

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051515\
 Data File : BF079299.D
 Acq On : 16 May 2015 7:43
 Operator : TP/IZ
 Sample : G2249-02
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SR-2

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-BF043015.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	1.324	10	12	15	rBV	4065032	5414382	35.71%	10.934%
2	1.473	20	25	27	rBV	7107438	15163478	100.00%	30.621%
3	1.587	31	35	36	rBV2	115851	175368	1.16%	0.354%
4	1.621	36	38	41	rVB	213188	253915	1.67%	0.513%
5	1.793	49	53	61	rVB2	102752	194110	1.28%	0.392%
6	4.947	326	329	338	rVB	427315	537136	3.54%	1.085%
7	5.473	372	375	378	rBV	1732552	1823943	12.03%	3.683%
8	6.753	484	487	489	rBV	2016757	1905410	12.57%	3.848%
9	6.890	497	499	502	rBV	2012646	1934661	12.76%	3.907%
10	7.176	522	524	527	rBV	498142	461604	3.04%	0.932%
11	7.359	538	540	543	rBV	1122683	1531666	10.10%	3.093%
12	7.873	582	585	587	rBV	1378684	1227267	8.09%	2.478%
13	8.753	660	662	667	rVB	720168	641520	4.23%	1.295%
14	10.102	778	780	782	rBV	2624662	2305315	15.20%	4.655%
15	10.925	850	852	854	rBV	698992	667212	4.40%	1.347%
16	11.908	935	938	941	rBV	1508769	1544509	10.19%	3.119%
17	12.765	1011	1013	1014	rBV	692692	679025	4.48%	1.371%
18	12.799	1014	1016	1018	rVV	559804	601244	3.97%	1.214%
19	12.857	1019	1021	1023	rVB	194339	196478	1.30%	0.397%
20	13.485	1074	1076	1078	rBV	122665	169799	1.12%	0.343%
21	13.531	1078	1080	1081	rVV	169042	226292	1.49%	0.457%
22	14.274	1142	1145	1148	rBV	1119456	1082311	7.14%	2.186%
23	14.548	1166	1169	1171	rVV	906075	1089079	7.18%	2.199%
24	14.765	1185	1188	1190	rVV	2467244	2530516	16.69%	5.110%
25	14.971	1204	1206	1208	rVB	154961	178139	1.17%	0.360%
26	15.097	1215	1217	1219	rBV	141833	154070	1.02%	0.311%
27	15.943	1288	1291	1296	rBV3	357216	611698	4.03%	1.235%
28	16.057	1297	1301	1302	rVV2	820737	1429405	9.43%	2.887%
29	16.091	1302	1304	1309	rVB2	466707	878305	5.79%	1.774%
30	16.766	1361	1363	1366	rVB2	129209	226470	1.49%	0.457%
31	17.360	1413	1415	1420	rVB	447378	966904	6.38%	1.953%
32	17.680	1441	1443	1446	rBV	277383	356479	2.35%	0.720%
33	17.749	1446	1449	1452	rVB	374686	499468	3.29%	1.009%
34	17.806	1452	1454	1456	rBV	473933	742401	4.90%	1.499%

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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-BF043015.M
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35	18.366	1501	1503	1506	rVB2	119568	181524	1.20%	0.367%
36	19.212	1574	1577	1583	rVB2	193116	406952	2.68%	0.822%
37	19.394	1590	1593	1595	rBV	61664	161934	1.07%	0.327%
38	19.623	1610	1613	1621	rVB	166502	369387	2.44%	0.746%

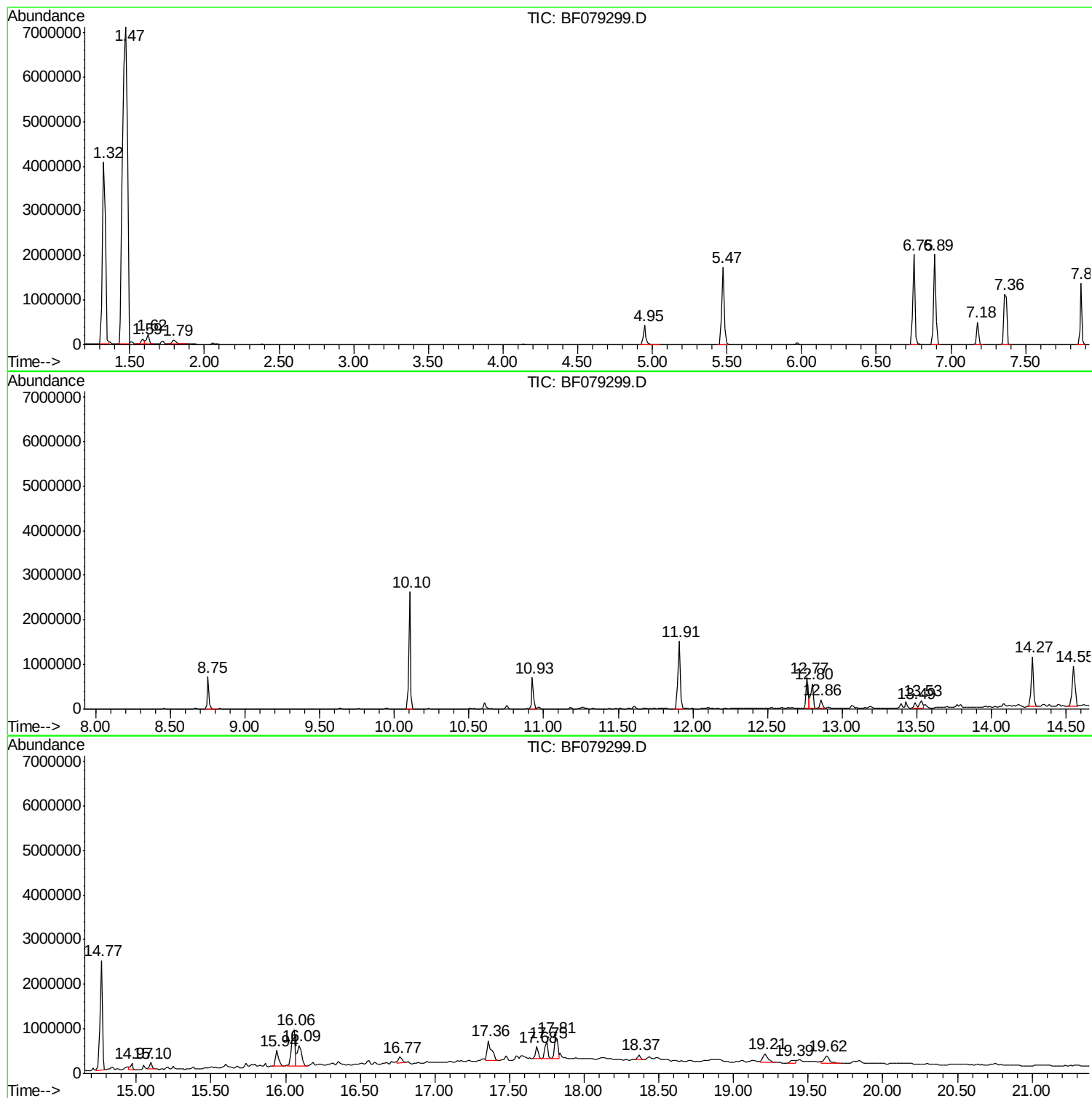
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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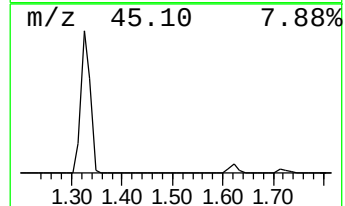
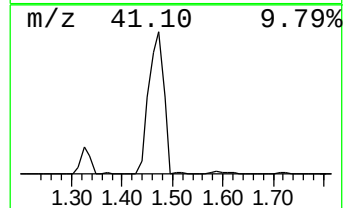
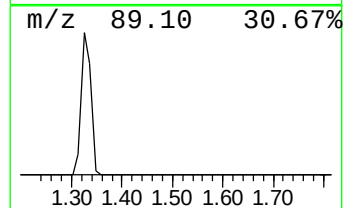
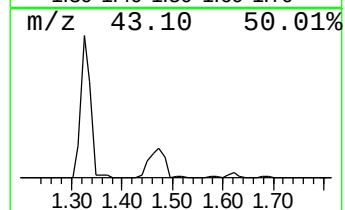
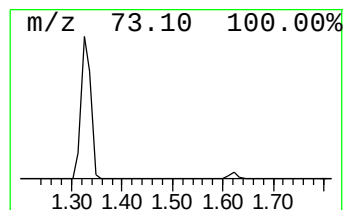
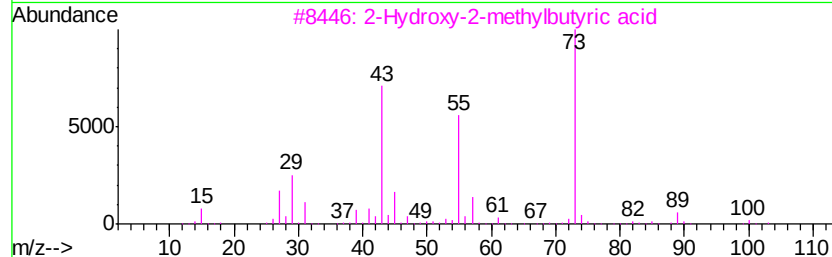
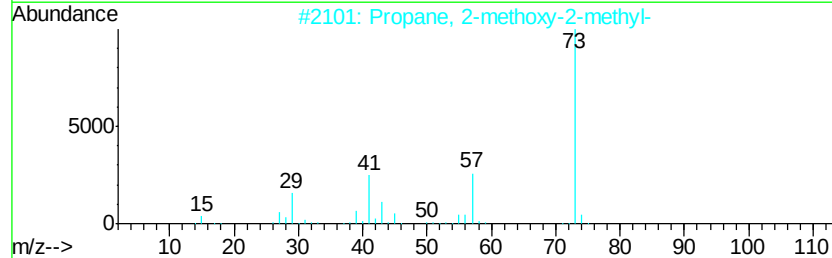
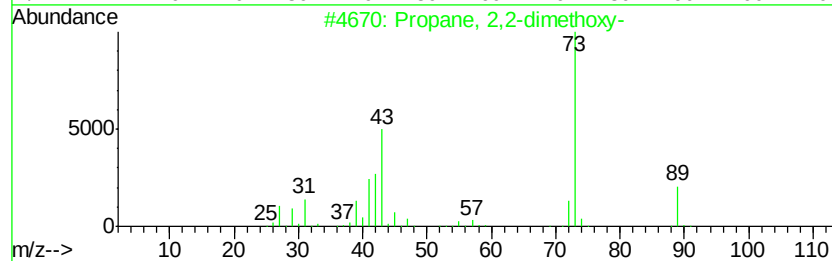
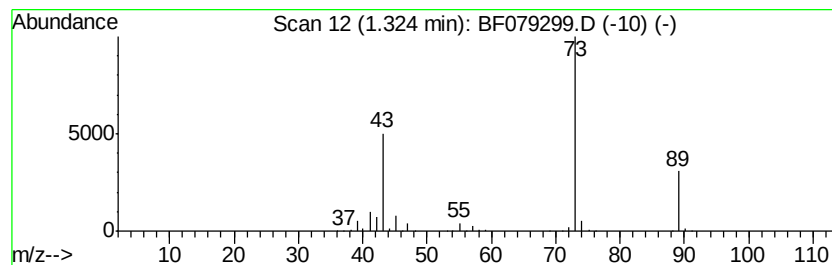
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown1.32 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.32	234.59 ng	5414380	1,4-Dichlorobenzene-d4	7.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	38
2		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
3		2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9
4		1-Hexen-4-ol, 1-chloro-3,5-dimet...	162	C8H15ClO	100244-09-5	9
5		N-Ethylformamide	73	C3H7NO	000627-45-2	5



