

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030915\
 Data File : BF077650.D
 Acq On : 9 Mar 2015 20:58
 Operator : TP/IZ
 Sample : G1453-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SEP4

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: CENTROID

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.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	1.344	37	40	43	rVB	4374270	4564109	100.00%	24.614%
2	1.470	48	51	60	rVB	115395	211447	4.63%	1.140%
3	1.641	64	66	69	rVB	55963	67387	1.48%	0.363%
4	1.687	69	70	73	rBV	61355	80082	1.75%	0.432%
5	1.801	79	80	101	rVB	596422	1053316	23.08%	5.681%
6	4.956	354	356	361	rBV	106330	123115	2.70%	0.664%
7	5.481	399	402	405	rBV	753362	777517	17.04%	4.193%
8	6.750	511	513	516	rBV	628544	785840	17.22%	4.238%
9	6.899	524	526	529	rBV	893933	838395	18.37%	4.522%
10	7.196	549	552	554	rBV	230562	307085	6.73%	1.656%
11	7.379	565	568	570	rBV	709404	663646	14.54%	3.579%
12	7.882	610	612	614	rBV	590267	508061	11.13%	2.740%
13	8.762	687	689	692	rBV	322787	429133	9.40%	2.314%
14	10.110	805	807	810	rBV	1073491	992545	21.75%	5.353%
15	10.933	877	879	882	rBV	444882	507696	11.12%	2.738%
16	11.916	962	965	967	rBV	706051	725199	15.89%	3.911%
17	12.774	1037	1040	1042	rBV	623424	706145	15.47%	3.808%
18	12.808	1042	1043	1045	rVV	245281	180179	3.95%	0.972%
19	12.865	1045	1048	1050	rVB	56189	61398	1.35%	0.331%
20	13.402	1093	1095	1097	rBV	90266	107980	2.37%	0.582%
21	13.436	1097	1098	1101	rVV	112750	101168	2.22%	0.546%
22	13.494	1101	1103	1105	rVV	40398	52781	1.16%	0.285%
23	13.539	1105	1107	1108	rVV2	113394	153148	3.36%	0.826%
24	13.562	1108	1109	1113	rVB	60424	86328	1.89%	0.466%
25	13.779	1126	1128	1130	rBV	41697	54108	1.19%	0.292%
26	14.088	1153	1155	1157	rBV	43228	60770	1.33%	0.328%
27	14.282	1170	1172	1174	rBV	165324	192156	4.21%	1.036%
28	14.557	1194	1196	1198	rVB	308246	324288	7.11%	1.749%
29	14.591	1198	1199	1208	rBV4	18312	78404	1.72%	0.423%
30	14.774	1212	1215	1217	rVV	1102939	1204714	26.40%	6.497%
31	14.842	1218	1221	1224	rBV2	40670	76236	1.67%	0.411%
32	14.979	1231	1233	1237	rVB	79147	95445	2.09%	0.515%
33	15.059	1238	1240	1242	rVB	51597	51121	1.12%	0.276%
34	15.105	1242	1244	1245	rBV	53348	60566	1.33%	0.327%

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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: CENTROID

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
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35	15.220	1252	1254	1255	rBV	31330	46871	1.03%	0.253%
36	15.940	1315	1317	1321	rBV2	125559	176401	3.86%	0.951%
37	16.065	1325	1328	1329	rVV2	478128	781365	17.12%	4.214%
38	16.088	1329	1330	1332	rVB	120980	120137	2.63%	0.648%
39	16.717	1383	1385	1388	rBV2	56387	106596	2.34%	0.575%
40	17.357	1439	1441	1445	rBV	71646	174780	3.83%	0.943%
41	17.677	1467	1469	1473	rVB	51431	70801	1.55%	0.382%
42	17.745	1473	1475	1477	rBV	78606	102900	2.25%	0.555%
43	17.803	1478	1480	1483	rBV	428167	621250	13.61%	3.350%
44	19.620	1636	1639	1642	rVB	30119	59774	1.31%	0.322%

