

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092518\
 Data File : BF109294.D
 Acq On : 25 Sep 2018 19:03
 Operator : JU/SJ
 Sample : J5055-16
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-D

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.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	4.916	477	481	489	rBV	80292	114293	3.55%	0.391%
2	5.328	545	551	554	rBV	2747414	2827143	87.86%	9.665%
3	6.351	720	725	737	rVB	2623952	3102340	96.41%	10.606%
4	6.481	742	747	750	rBV	3014876	2980137	92.62%	10.188%
5	6.710	782	786	789	rBV	958079	737758	22.93%	2.522%
6	6.869	809	813	816	rBV	2708616	2309900	71.79%	7.897%
7	7.275	876	882	885	rBV	1967105	1883441	58.53%	6.439%
8	7.992	1000	1004	1007	rBV	1156591	912344	28.35%	3.119%
9	9.069	1182	1187	1190	rBV	3691104	3217762	100.00%	11.001%
10	9.151	1197	1201	1205	rBV5	65781	82239	2.56%	0.281%
11	9.192	1205	1208	1211	rBV3	40879	45704	1.42%	0.156%
12	9.404	1242	1244	1250	rBV3	54428	75621	2.35%	0.259%
13	9.463	1250	1254	1257	rBV	1122504	966244	30.03%	3.303%
14	9.516	1260	1263	1267	rVV2	105381	115563	3.59%	0.395%
15	9.616	1277	1280	1283	rBV3	211264	249406	7.75%	0.853%
16	9.663	1285	1288	1291	rVB	426071	383067	11.90%	1.310%
17	9.698	1291	1294	1296	rBV2	136721	157976	4.91%	0.540%
18	9.745	1296	1302	1306	rVB	1087521	1198537	37.25%	4.097%
19	9.792	1307	1310	1315	rBV2	184388	276801	8.60%	0.946%
20	9.910	1328	1330	1332	rBV2	68343	73433	2.28%	0.251%
21	10.016	1346	1348	1349	rBV2	79062	62946	1.96%	0.215%
22	10.086	1355	1360	1363	rBV	297797	377919	11.74%	1.292%
23	10.121	1363	1366	1369	rBV3	242441	273502	8.50%	0.935%
24	10.157	1369	1372	1375	rBV	616558	493555	15.34%	1.687%
25	10.210	1378	1381	1382	rBV	174603	137725	4.28%	0.471%
26	10.527	1432	1435	1439	rBV	1533403	1651121	51.31%	5.645%
27	11.221	1551	1553	1556	rVB	759215	630411	19.59%	2.155%
28	12.804	1818	1822	1826	rBV	2403521	2281350	70.90%	7.799%
29	13.715	1975	1977	1987	rVB2	113751	186518	5.80%	0.638%
30	13.845	1996	1999	2003	rVB	732669	692808	21.53%	2.369%
31	14.351	2083	2085	2087	rBV	50109	39065	1.21%	0.134%
32	14.692	2140	2143	2147	rVB	154450	147176	4.57%	0.503%
33	14.939	2183	2185	2190	rVB	70062	67467	2.10%	0.231%
34	15.251	2233	2238	2242	rBV	441078	499207	15.51%	1.707%

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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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

