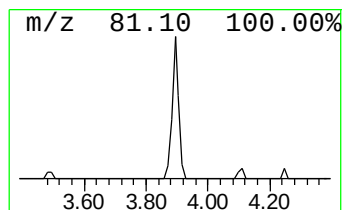
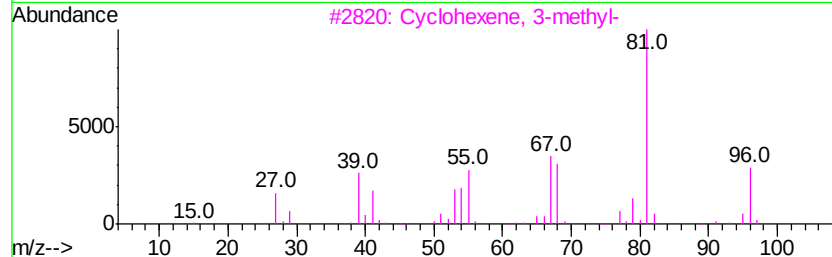
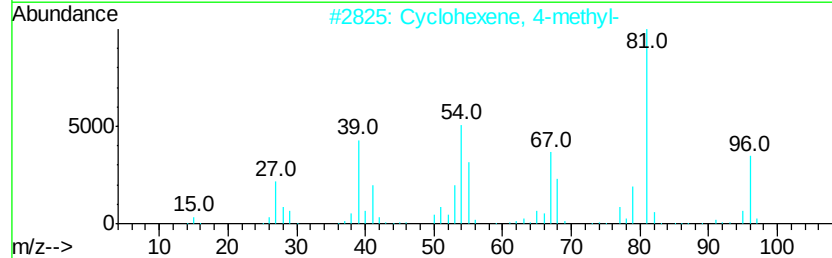
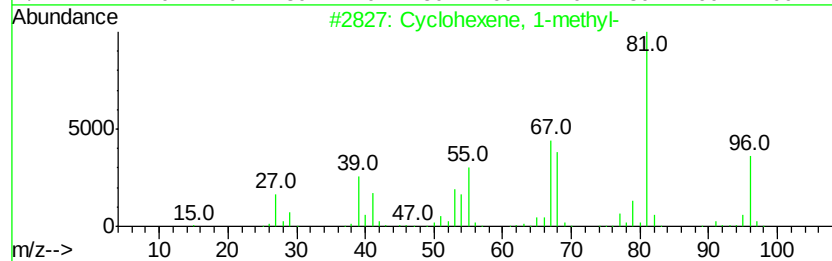
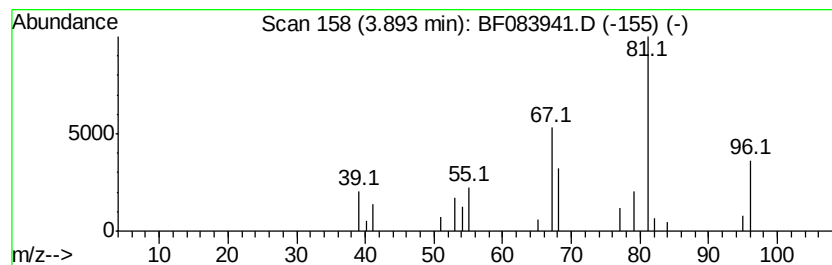
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclohexene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.89	2.22 ng	38586	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexene, 1-methyl-	96	C7H12	000591-49-1	90
2		Cyclohexene, 4-methyl-	96	C7H12	000591-47-9	80
3		Cyclohexene, 3-methyl-	96	C7H12	000591-48-0	72
4		Norbornane	96	C7H12	000279-23-2	72
5		Cyclopentene, 4,4-dimethyl-	96	C7H12	019037-72-0	59



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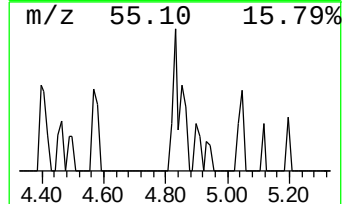
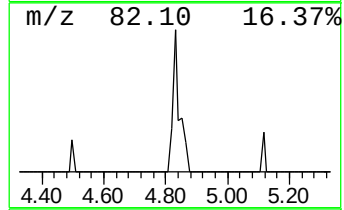
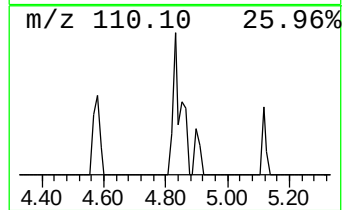
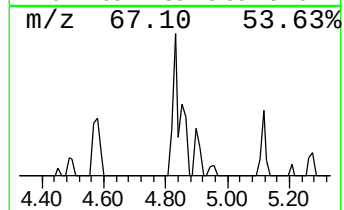
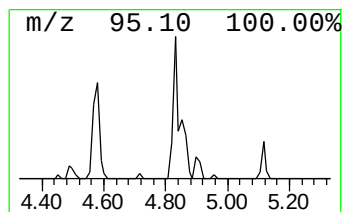
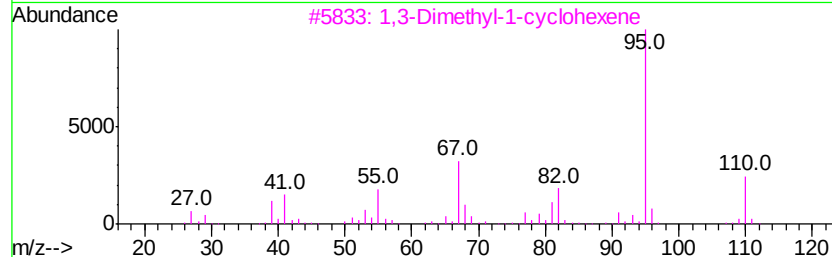
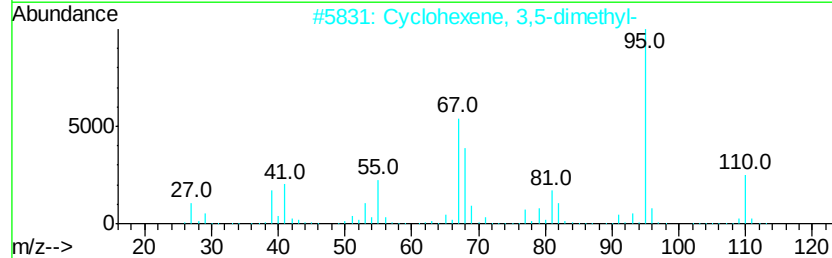
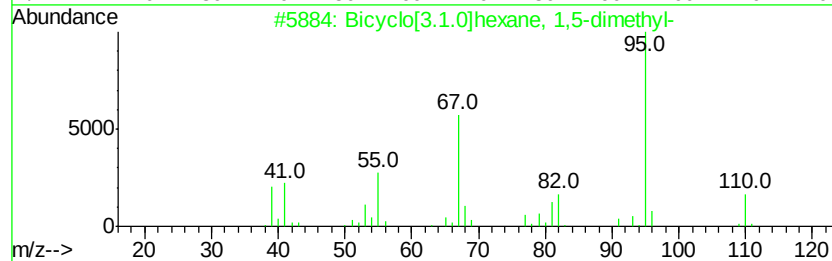
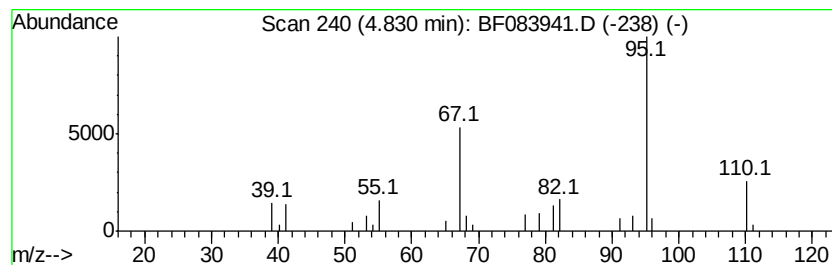
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Bicyclo[3.1.0]hexane, 1,5-d... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.83	2.07 ng	36072	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[3.1.0]hexane, 1,5-dimethyl-	110	C8H14	1000142-17-5	91
