

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032415\
 Data File : BF077943.D
 Acq On : 25 Mar 2015 8:22
 Operator : TP/IZ
 Sample : G1615-11
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-22

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-BF031115.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	1.047	11	14	17	rVB	568632	861145	22.25%	4.839%
2	1.264	31	33	36	rBV	3161786	3870061	100.00%	21.748%
3	1.390	40	44	47	rVB	529723	759998	19.64%	4.271%
4	1.550	56	58	61	rVB	47436	57909	1.50%	0.325%
5	1.767	74	77	83	rBV	55301	82605	2.13%	0.464%
6	4.864	346	348	353	rVB	102749	127086	3.28%	0.714%
7	5.401	393	395	398	rBV	678935	821246	21.22%	4.615%
8	6.693	506	508	511	rBV	641953	874639	22.60%	4.915%
9	6.830	518	520	523	rBV	973651	933388	24.12%	5.245%
10	7.116	543	545	548	rBV	262930	238706	6.17%	1.341%
11	7.299	559	561	564	rBV	617318	739791	19.12%	4.157%
12	7.813	604	606	608	rBV	650884	566027	14.63%	3.181%
13	8.693	681	683	686	rBV	371767	332971	8.60%	1.871%
14	10.042	799	801	804	rBV	1357338	1197211	30.94%	6.728%
15	10.556	843	846	848	rBV	144053	189276	4.89%	1.064%
16	10.865	871	873	875	rVB	448255	400945	10.36%	2.253%
17	11.848	956	959	961	rBV2	462784	643700	16.63%	3.617%
18	12.042	974	976	980	rBV	42223	55398	1.43%	0.311%
19	12.602	1023	1025	1027	rBV	68724	67902	1.75%	0.382%
20	12.705	1031	1034	1035	rVV	354662	431211	11.14%	2.423%
21	12.728	1035	1036	1039	rVB	74345	66405	1.72%	0.373%
22	13.139	1070	1072	1074	rVB	32002	42281	1.09%	0.238%
23	14.202	1162	1165	1167	rBV	129326	128369	3.32%	0.721%
24	14.477	1187	1189	1191	rBV	108608	129919	3.36%	0.730%
25	14.694	1205	1208	1211	rBV	1379198	1390165	35.92%	7.812%
26	15.037	1235	1238	1240	rBV3	22378	39684	1.03%	0.223%
27	15.334	1260	1264	1267	rBV2	38667	67661	1.75%	0.380%
28	15.882	1309	1312	1317	rBV2	85759	173635	4.49%	0.976%
29	15.974	1317	1320	1322	rBV	377049	568504	14.69%	3.195%
30	16.008	1322	1323	1325	rVB	46934	46337	1.20%	0.260%
31	16.637	1376	1378	1379	rBV	27172	40622	1.05%	0.228%
32	17.037	1411	1413	1414	rBV2	26968	40936	1.06%	0.230%
33	17.151	1421	1423	1428	rBV	29207	47849	1.24%	0.269%
34	17.265	1431	1433	1438	rVB	55458	142306	3.68%	0.800%

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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-BF031115.M
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35	17.471	1449	1451	1453	rBV2	29131	54927	1.42%	0.309%
36	17.586	1459	1461	1464	rBV	29576	53022	1.37%	0.298%
37	17.654	1464	1467	1470	rVB	63417	90180	2.33%	0.507%
38	17.711	1470	1472	1476	rBV	339588	538025	13.90%	3.024%
39	19.266	1604	1608	1612	rBV	378034	784571	20.27%	4.409%
40	19.574	1632	1635	1639	rVB4	46636	98115	2.54%	0.551%

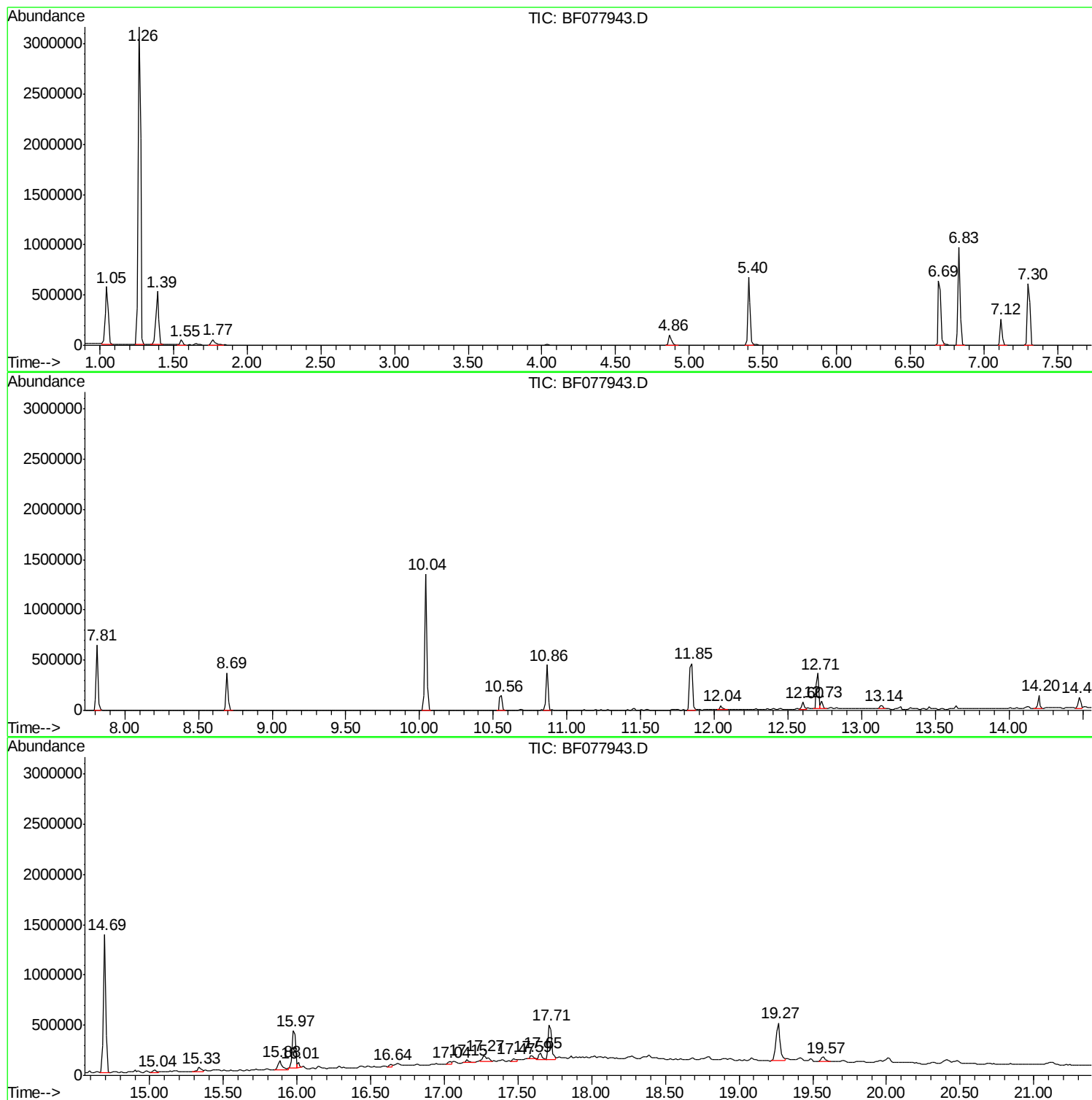
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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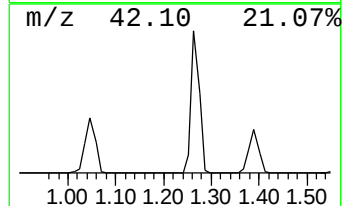
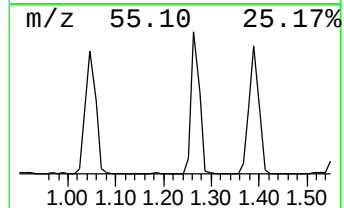
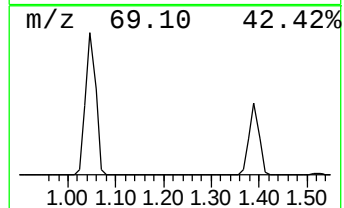
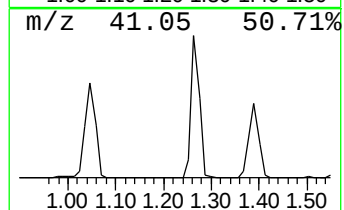
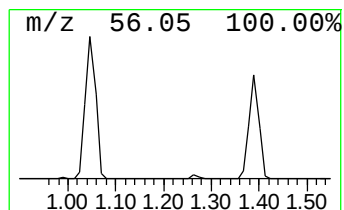
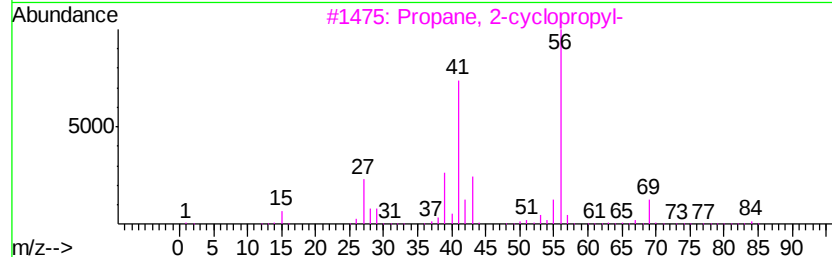
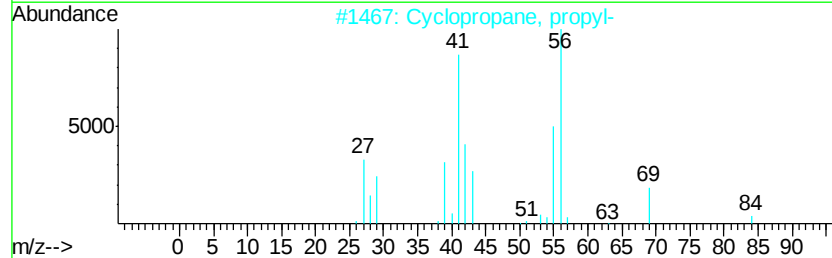
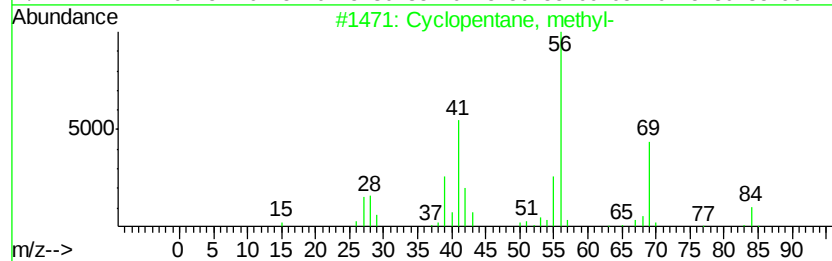
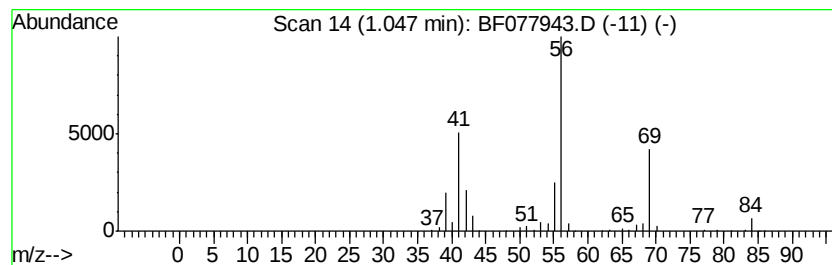
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Cyclopentane, methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.05	72.15 ng	861145	1,4-Dichlorobenzene-d4	7.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	87
2		Cyclopropane, propyl-	84	C6H12	002415-72-7	78
3		Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	64
4		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
5		Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	56