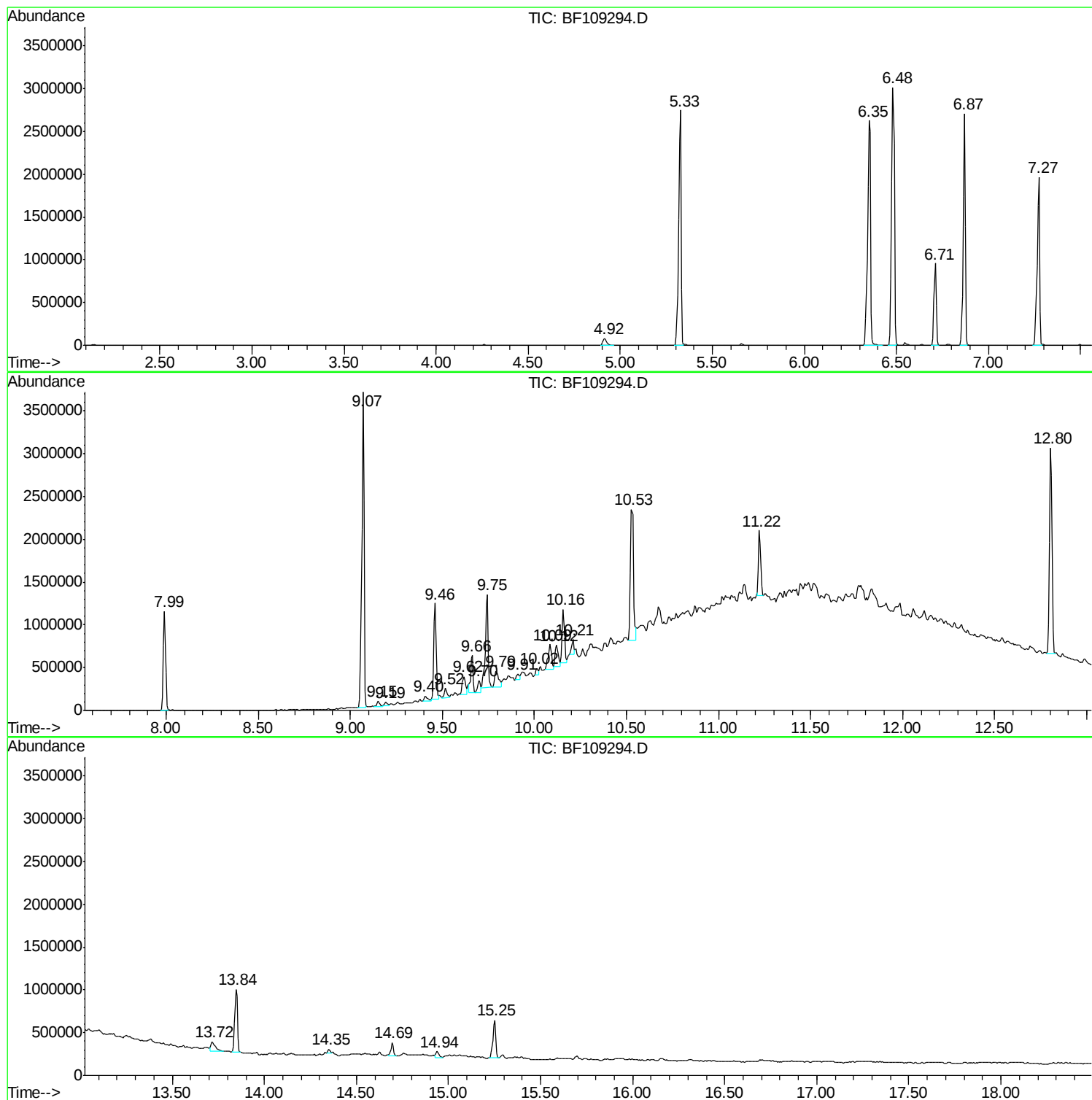
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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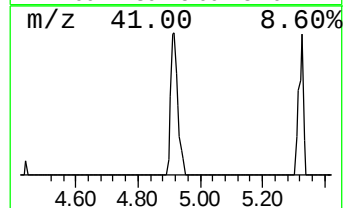
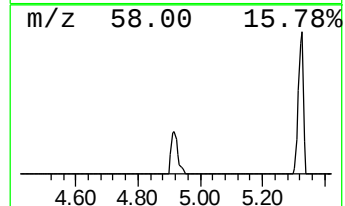
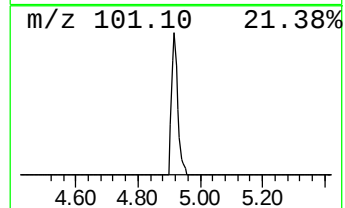
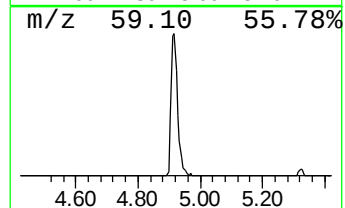
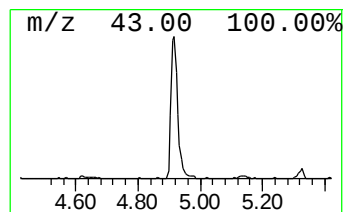
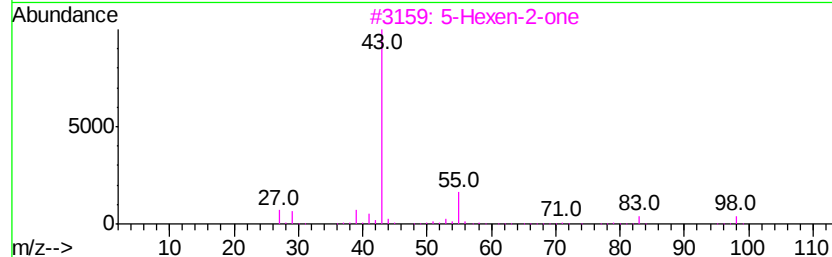
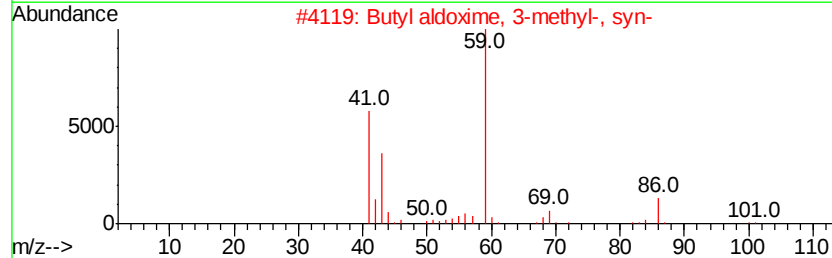
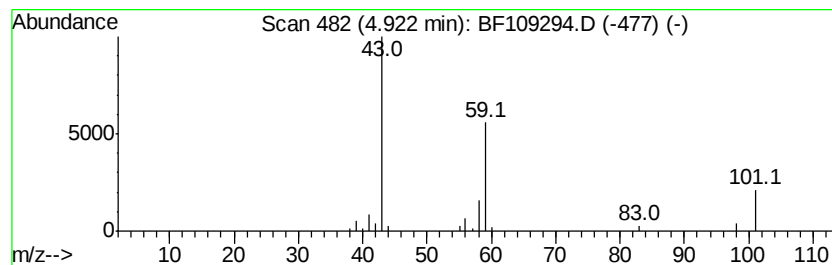
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.92	3.10 ng	114293	1,4-Dichlorobenzene-d4	6.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17
3			5-Hexen-2-one	98	C6H10O	000109-49-9	9
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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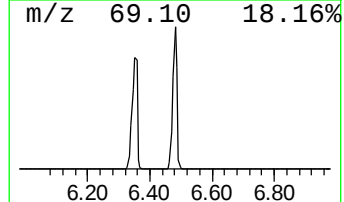
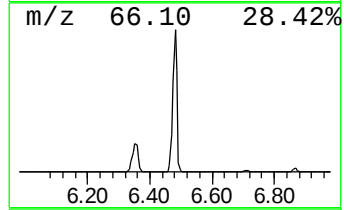
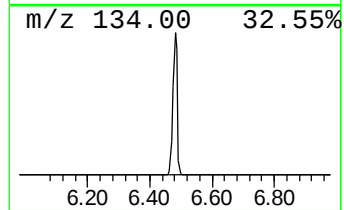
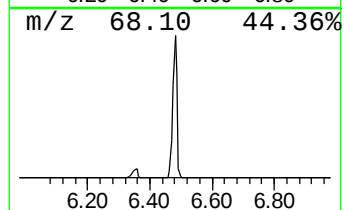
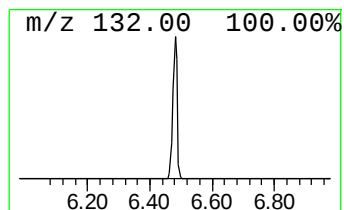
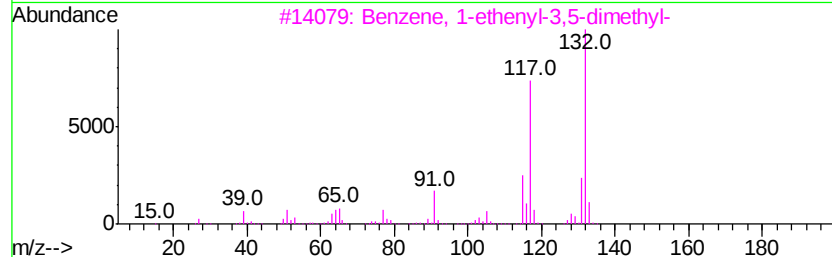
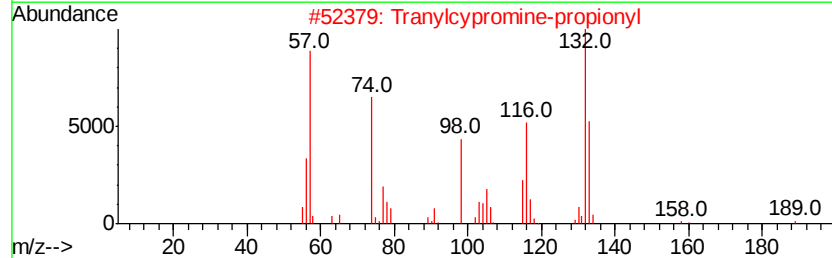
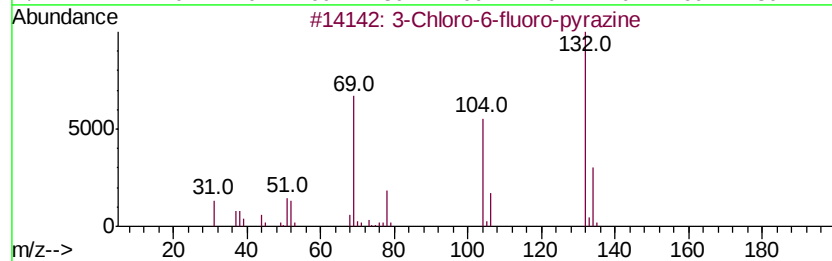
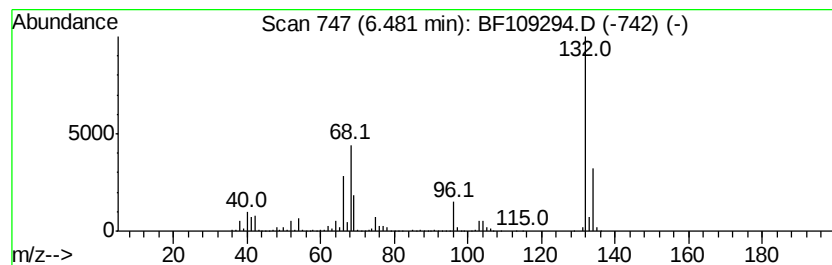
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.48 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.48	80.79 ng	2980140	1,4-Dichlorobenzene-d4	6.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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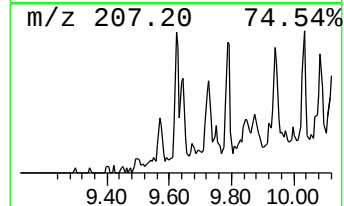
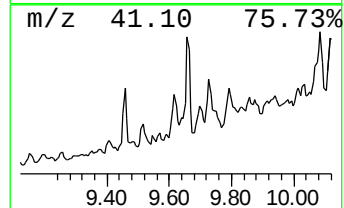
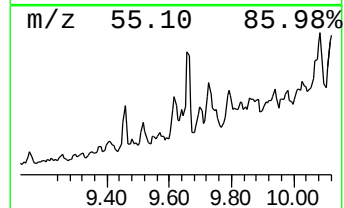
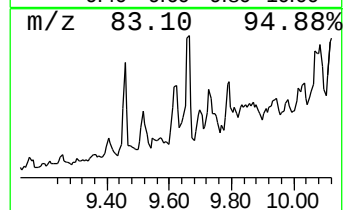
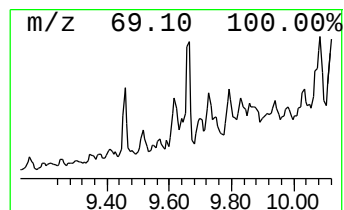
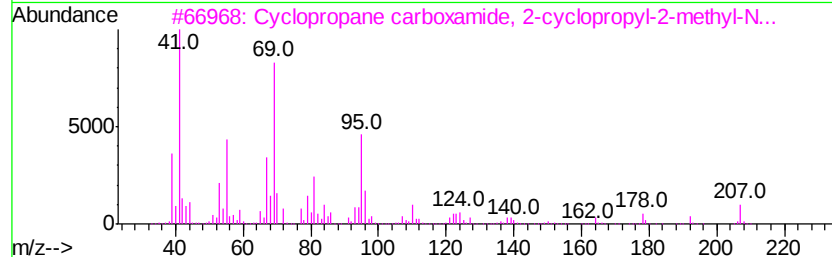
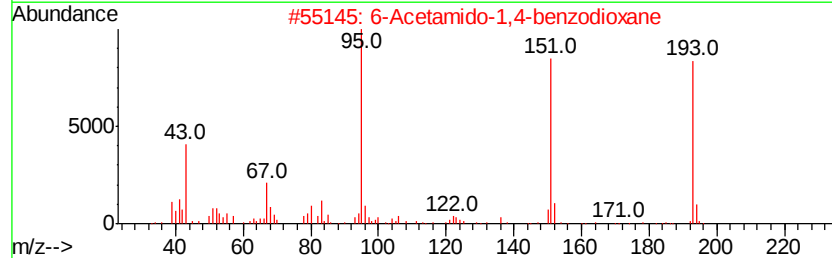
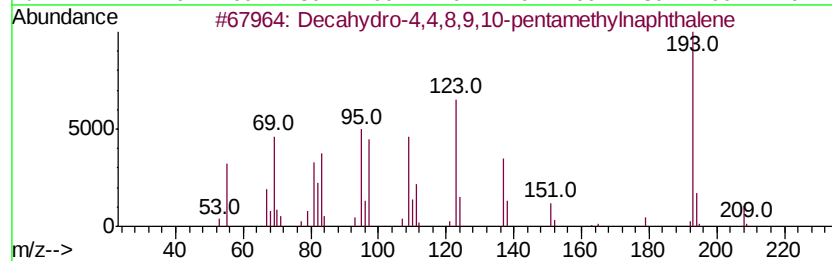
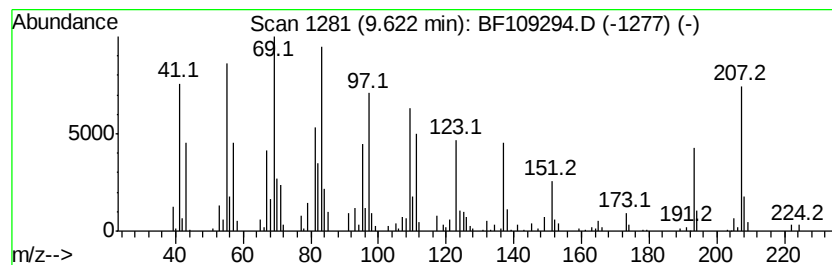
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Decahydro-4,4,8,9,10-pentam... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.62	4.16 ng	249406	Acenaphthene-d10	9.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	83
2		6-Acetamido-1,4-benzodioxane	193	C10H11NO3	063546-19-0	25
3		Cyclopropane carboxamide, 2-cycl...	207	C13H21NO	331416-19-4	20
4		2,6-Heptadienal, 2,4-dimethyl-	138	C9H14O	085136-08-9	18
5		Ethanone, 1-(1,4-dimethyl-3-cycl...	152	C10H16O	043219-68-7	15



