

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042915\
 Data File : BF078830.D
 Acq On : 30 Apr 2015 1:21
 Operator : TP/IZ
 Sample : G2002-06
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-03

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_F\METHODS\8270-BF042815.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	1.541	28	31	34	rVB	6445198	7435094	100.00%	19.894%
2	1.678	40	43	51	rVB	212690	445339	5.99%	1.192%
3	1.781	51	52	57	rVV	532328	809877	10.89%	2.167%
4	1.861	57	59	62	rVV	153876	274651	3.69%	0.735%
5	1.918	62	64	99	rVB	2894660	5396454	72.58%	14.439%
6	5.187	348	350	357	rVB	252710	295762	3.98%	0.791%
7	5.679	390	393	395	rBV	1484309	1544785	20.78%	4.133%
8	6.924	499	502	504	rBV	1889853	1694822	22.79%	4.535%
9	7.084	514	516	519	rBV	1362644	1642953	22.10%	4.396%
10	7.382	540	542	544	rBV	495626	430899	5.80%	1.153%
11	7.564	556	558	561	rVB	1131452	1226298	16.49%	3.281%
12	8.067	600	602	605	rBV	1041487	938595	12.62%	2.511%
13	8.959	678	680	682	rBV	689452	592109	7.96%	1.584%
14	10.308	795	798	800	rBV	1870027	1842548	24.78%	4.930%
15	10.811	839	842	844	rVB2	59414	74634	1.00%	0.200%
16	10.959	853	855	858	rVB	133353	129665	1.74%	0.347%
17	11.142	868	871	873	rBV	503380	632769	8.51%	1.693%
18	12.034	947	949	952	rVB	90852	108794	1.46%	0.291%
19	12.114	954	956	959	rBV	1217371	1191361	16.02%	3.188%
20	12.982	1030	1032	1034	rBV	732013	749566	10.08%	2.006%
21	13.016	1034	1035	1037	rVV	139353	101180	1.36%	0.271%
22	13.074	1038	1040	1043	rVB	70300	75356	1.01%	0.202%
23	13.645	1089	1090	1092	rVB	96193	83932	1.13%	0.225%
24	13.691	1092	1094	1097	rBV2	59834	98606	1.33%	0.264%
25	13.748	1097	1099	1100	rBV	130798	134494	1.81%	0.360%
26	13.988	1116	1120	1127	rVB3	35948	90513	1.22%	0.242%
27	14.297	1144	1147	1149	rBV	77333	107360	1.44%	0.287%
28	14.491	1162	1164	1166	rBV	533020	529678	7.12%	1.417%
29	14.674	1177	1180	1186	rBV5	47023	101058	1.36%	0.270%
30	14.777	1186	1189	1191	rBV	625606	678232	9.12%	1.815%
31	14.982	1204	1207	1210	rVB	2221521	2138838	28.77%	5.723%
32	15.051	1210	1213	1218	rBV2	54311	95555	1.29%	0.256%
33	15.188	1220	1225	1228	rVB	105692	214258	2.88%	0.573%
34	15.280	1231	1233	1235	rBV	74706	82564	1.11%	0.221%

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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_F\METHODS\8270-BF042815.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.314	1235	1236	1239	rVB	84549	97927	1.32%	0.262%
36	15.702	1265	1270	1277	rBV6	27415	122048	1.64%	0.327%
37	15.817	1277	1280	1284	rVB2	39108	87616	1.18%	0.234%
38	15.965	1289	1293	1295	rBV2	60834	98261	1.32%	0.263%
39	16.000	1295	1296	1299	rBV2	51183	87506	1.18%	0.234%
40	16.148	1306	1309	1315	rBV3	60258	153926	2.07%	0.412%
41	16.274	1316	1320	1322	rBV2	888975	1333395	17.93%	3.568%
42	16.308	1322	1323	1325	rVB	316560	252443	3.40%	0.675%
43	16.765	1360	1363	1365	rBV	86609	114797	1.54%	0.307%
44	16.948	1375	1379	1383	rVB5	44388	130921	1.76%	0.350%
45	17.531	1427	1430	1438	rBV	289405	696813	9.37%	1.864%
46	17.645	1438	1440	1442	rVB	67903	79803	1.07%	0.214%
47	17.851	1455	1458	1460	rBV	178714	236486	3.18%	0.633%
48	17.908	1460	1463	1467	rVB	274923	345169	4.64%	0.924%
49	17.977	1467	1469	1471	rBV	588442	792366	10.66%	2.120%
50	18.560	1517	1520	1523	rBV3	67862	109635	1.47%	0.293%
51	19.486	1598	1601	1608	rBV2	110720	276684	3.72%	0.740%
52	19.931	1636	1640	1647	rBV	112818	247093	3.32%	0.661%
53	20.149	1655	1659	1665	rVB4	33808	122856	1.65%	0.329%