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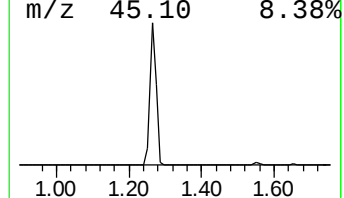
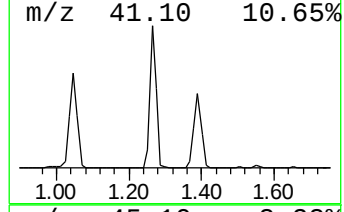
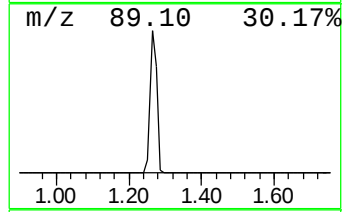
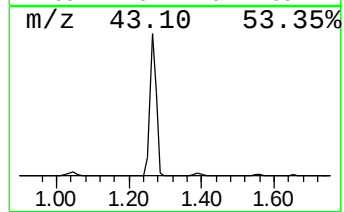
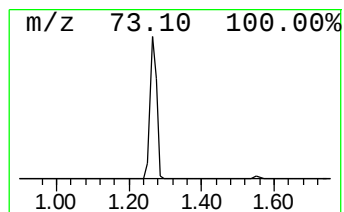
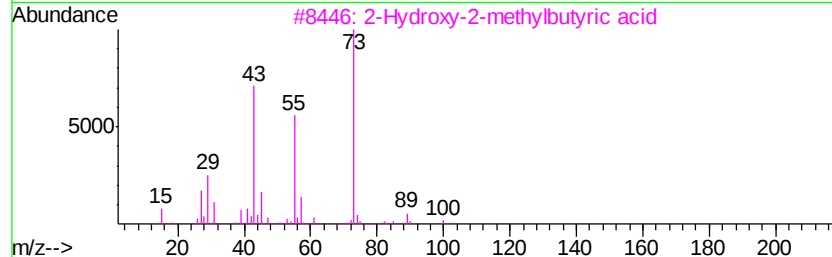
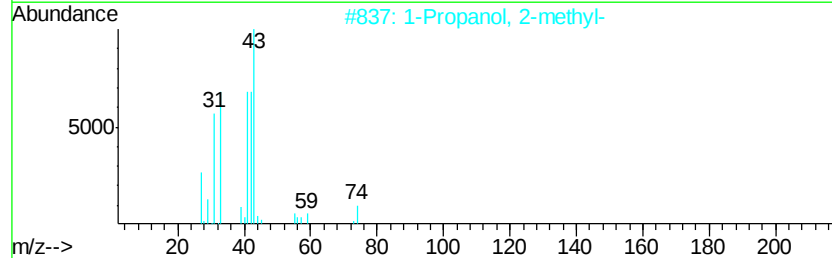
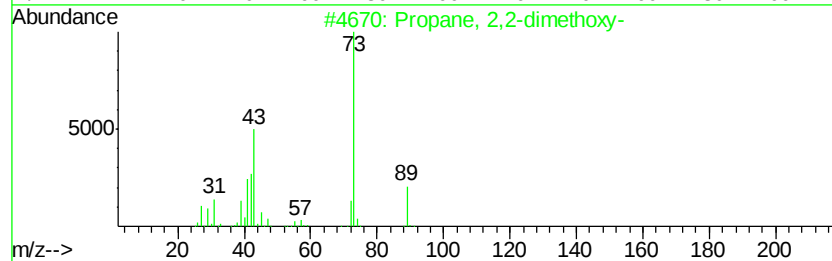
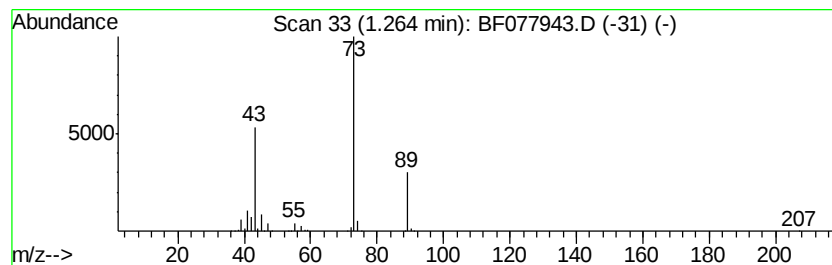
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.26	324.25 ng	3870060	1,4-Dichlorobenzene-d4	7.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	72
2		1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	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		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9