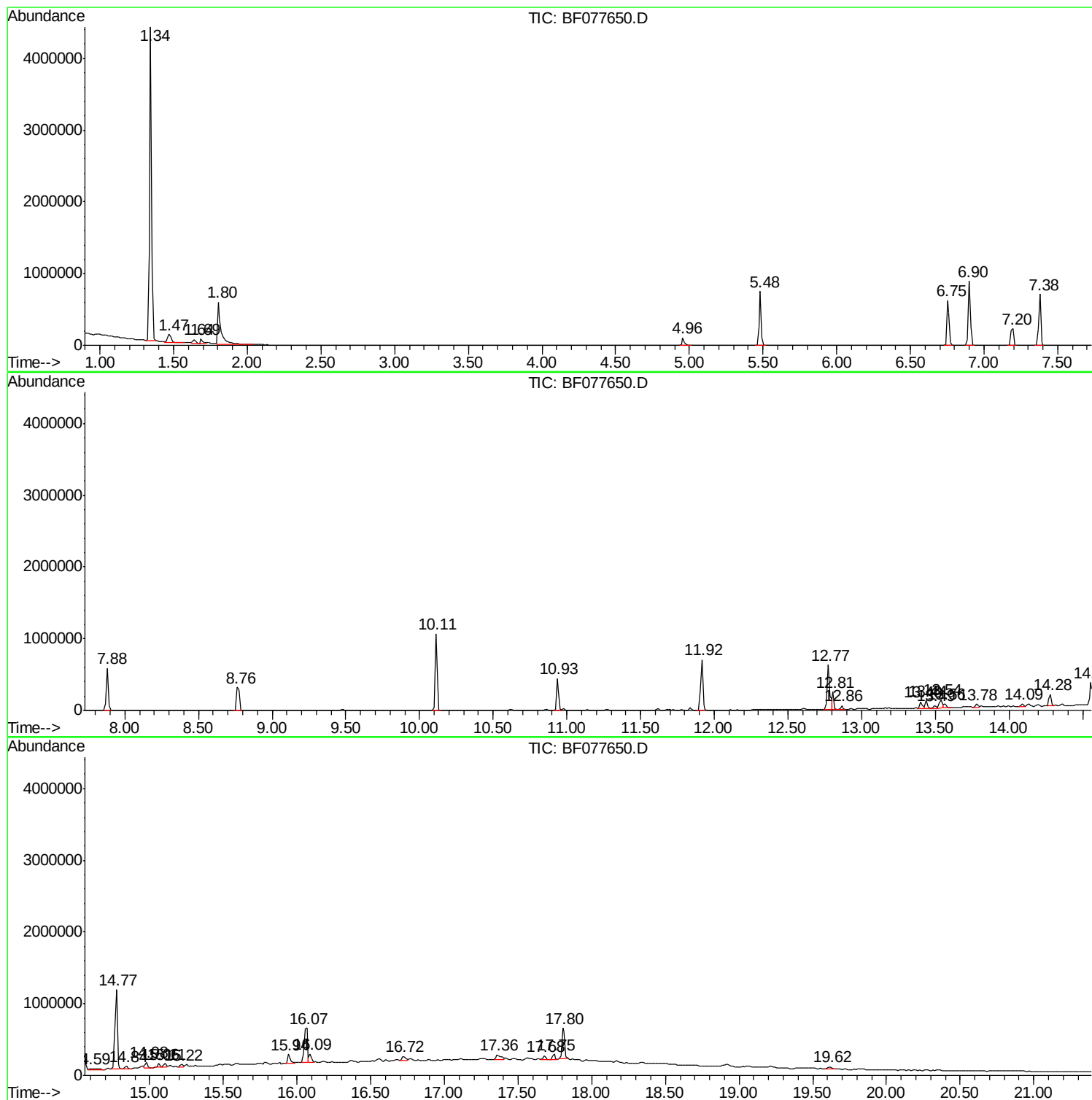
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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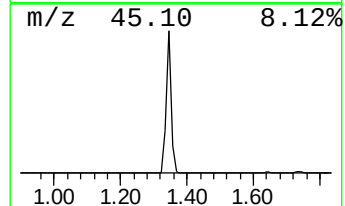
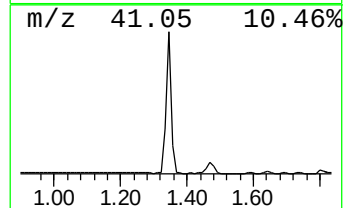
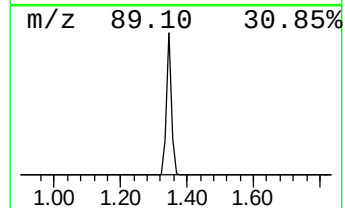
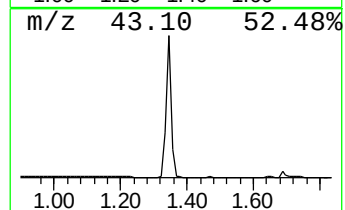
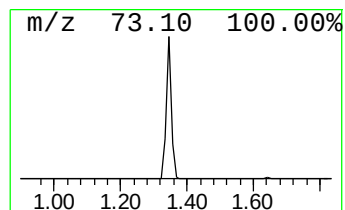
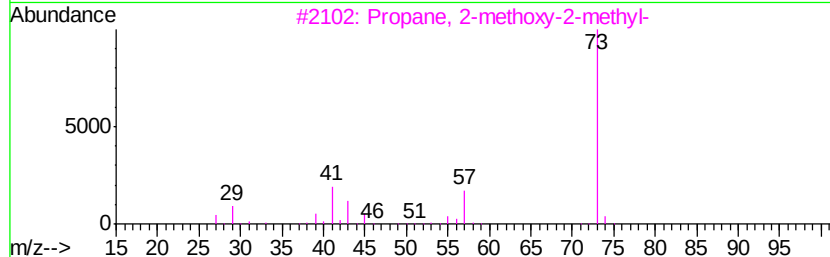
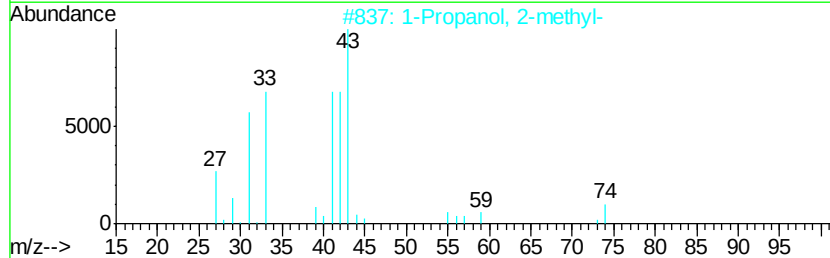
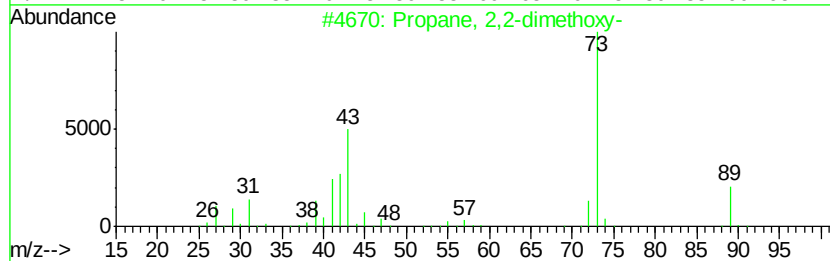
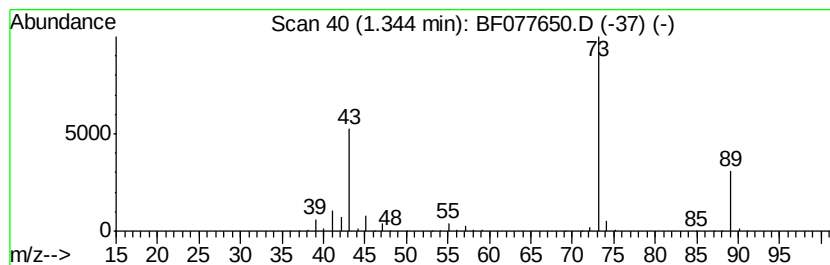
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown1.34 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.34	297.25 ng	4564110	1,4-Dichlorobenzene-d4	7.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	38
2		1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	9
3		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
4		1-Hexen-4-ol, 1-chloro-3,5-dimet...	162	C8H15ClO	100244-09-5	9
5		1,3-Dioxolane, 2-(3-bromo-5,5,5-...	352	C10H16BrCl3O2	1000115-31-4	9



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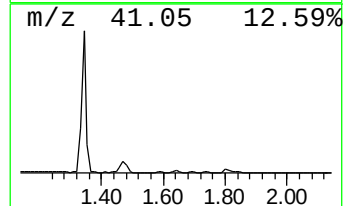
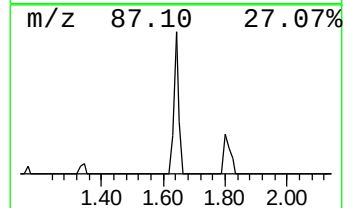
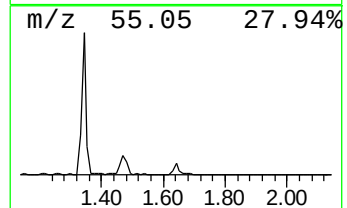
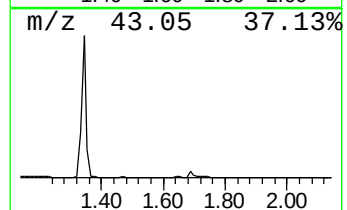
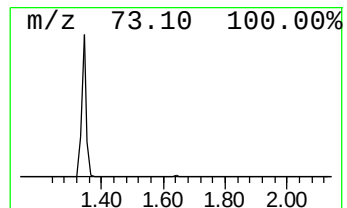
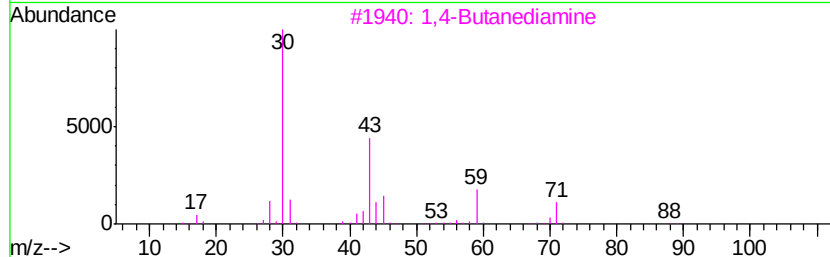
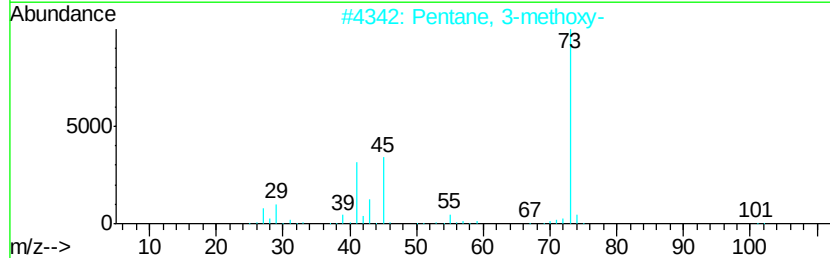
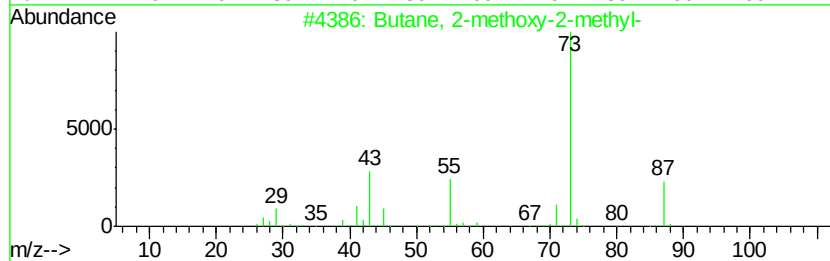
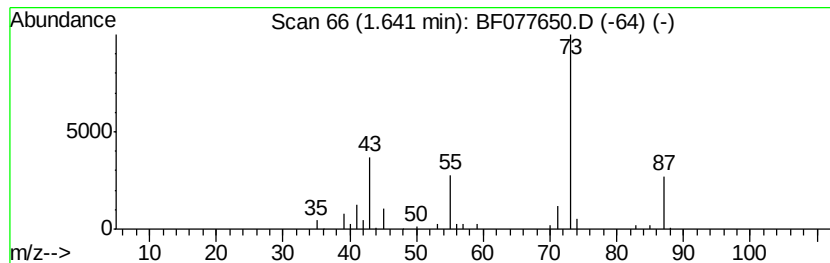
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.64	4.39 ng	67387	1,4-Dichlorobenzene-d4	7.19