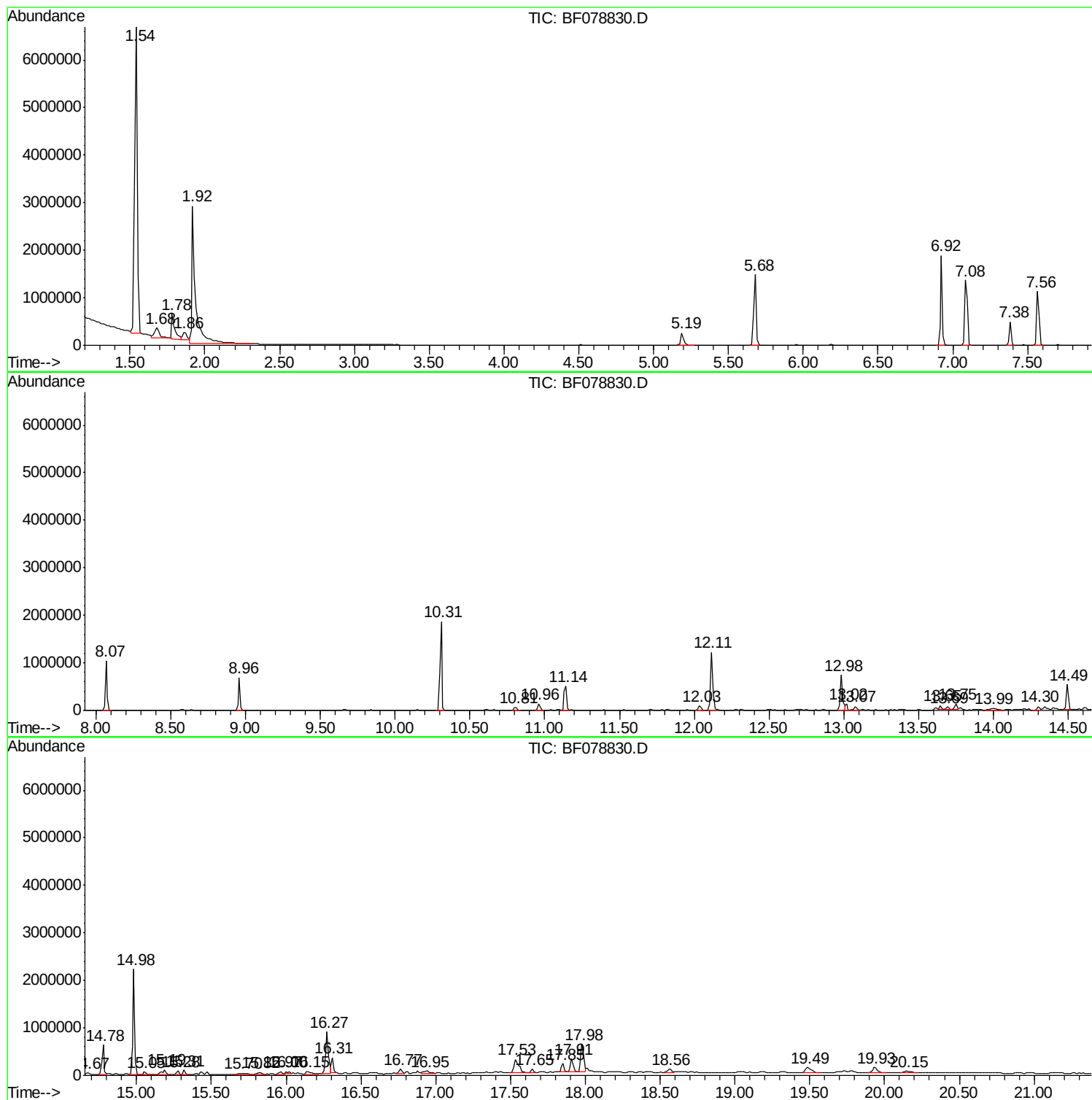
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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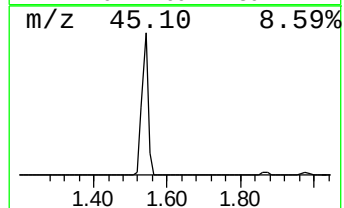
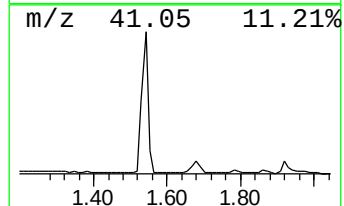
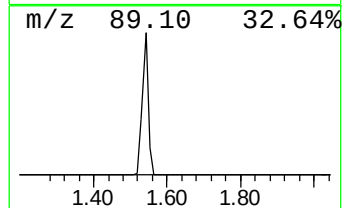
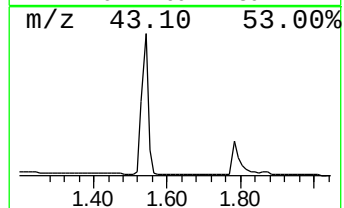
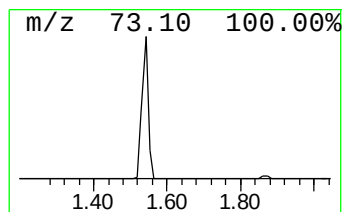
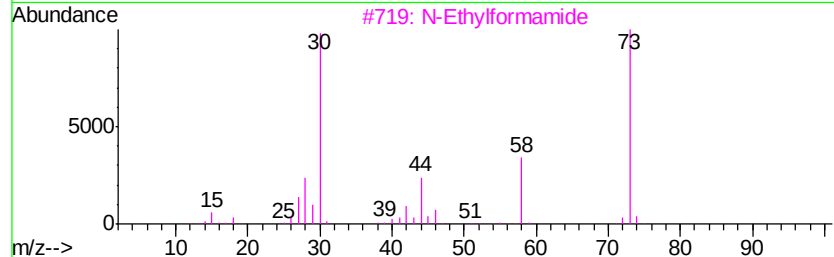
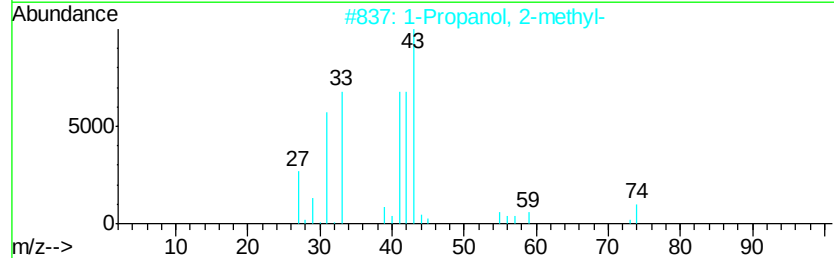
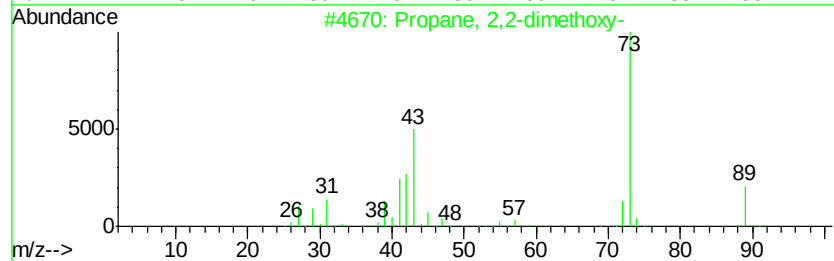
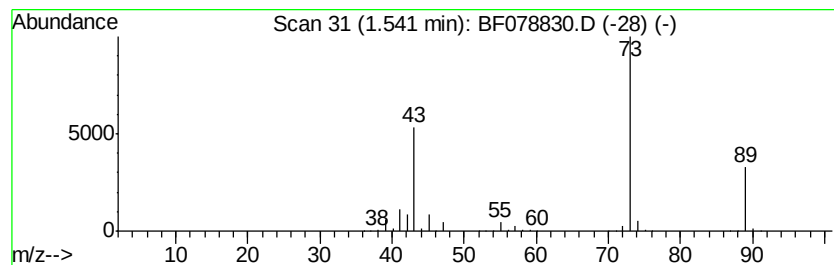
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Propane, 2,2-dimethoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
1.54	345.10 ng	7435090	1,4-Dichlorobenzene-d4	7.38	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	56
2	1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	9
3	N-Ethylformamide	73	C3H7NO	000627-45-2	9
4	1,3-Diaminoquanidine	89	CH7N5	004364-78-7	9
5	2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9