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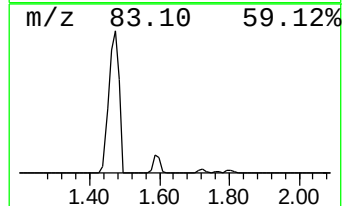
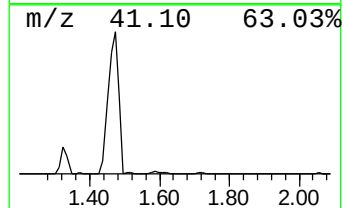
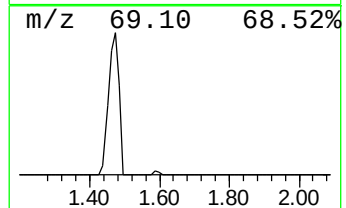
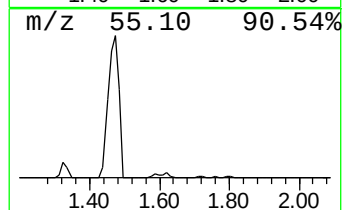
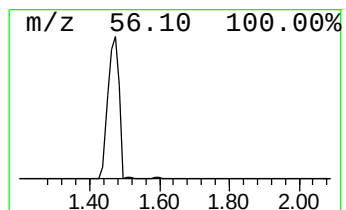
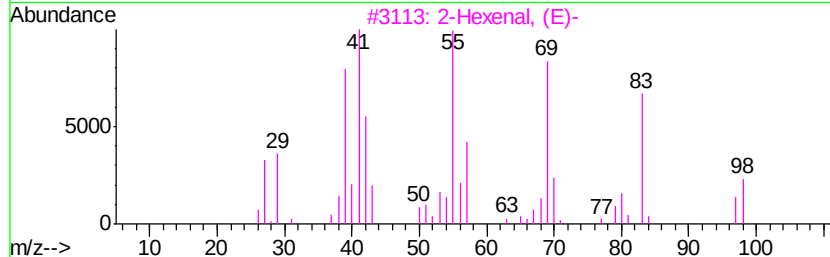
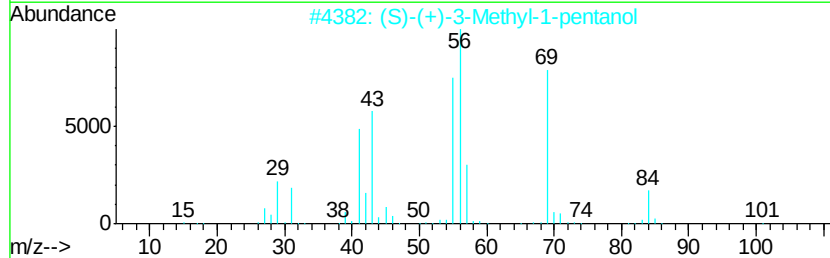
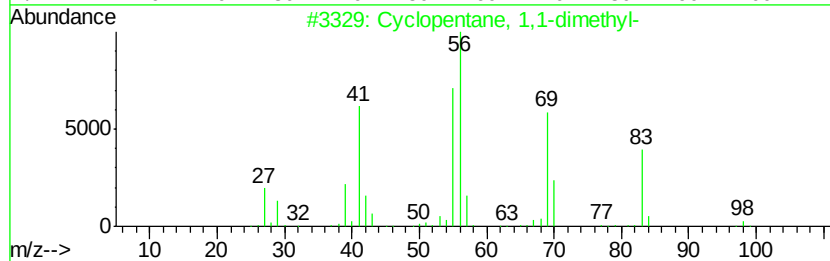
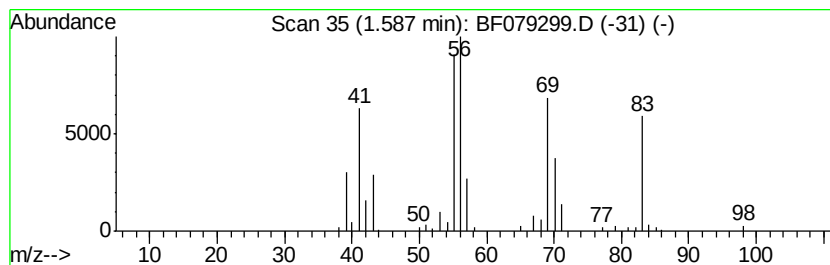
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Cyclopentane, 1,1-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.59	7.60 ng	175368	1,4-Dichlorobenzene-d4	7.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	86
2		(S)-(+)-3-Methyl-1-pentanol	102	C6H14O	042072-39-9	50
3		2-Hexenal, (E)-	98	C6H10O	006728-26-3	38
4		Cyclobutane, ethyl-	84	C6H12	004806-61-5	38
5		1-Butene, 2,3,3-trimethyl-	98	C7H14	000594-56-9	27