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
3			1,4-Butanediamine	88	C4H12N2	000110-60-1	10
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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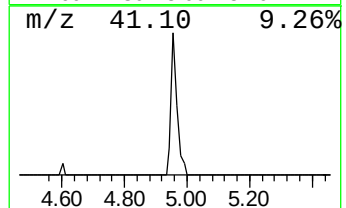
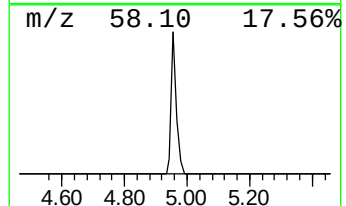
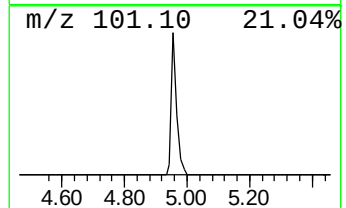
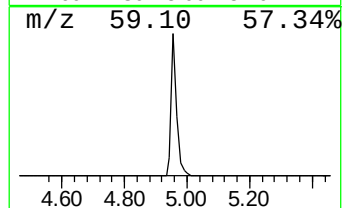
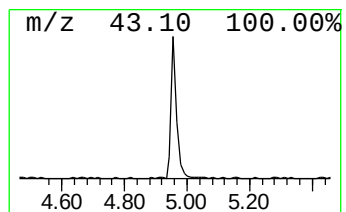
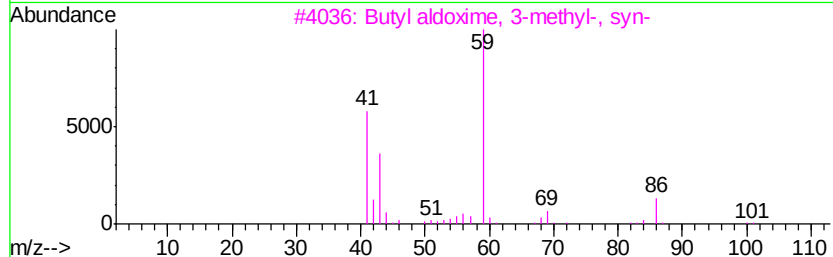
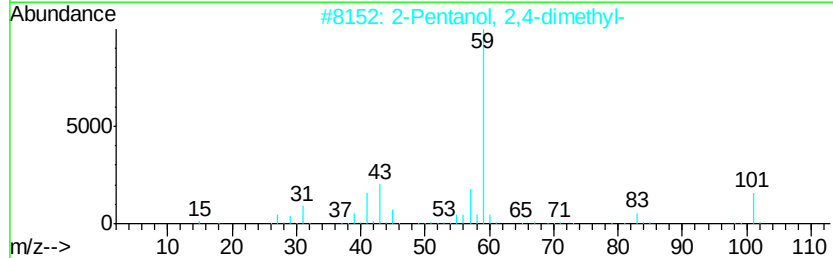
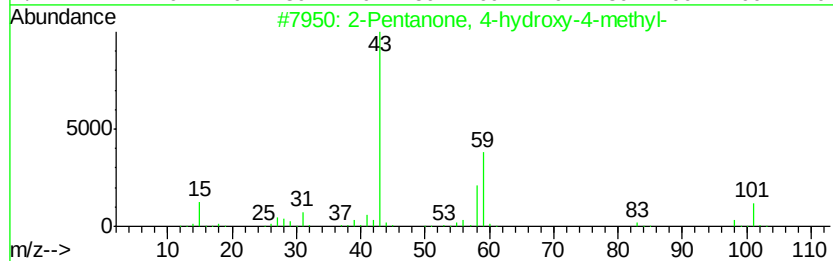
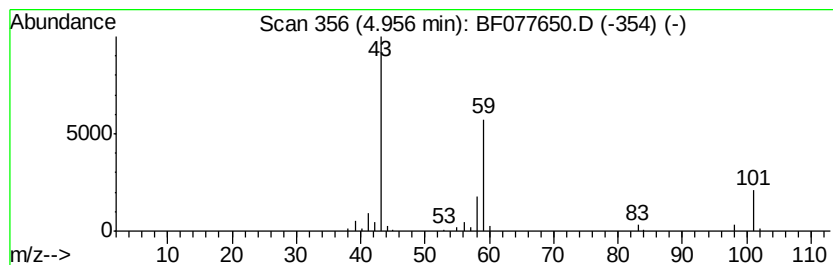
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	8.02 ng	123115	1,4-Dichlorobenzene-d4	7.19