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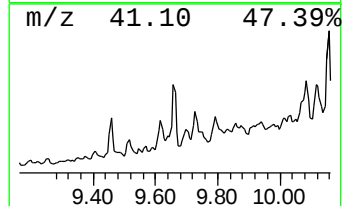
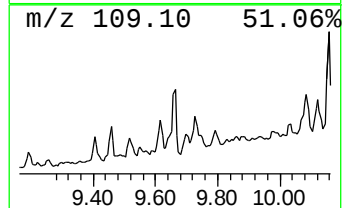
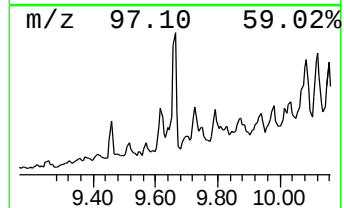
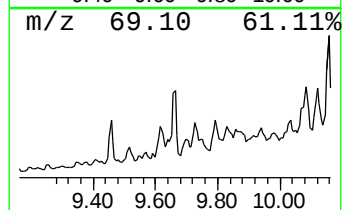
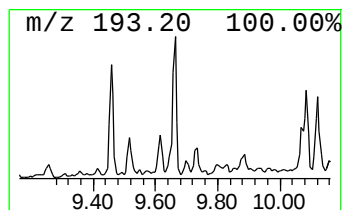
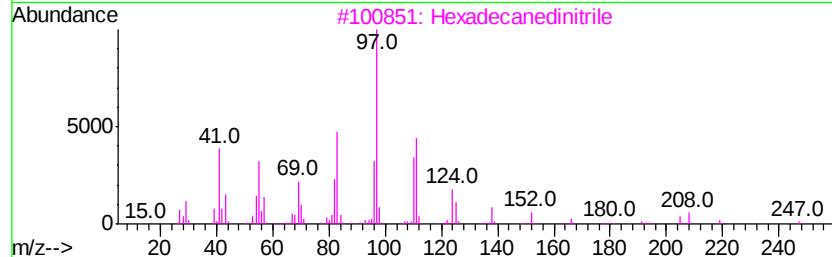
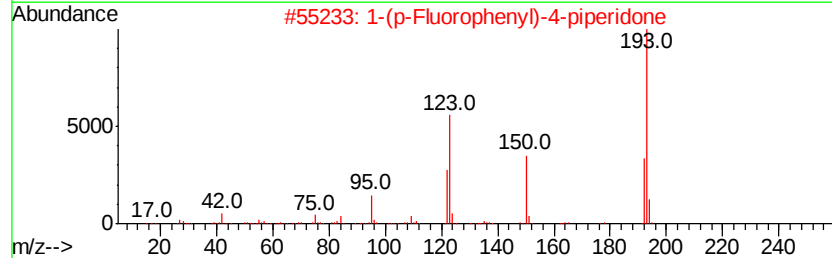
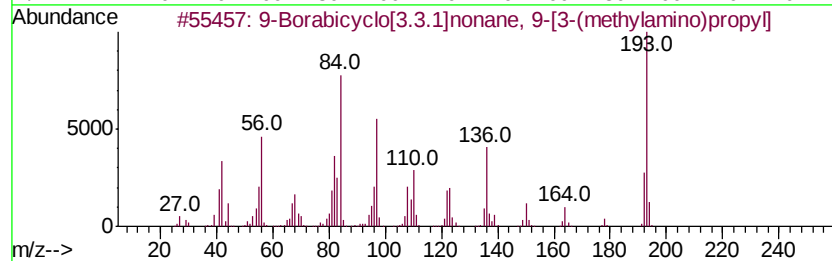
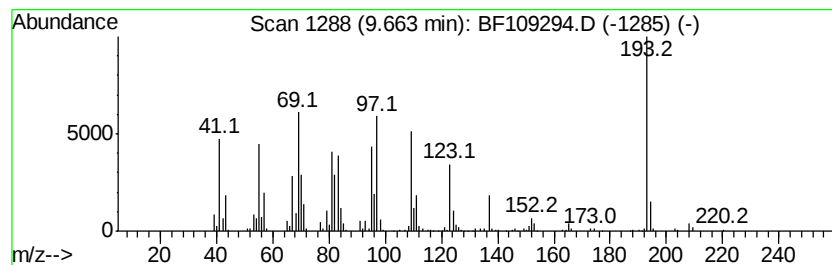
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 unknown9.66 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.		
9.66	6.39 ng	383067	Acenaphthene-d10		9.75		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Borabicyclo[3.3.1]nonane, 9-[3...	193	C12H24BN	1000162-56-0	43		
2	1-(p-Fluorophenyl)-4-piperidone	193	C11H12FN0	1000238-56-7	35		
3	Hexadecanedinitrile	248	C16H28N2	006812-44-8	22		
4	2-Anthracenamine	193	C14H11N	000613-13-8	14		
5	9-Vinylcarbazole	193	C14H11N	001484-13-5	14		