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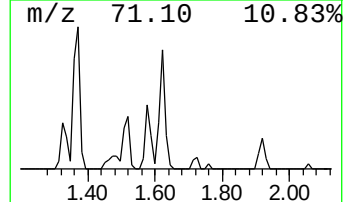
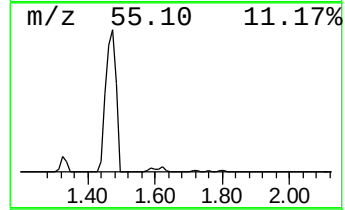
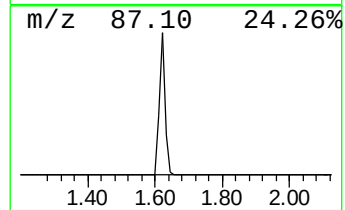
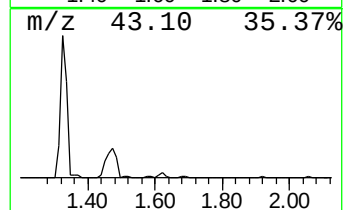
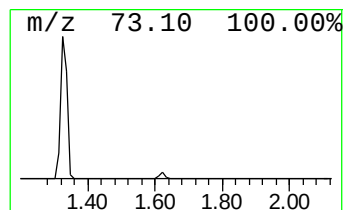
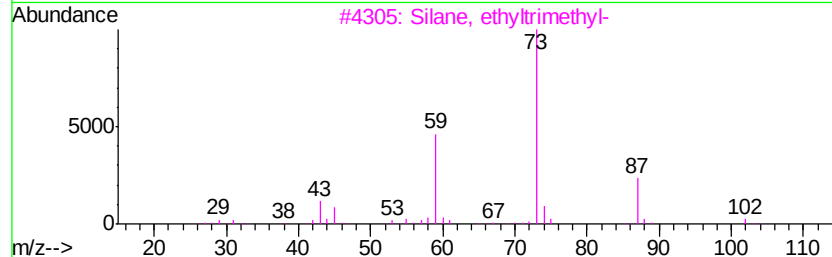
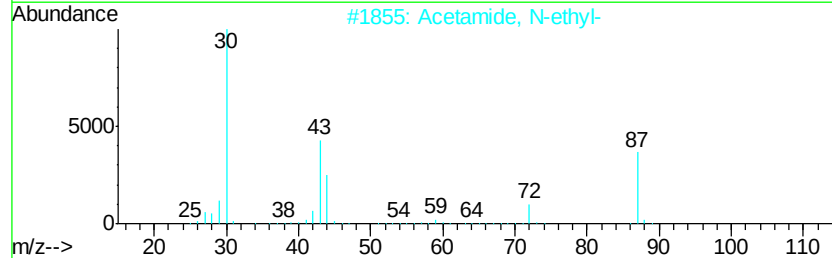
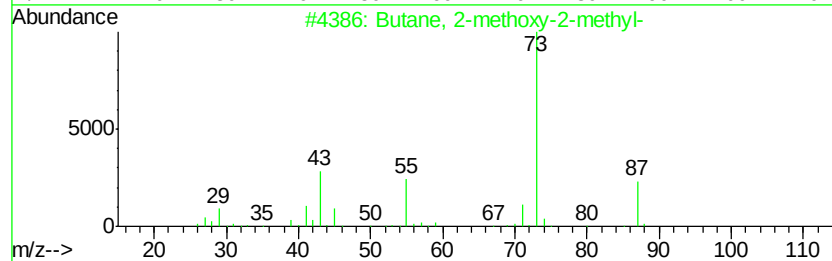
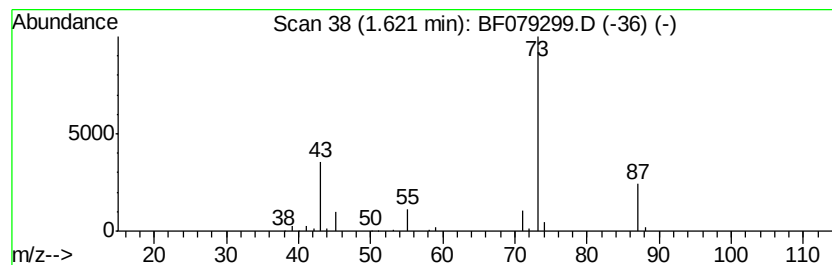
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Butane, 2-methoxy-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.62	11.00 ng	253915	1,4-Dichlorobenzene-d4	7.18

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
3			Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	9
4			2-Propanone, oxime	73	C3H7NO	000127-06-0	7
5			N-Ethylformamide	73	C3H7NO	000627-45-2	7



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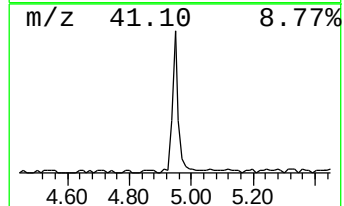
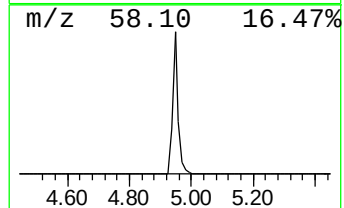
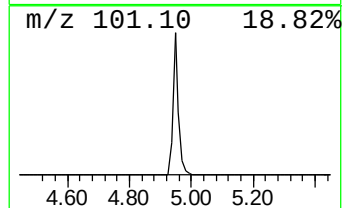
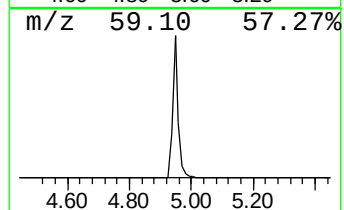
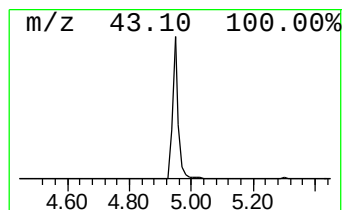
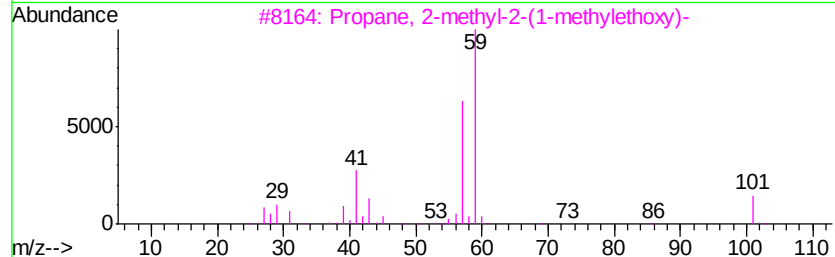
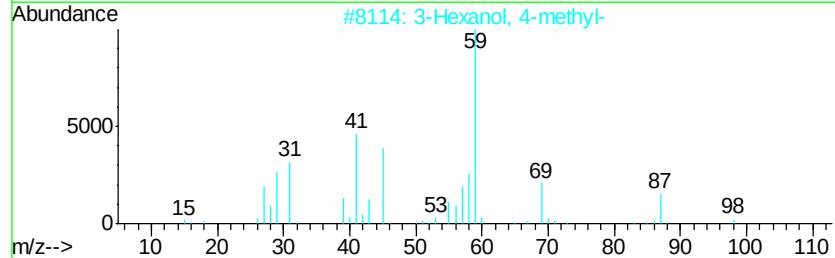
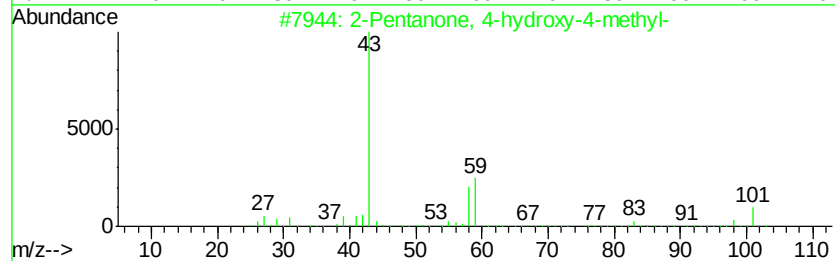
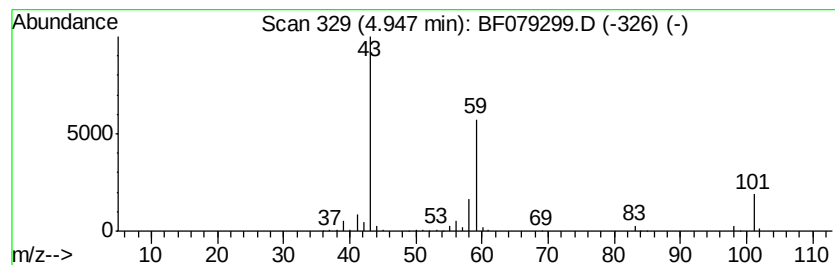
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.95	23.27 ng	537136	1,4-Dichlorobenzene-d4	7.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3		Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	33
4		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	25
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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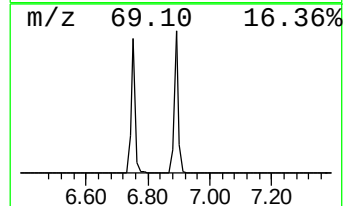
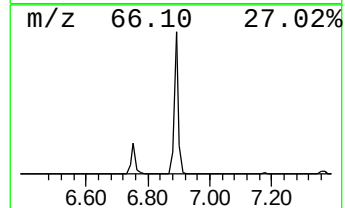
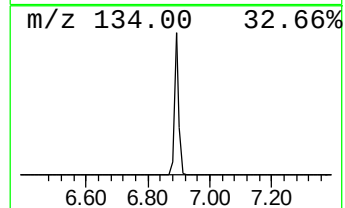
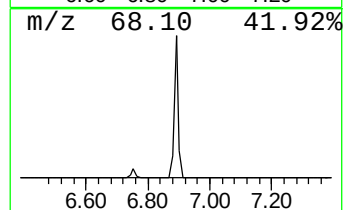
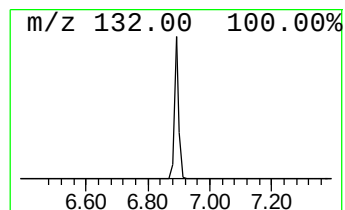
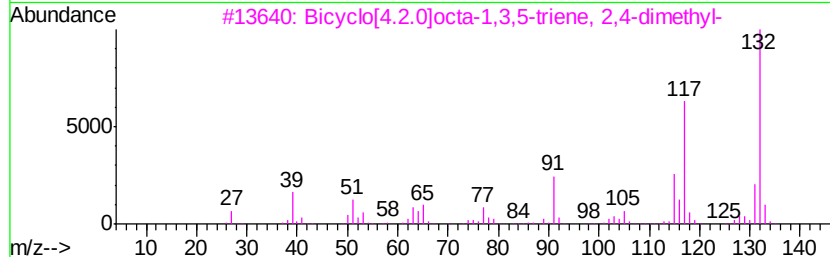
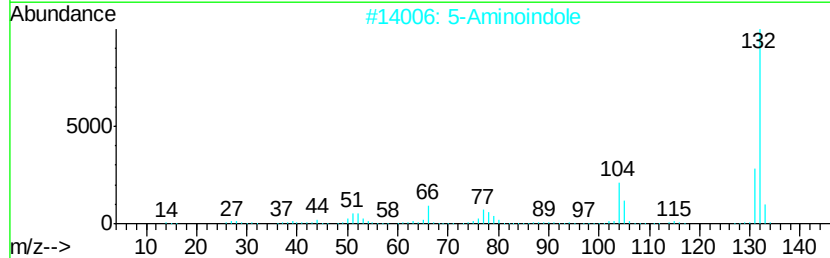
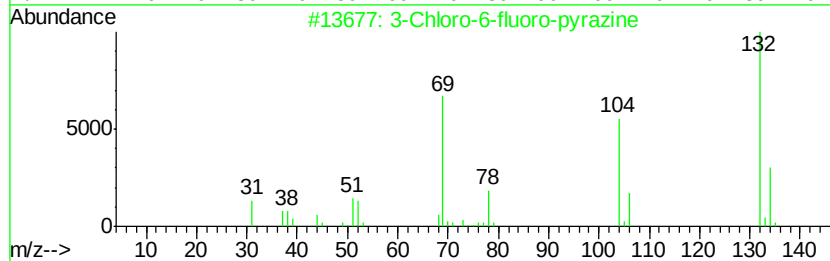
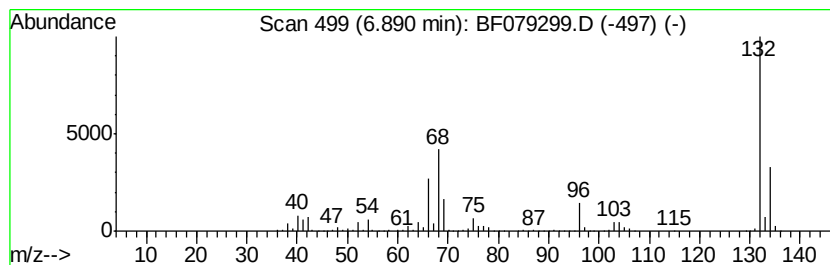
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 7 unknown6.89 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.89	83.82 ng	1934660	1,4-Dichlorobenzene-d4	7.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051515\
Data File : BF079299.D
Acq On : 16 May 2015 7:43
Operator : TP/IZ
Sample : G2249-02
Misc :
ALS Vial : 32 Sample Multiplier: 1