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	42
3	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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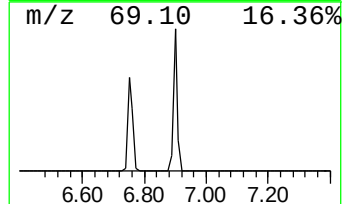
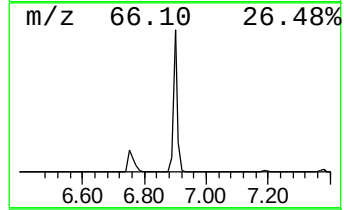
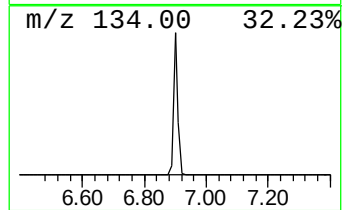
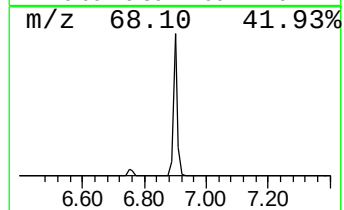
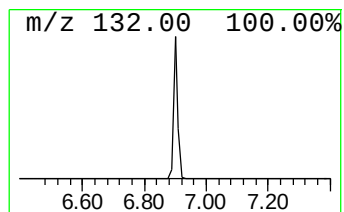
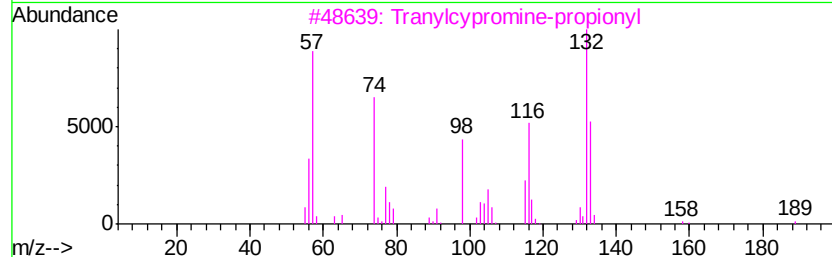
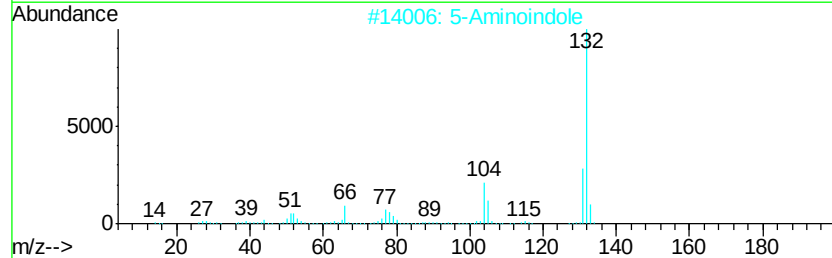
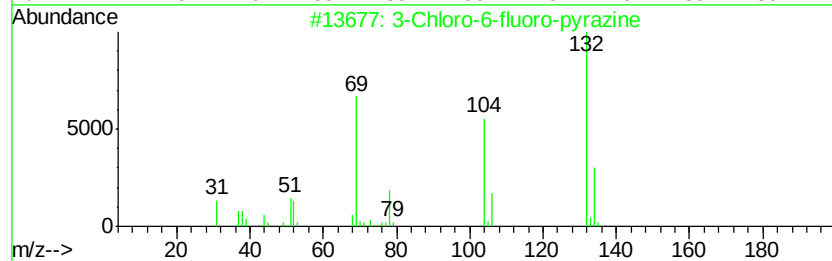
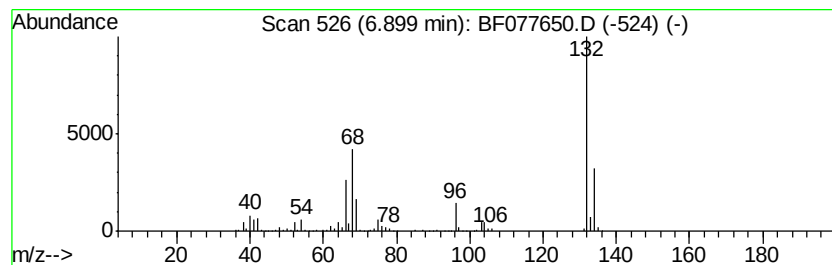
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 7 unknown6.90 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.90	54.60 ng	838395	1,4-Dichlorobenzene-d4	7.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3	Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2	5	Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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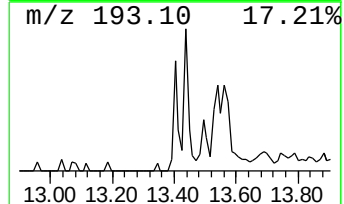
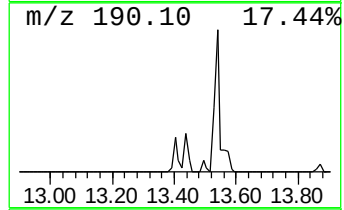
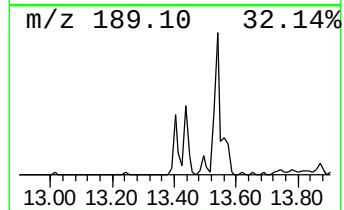
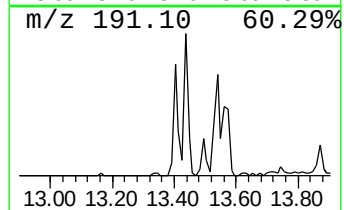
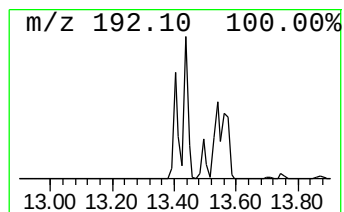
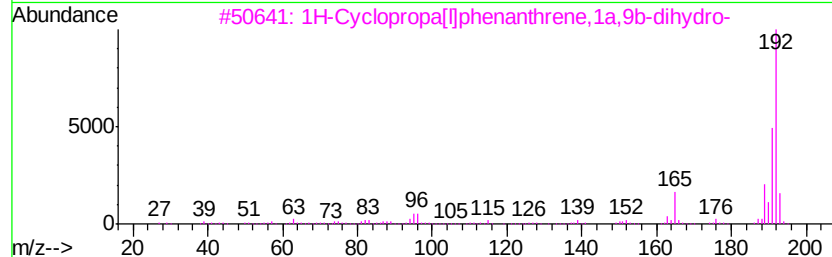
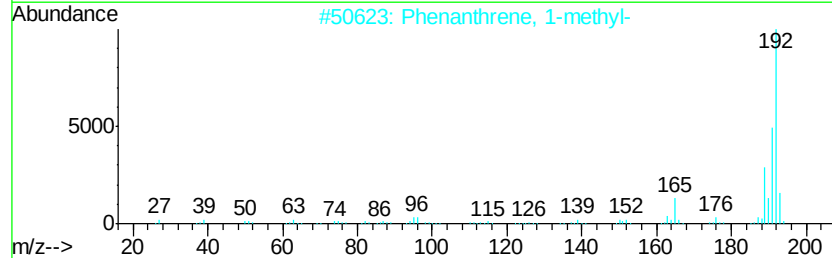
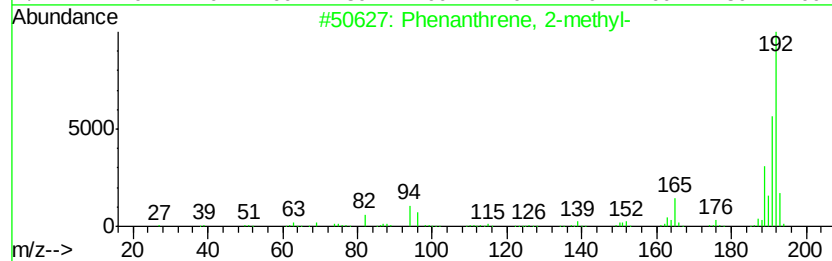
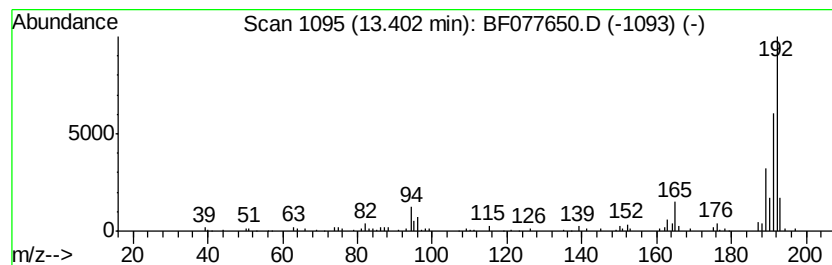
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TIC Integration Parameters: LSCINT.P