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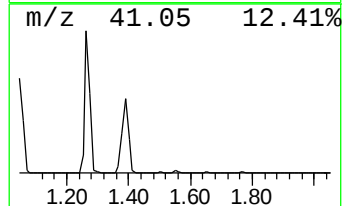
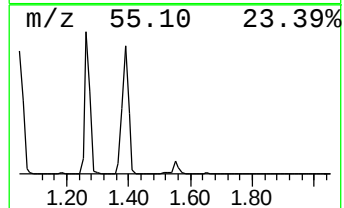
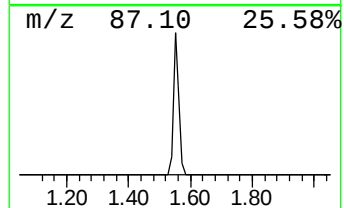
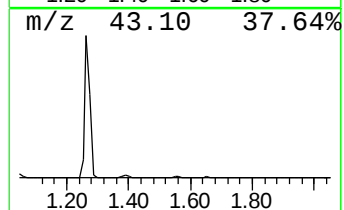
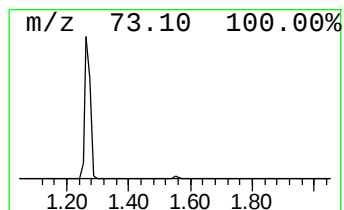
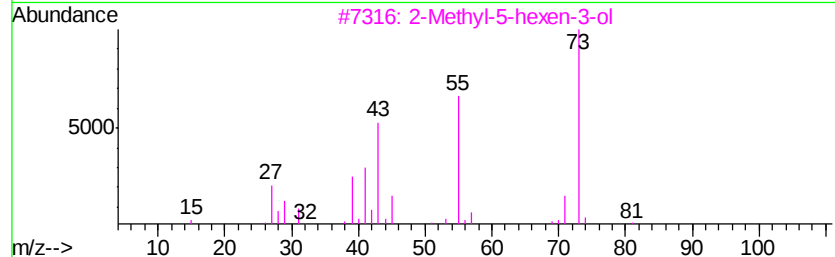
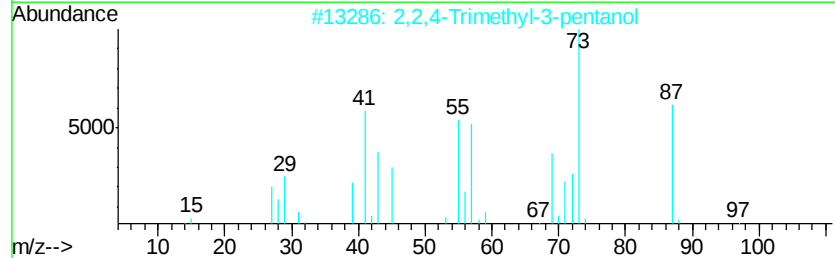
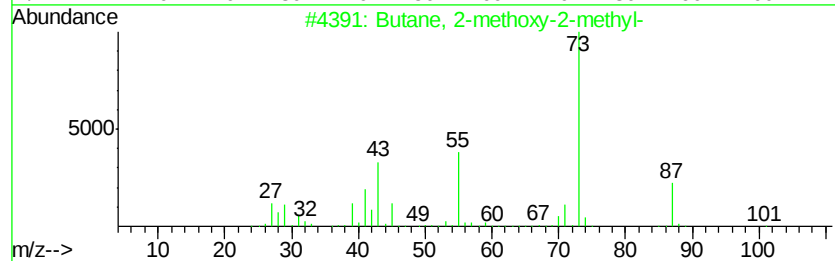
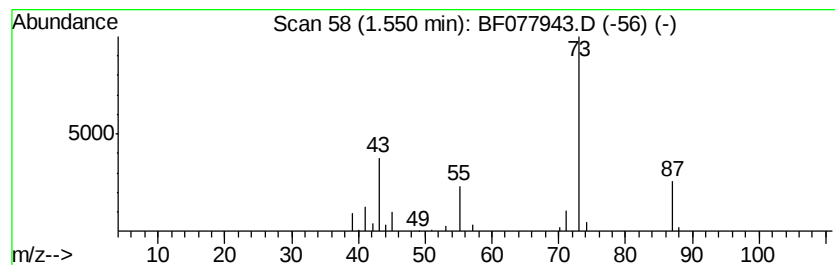
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Butane, 2-methoxy-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.55	4.85 ng	57909	1,4-Dichlorobenzene-d4	7.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	32
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5		Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	9



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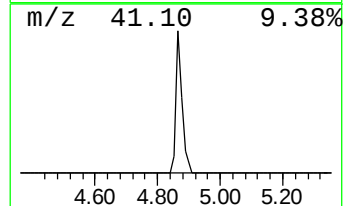
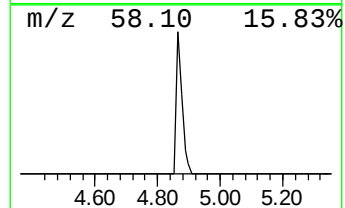
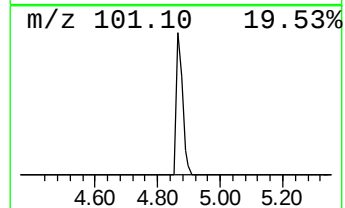
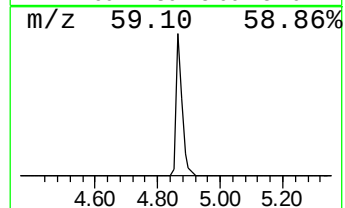
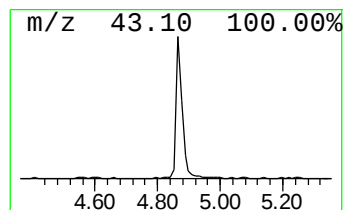
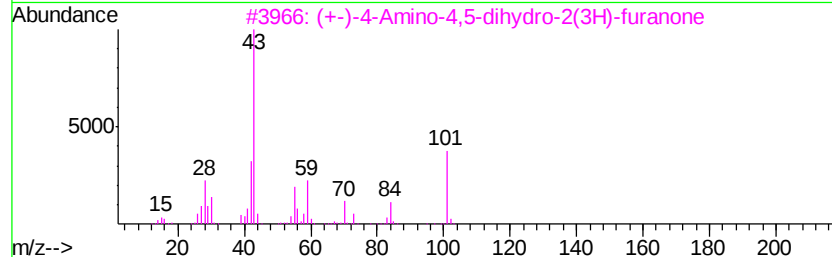
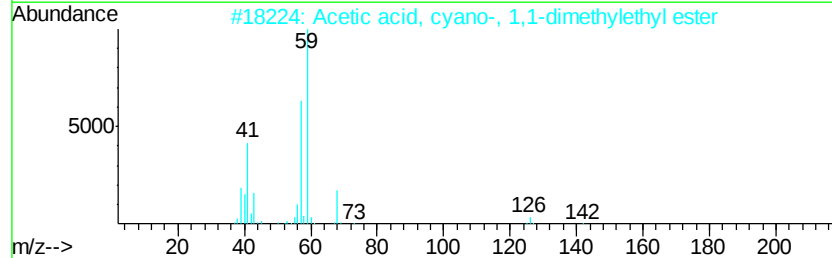
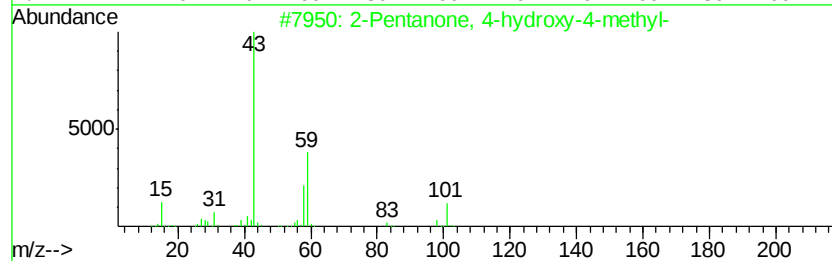
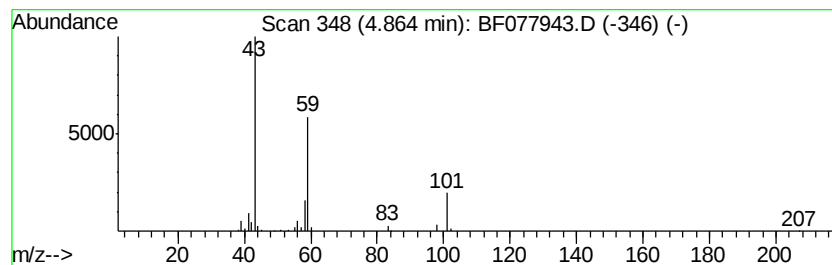
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.86	10.65 ng	127086	1,4-Dichlorobenzene-d4	7.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
3		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