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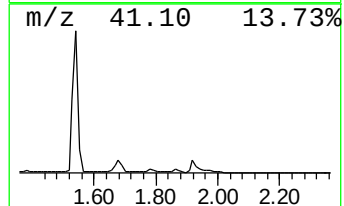
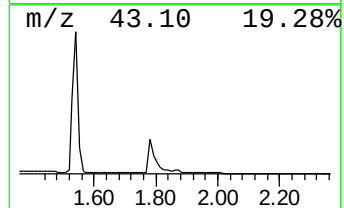
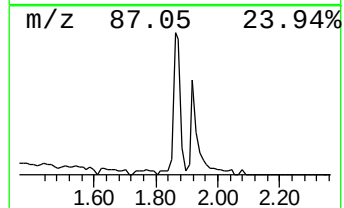
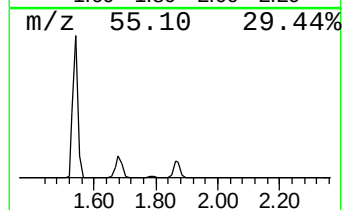
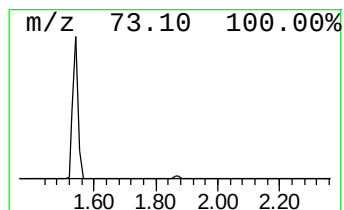
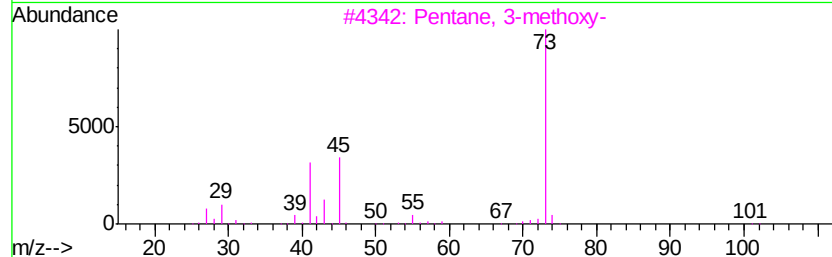
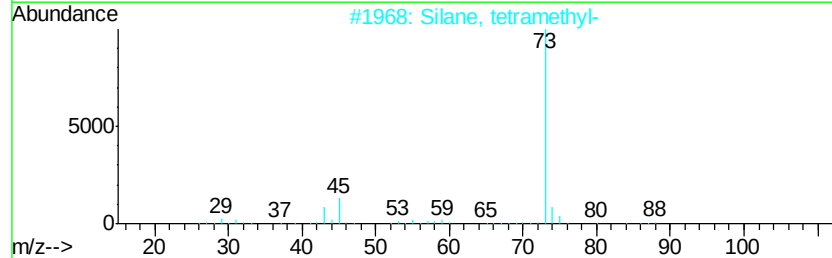
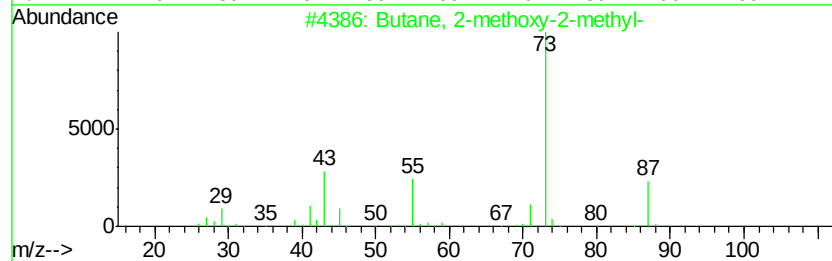
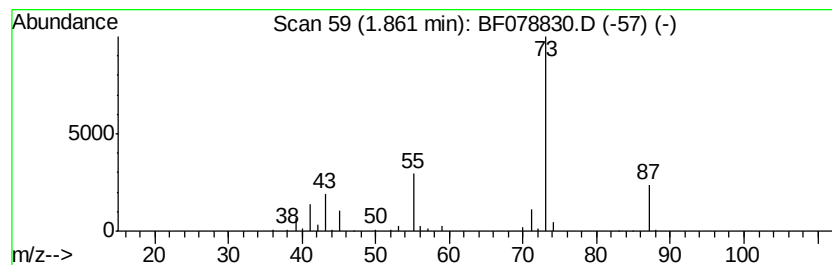
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042815.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Butane, 2-methoxy-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.86	12.75 ng	274651	1,4-Dichlorobenzene-d4	7.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	59
2		Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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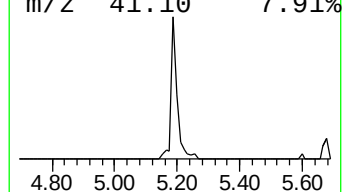
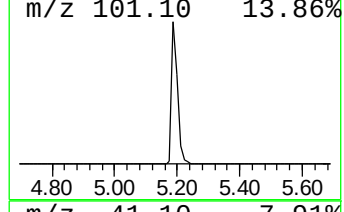
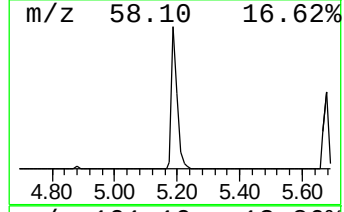
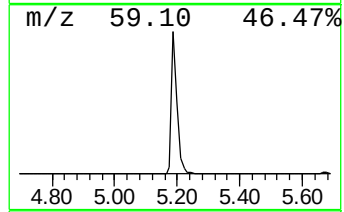
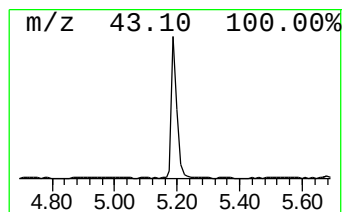
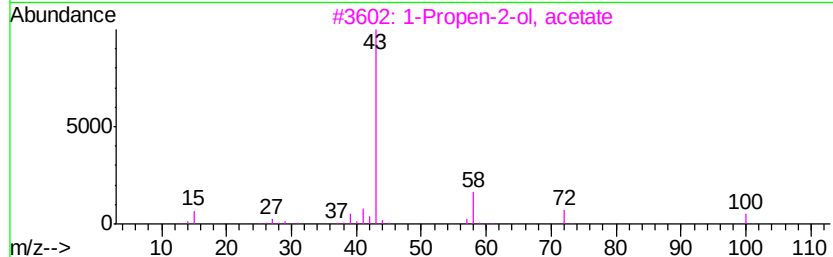
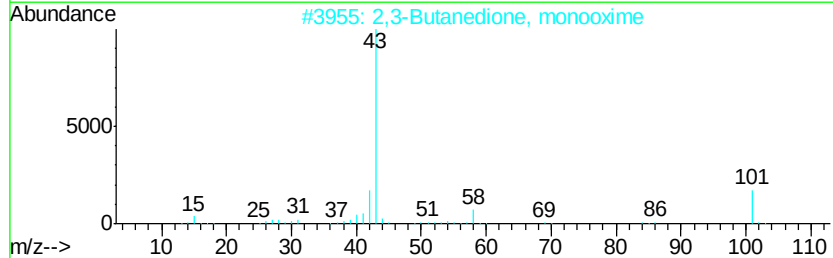
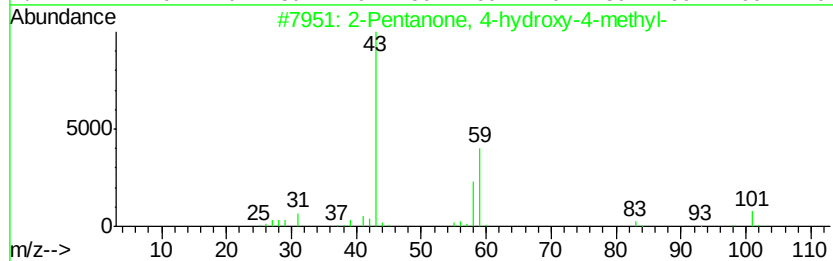
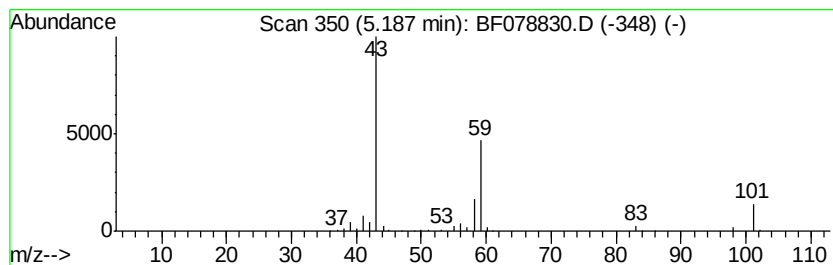
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	13.73 ng	295762	1,4-Dichlorobenzene-d4	7.38

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50	
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27	
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10	
4	2-Heptanone, 7,7-dichloro-	182	C7H12Cl2O	066241-43-8	9	
5	2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9	