2		Cyclohexene, 3,5-dimethyl-	110	C8H14	000823-17-6	91
3		1,3-Dimethyl-1-cyclohexene	110	C8H14	002808-76-6	90
4		Cyclopentane, 1,2-dimethyl-3-met...	110	C8H14	091884-67-2	87
5		5,5-Dimethyl-1,3-hexadiene	110	C8H14	001515-79-3	87



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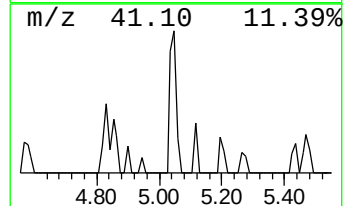
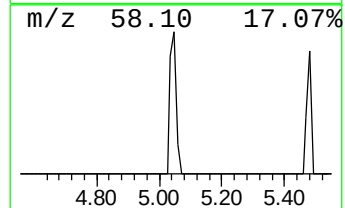
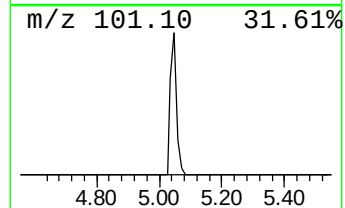
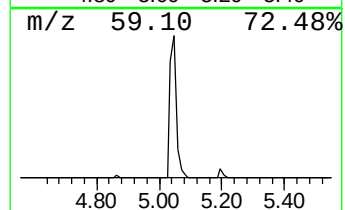
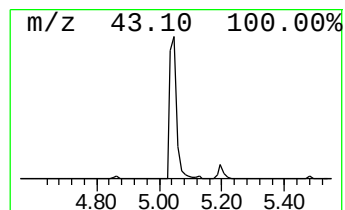
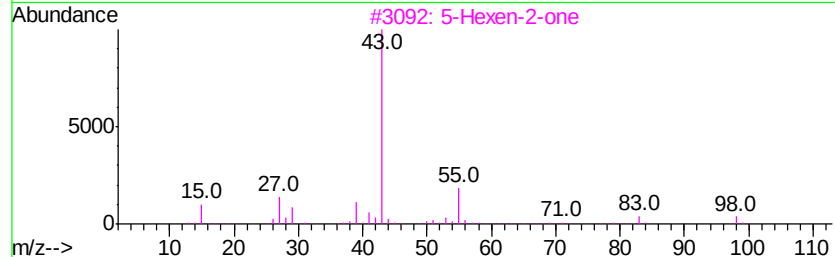
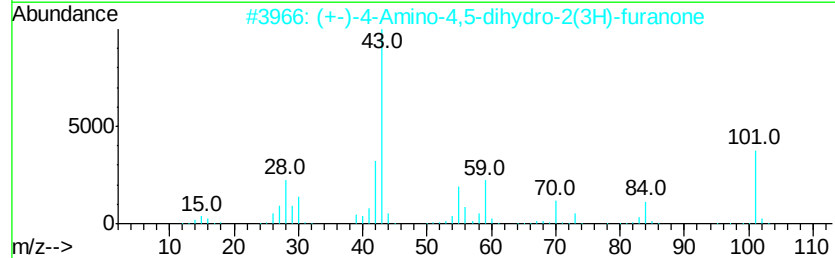
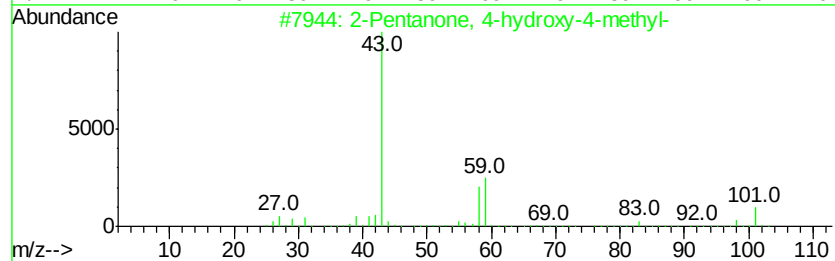
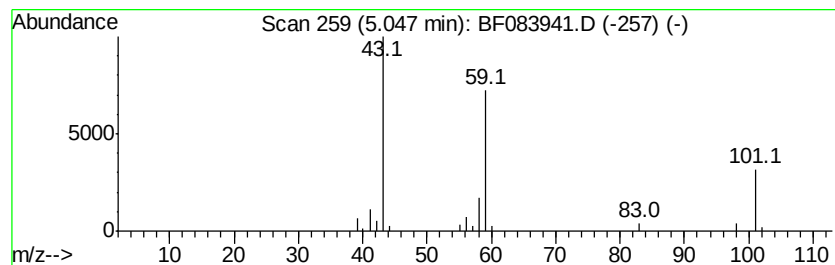
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	5.95 ng	103581	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	47
3		5-Hexen-2-one	98	C6H10O	000109-49-9	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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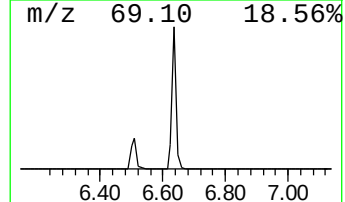
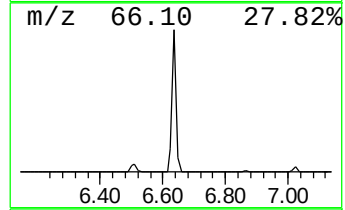
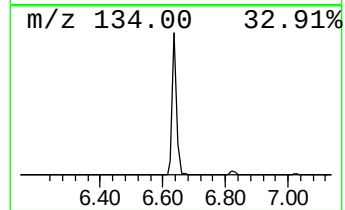
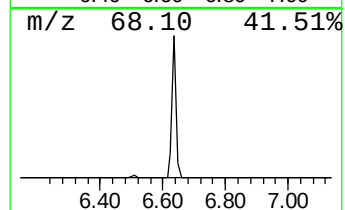
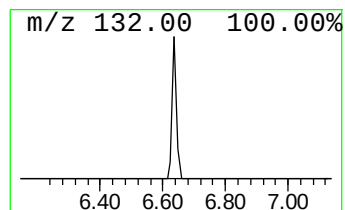
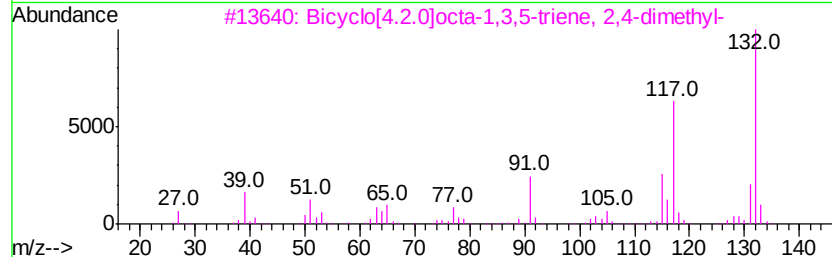