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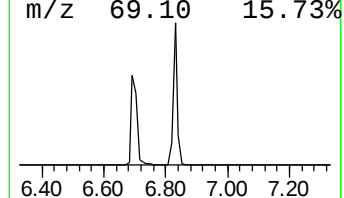
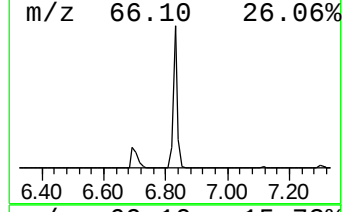
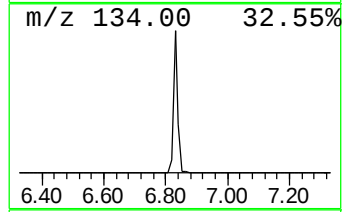
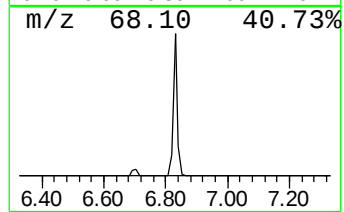
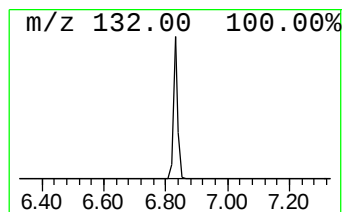
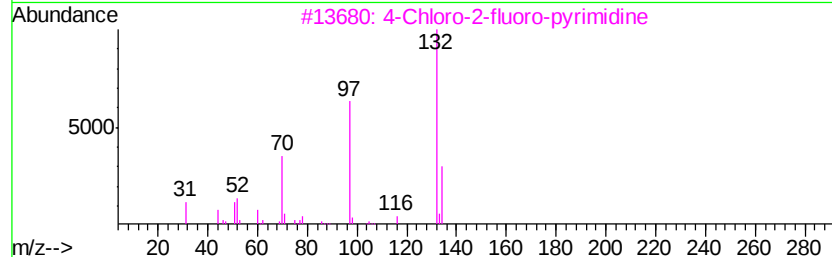
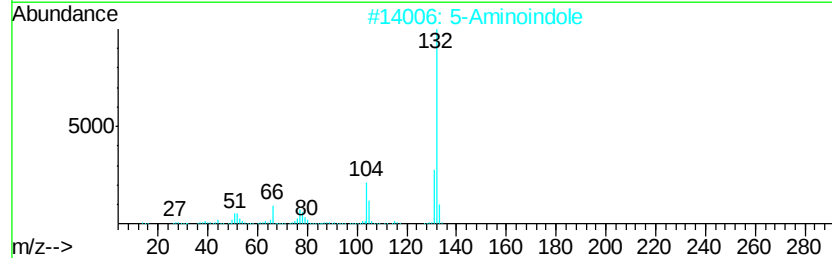
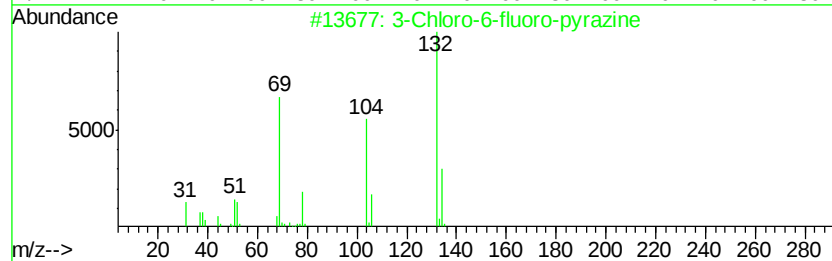
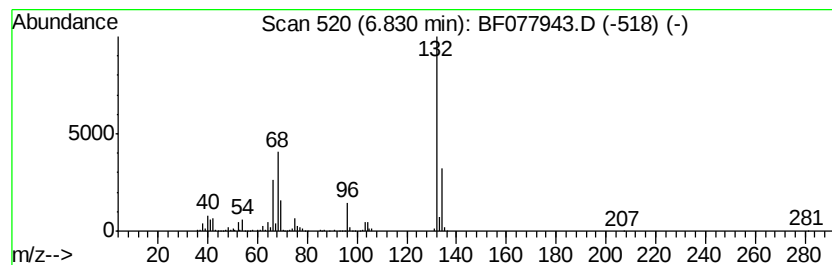
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown6.83 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.83	78.20 ng	933388	1,4-Dichlorobenzene-d4	7.12

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		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
4		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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Sample : G1615-11
Misc :
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ClientSampleId :
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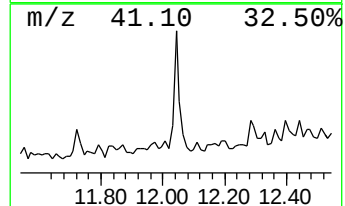
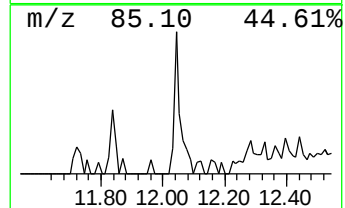
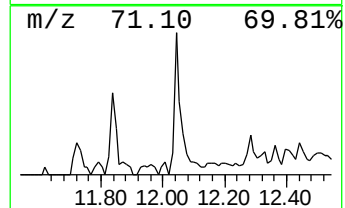
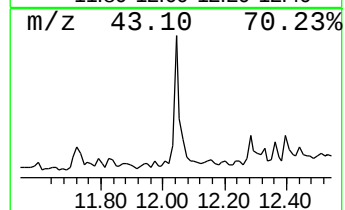
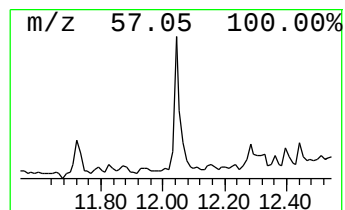
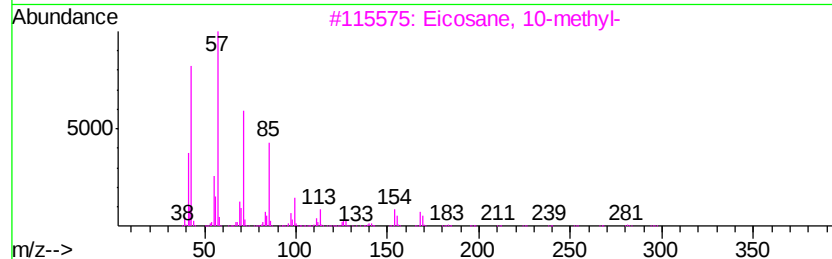
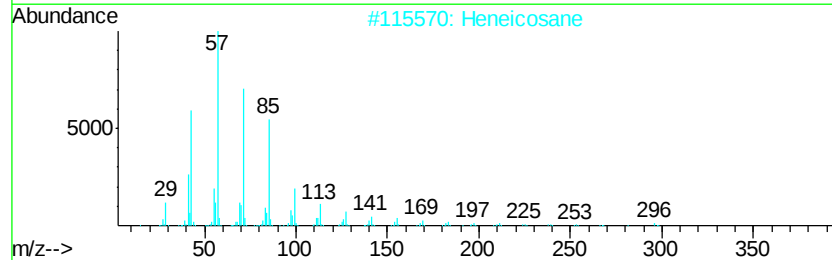
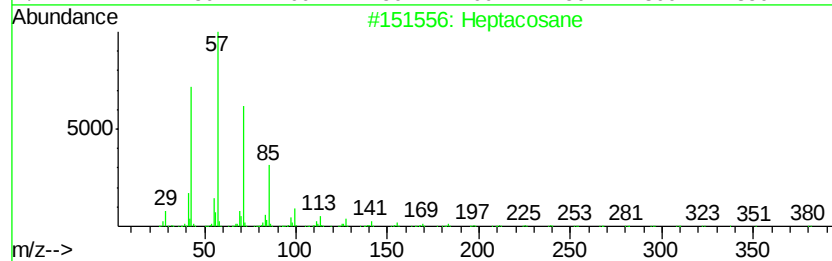
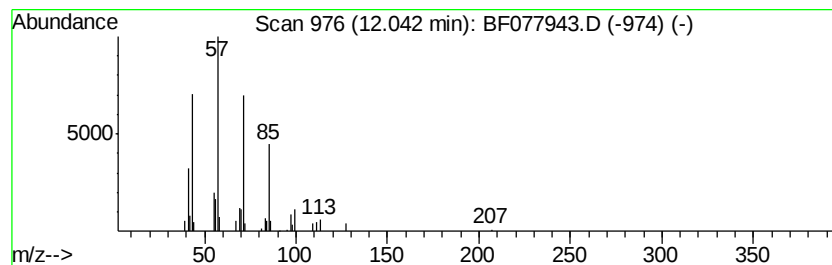
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptacosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.04	2.57 ng	55398	Phenanthrene-d10	12.71