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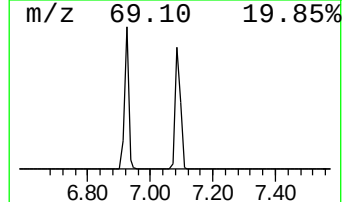
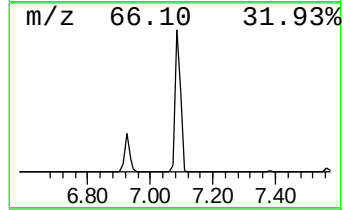
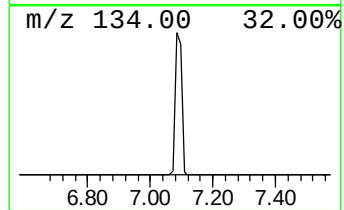
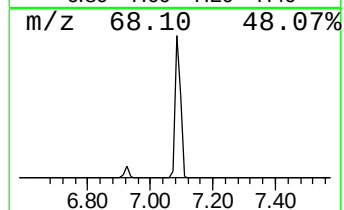
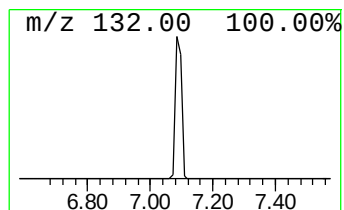
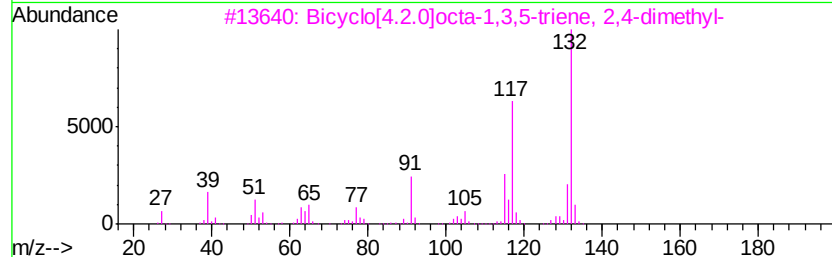
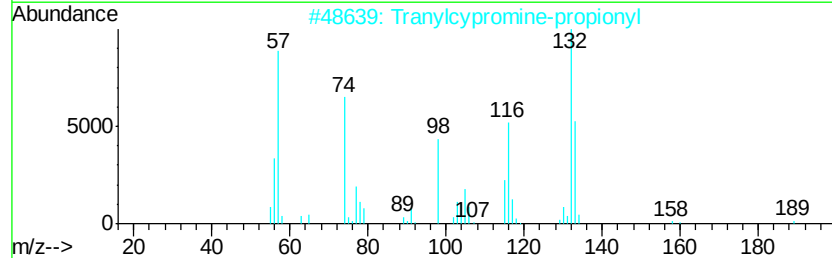
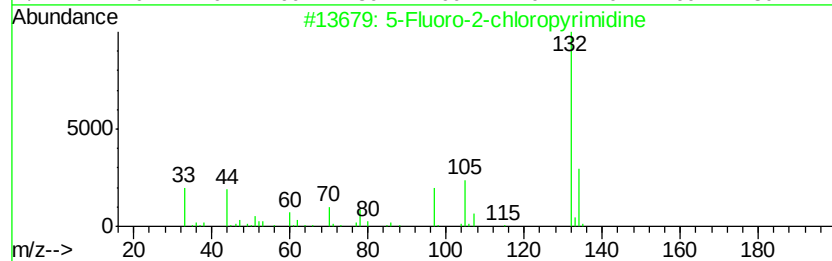
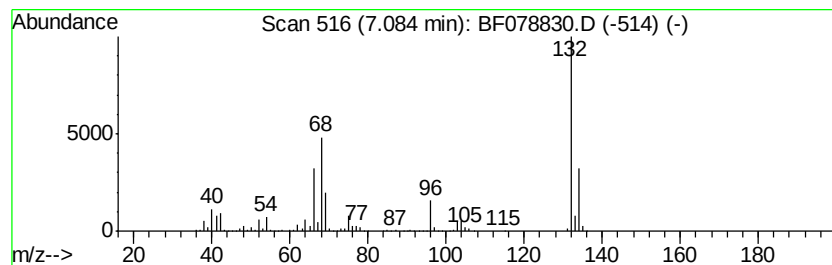
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown7.08 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.08	76.26 ng	1642950	1,4-Dichlorobenzene-d4	7.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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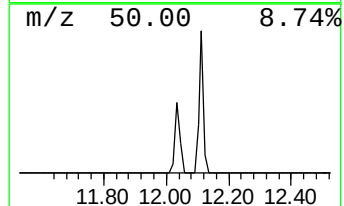
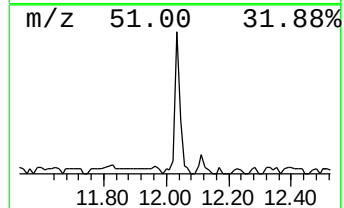
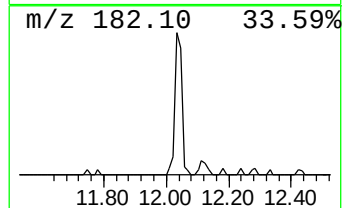
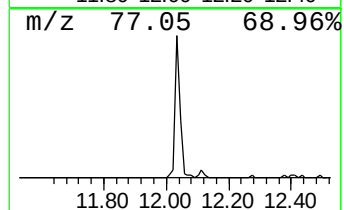
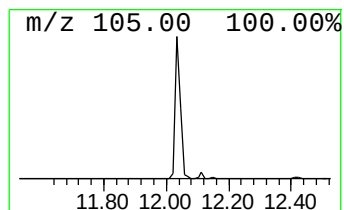
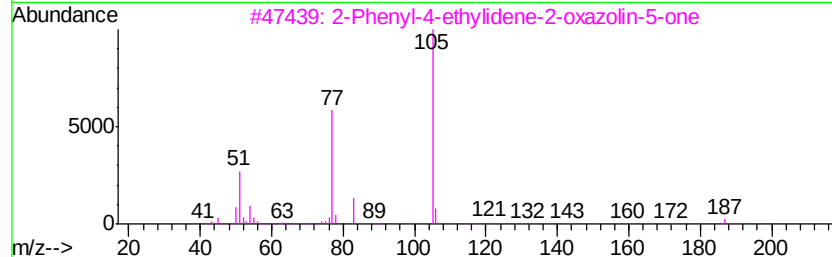
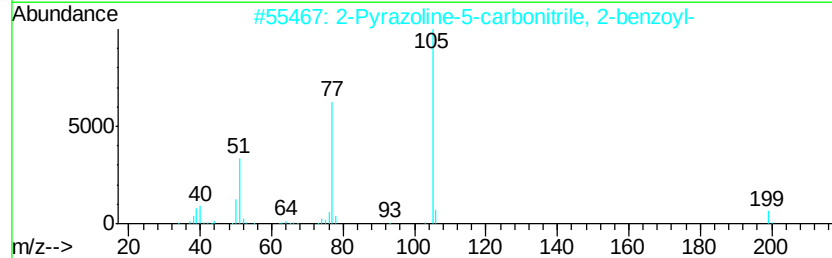
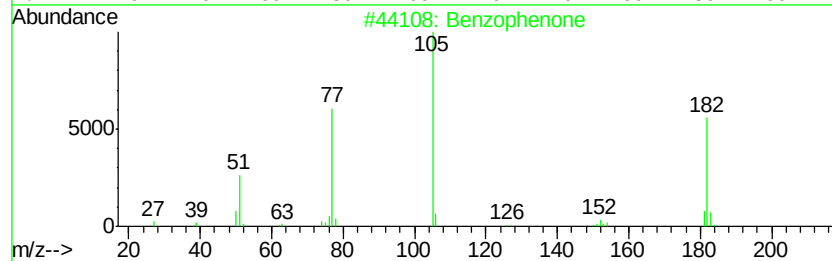
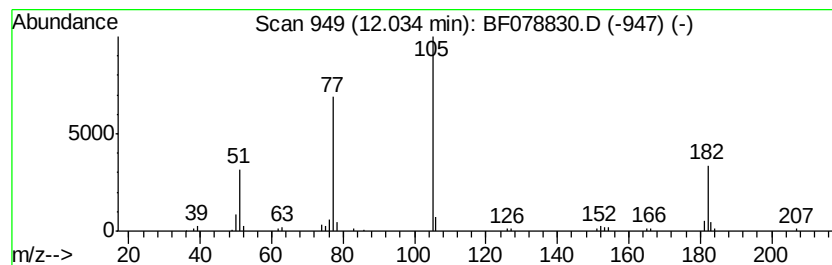
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzophenone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.03	3.44 ng	108794	Acenaphthene-d10	11.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzophenone	182	C13H10O	000119-61-9	96
2		2-Pyrazoline-5-carbonitrile, 2-b...	199	C11H9N3O	067872-68-8	53
3		2-Phenyl-4-ethylidene-2-oxazolin...	187	C11H9N02	037791-41-6	53
4		1-Propanone, 2-bromo-1-phenyl-	212	C9H9BrO	002114-00-3	53
5		Benzoic acid, phenyl ester	198	C13H10O2	000093-99-2	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042915\
Data File : BF078830.D
Acq On : 30 Apr 2015 1:21
Operator : TP/IZ
Sample : G2002-06
Misc :
ALS Vial : 24 Sample Multiplier: 1

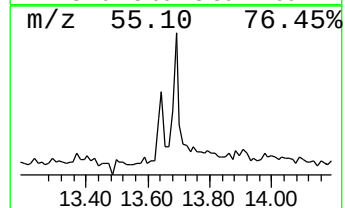
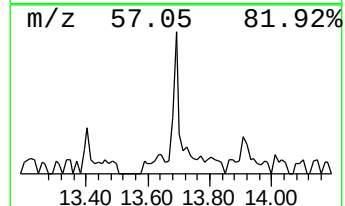
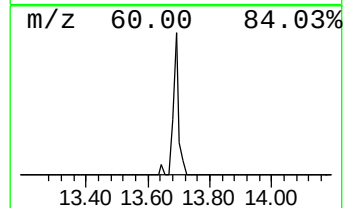
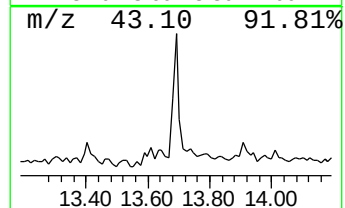
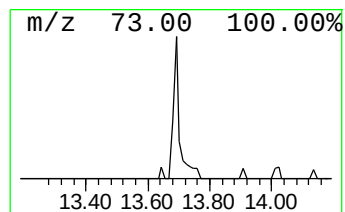
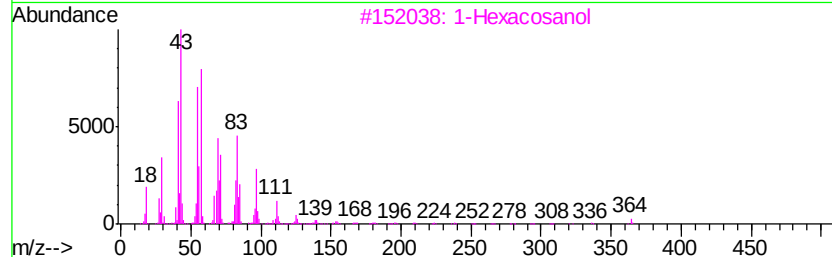
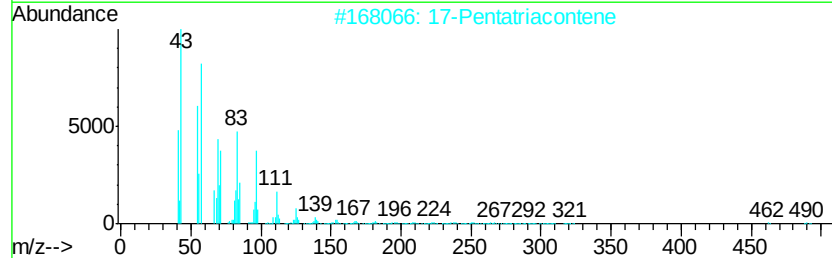
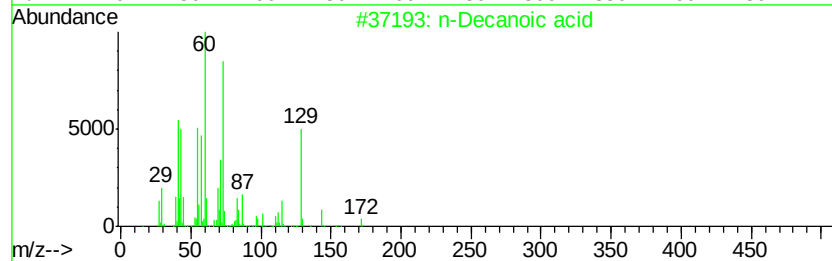
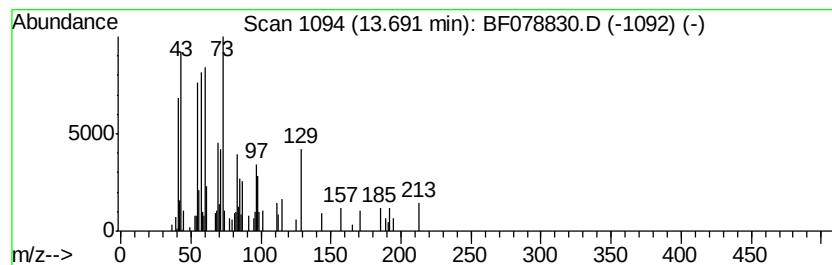
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 n-Decanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD		R.T.		
13.69	2.63 ng	98606	Phenanthrene-d10		12.98		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Decanoic acid	172	C10H20O2	000334-48-5	50
2			17-Pentatriacontene	491	C35H70	006971-40-0	25
3			1-Hexacosanol	382	C26H54O	000506-52-5	25
4			Chloroacetic acid, 3-pentadecyl ...	304	C17H33ClO2	1000280-08-8	22
5			.alpha.-D-Glucopyranoside, 0-.al...	504	C18H32O16	000597-12-6	14