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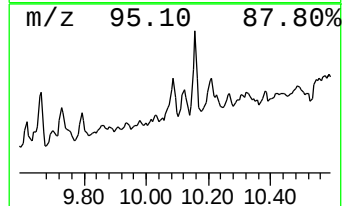
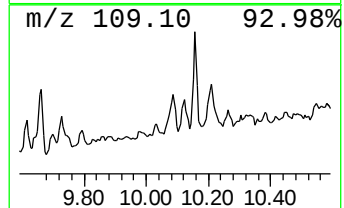
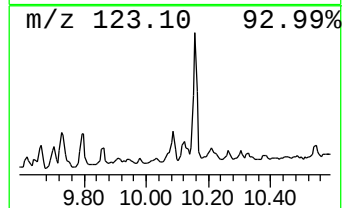
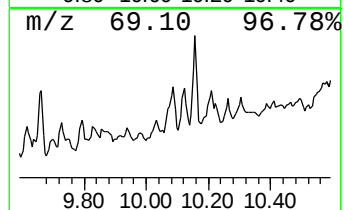
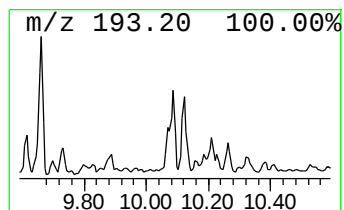
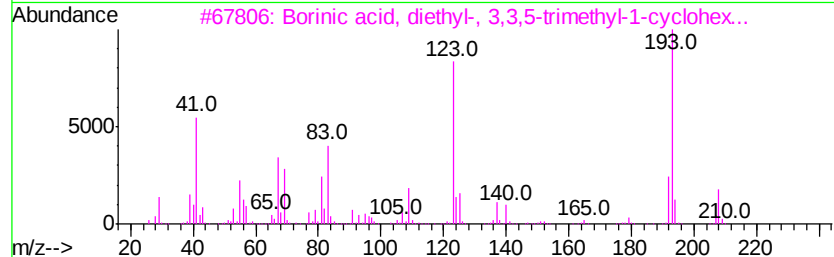
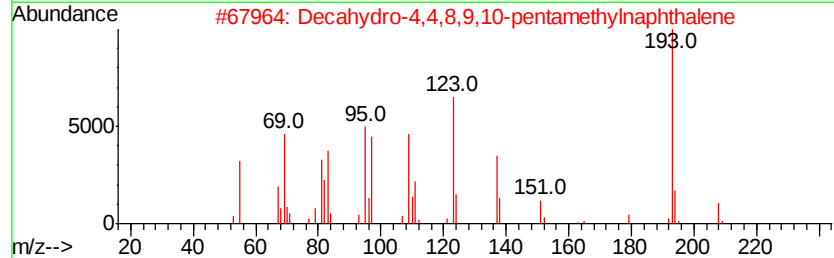
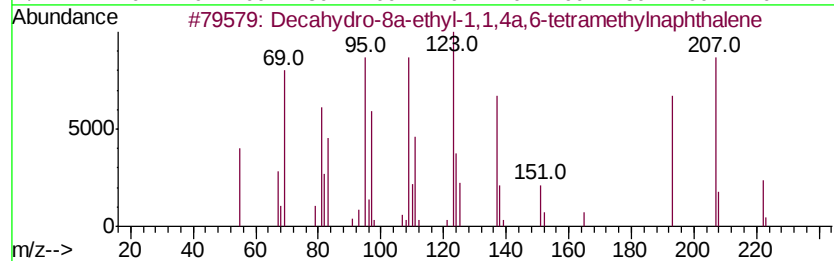
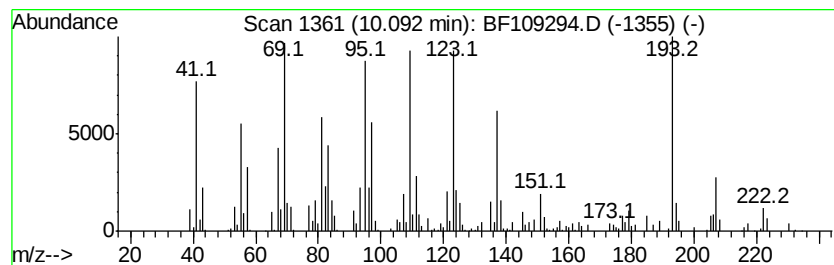
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Peak Number 6 Decahydro-8a-ethyl-1,1,4a,6... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.09	6.31 ng	377919	Acenaphthene-d10	9.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decahydro-8a-ethyl-1,1,4a,6-tetr...	222	C16H30	1000100-23-6	83
2		Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	64
3		Borinic acid, diethyl-, 3,3,5-tr...	208	C13H25BO	057387-76-5	27
4		Naphthalene, 2-butyldecahydro-	194	C14H26	006305-52-8	25
5		Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	020536-40-7	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092518\
Data File : BF109294.D
Acq On : 25 Sep 2018 19:03
Operator : JU/SJ
Sample : J5055-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-D