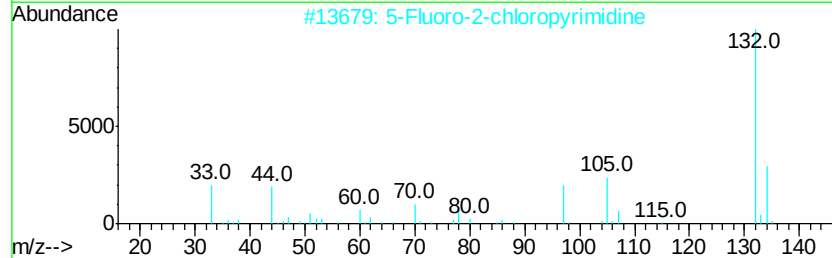
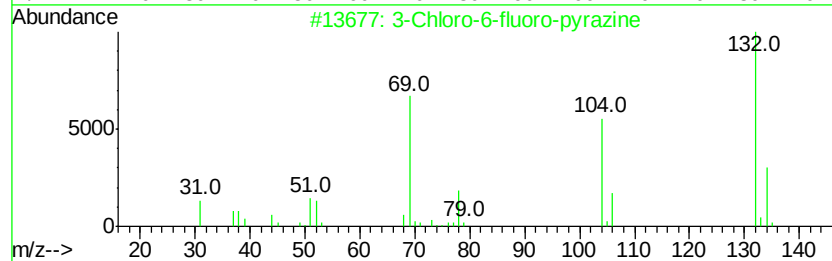
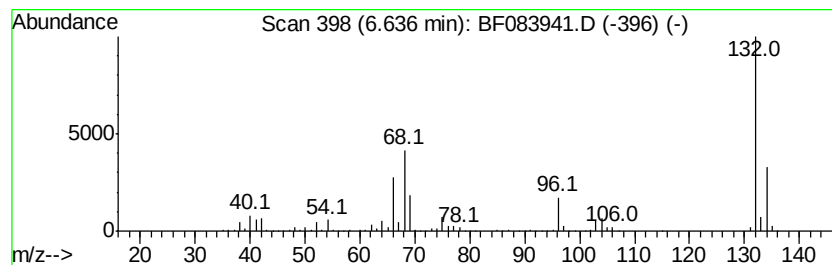
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Peak Number 5 unknown6.64 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.64	70.12 ng	1219720	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		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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Sample : G4840-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

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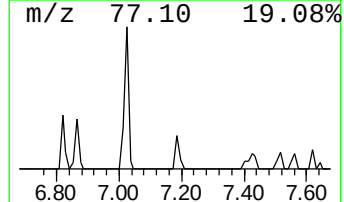
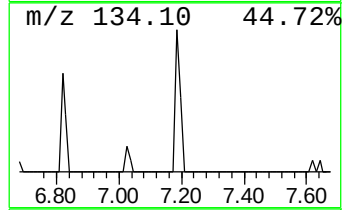
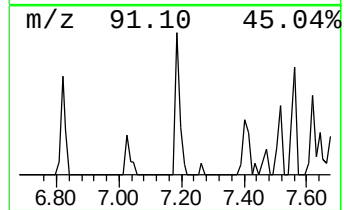
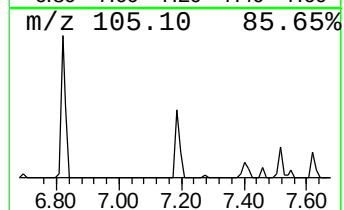
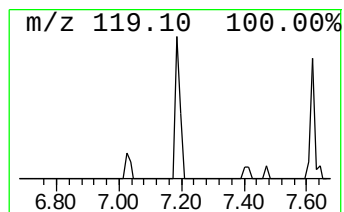
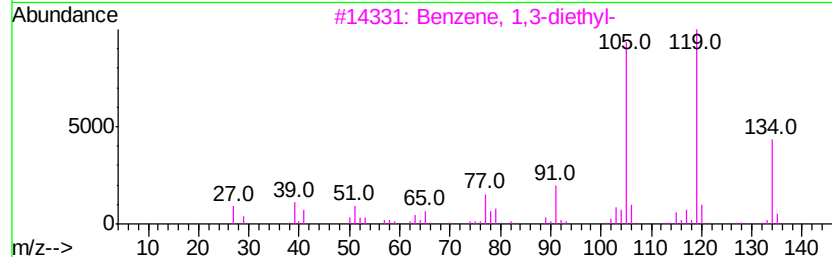
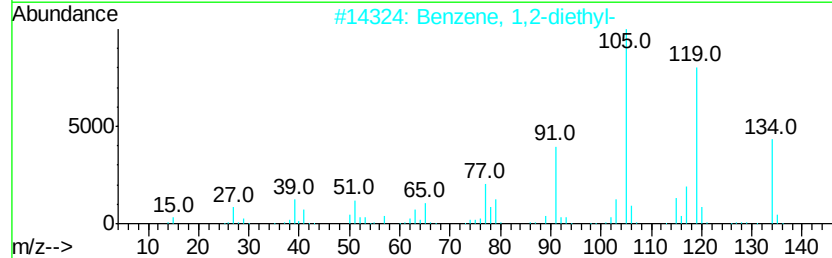
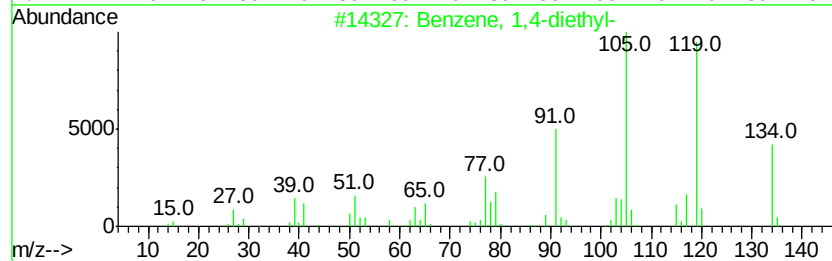
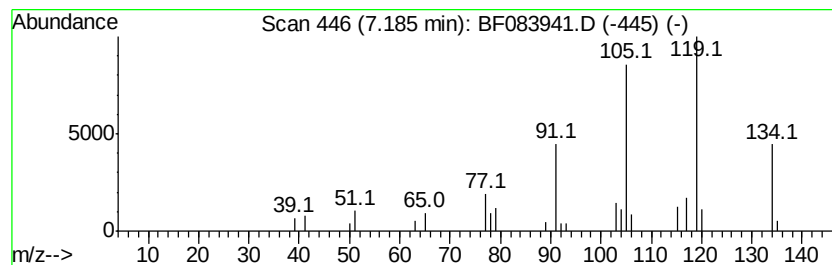
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Benzene, 1,4-diethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.18	2.39 ng	41656	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	96
2		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	91
4		Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	87