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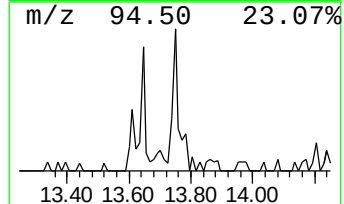
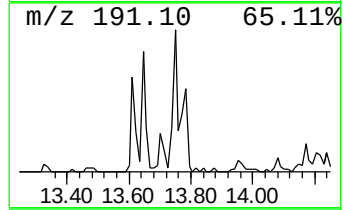
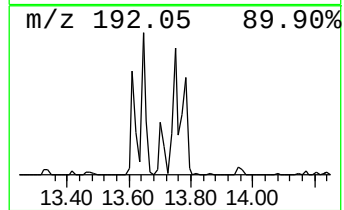
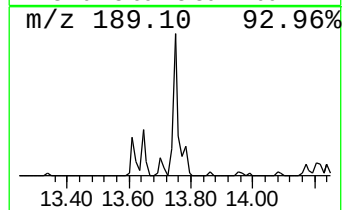
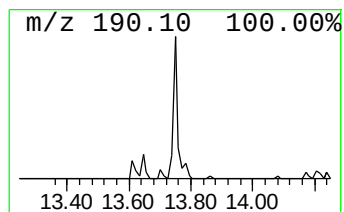
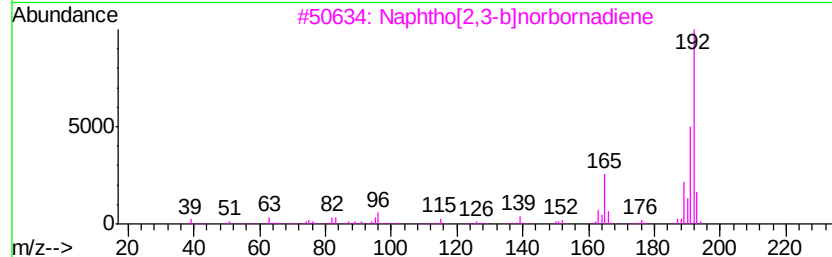
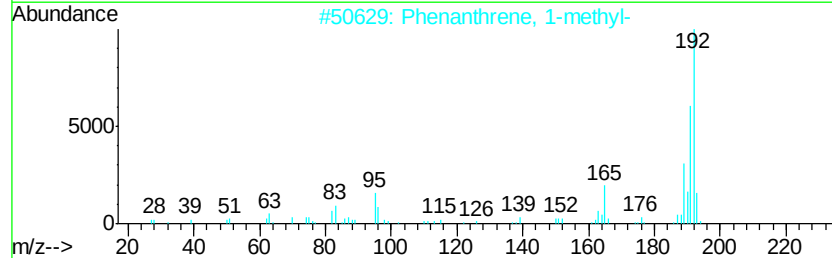
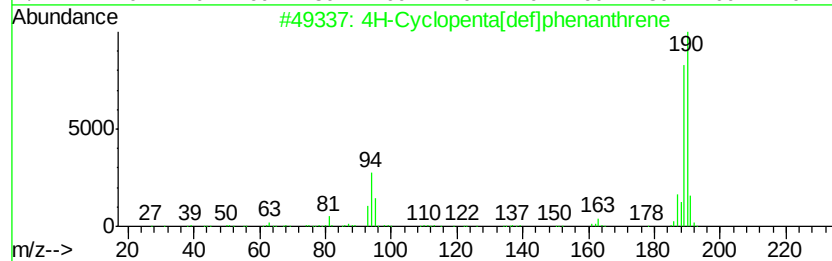
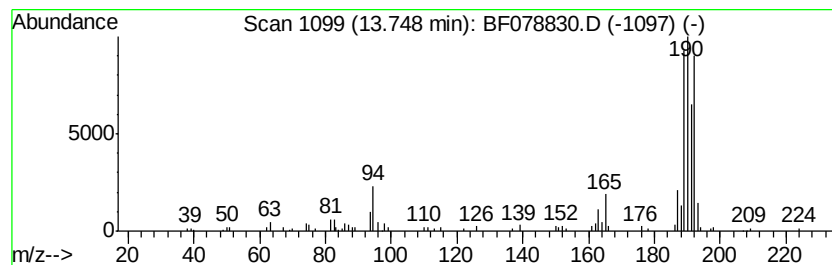
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 4H-Cyclopenta[def]phenanthrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.75	3.59 ng	134494	Phenanthrene-d10	12.98

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	53
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	46
3			Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	42
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	42
5			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	42



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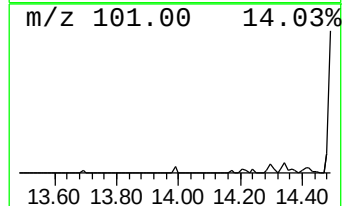
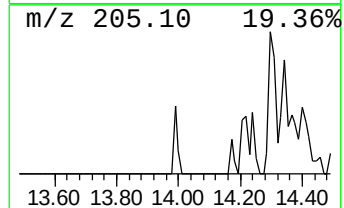
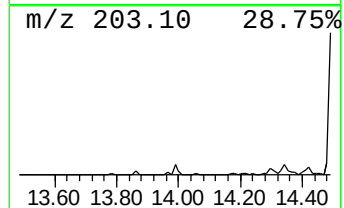
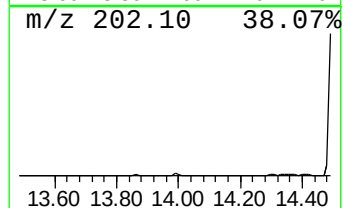
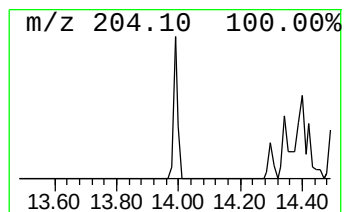
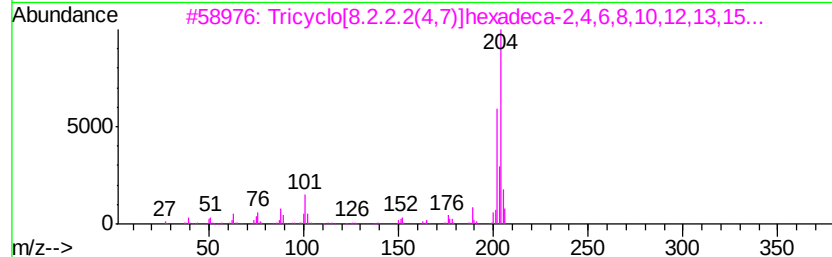
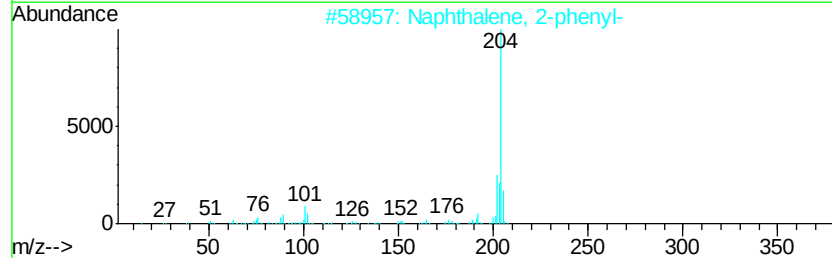
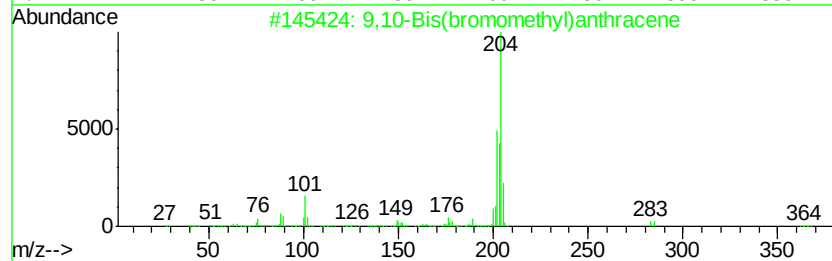
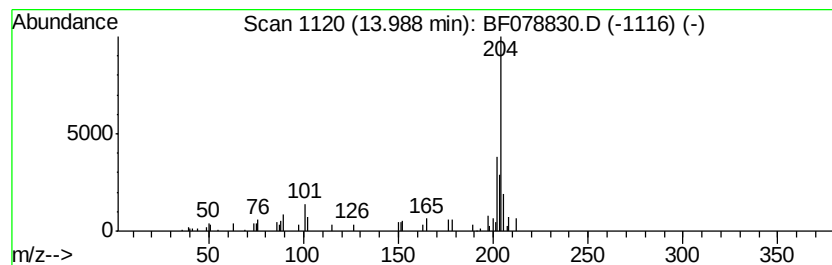
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 9,10-Bis(bromomethyl)anthra... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	2.42 ng	90513	Phenanthrene-d10	12.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	78
2		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	70
3		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	70
4		5H-Dibenzo[<i>a,d</i>]cycloheptene, 5-m...	204	C16H12	002975-79-3	49
5		6-Cyanophenanthridine	204	C14H8N2	002176-28-5	49