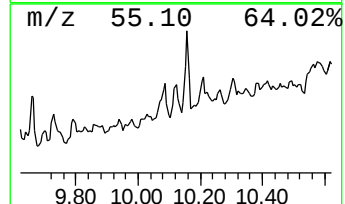
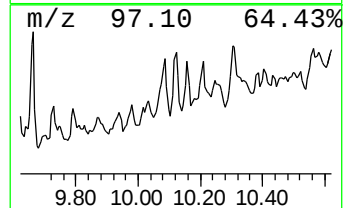
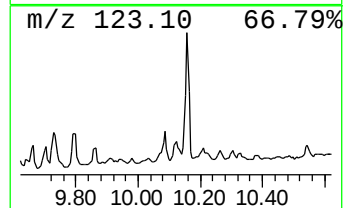
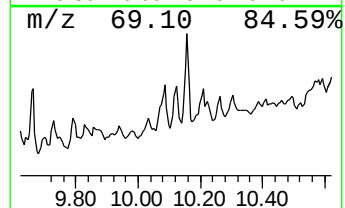
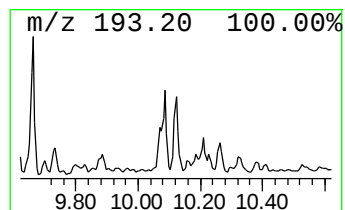
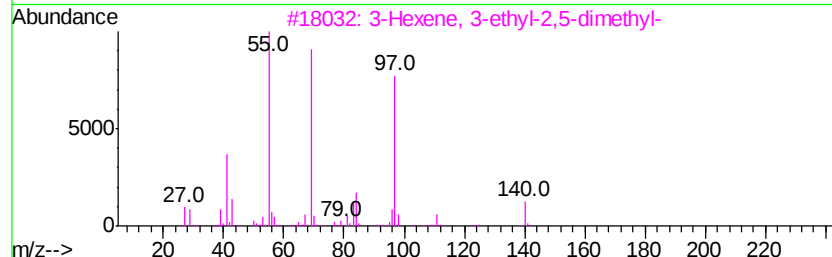
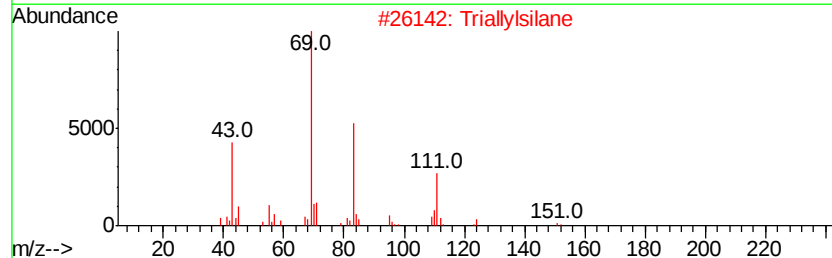
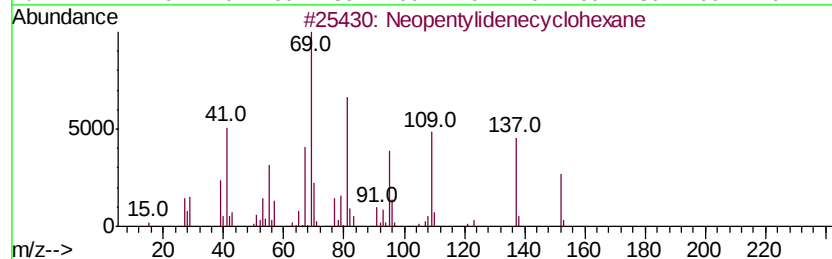
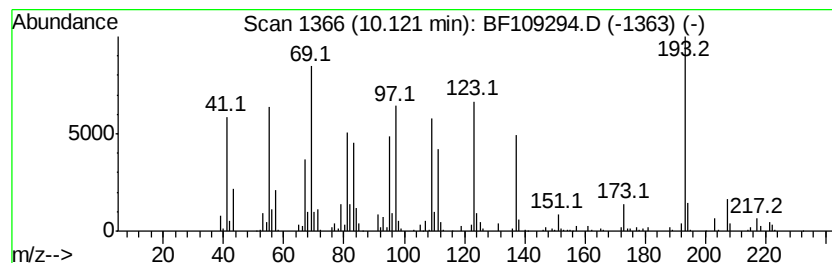
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown10.12 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.12	4.56 ng	273502	Acenaphthene-d10	9.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Neopentylidenecyclohexane	152	C11H20	039546-80-0	35
2		Triallylsilane	152	C9H16Si	001116-62-7	18
3		3-Hexene, 3-ethyl-2,5-dimethyl-	140	C10H20	062338-08-3	14
4		Cyclohexane, 1-ethyl-1,4-dimethyl...	140	C10H20	062238-30-6	14
5		2-Anthracenamine	193	C14H11N	000613-13-8	11



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092518\
Data File : BF109294.D
Acq On : 25 Sep 2018 19:03
Operator : JU/SJ
Sample : J5055-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

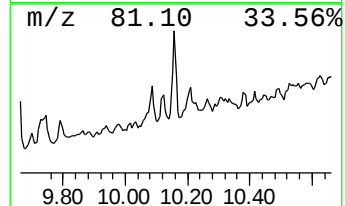
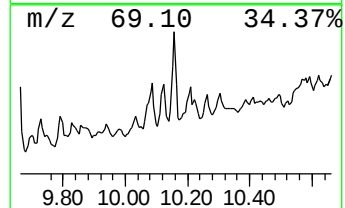
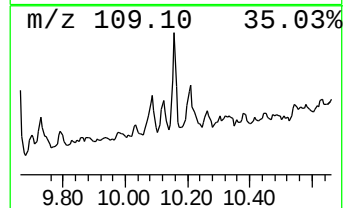
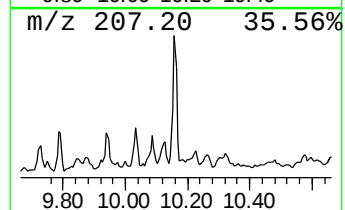
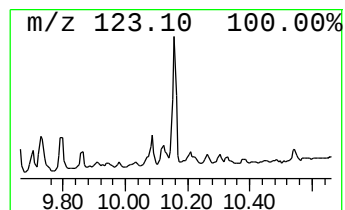
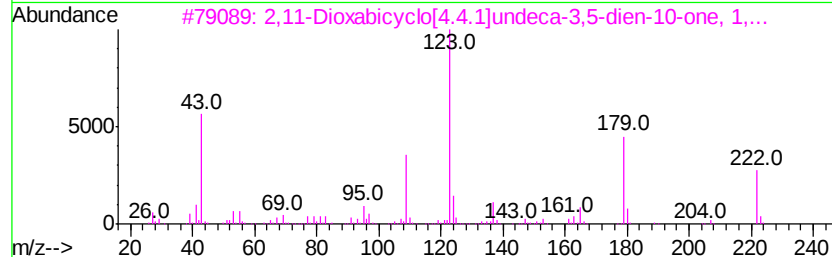
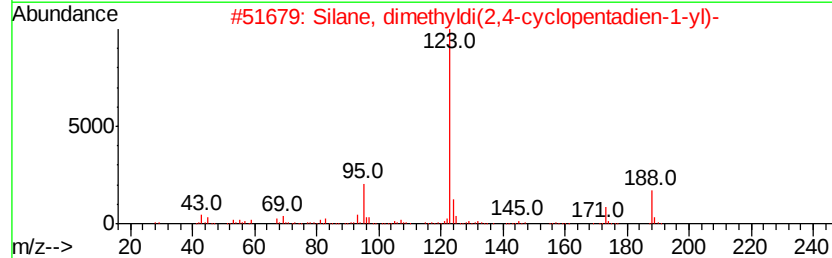
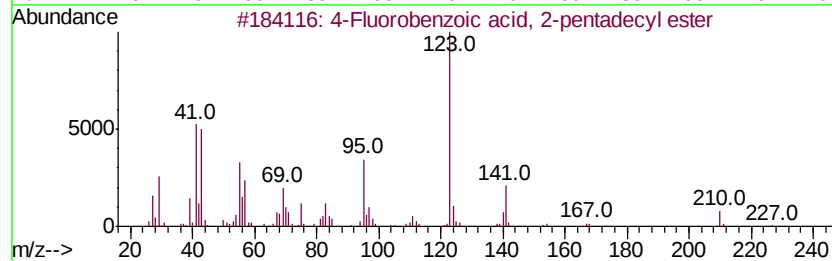
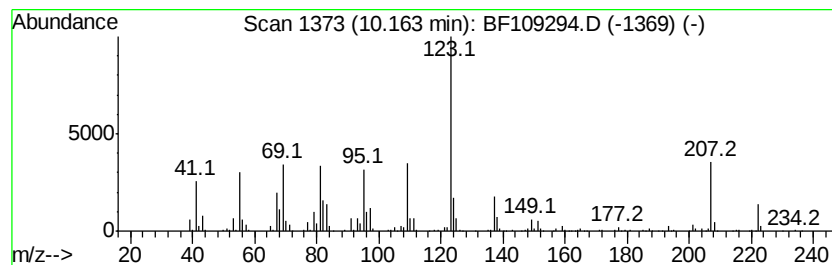
Instrument :
BNA_F
ClientSampleId :
TP-D