5		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121915\
Data File : BF083941.D
Acq On : 19 Dec 2015 4:43
Operator : UM/SJ
Sample : G4840-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
W-48

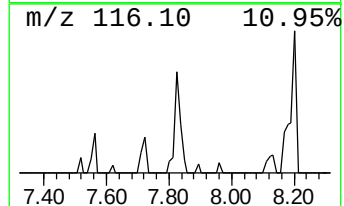
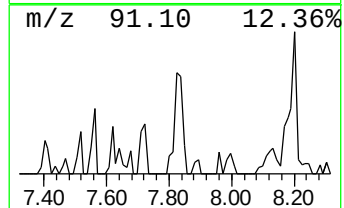
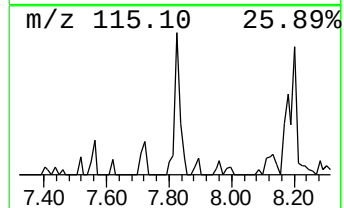
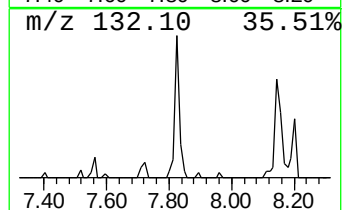
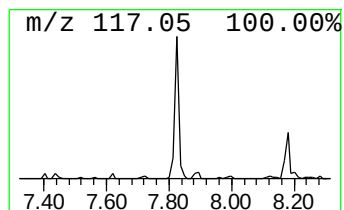
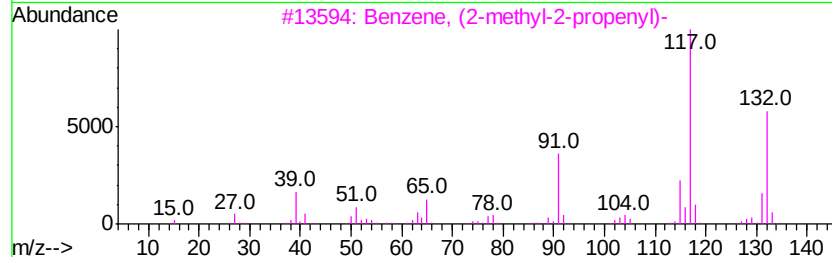
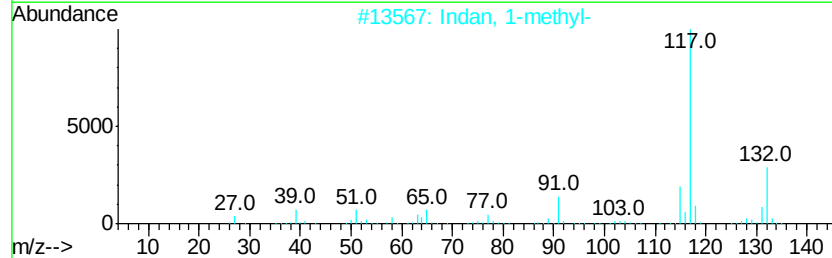
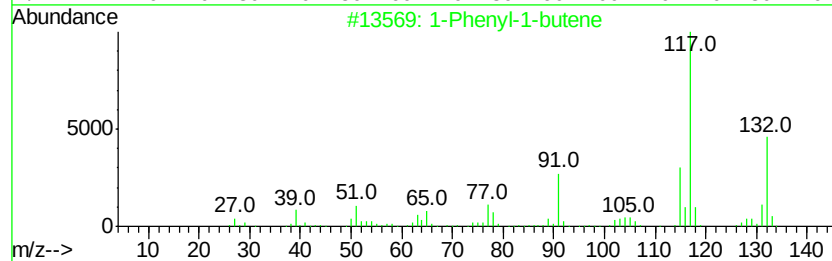
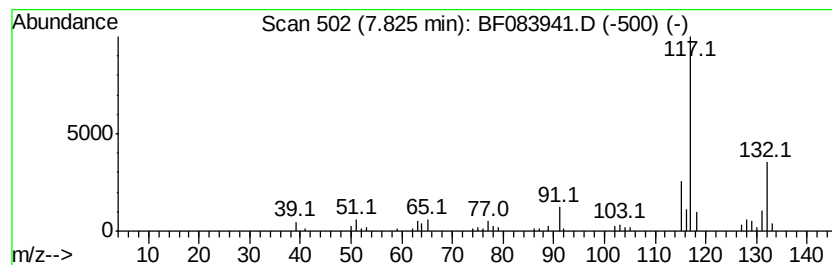
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1-Phenyl-1-butene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.82	6.16 ng	135516	Napthalene-d8	8.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenyl-1-butene	132	C10H12	000824-90-8	94
2		Indan, 1-methyl-	132	C10H12	000767-58-8	91
3		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	91
4		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	91
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121915\
Data File : BF083941.D
Acq On : 19 Dec 2015 4:43
Operator : UM/SJ
Sample : G4840-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