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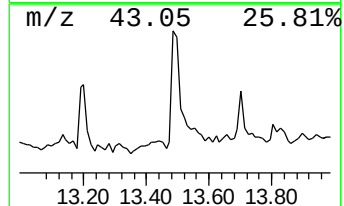
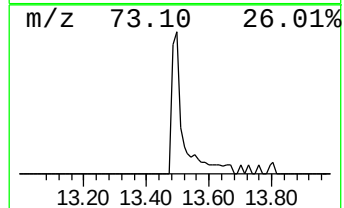
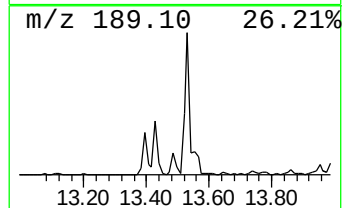
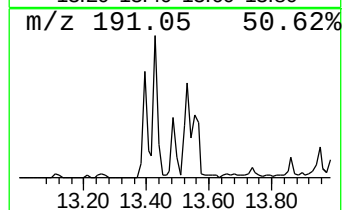
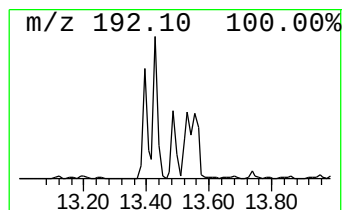
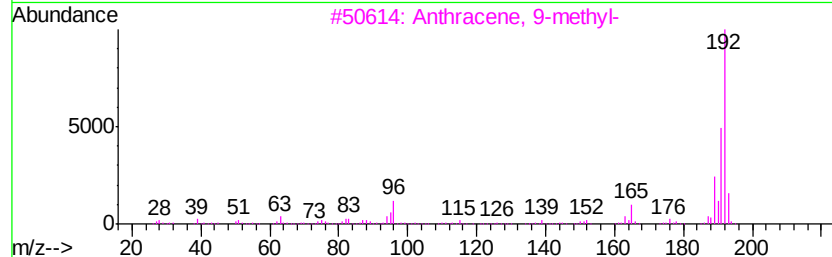
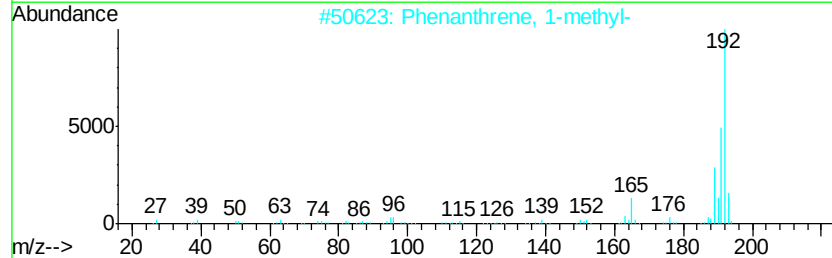
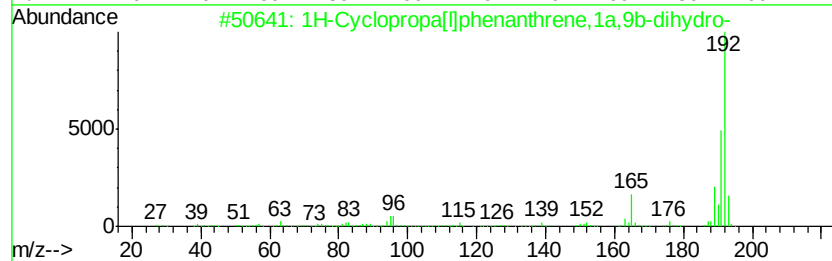
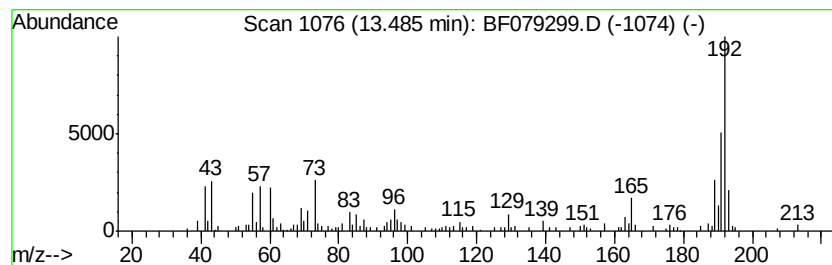
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1H-Cyclopropa[l]phenanthren... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.49	5.00 ng	169799	Phenanthrene-d10	12.77