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane		380	C27H56	000593-49-7	90
2	Heneicosane		296	C21H44	000629-94-7	90
3	Eicosane, 10-methyl-		296	C21H44	054833-23-7	86
4	Octacosane		394	C28H58	000630-02-4	86
5	Pentadecane		212	C15H32	000629-62-9	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032415\
Data File : BF077943.D
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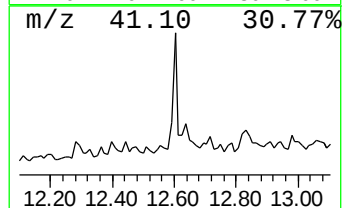
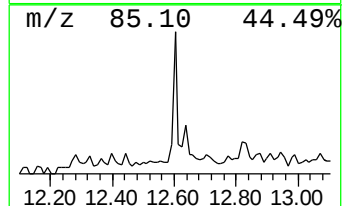
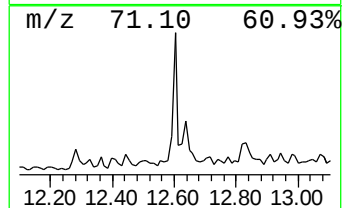
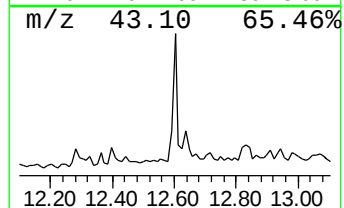
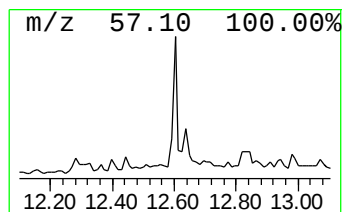
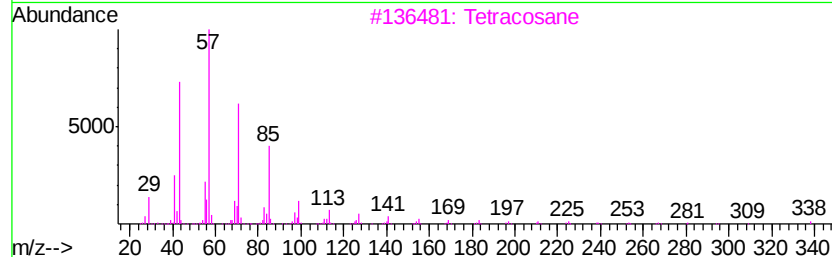
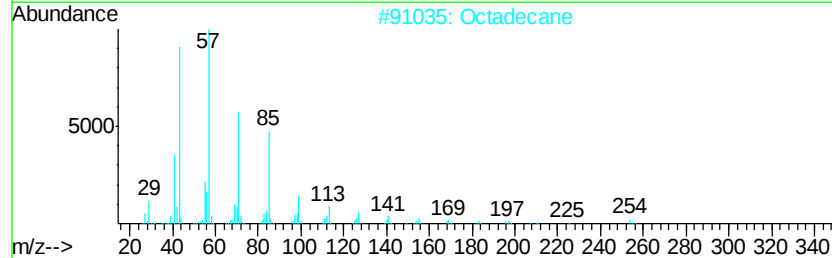
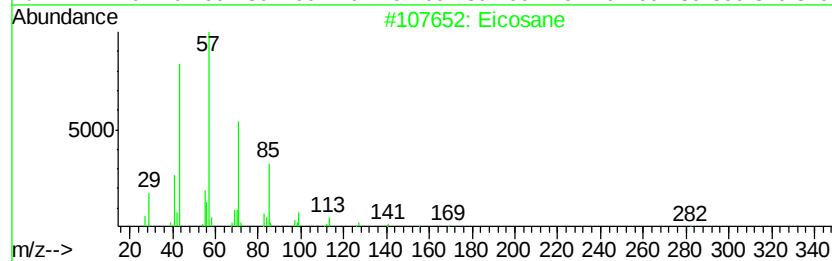
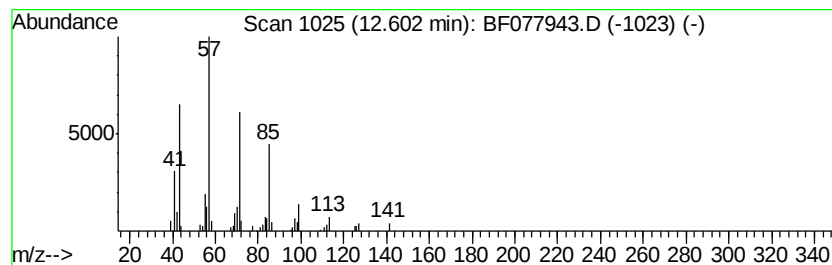
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Eicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.60	3.15 ng	67902	Phenanthrene-d10	12.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane			282	C20H42	000112-95-8	91
2	Octadecane			254	C18H38	000593-45-3	91
3	Tetracosane			338	C24H50	000646-31-1	91
4	Tetradecane			198	C14H30	000629-59-4	90
5	Pentadecane			212	C15H32	000629-62-9	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032415\
Data File : BF077943.D
Acq On : 25 Mar 2015 8:22
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Misc :
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ClientSampleId :
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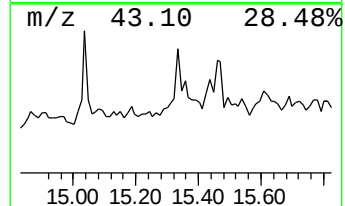
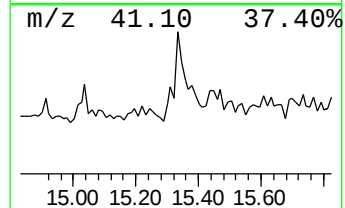
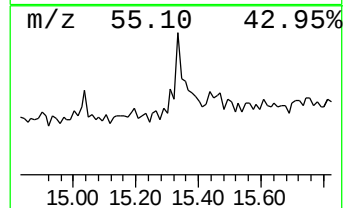
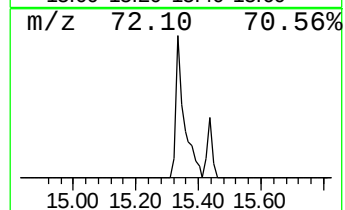
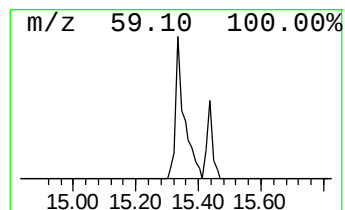
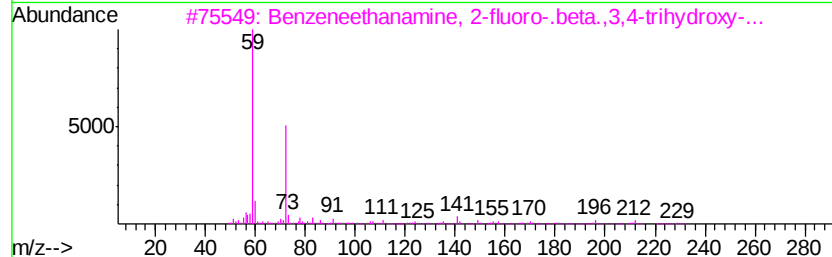
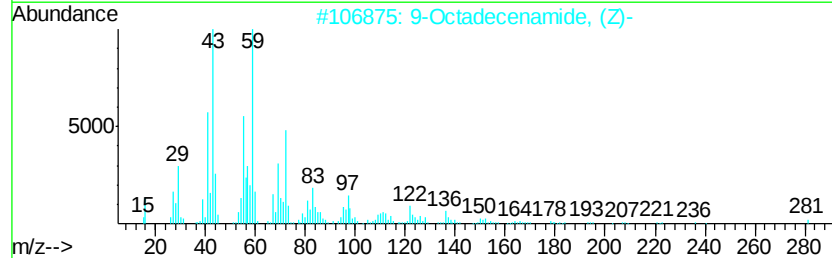
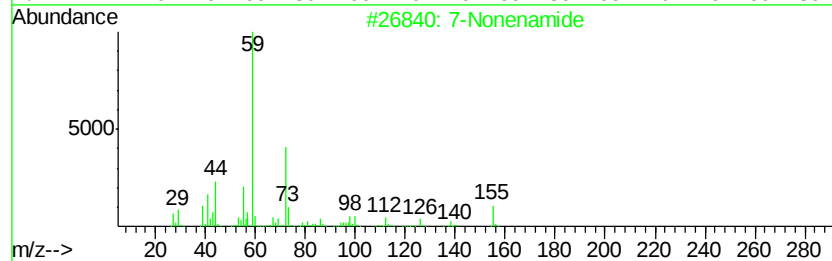
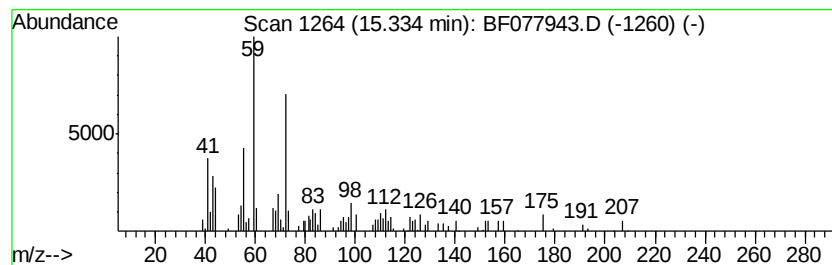
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 7-Nonenamide Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.33	2.38 ng	67661	Chrysene-d12	15.99