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 unknown10.16 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.16	8.24 ng	493555	Acenaphthene-d10	9.75	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Fluorobenzoic acid, 2-pentadec...	350	C22H35FO2	1000280-68-3	47
2	Silane, dimethyldi(2,4-cyclopent...	188	C12H16Si	018053-74-2	46
3	2,11-Dioxabicyclo[4.4.1]undeca-3...	222	C13H18O3	070412-52-1	45
4	3-Fluorobenzoic acid, 1-cyclopent...	236	C14H17FO2	1000279-04-7	43
5	4-Fluorobenzoic acid, 3-tetradec...	336	C21H33FO2	1000280-68-1	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092518\
Data File : BF109294.D
Acq On : 25 Sep 2018 19:03
Operator : JU/SJ
Sample : J5055-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

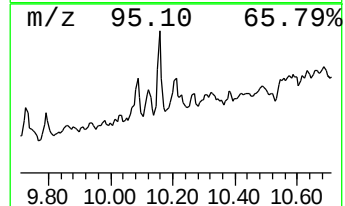
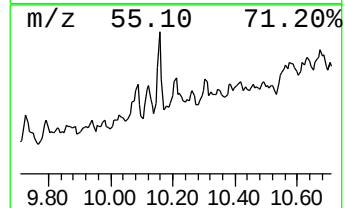
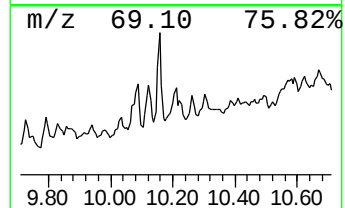
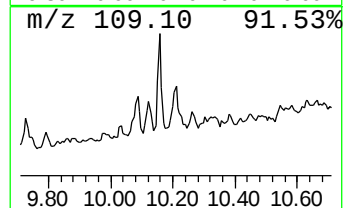
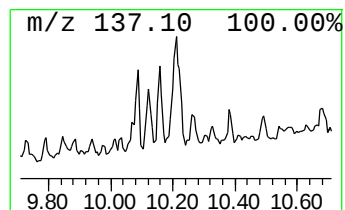
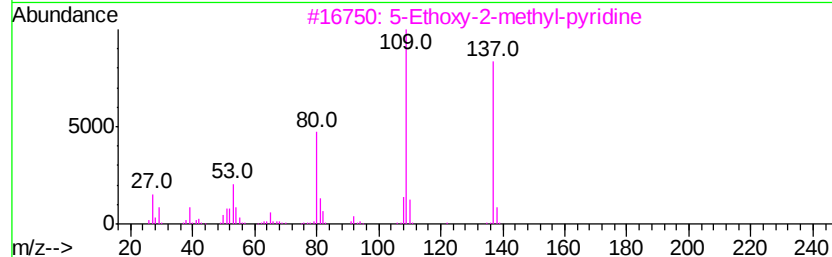
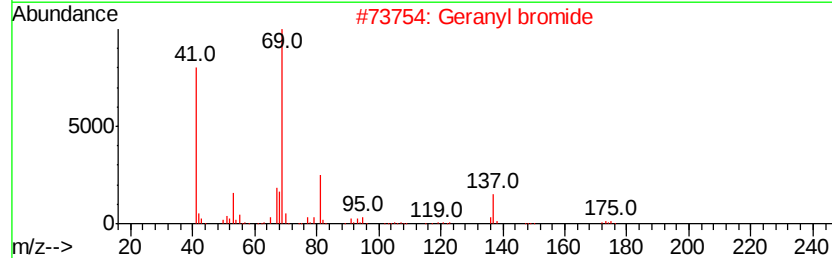
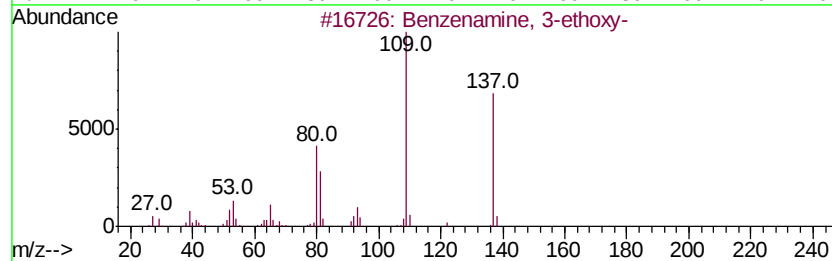
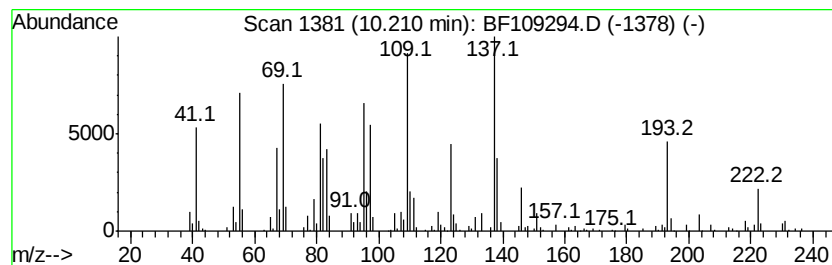
Instrument :
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ClientSampleId :
TP-D