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	98
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
3		Anthracene, 9-methyl-	192	C15H12	000779-02-2	95
4		Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051515\
Data File : BF079299.D
Acq On : 16 May 2015 7:43
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Sample : G2249-02
Misc :
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Instrument :
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ClientSampled :
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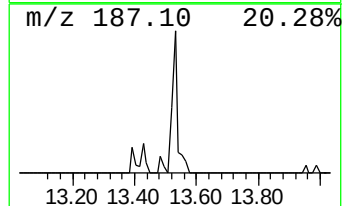
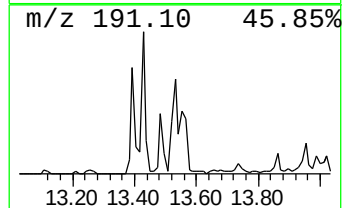
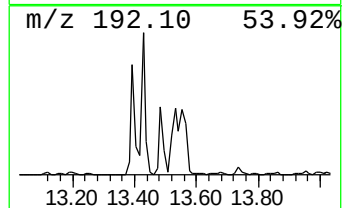
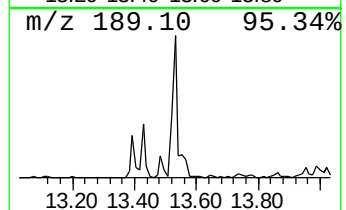
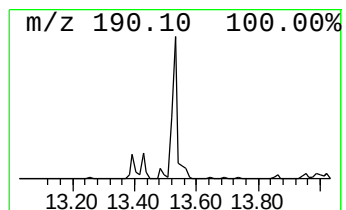
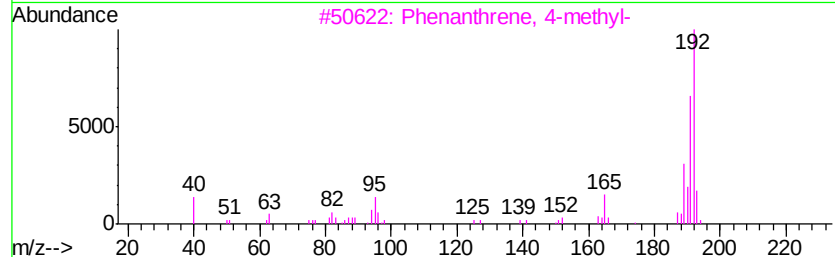
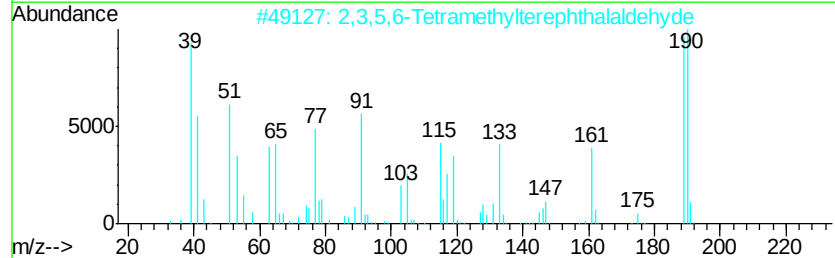
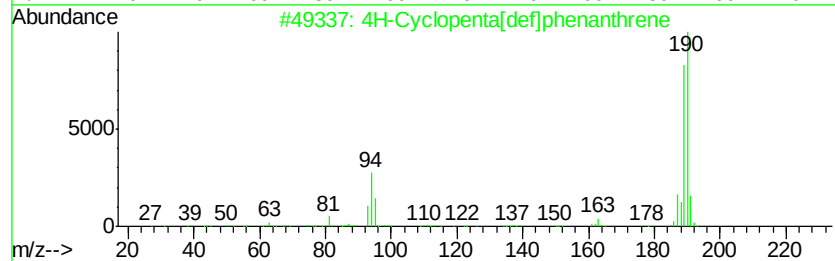
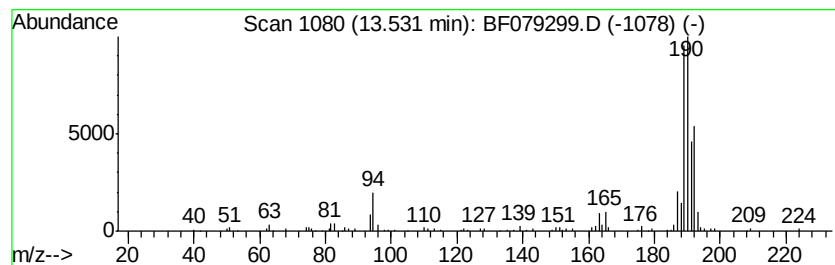
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.53	6.67 ng	226292	Phenanthrene-d10	12.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
2		2,3,5,6-Tetramethylterephthalaldehyde	190	C12H14O2	007072-01-7	38
3		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	38
4		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	30
5		5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	25



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051515\
Data File : BF079299.D
Acq On : 16 May 2015 7:43
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Sample : G2249-02
Misc :
ALS Vial : 32 Sample Multiplier: 1

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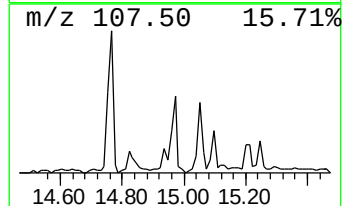