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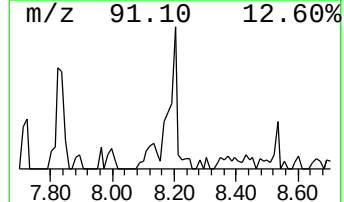
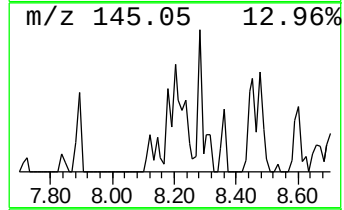
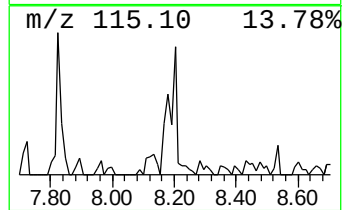
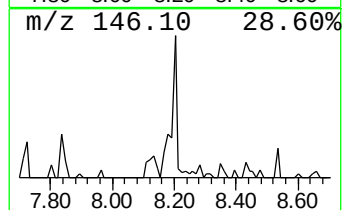
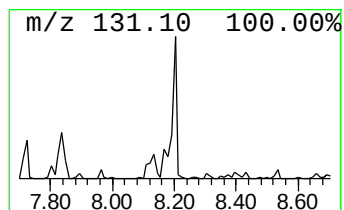
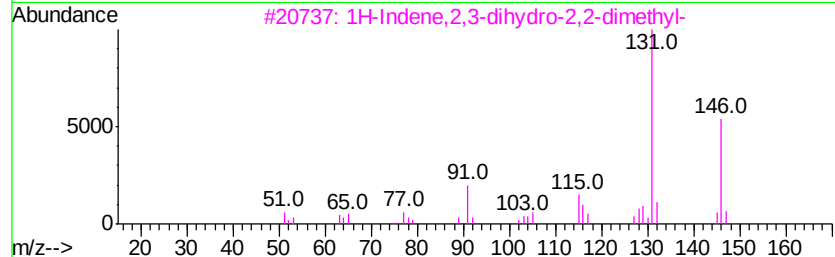
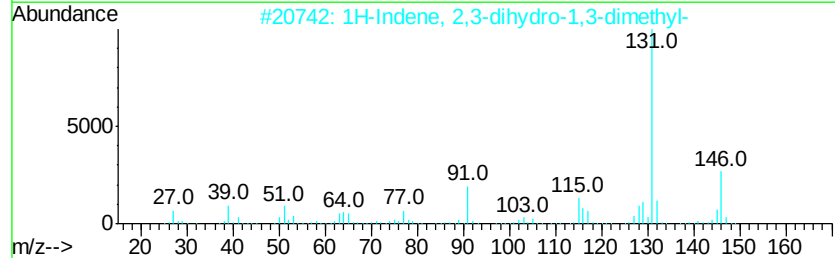
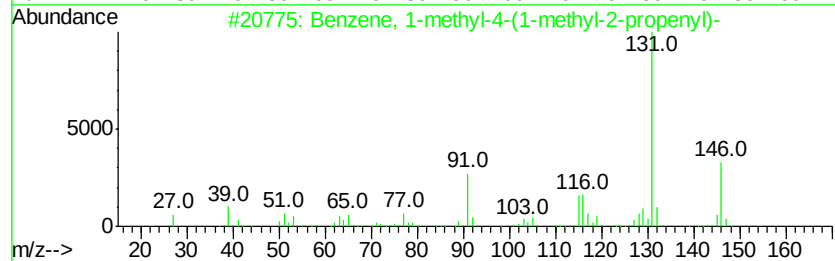
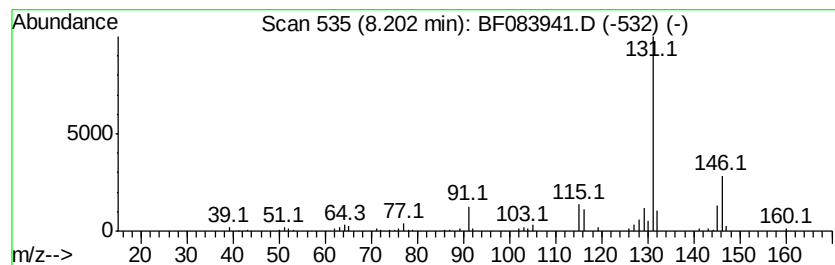
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzene, 1-methyl-4-(1-meth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.20	7.25 ng	159507	Napthalene-d8	8.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	91
2		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
3		1H-Indene, 2,3-dihydro-2,2-dimethyl-	146	C11H14	020836-11-7	91
4		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
5		Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121915\
Data File : BF083941.D
Acq On : 19 Dec 2015 4:43
Operator : UM/SJ
Sample : G4840-06
Misc :
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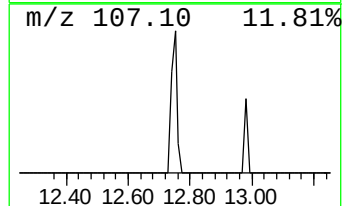
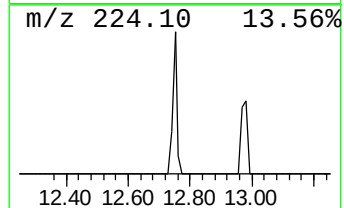
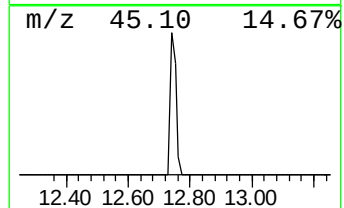
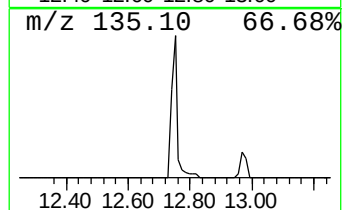
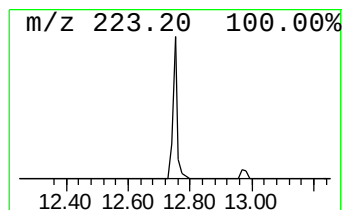
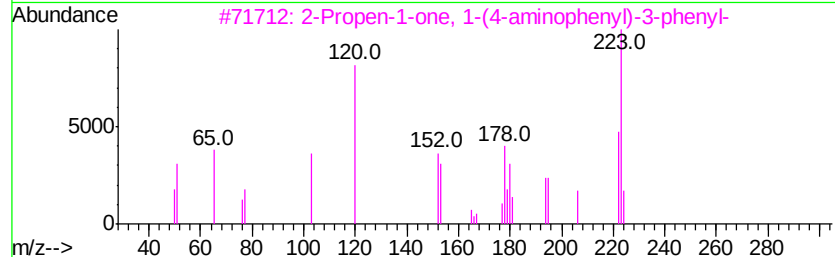