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown10.21 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.21	2.30 ng	137725	Acenaphthene-d10		9.75
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-ethoxy-	137	C8H11NO	000621-33-0	22
2	Geranyl bromide	216	C10H17Br	006138-90-5	22
3	5-Ethoxy-2-methyl-pyridine	137	C8H11NO	055270-48-9	22
4	cis,trans-2-Ethylbicyclo[4.4.0]d...	166	C12H22	066660-38-6	14
5	Naphthalene, 2-butyldecahydro-	194	C14H26	006305-52-8	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092518\
Data File : BF109294.D
Acq On : 25 Sep 2018 19:03
Operator : JU/SJ
Sample : J5055-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

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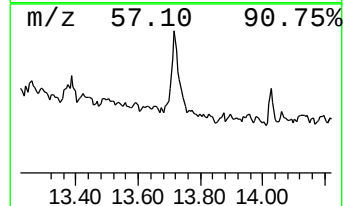
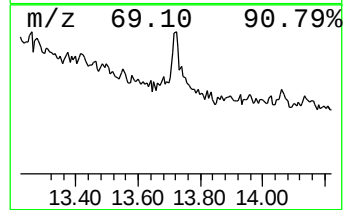
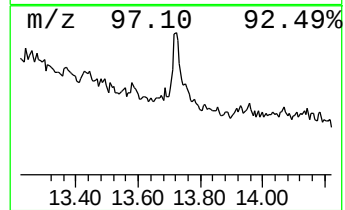
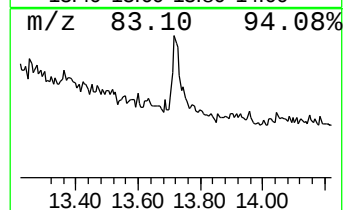
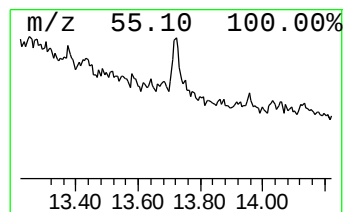
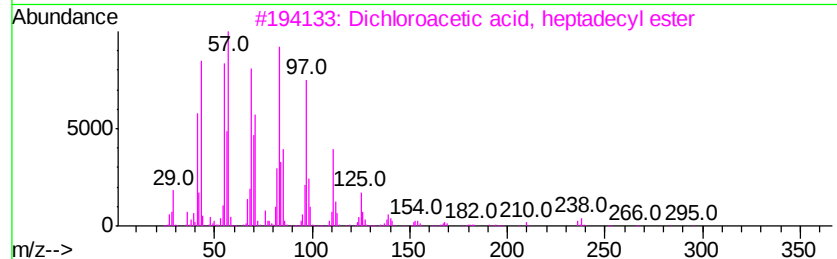
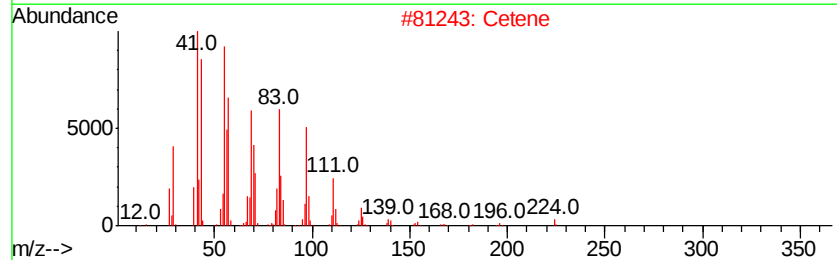
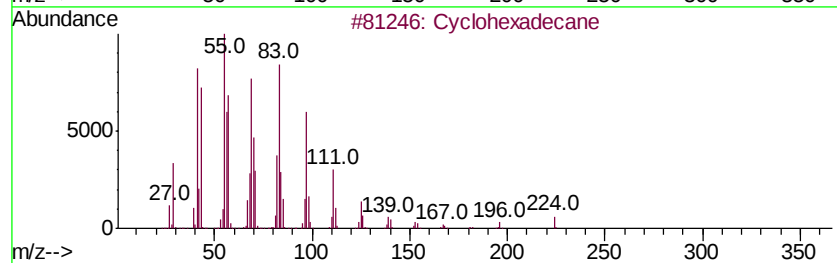
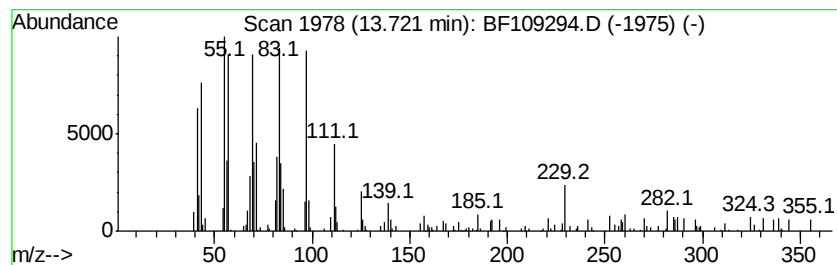
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Cyclohexadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	5.38 ng	186518	Chrysene-d12	13.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexadecane	224	C16H32	000295-65-8	97
2		Cetene	224	C16H32	000629-73-2	95
3		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	93
4		Carbonic acid, octadecyl 2,2,2-t...	444	C21H39Cl3O3	1000314-56-3	91
5		Bacchotricuneatin c	342	C20H22O5	066563-30-2	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092518\
Data File : BF109294.D
Acq On : 25 Sep 2018 19:03
Operator : JU/SJ
Sample : J5055-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-D