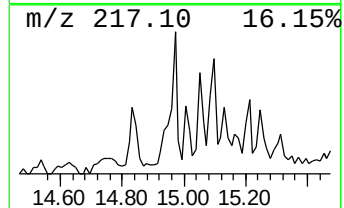
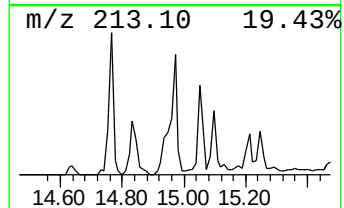
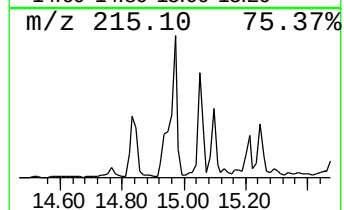
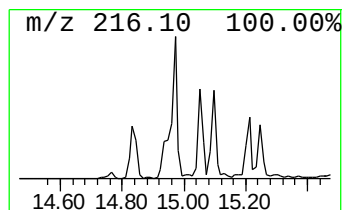
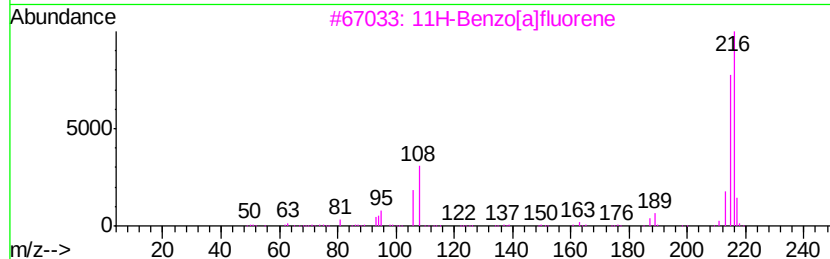
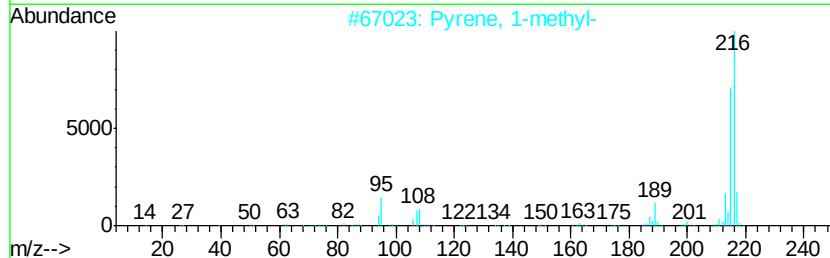
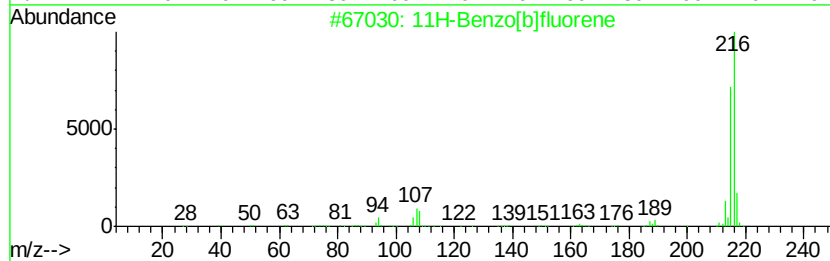
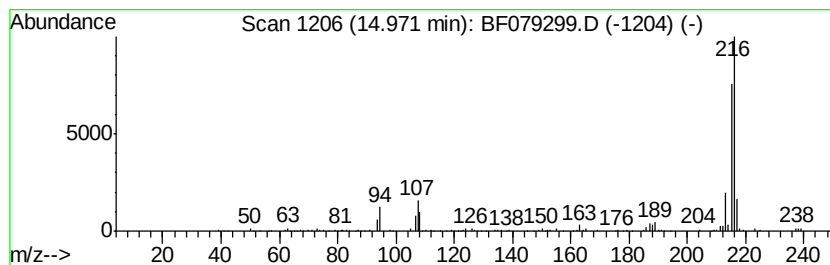
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 11H-Benzo[b]fluorene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.97	2.49 ng	178139	Chrysene-d12	16.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	81
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051515\
Data File : BF079299.D
Acq On : 16 May 2015 7:43
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Sample : G2249-02
Misc :
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ClientSampleId :
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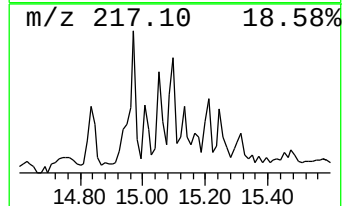
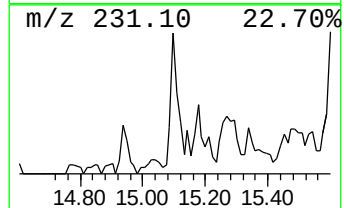
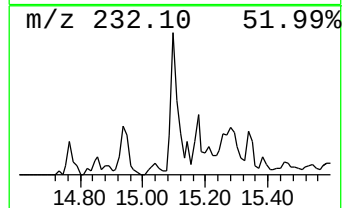
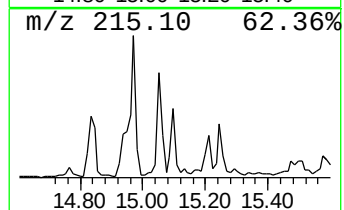
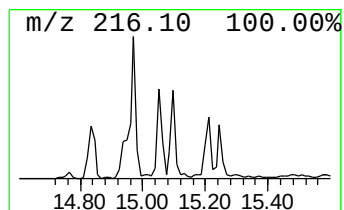
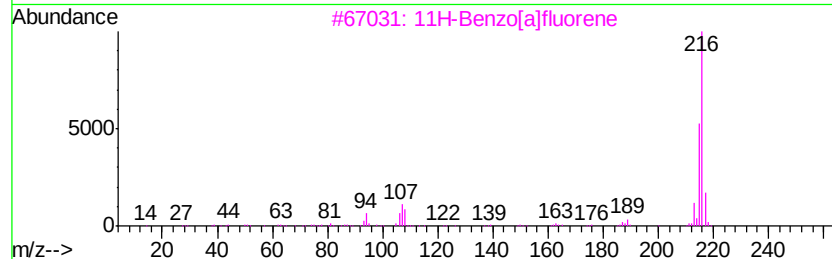
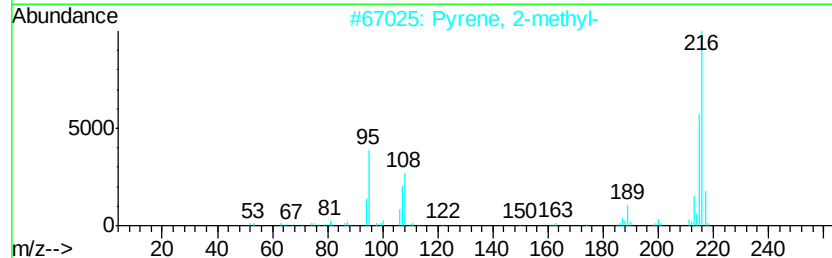
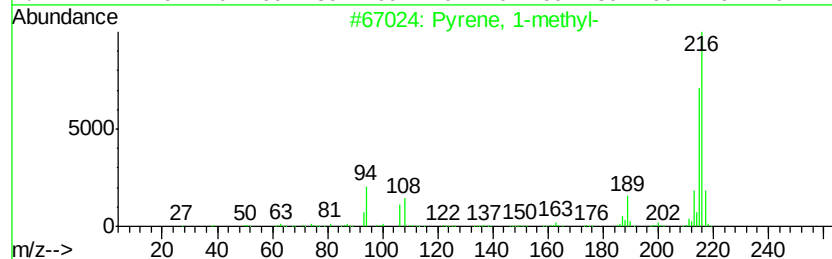
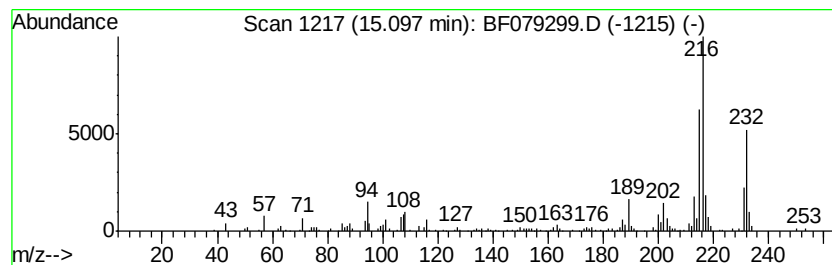
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Pyrene, 1-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.10	2.16 ng	154070	Chrysene-d12	16.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	92
2		Pyrene, 2-methyl-	216	C17H12	003442-78-2	83
3		11H-Benzofluorene	216	C17H12	000238-84-6	64
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	60
5		Pyrene, 4-methyl-	216	C17H12	003353-12-6	55