Peak Number 9 Phenanthrene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.40	3.06 ng	107980	Phenanthrene-d10	12.78	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	94
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030915\
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Acq On : 9 Mar 2015 20:58
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Sample : G1453-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

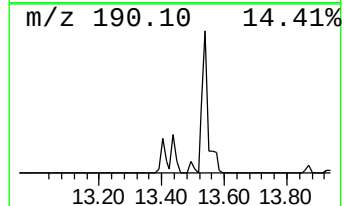
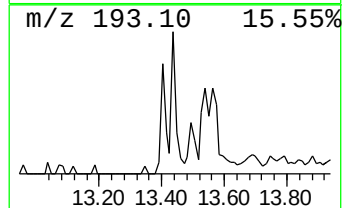
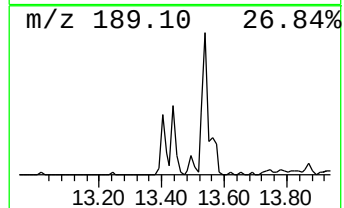
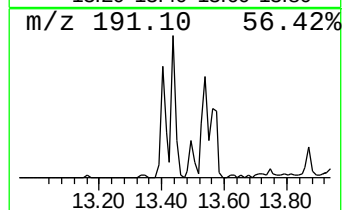
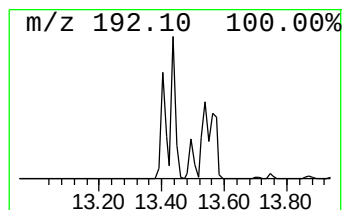
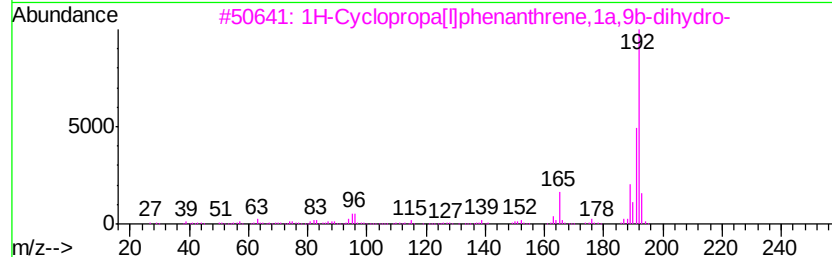
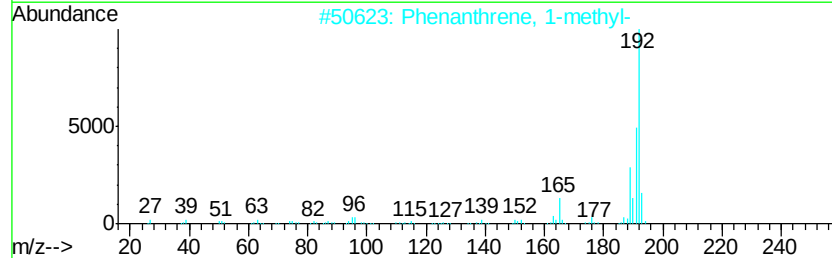
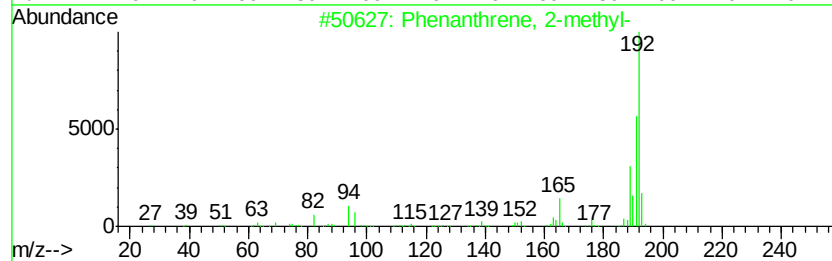
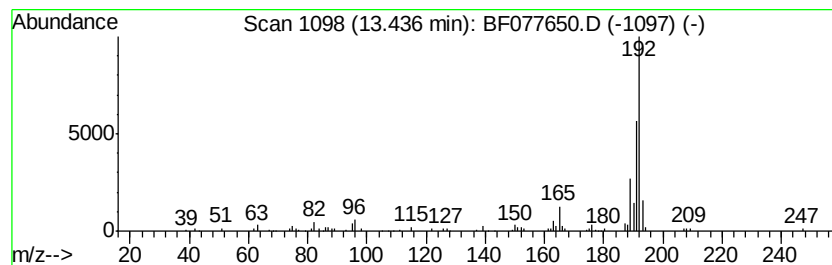
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Phenanthrene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.44	2.87 ng	101168	Phenanthrene-d10	12.78	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3	1H-Cyclopropa[1,1a]phenanthrene, 1a, ...	192	C15H12	000949-41-7	95
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	94



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030915\
Data File : BF077650.D
Acq On : 9 Mar 2015 20:58
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Sample : G1453-03
Misc :
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Instrument :
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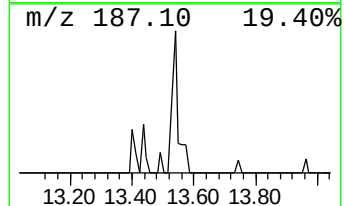
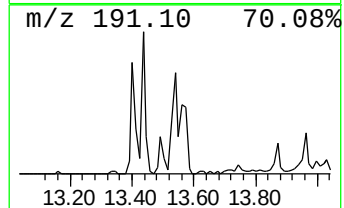
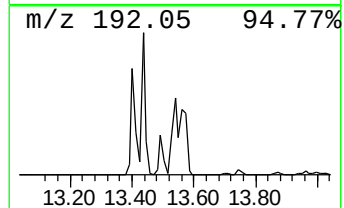
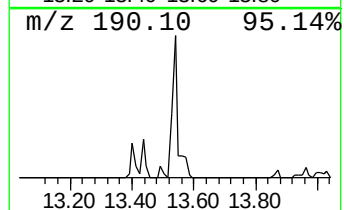
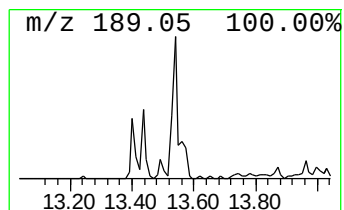
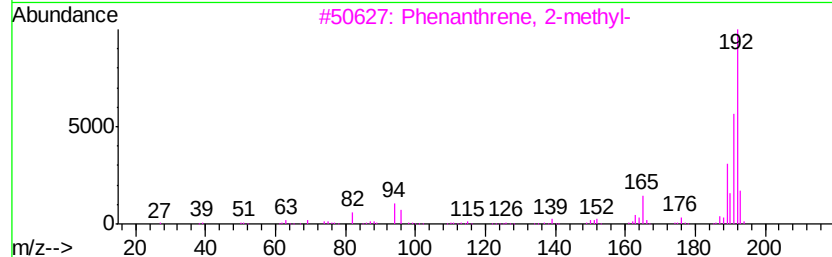
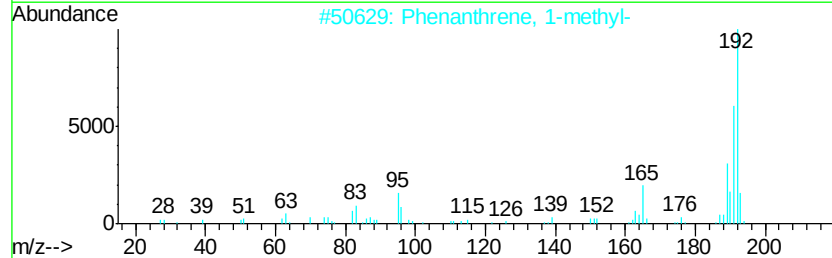
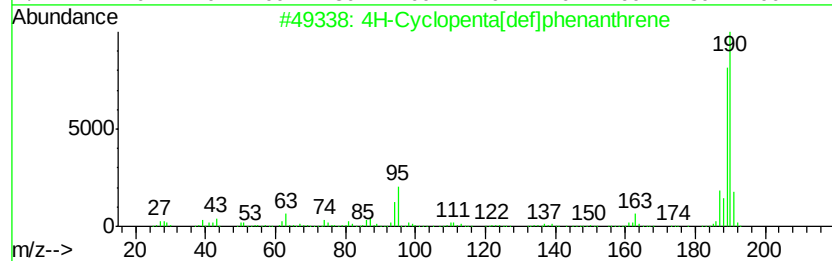
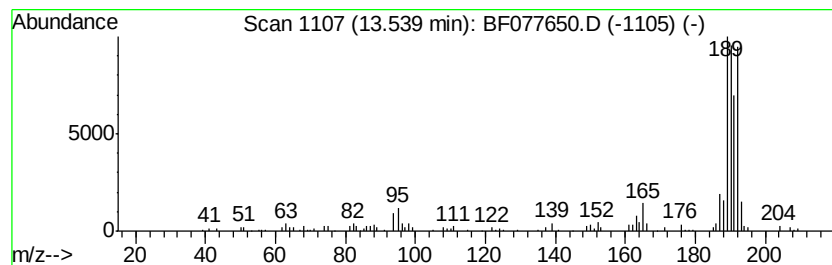
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 4H-Cyclopenta[def]phenanthrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.54	4.34 ng	153148	Phenanthrene-d10	12.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	55
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	46
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	46
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	46
5			Anthracene, 9-methyl-	192	C15H12	000779-02-2	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF030915\
Data File : BF077650.D
Acq On : 9 Mar 2015 20:58
Operator : TP/IZ
Sample : G1453-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