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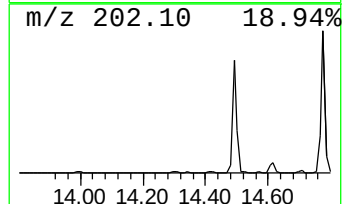
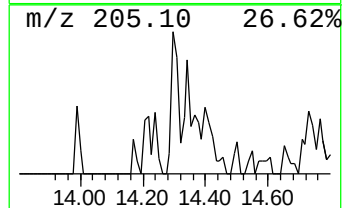
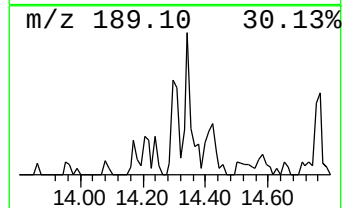
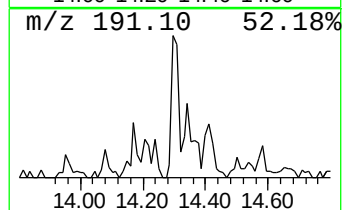
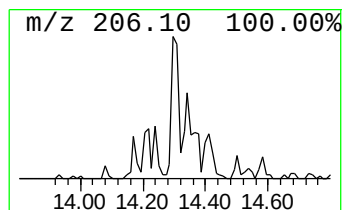
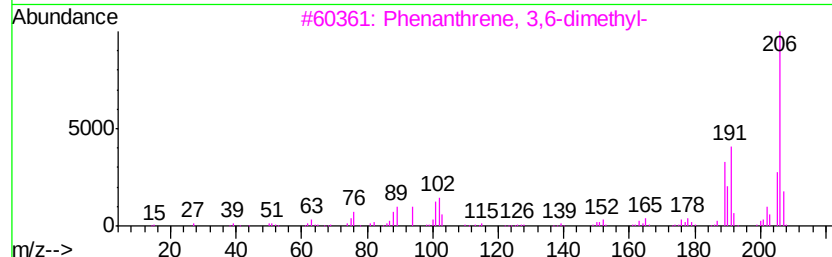
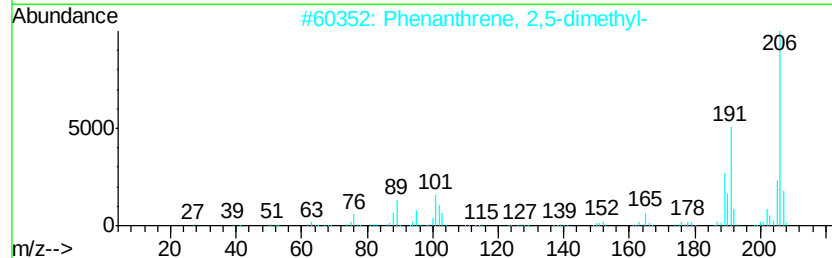
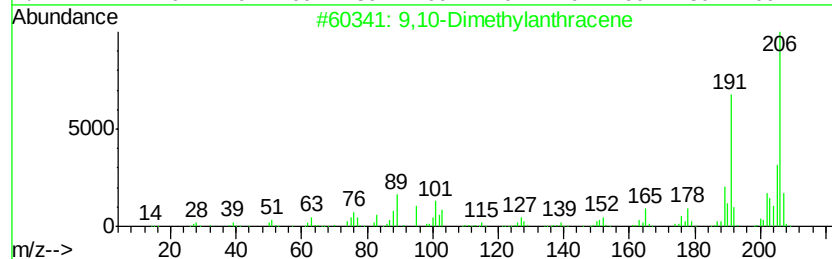
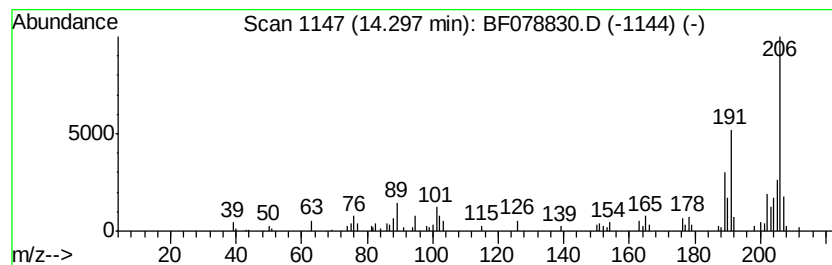
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 9,10-Dimethylantracene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.30	2.86 ng	107360	Phenanthrene-d10	12.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Dimethylantracene	206	C16H14	000781-43-1	98
2		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	97
3		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	90
4		Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	90
5		Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	86



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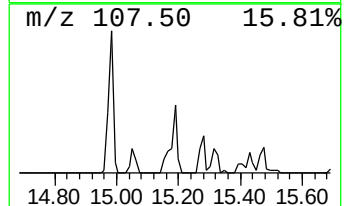
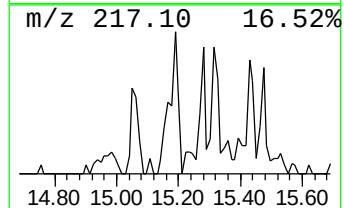
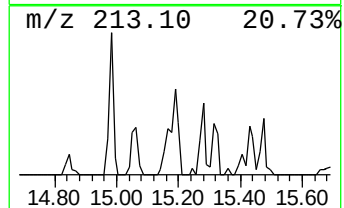
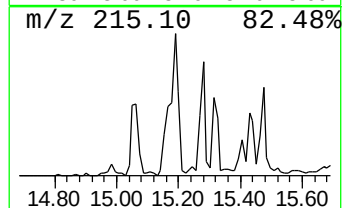
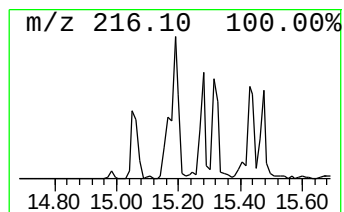
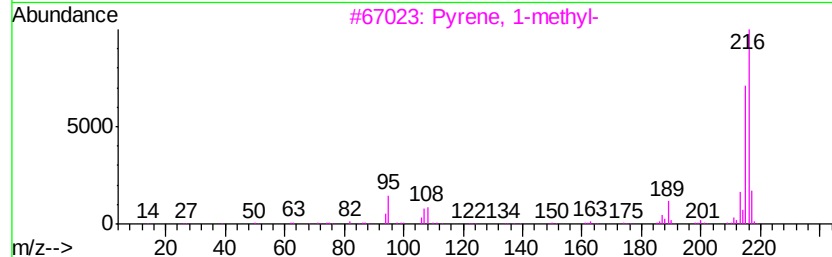
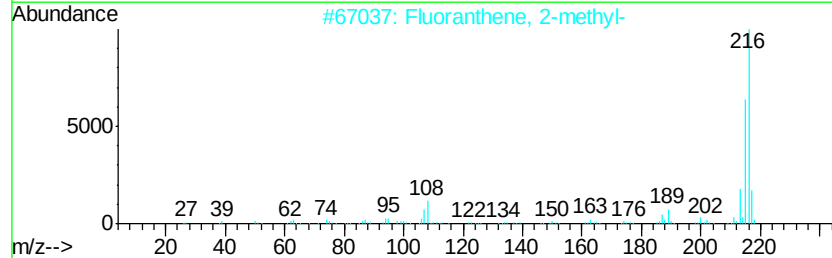
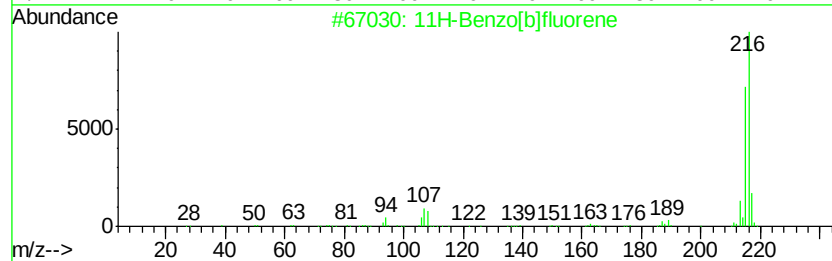
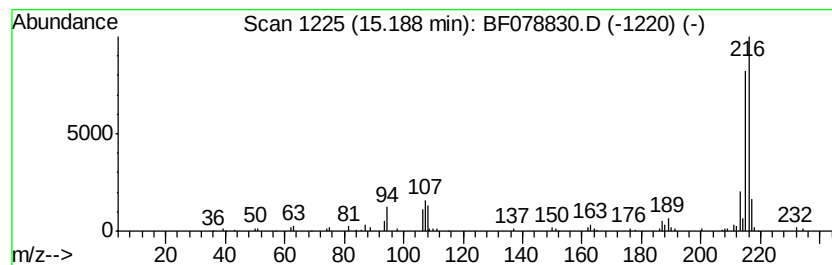
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 11H-Benzo[b]fluorene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.19	3.21 ng	214258	Chrysene-d12	16.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	95
2		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	95
3		Pvrene, 1-methyl-	216	C17H12	002381-21-7	93
4		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
5		7H-Benzo[c]fluorene	216	C17H12	000205-12-9	81