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051515\
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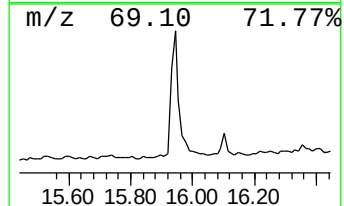
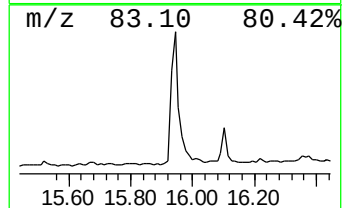
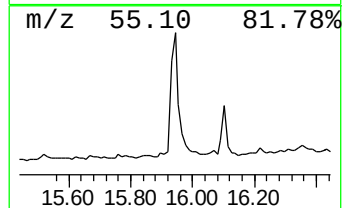
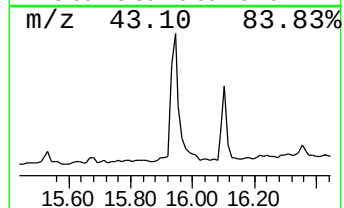
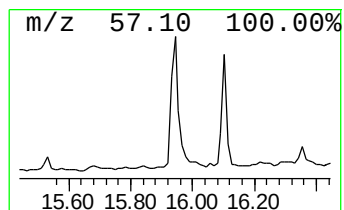
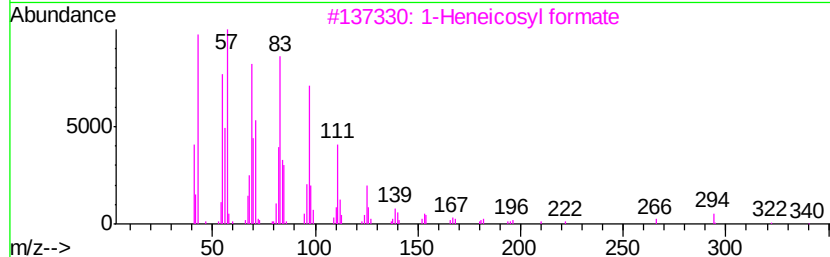
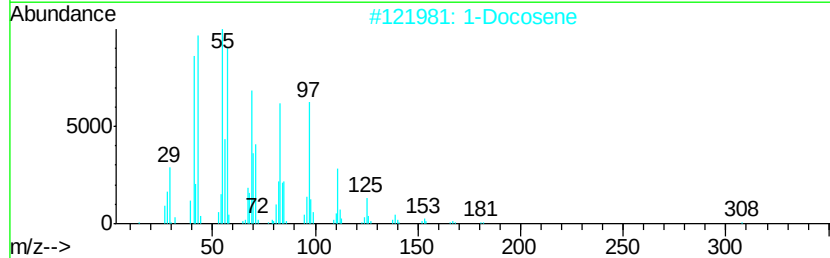
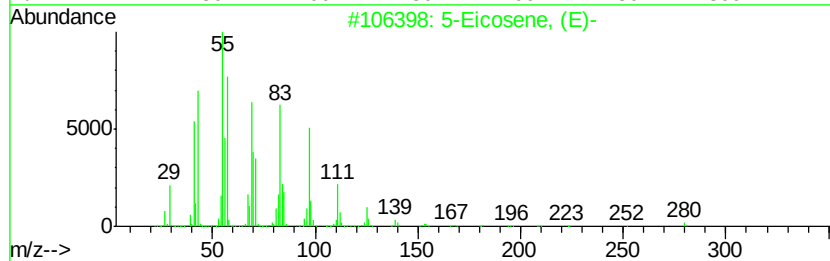
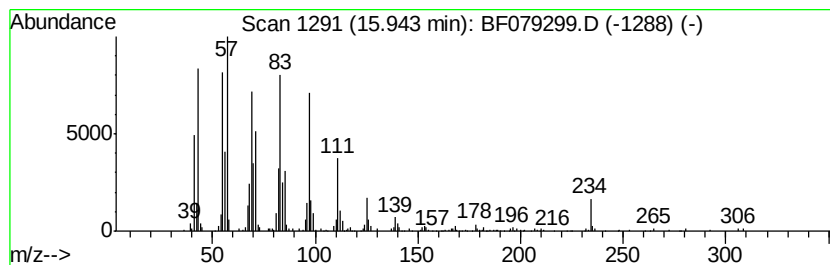
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 5-Eicosene, (E)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.94	8.56 ng	611698	Chrysene-d12	16.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Eicosene, (E)-	280	C20H40	074685-30-6	97
2		1-Docosene	308	C22H44	001599-67-3	96
3		1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
4		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	94
5		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051515\
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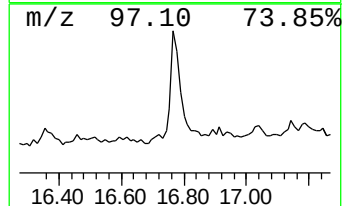
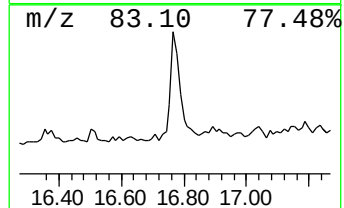
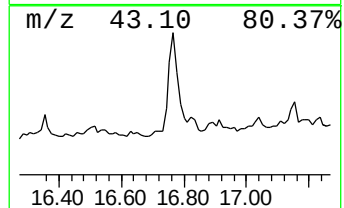
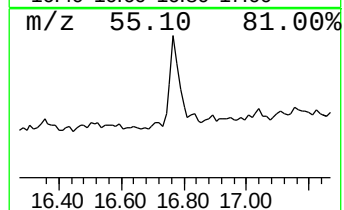
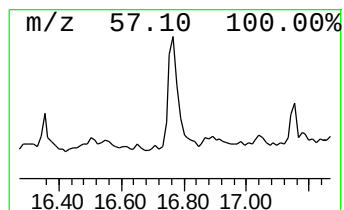
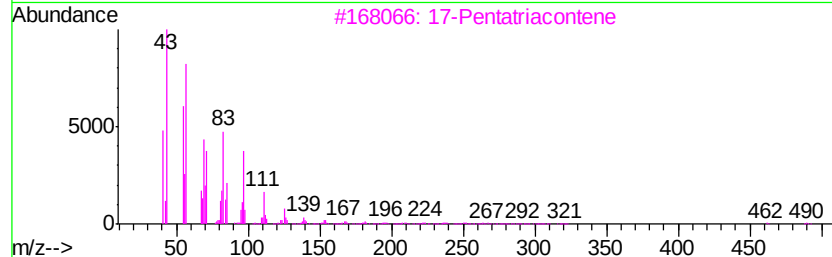
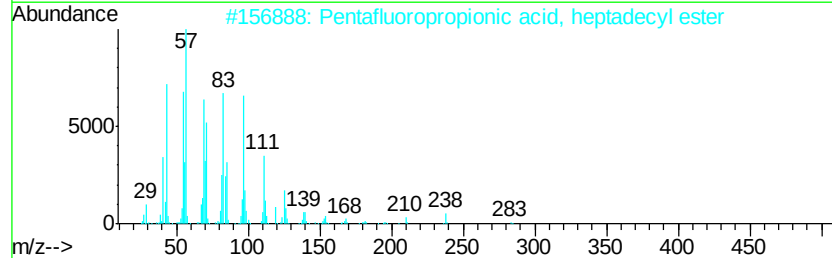
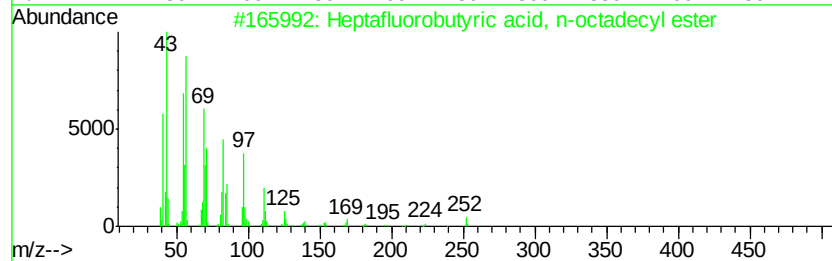
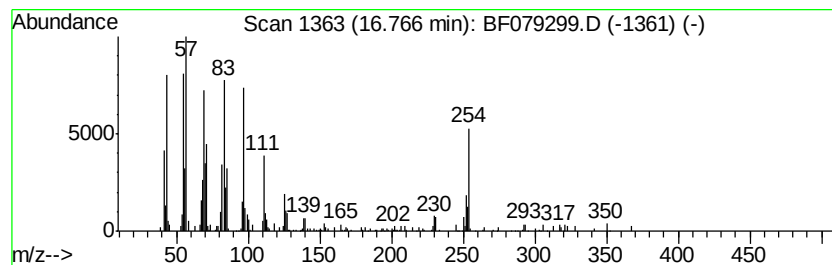
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Heptafluorobutyric acid, n-... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.77	3.17 ng	226470	Chrysene-d12	16.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptafluorobutyric acid, n-octadecyl...	466	C22H37F7O2	000400-57-7	80
2		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	74
3		17-Pentatriacontene	491	C35H70	006971-40-0	74
4		1-Docosene	308	C22H44	001599-67-3	74
5		Acetic acid, chloro-, octadecyl ...	346	C20H39ClO2	005348-82-3	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051515\
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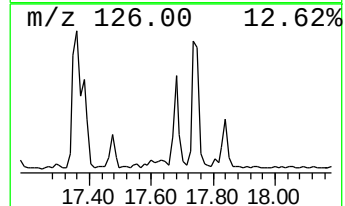
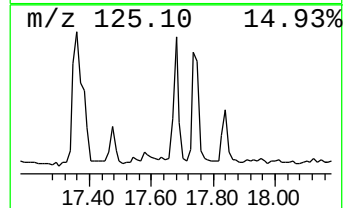
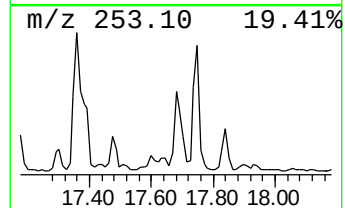
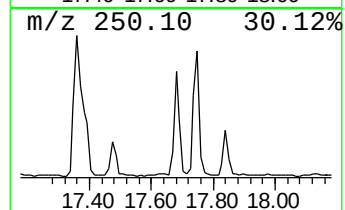
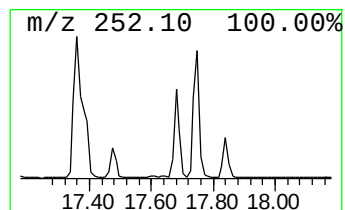
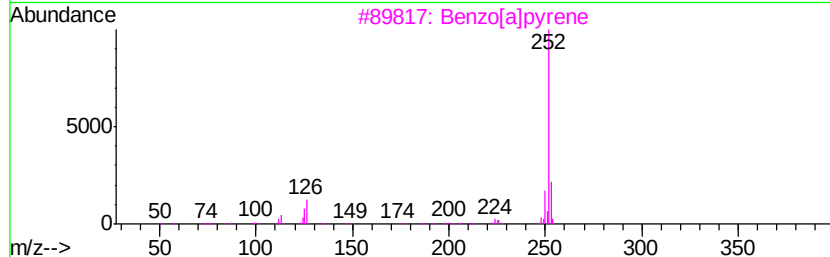
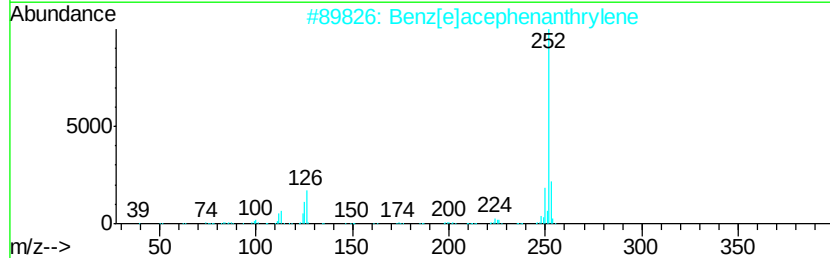
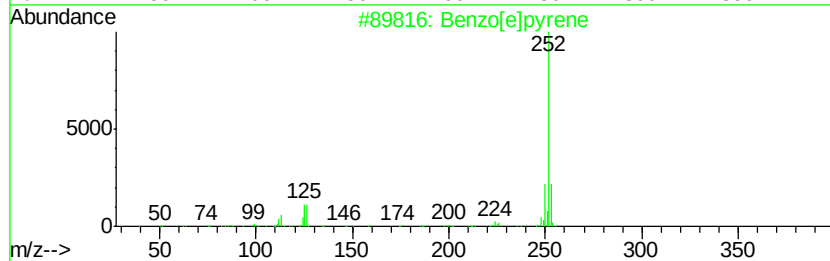
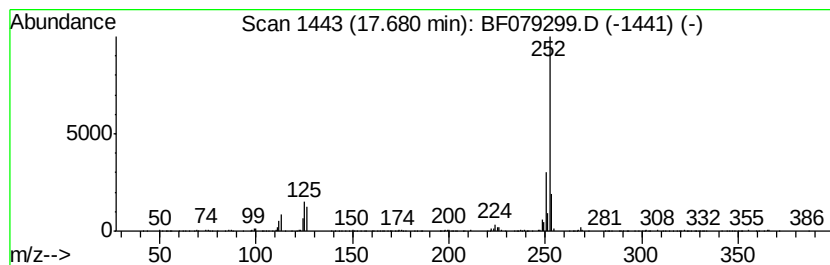
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Benzo[e]pyrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.68	9.60 ng	356479	Perylene-d12	17.82