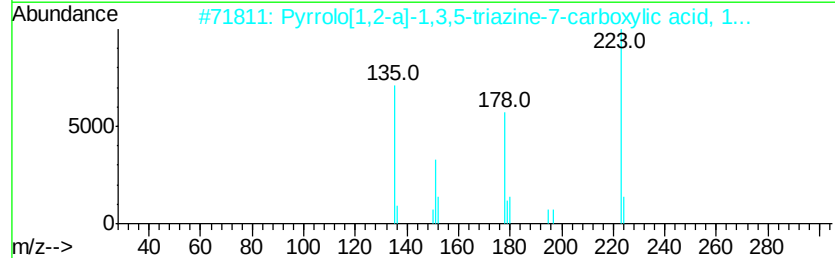
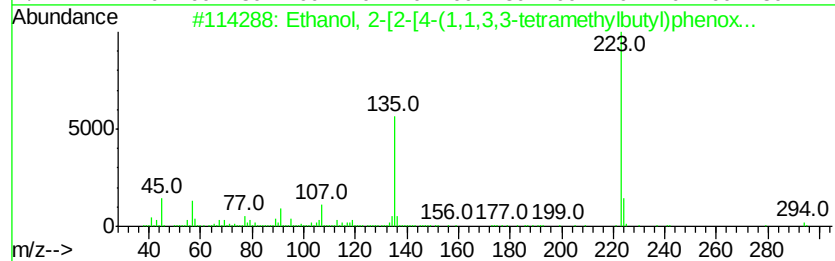
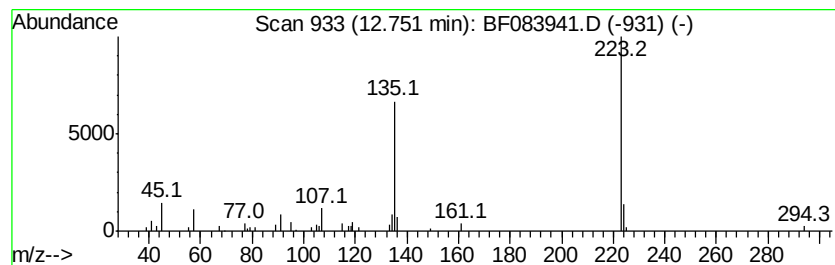
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Ethanol, 2-[2-[4-(1,1,3,3-t... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.75	3.06 ng	72608	Chrysene-d12	14.02		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol,	2-[2-[4-(1,1,3,3-tetram...	294	C18H30O3	002315-61-9	98
2	Pyrrolo[1,2-a]-1,3,5-triazine-7-	...	223	C9H9N3O4	054449-89-7	90
3	2-Propen-1-one,	1-(4-aminophenyl...	223	C15H13NO	002403-30-7	38
4	Tetrazole,	1-(4-ethylphenyl)-5-(...	410	C25H26N6	125038-06-4	37
5	4,4'-Trimethylenebis(1-methylpip...		238	C15H30N2	064168-11-2	32



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121915\
Data File : BF083941.D
Acq On : 19 Dec 2015 4:43
Operator : UM/SJ
Sample : G4840-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
W-48

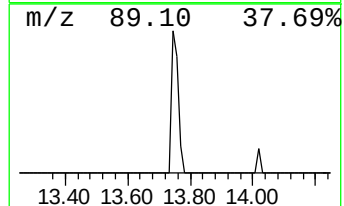
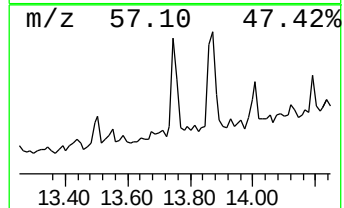
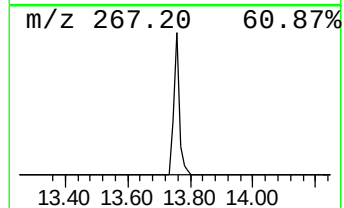
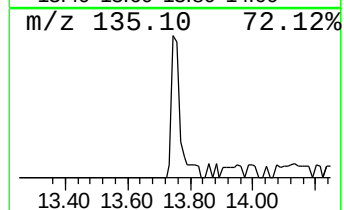
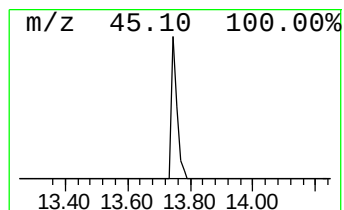
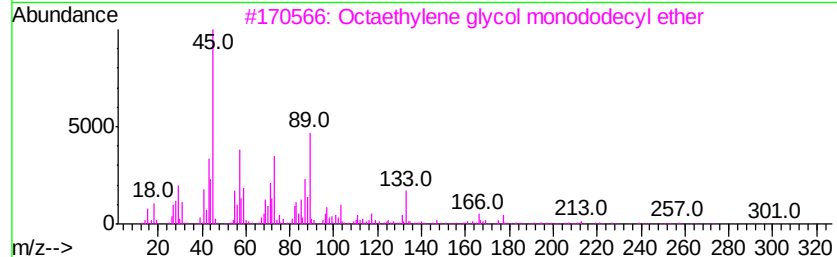
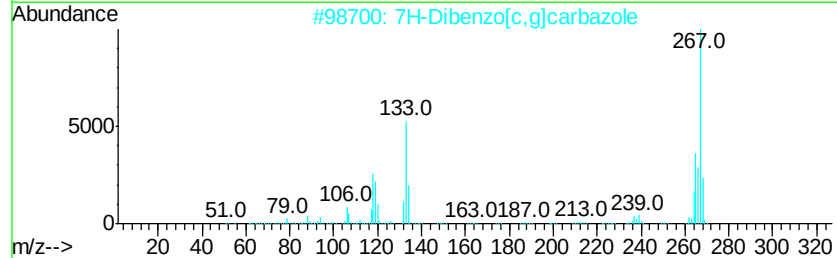
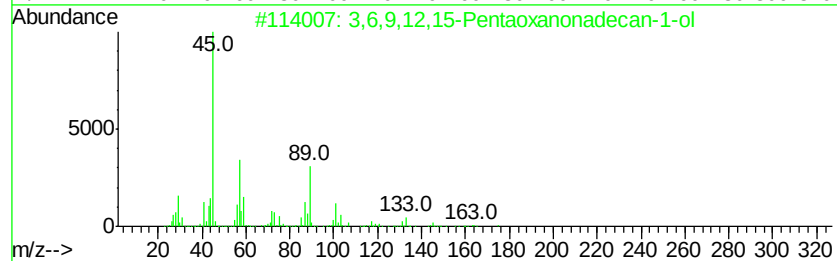