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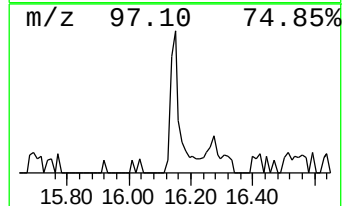
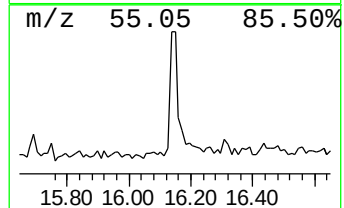
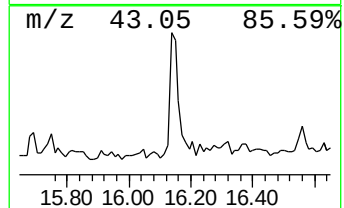
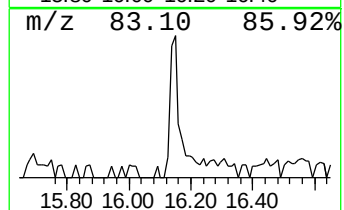
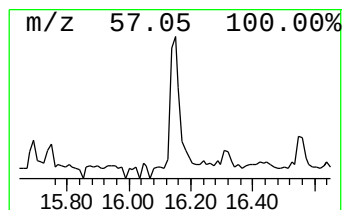
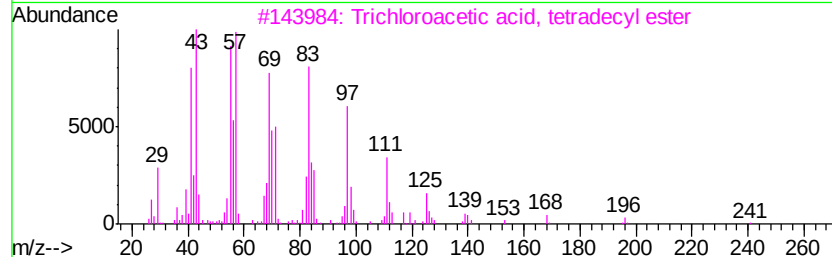
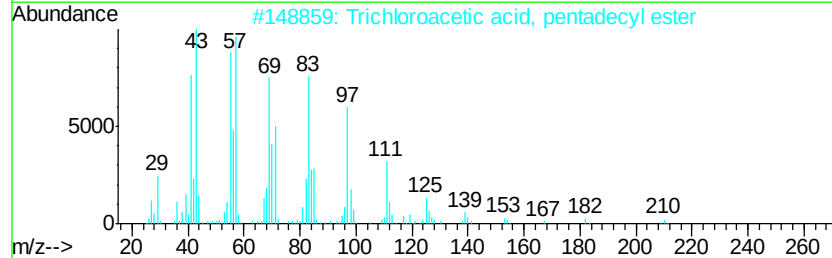
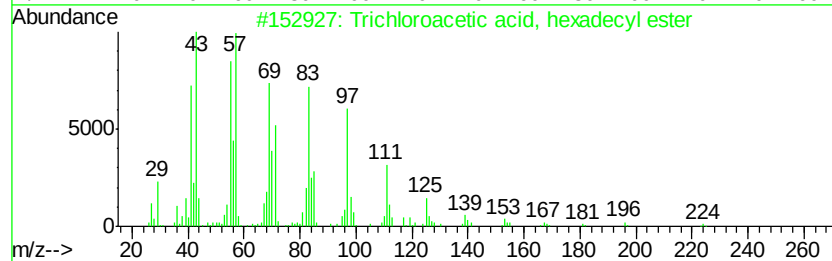
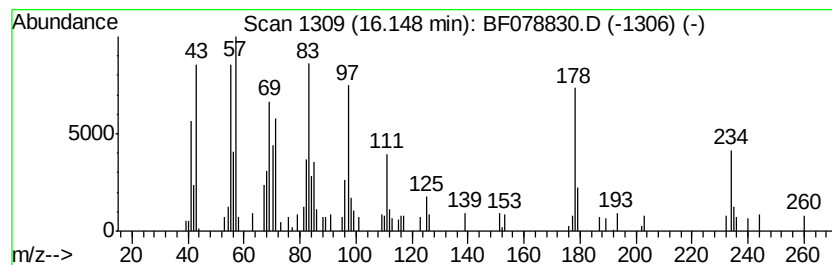
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Trichloroacetic acid, hexad... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.15	2.31 ng	153926	Chrysene-d12	16.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	89
2			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	87
3			Trichloroacetic acid, tetradecyl...	358	C16H29Cl3O2	074339-52-9	60
4			Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	60
5			Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	60