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7-Nonenamide	155	C9H17NO	090949-53-4	64
2			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	64
3			Benzeneethanamine, 2-fluoro-.bet...	229	C11H16FN03	061338-98-5	42
4			Nonanamide	157	C9H19NO	001120-07-6	40
5			Pentanamide, 4-methyl-	115	C6H13NO	001119-29-5	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032415\
Data File : BF077943.D
Acq On : 25 Mar 2015 8:22
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Sample : G1615-11
Misc :
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ClientSampleId :
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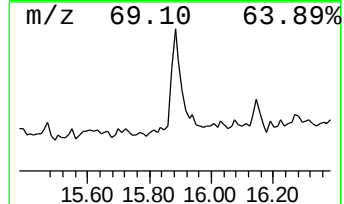
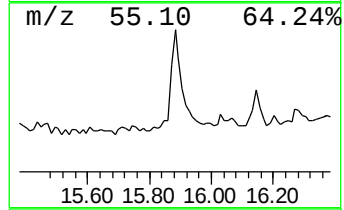
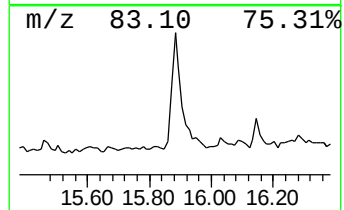
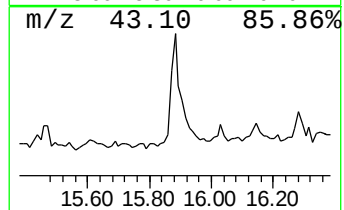
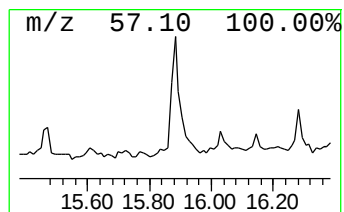
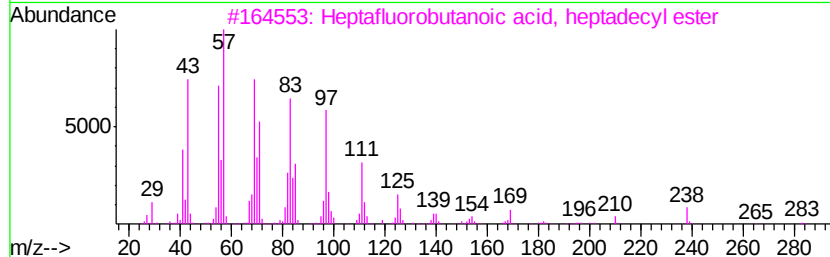
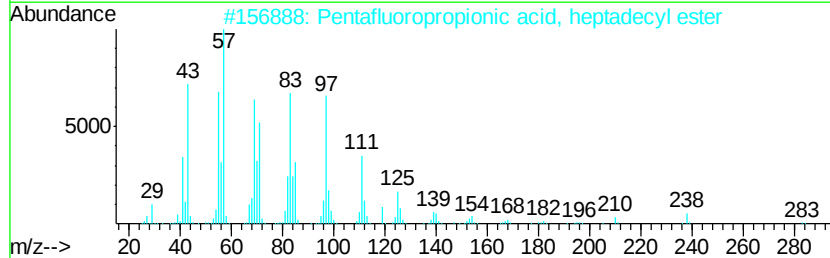
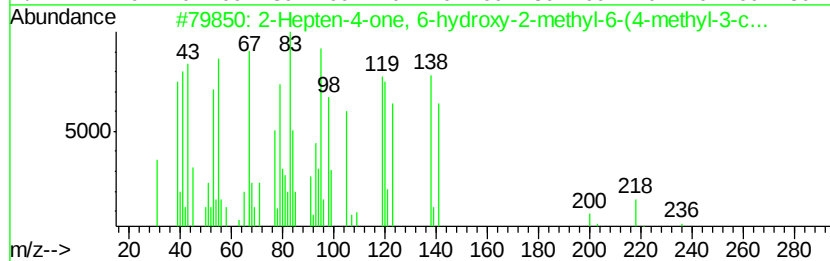
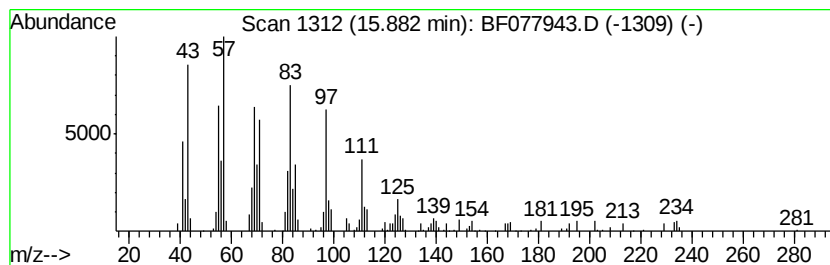
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 2-Hepten-4-one, 6-hydroxy-2... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.88	6.11 ng	173635	Chrysene-d12	15.99