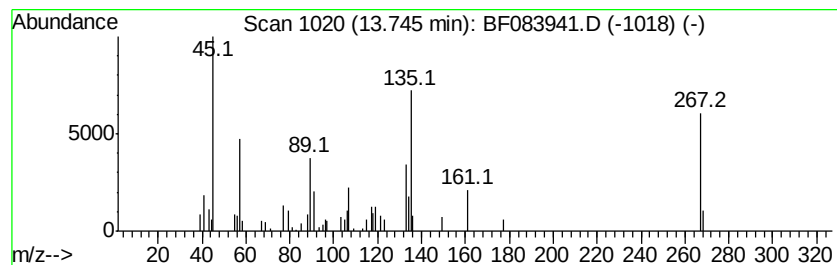
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown13.75 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.75	2.46 ng	58203	Chrysene-d12	14.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,6,9,12,15-Pentaoxanonadecan-1-ol	294	C14H30O6	001786-94-3	35		
2	7H-Dibenzo[c,q]carbazole	267	C20H13N	000194-59-2	32		
3	Octaethylene glycol monododecyl ...	538	C28H58O9	003055-98-9	25		
4	Ethanol, 2-[2-[2-[4-(1,1,3,3-tet...	338	C20H34O4	002315-62-0	25		
5	Phenol, 4,4'-(1,2-diethyl-1,2-et...	270	C18H22O2	000084-16-2	22		



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121915\
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Operator : UM/SJ
Sample : G4840-06
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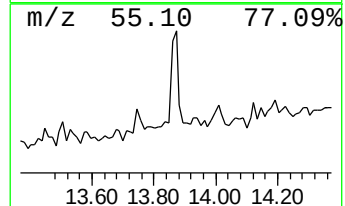
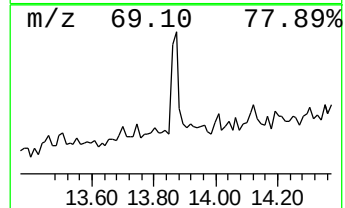
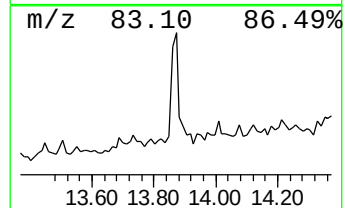
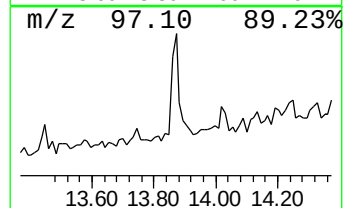
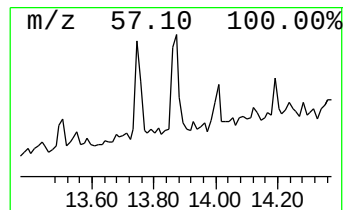
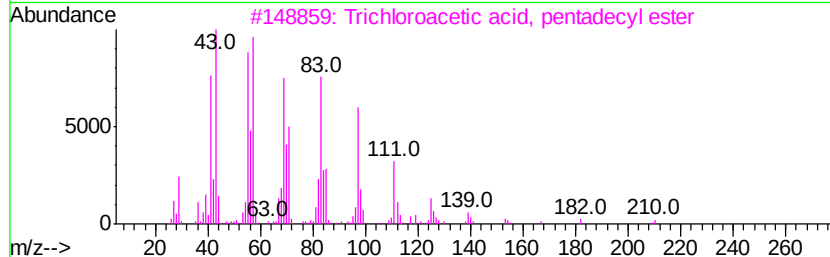
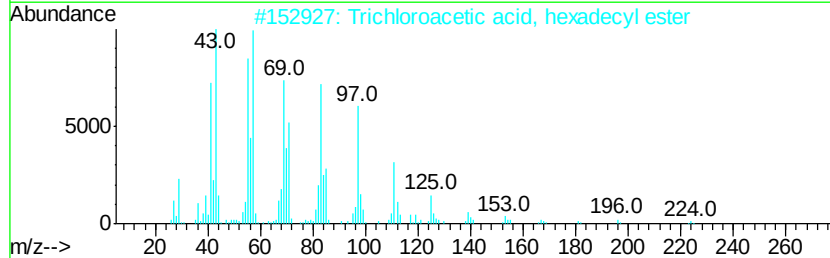
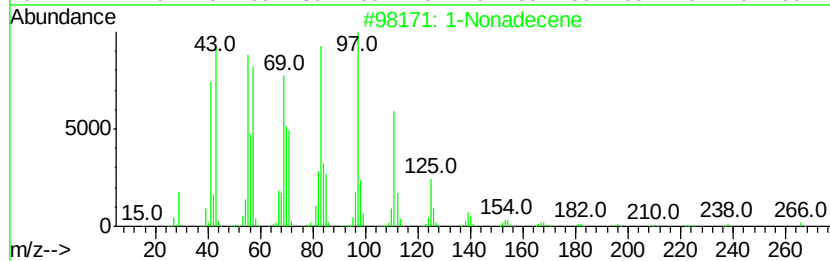
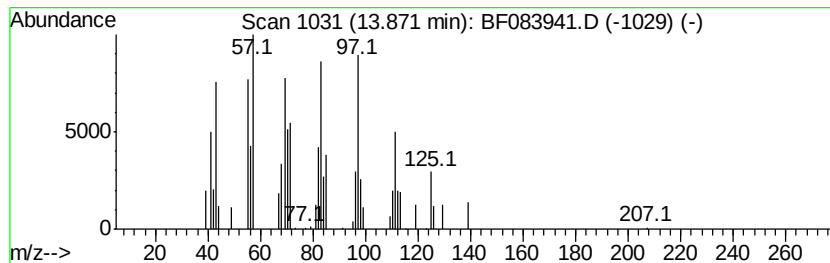