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[e]pyrene	252	C20H12	000192-97-2	94
2		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	94
3		Benzo[a]pyrene	252	C20H12	000050-32-8	93
4		Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
5		Benzo[j]fluoranthene	252	C20H12	000205-82-3	90



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051515\
Data File : BF079299.D
Acq On : 16 May 2015 7:43
Operator : TP/IZ
Sample : G2249-02
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SR-2

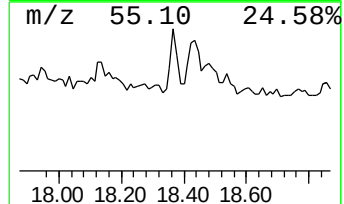
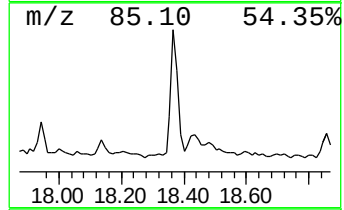
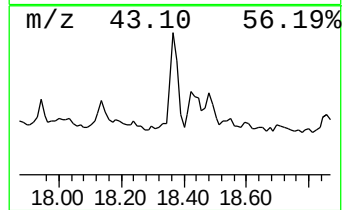
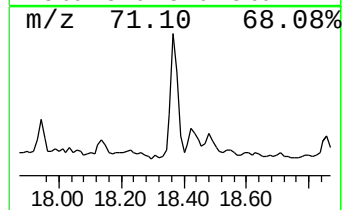
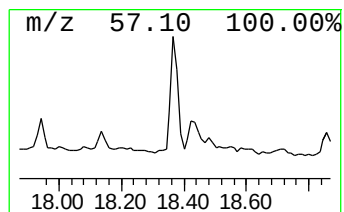
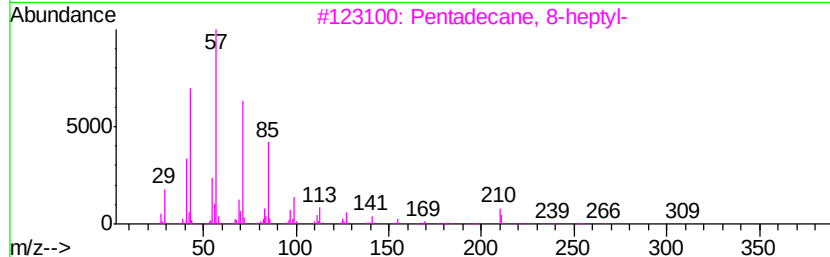
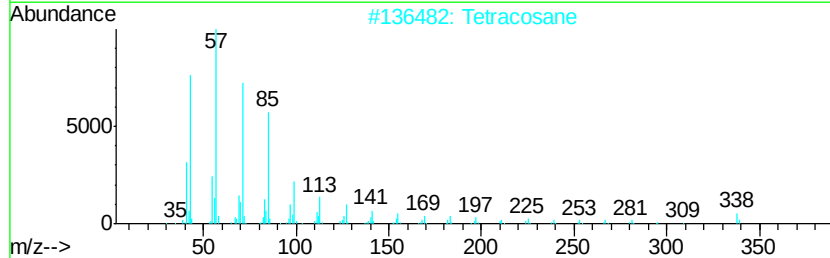
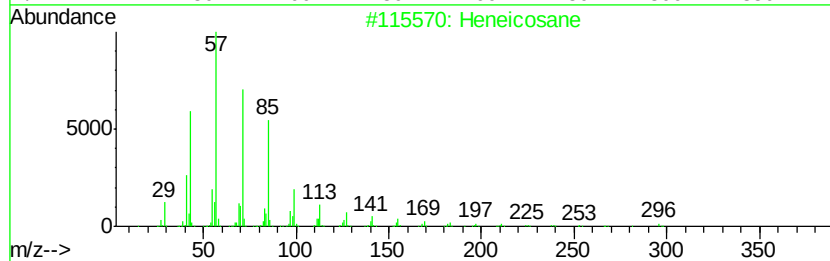
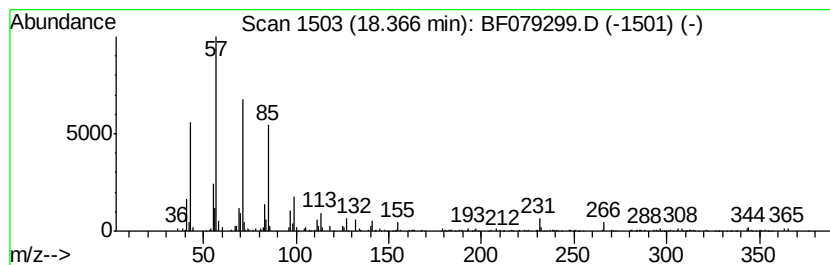
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Heneicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.37	4.89 ng	181524	Perylene-d12	17.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	97
2			Tetracosane	338	C24H50	000646-31-1	93
3			Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	87
4			Docosane, 11-butyl-	366	C26H54	013475-76-8	86
5			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051515\
Data File : BF079299.D
Acq On : 16 May 2015 7:43
Operator : TP/IZ
Sample : G2249-02
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SR-2

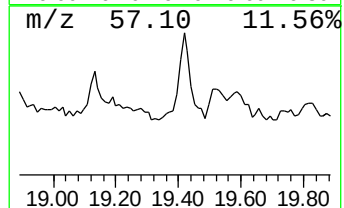
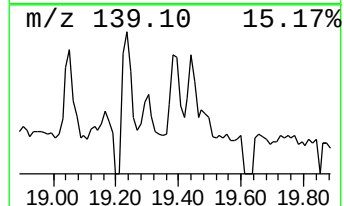
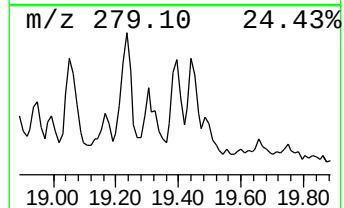
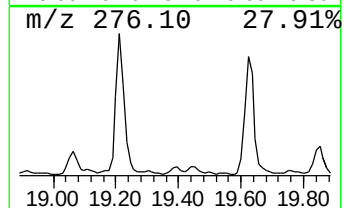
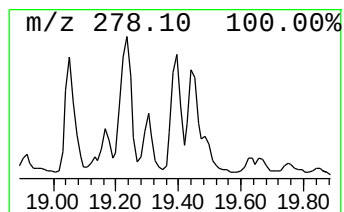
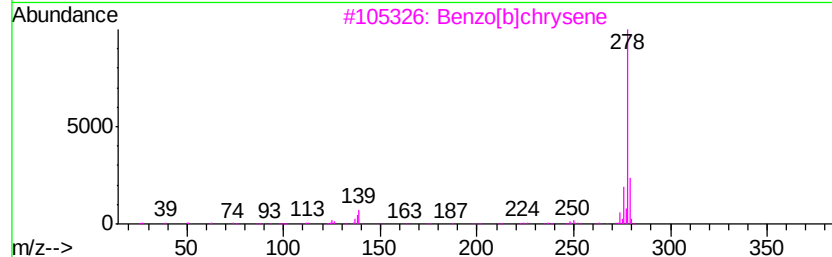
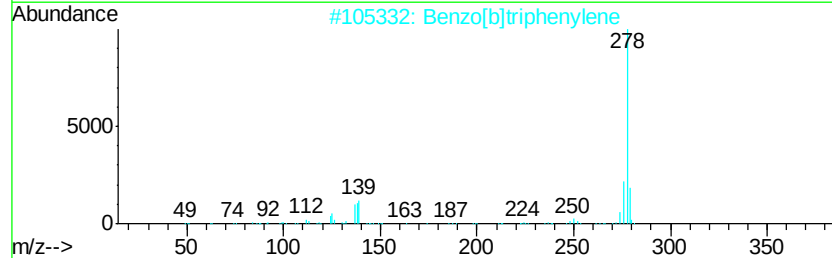
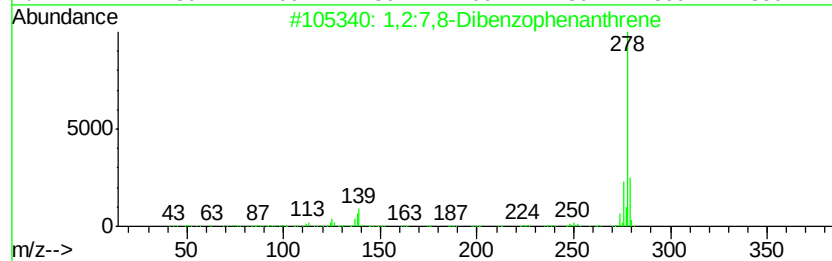
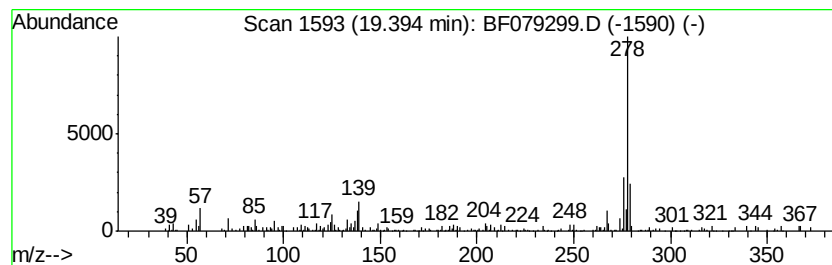
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 1,2:7,8-Dibenzophenanthrene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.39	4.36 ng	161934	Perylene-d12	17.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2:7,8-Dibenzophenanthrene	278	C22H14	000213-46-7	89
2		Benzo[b]triphenylene	278	C22H14	000215-58-7	86
3		Benzo[b]chrysene	278	C22H14	000214-17-5	76
4		Dibenz[a,h]anthracene	278	C22H14	000053-70-3	76
5		Pentaphene	278	C22H14	000222-93-5	68



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051515\
 Data File : BF079299.D
 Acq On : 16 May 2015 7:43
 Operator : TP/IZ
 Sample : G2249-02
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SR-2

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.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
unknown1.32	1.32	234.6	ng	5414380	1	7.18	461604	20.0
Cyclopentane, 1,1...	1.59	7.6	ng	175368	1	7.18	461604	20.0
Butane, 2-methoxy...	1.62	11.0	ng	253915	1	7.18	461604	20.0
2-Pentanone, 4-hy...	4.95	23.3	ng	537136	1	7.18	461604	20.0
unknown6.89	6.89	83.8	ng	1934660	1	7.18	461604	20.0
1H-Cyclopropa[l]p...	13.49	5.0	ng	169799	4	12.77	679025	20.0
4H-Cyclopenta[def...	13.53	6.7	ng	226292	4	12.77	679025	20.0
11H-Benzo[b]fluorene	14.97	2.5	ng	178139	5	16.06	1429410	20.0
Pyrene, 1-methyl-	15.10	2.2	ng	154070	5	16.06	1429410	20.0
5-Eicosene, (E)-	15.94	8.6	ng	611698	5	16.06	1429410	20.0
Heptafluorobutyri...	16.77	3.2	ng	226470	5	16.06	1429410	20.0
Benzo[e]pyrene	17.68	9.6	ng	356479	6	17.82	742401	20.0
Heneicosane	18.37	4.9	ng	181524	6	17.82	742401	20.0
1,2:7,8-Dibenzoph...	19.39	4.4	ng	161934	6	17.82	742401	20.0