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Hepten-4-one, 6-hydroxy-2-meth...	236	C15H24O2	038043-98-0	93
2			Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	93
3			Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	90
4			2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	90
5			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032415\
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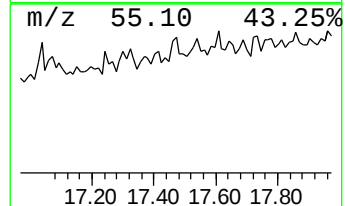
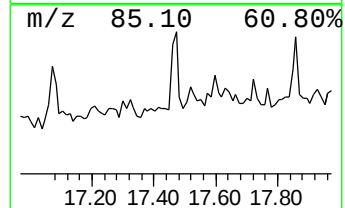
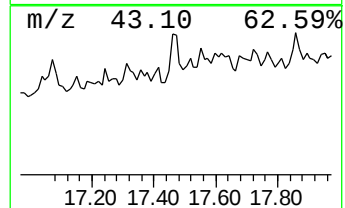
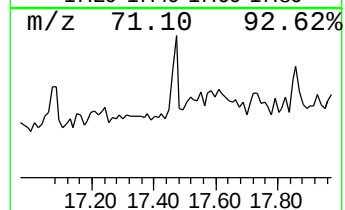
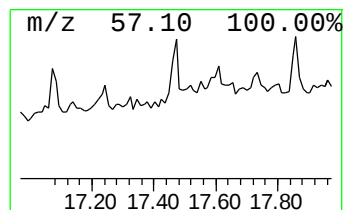
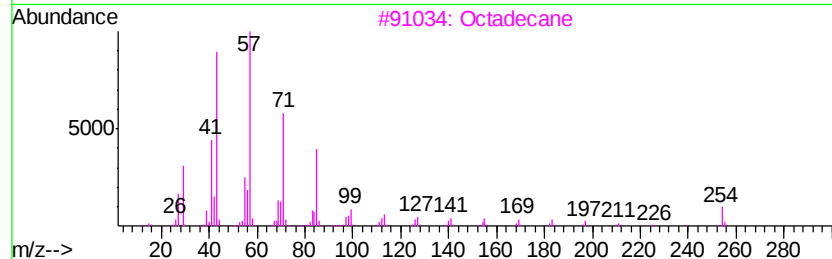
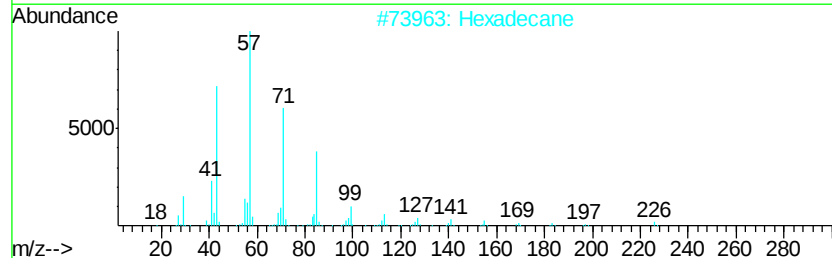
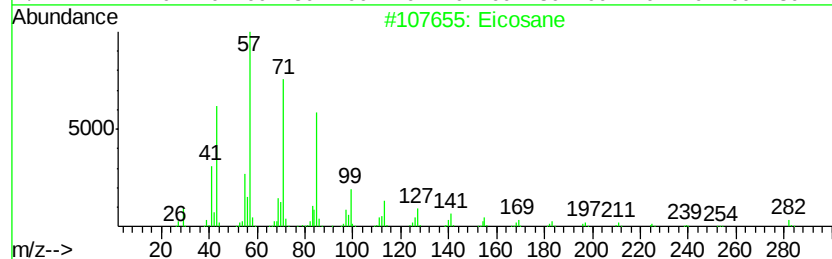
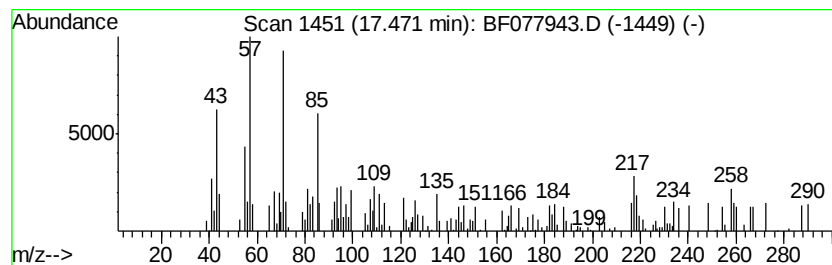
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hexadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.47	2.04 ng	54927	Perylene-d12	17.71
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane	282 C20H42	000112-95-8	55
2	Hexadecane	226 C16H34	000544-76-3	55
3	Octadecane	254 C18H38	000593-45-3	53
4	Tridecane	184 C13H28	000629-50-5	49
5	Tetratriacontane	479 C34H70	014167-59-0	49



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032415\
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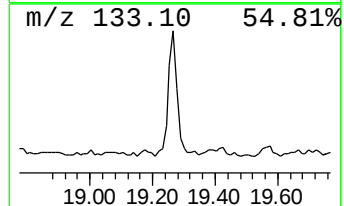
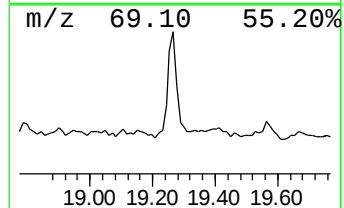
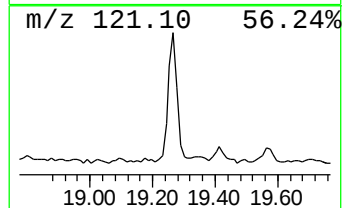
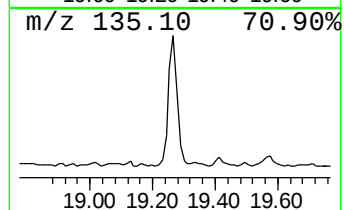
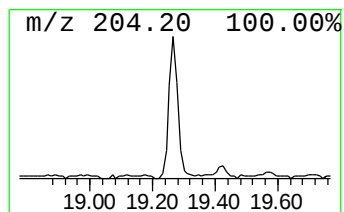
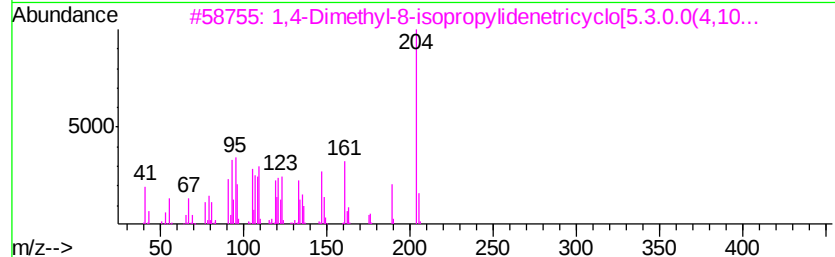
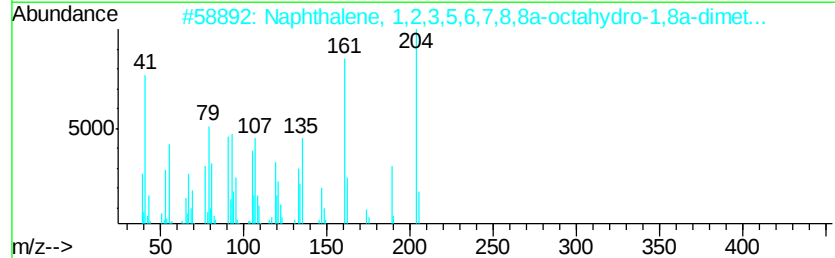
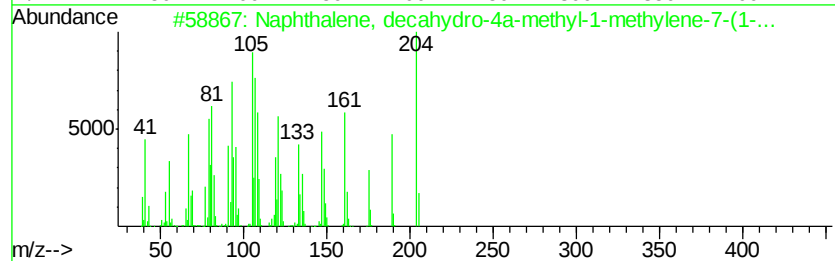
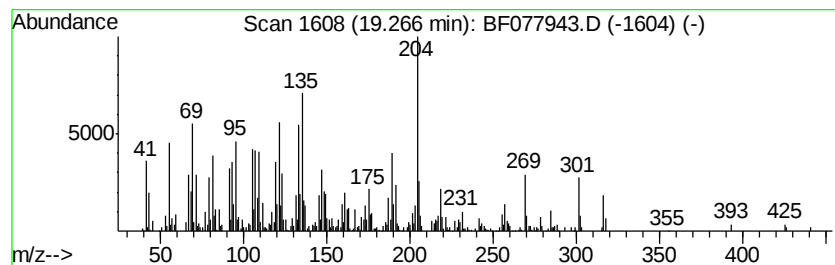
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Naphthalene, decahydro-4a-m... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.27	29.16 ng	784571	Perylene-d12	17.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, decahydro-4a-methyl...	204	C15H24	017066-67-0	55
2		Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204	C15H24	004630-07-3	53
3		1,4-Dimethyl-8-isopropylidenetri...	204	C15H24	1000140-07-7	53
4		1,1,4a-Trimethyl-5,6-dimethylene...	204	C15H24	1000193-60-8	50
5		1H-3a,7-Methanoazulene, 2,3,6,7,...	204	C15H24	000560-32-7	49