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-Nonadecene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.87	2.12 ng	50281	Chrysene-d12	14.02	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene	266	C19H38	018435-45-5	90
2	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	90
3	Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	87
4	Trichloroacetic acid, tetradecyl...	358	C16H29Cl3O2	074339-52-9	87
5	17-Pentatriacontene	491	C35H70	006971-40-0	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121915\
Data File : BF083941.D
Acq On : 19 Dec 2015 4:43
Operator : UM/SJ
Sample : G4840-06
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121815.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
Cyclohexene, 3-me...	2.37	2.9	ng	50436	1	6.86	347880	20.0
Cyclohexene, 1-me...	3.89	2.2	ng	38586	1	6.86	347880	20.0
Bicyclo[3.1.0]hex...	4.83	2.1	ng	36072	1	6.86	347880	20.0
2-Pentanone, 4-hy...	5.05	6.0	ng	103581	1	6.86	347880	20.0
unknown6.64	6.64	70.1	ng	1219720	1	6.86	347880	20.0
Benzene, 1,4-diet...	7.18	2.4	ng	41656	1	6.86	347880	20.0
1-Phenyl-1-butene	7.82	6.2	ng	135516	2	8.14	440165	20.0
Benzene, 1-methyl...	8.20	7.3	ng	159507	2	8.14	440165	20.0
Ethanol, 2-[2-[4-...	12.75	3.1	ng	72608	5	14.02	473859	20.0
unknown13.75	13.75	2.5	ng	58203	5	14.02	473859	20.0
1-Nonadecene	13.87	2.1	ng	50281	5	14.02	473859	20.0