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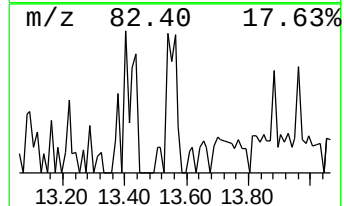
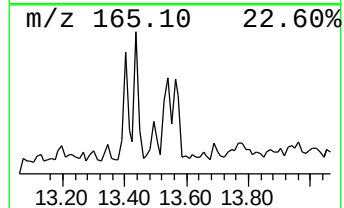
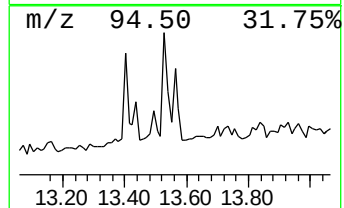
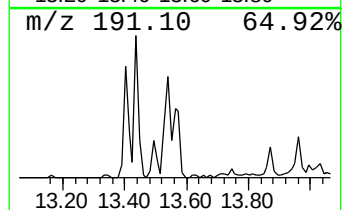
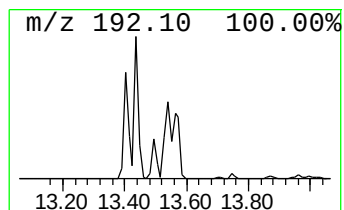
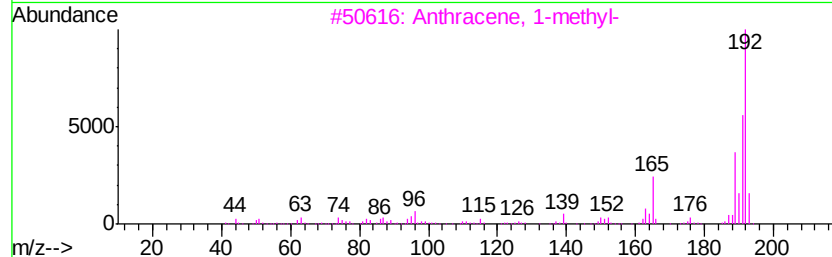
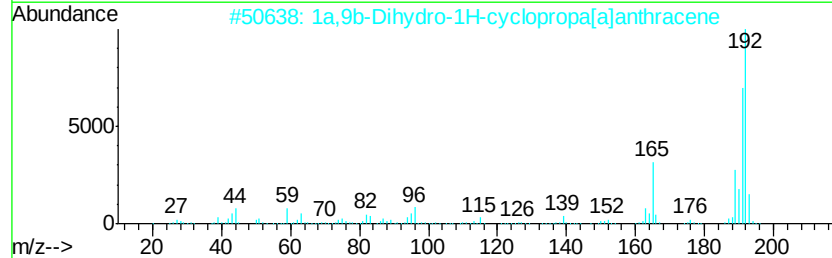
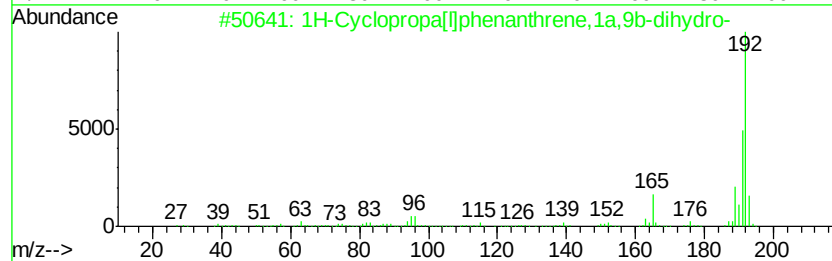
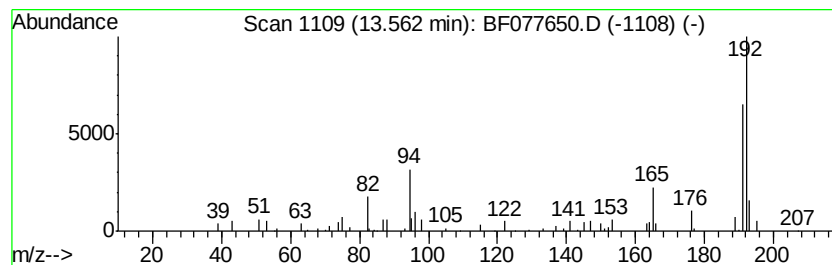
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1H-Cyclopropa[l]phenanthren... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.56	2.45 ng	86328	Phenanthrene-d10	12.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Cyclopropa[l]phenanthrene, 1a,...	192	C15H12	000949-41-7	81
2		1a,9b-Dihydro-1H-cyclopropa[al]an...	192	C15H12	006109-24-6	74
3		Anthracene, 1-methyl-	192	C15H12	000610-48-0	74
4		Naphtho[2,3-b]norborene	192	C15H12	107426-38-0	74
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030915\
Data File : BF077650.D
Acq On : 9 Mar 2015 20:58
Operator : TP/IZ
Sample : G1453-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