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032415\
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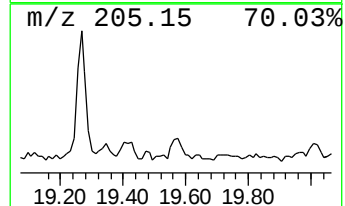
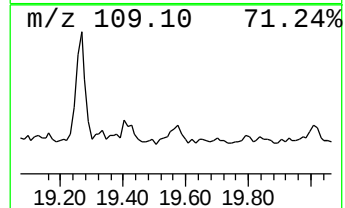
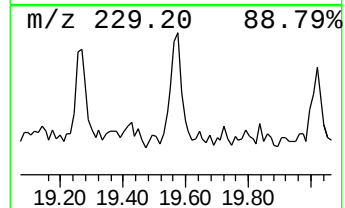
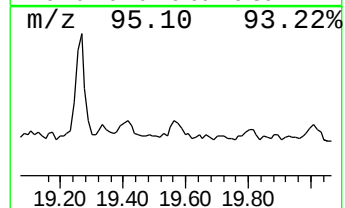
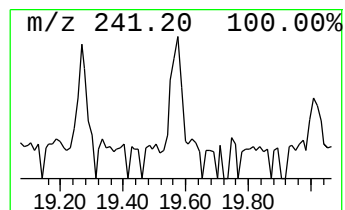
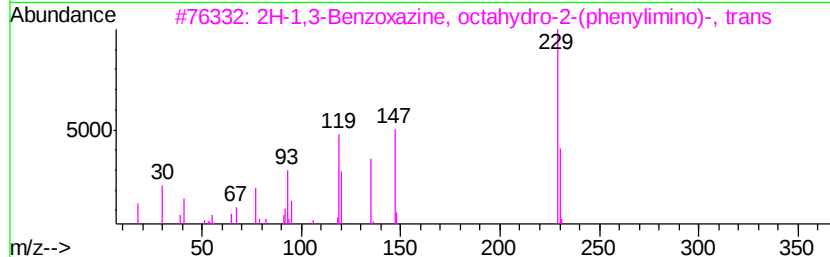
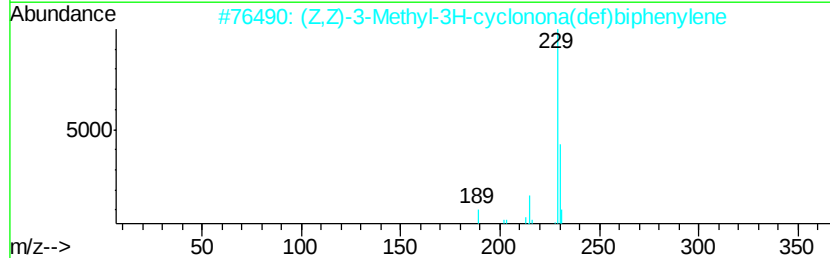
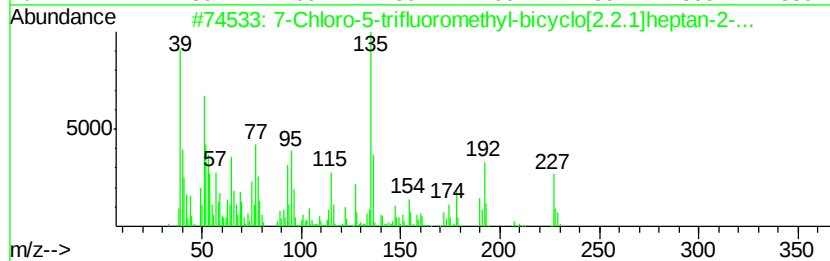
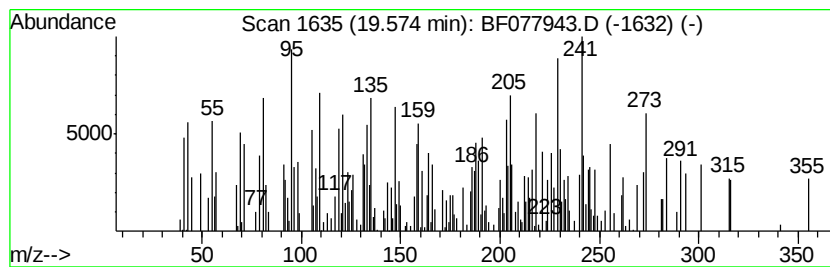
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 7-Chloro-5-trifluoromethyl-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.57	3.65 ng	98115	Perylene-d12	17.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	7-Chloro-5-trifluoromethyl-bicyc...	227	C8H9ClF3NO	1000278-15-7	90
2		(Z,Z)-3-Methyl-3H-cyclonona(def)...	230	C18H14	110823-80-8	27
3		2H-1,3-Benzoxazine, octahydro-2-...	230	C14H18N2O	1000138-91-8	15
4		5-Cyano-2-methyl-4-methylthio-6-...	241	C13H11N3S	068364-38-5	11
5		.beta.-Chlordene	336	C10H6Cl6	056534-03-3	11



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032415\
Data File : BF077943.D
Acq On : 25 Mar 2015 8:22
Operator : TP/IZ
Sample : G1615-11
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-22

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.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
Cyclopentane, met...	1.05	72.2	ng	861145	1	7.12	238706	20.0
Propane, 2,2-dime...	1.26	324.3	ng	3870060	1	7.12	238706	20.0
Butane, 2-methoxy...	1.55	4.8	ng	57909	1	7.12	238706	20.0
2-Pentanone, 4-hy...	4.86	10.7	ng	127086	1	7.12	238706	20.0
unknown6.83	6.83	78.2	ng	933388	1	7.12	238706	20.0
Heptacosane	12.04	2.6	ng	55398	4	12.71	431211	20.0
Eicosane	12.60	3.1	ng	67902	4	12.71	431211	20.0
7-Nonenamide	15.33	2.4	ng	67661	5	15.99	568504	20.0
2-Hepten-4-one, 6...	15.88	6.1	ng	173635	5	15.99	568504	20.0
Hexadecane	17.47	2.0	ng	54927	6	17.71	538025	20.0
Naphthalene, deca...	19.27	29.2	ng	784571	6	17.71	538025	20.0
7-Chloro-5-triflu...	19.57	3.6	ng	98115	6	17.71	538025	20.0