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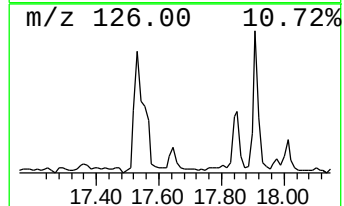
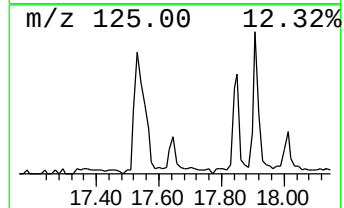
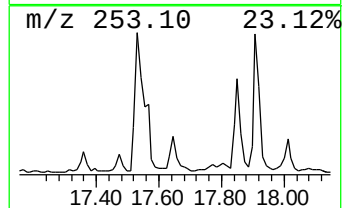
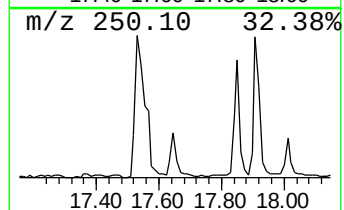
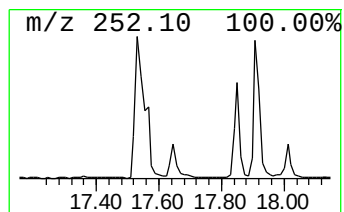
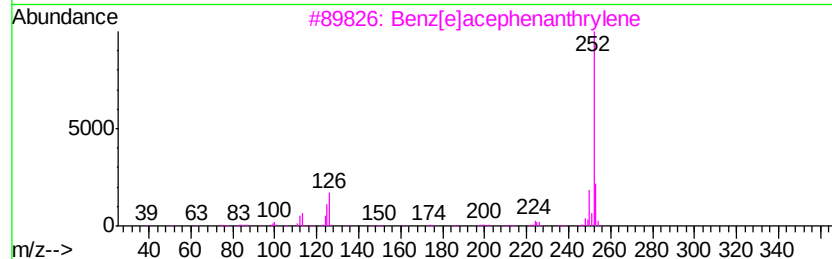
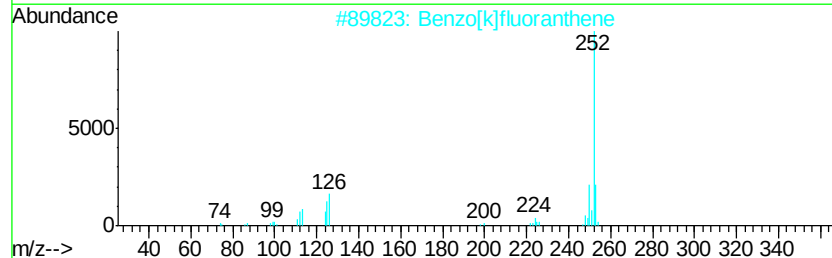
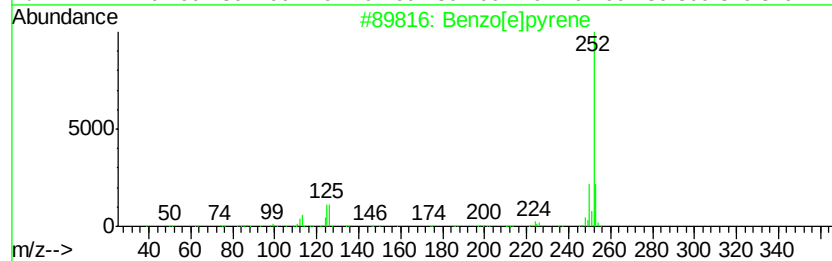
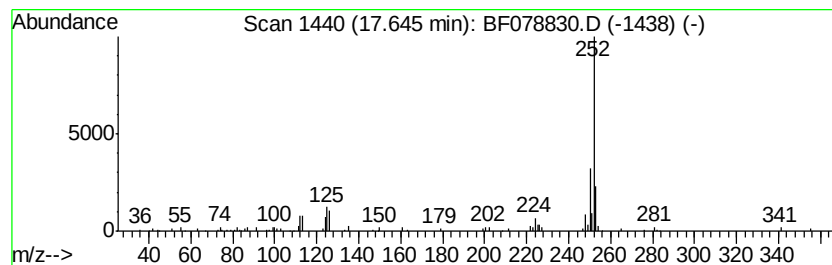
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Benzo[e]pyrene Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.65	2.01 ng	79803	Perylene-d12	17.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	96
2		Benzo[k]fluoranthene	252	C20H12	000207-08-9	94
3		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	93
4		Benzo[a]pyrene	252	C20H12	000050-32-8	81
5		Perylene	252	C20H12	000198-55-0	76



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042915\
Data File : BF078830.D
Acq On : 30 Apr 2015 1:21
Operator : TP/IZ
Sample : G2002-06
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TEC-03

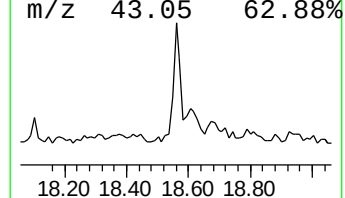
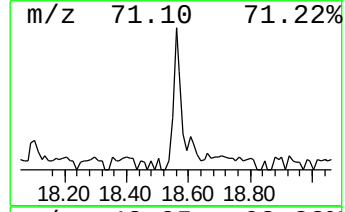
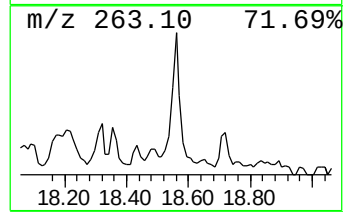
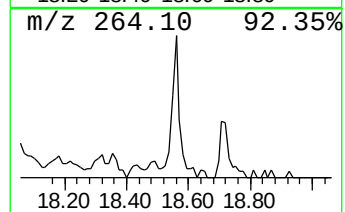
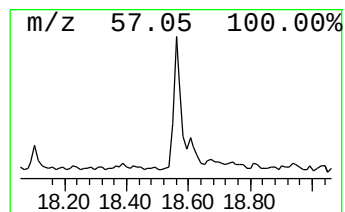
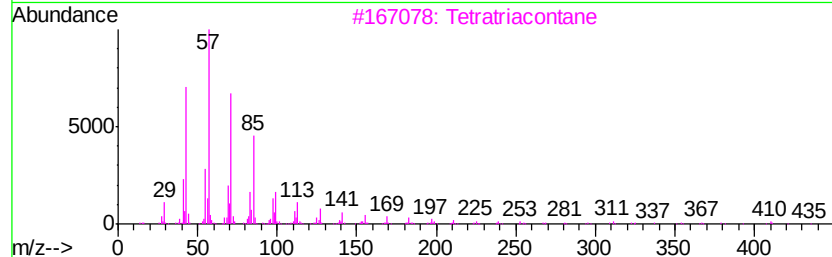
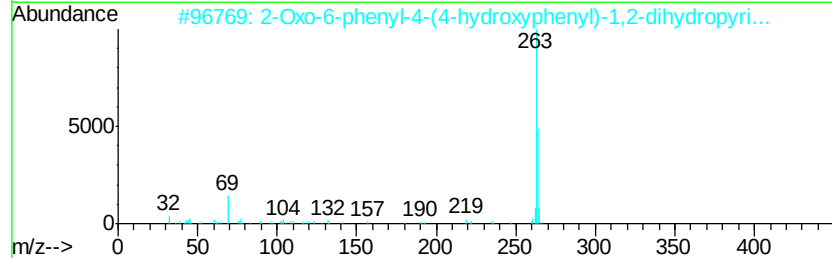
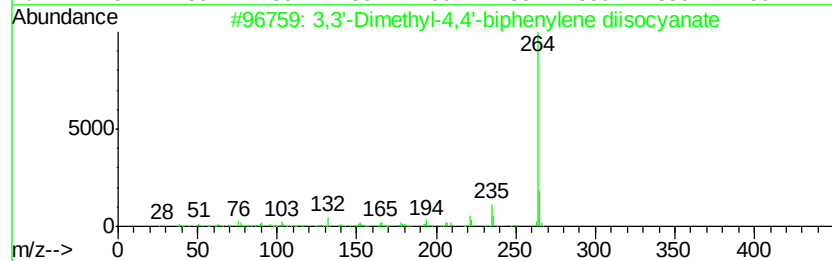
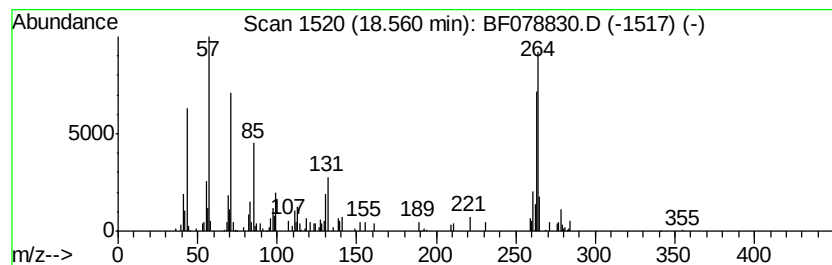
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 unknown18.56 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.56	2.77 ng	109635	Perylene-d12	17.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3,3'	-Dimethyl-4,4'-biphenylene d...	264	C16H12N2O2	000091-97-4	22
2	2-Oxo-6-phenyl-4-(4-hydroxypheny...		264	C16H12N2O2	161200-20-0	22
3	Tetratriacontane		479	C34H70	014167-59-0	15
4	Octacosane		394	C28H58	000630-02-4	15
5	Heneicosane		296	C21H44	000629-94-7	15



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042915\
Data File : BF078830.D
Acq On : 30 Apr 2015 1:21
Operator : TP/IZ
Sample : G2002-06
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TEC-03

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042815.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
Propane, 2,2-dime...	1.54	345.1	ng	7435090	1	7.38	430899	20.0
Butane, 2-methoxy...	1.86	12.8	ng	274651	1	7.38	430899	20.0
2-Pentanone, 4-hy...	5.19	13.7	ng	295762	1	7.38	430899	20.0
unknown7.08	7.08	76.3	ng	1642950	1	7.38	430899	20.0
Benzophenone	12.03	3.4	ng	108794	3	11.14	632769	20.0
n-Decanoic acid	13.69	2.6	ng	98606	4	12.98	749566	20.0
4H-Cyclopenta[def...	13.75	3.6	ng	134494	4	12.98	749566	20.0
9,10-Bis(bromomet...	13.99	2.4	ng	90513	4	12.98	749566	20.0
9,10-Dimethylanthe...	14.30	2.9	ng	107360	4	12.98	749566	20.0
11H-Benzo[b]fluorene	15.19	3.2	ng	214258	5	16.27	1333400	20.0
Trichloroacetic a...	16.15	2.3	ng	153926	5	16.27	1333400	20.0
Benzo[e]pyrene	17.65	2.0	ng	79803	6	17.98	792366	20.0
unknown18.56	18.56	2.8	ng	109635	6	17.98	792366	20.0