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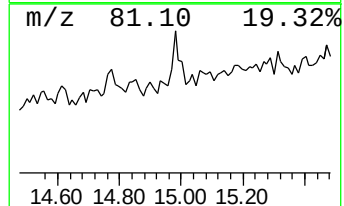
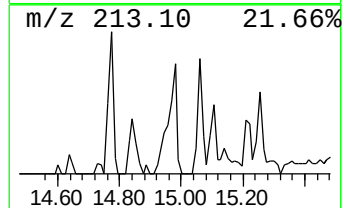
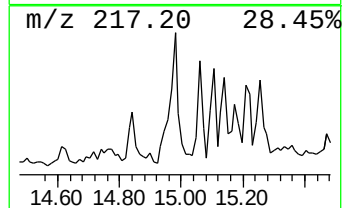
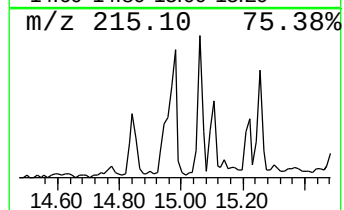
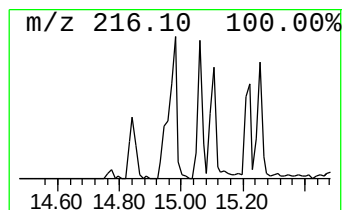
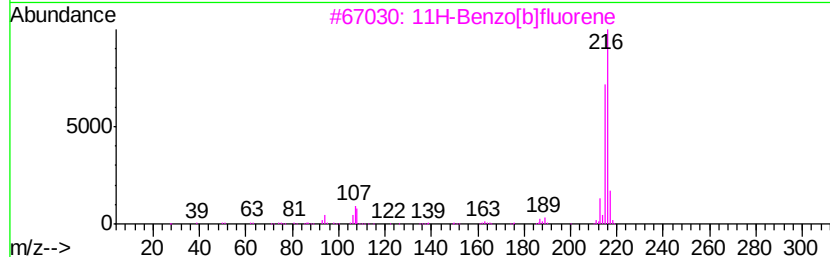
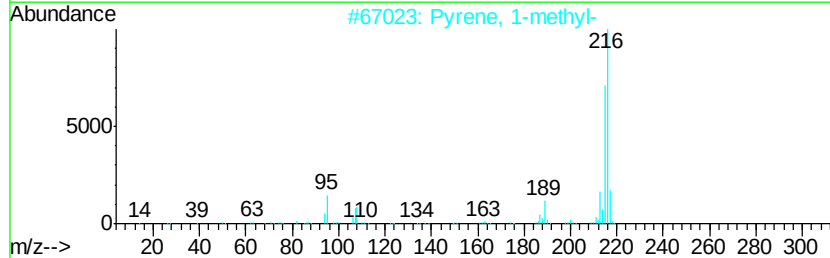
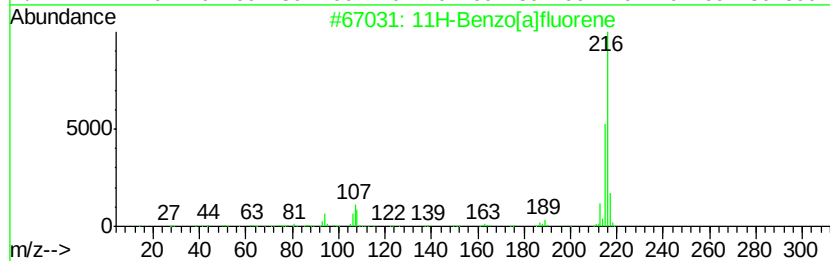
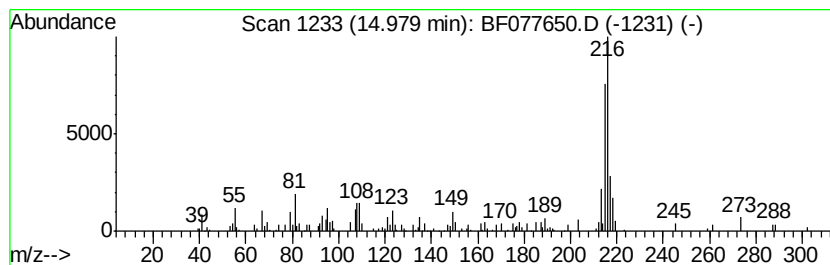
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 11H-Benzo[a]fluorene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.98	2.44 ng	95445	Chrysene-d12	16.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	81
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	80
3		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	74
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	68
5		7H-Benzo[c]fluorene	216	C17H12	000205-12-9	62



Data Path : Z:\HPCHEM1\BNA F\DATA\BF030915\
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Sample : G1453-03
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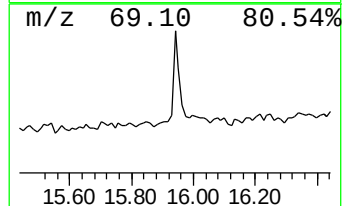
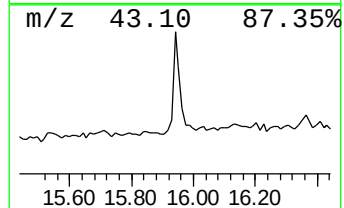
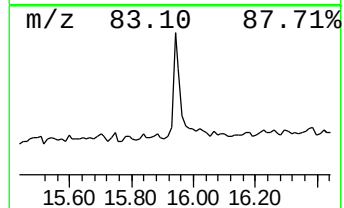
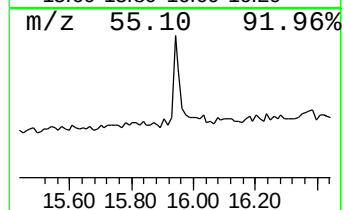
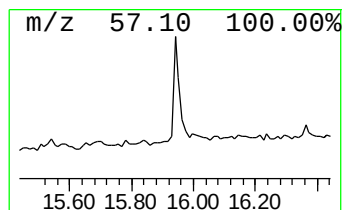
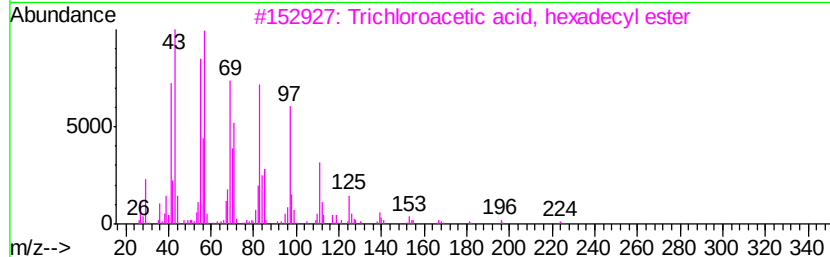
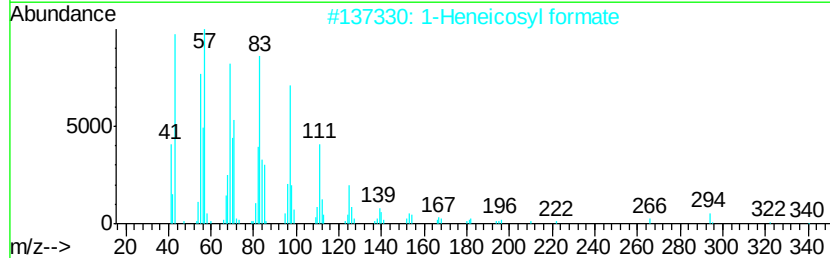
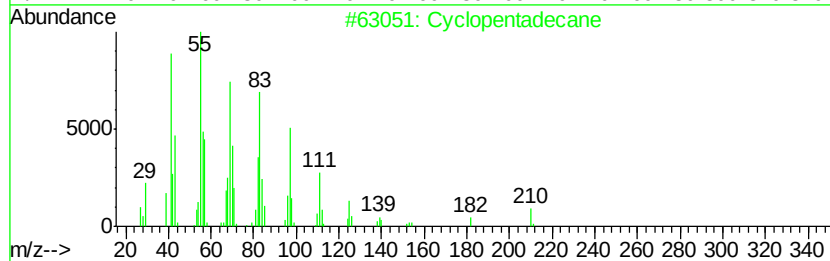
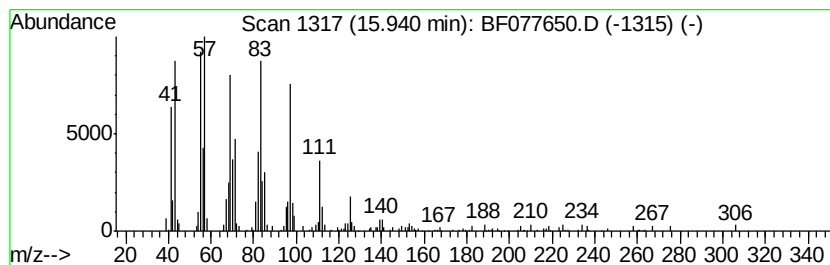
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Cyclopentadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.94	4.52 ng	176401	Chrysene-d12	16.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclopentadecane	210 C15H30	000295-48-7 97
2		1-Heneicosyl formate	340 C22H44O2	077899-03-7 94
3		Trichloroacetic acid, hexadecyl ...	386 C18H33Cl3O2	074339-54-1 93
4		Heptafluorobutyric acid, n-octad...	466 C22H37F7O2	000400-57-7 91
5		Bromoacetic acid, octadecyl ester	390 C20H39BrO2	018992-03-5 91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030915\
Data File : BF077650.D
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ClientSampleId :
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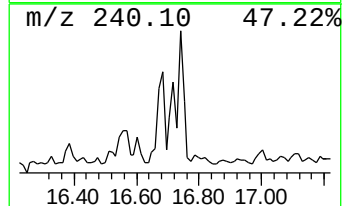
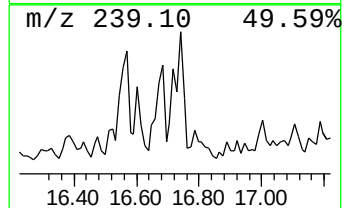
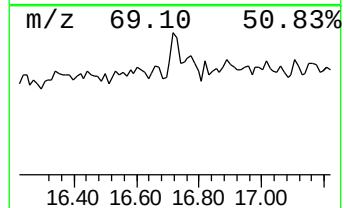
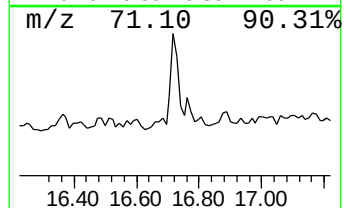
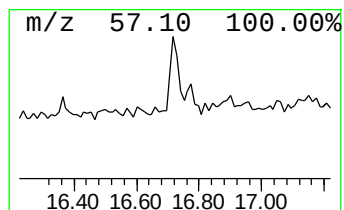
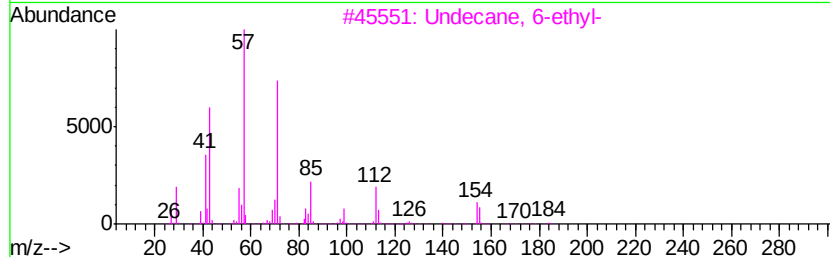
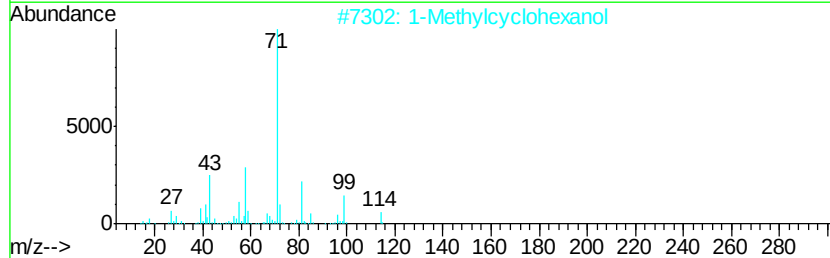
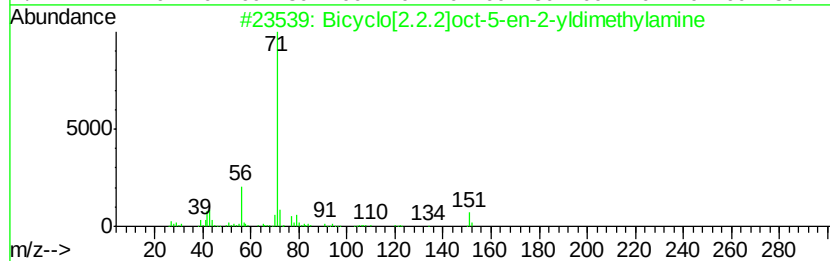
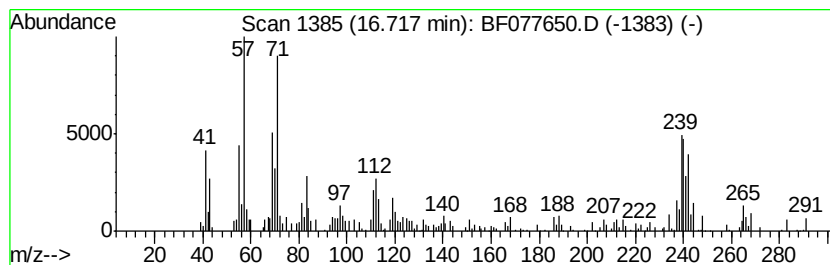
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown16.72 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.72	2.73 ng	106596	Chrysene-d12	16.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[2.2.2]oct-5-en-2-yl dimethylam...	151	C10H17N	033162-94-6	18
2		1-Methylcyclohexanol	114	C7H14O	000590-67-0	18
3		Undecane, 6-ethyl-	184	C13H28	017312-60-6	14
4		Octane, 3-ethyl-	142	C10H22	005881-17-4	14
5		Isophytol	296	C20H40O	000505-32-8	14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030915\
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Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.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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Butane, 2-methoxy...	1.64	4.4	ng	67387	1	7.19	307085	20.0
2-Pentanone, 4-hy...	4.96	8.0	ng	123115	1	7.19	307085	20.0
unknown6.90	6.90	54.6	ng	838395	1	7.19	307085	20.0
Phenanthrene, 2-m...	13.40	3.1	ng	107980	4	12.78	706145	20.0
Phenanthrene, 1-m...	13.44	2.9	ng	101168	4	12.78	706145	20.0
4H-Cyclopenta[def...	13.54	4.3	ng	153148	4	12.78	706145	20.0
1H-Cyclopropa[l]p...	13.56	2.5	ng	86328	4	12.78	706145	20.0
11H-Benzo[a]fluorene	14.98	2.4	ng	95445	5	16.06	781365	20.0
Cyclopentadecane	15.94	4.5	ng	176401	5	16.06	781365	20.0
unknown16.72	16.72	2.7	ng	106596	5	16.06	781365	20.0