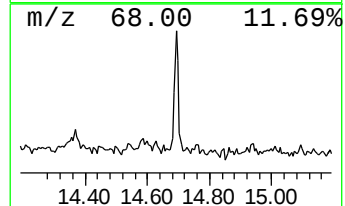
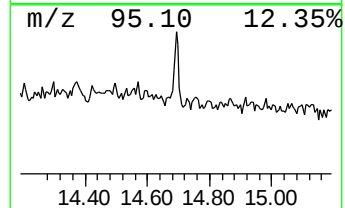
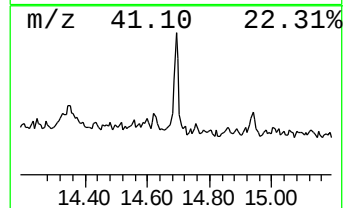
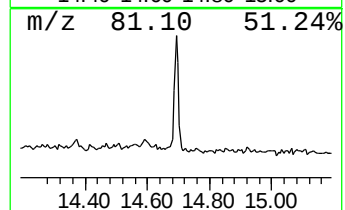
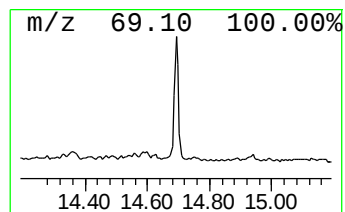
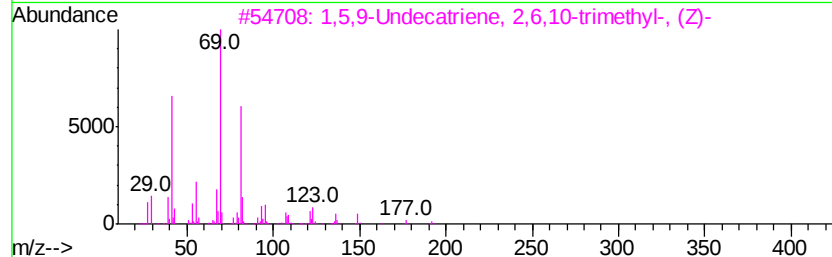
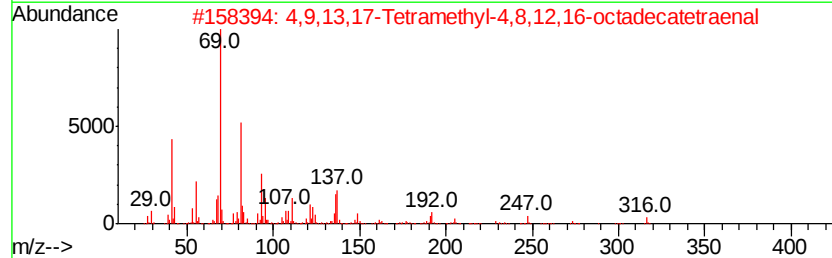
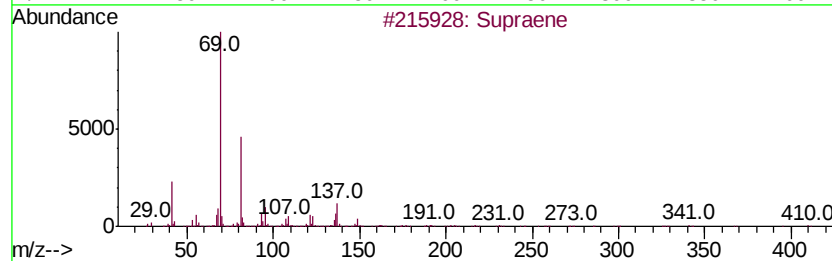
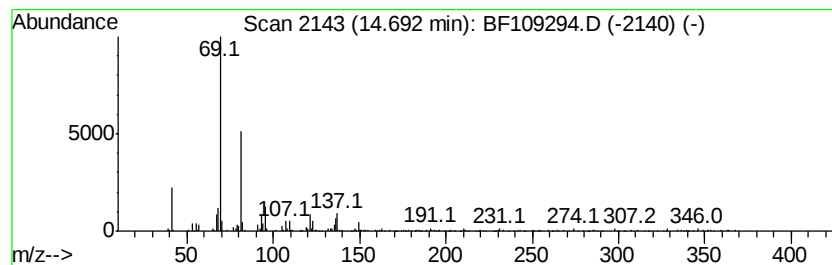
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Supraene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	5.90 ng	147176	Perylene-d12	15.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Supraene	410	C30H50	007683-64-9	87
2		4,9,13,17-Tetramethyl-4,8,12,16-...	316	C22H36O	056882-09-8	72
3		1,5,9-Undecatriene, 2,6,10-trime...	192	C14H24	062951-96-6	64
4		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	64
5		4,8,12-Tetradecatrienenitrile, 5...	245	C17H27N	005981-31-7	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092518\
Data File : BF109294.D
Acq On : 25 Sep 2018 19:03
Operator : JU/SJ
Sample : J5055-16
Misc :
ALS Vial : 14 Sample Multiplier: 1

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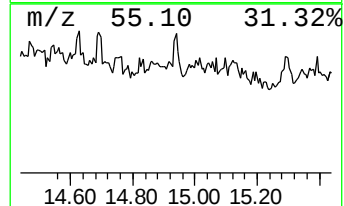
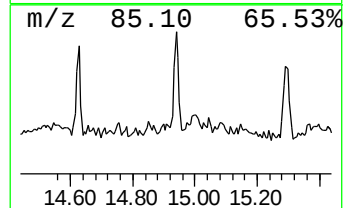
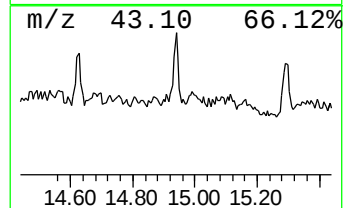
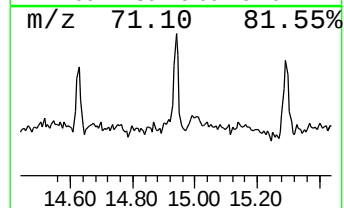
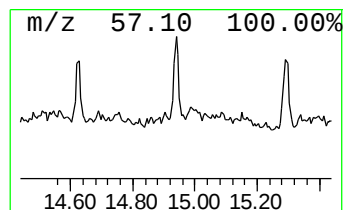
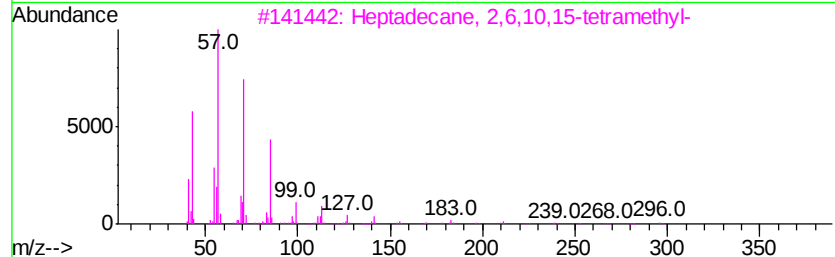
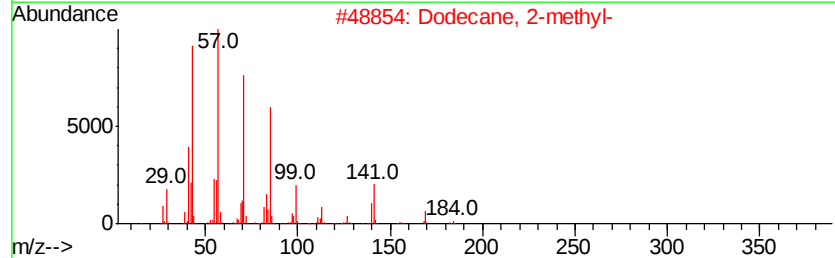
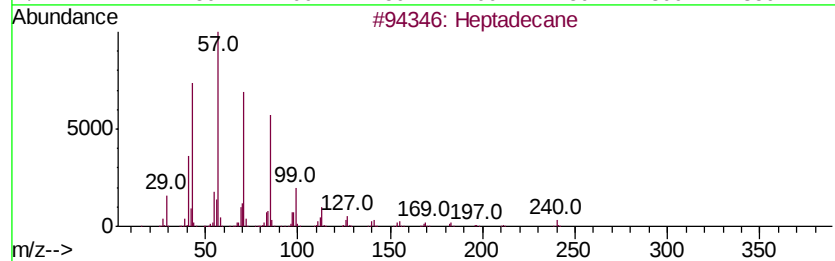
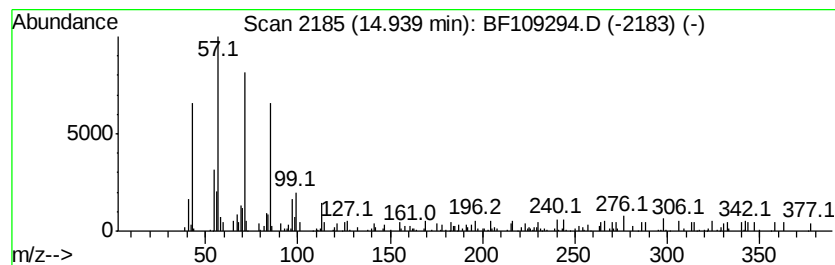
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Heptadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.94	2.70 ng	67467	Perylene-d12	15.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	90
2		Dodecane, 2-methyl-	184	C13H28	001560-97-0	74
3		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	74
4		Pentadecane, 2-methyl-	226	C16H34	001560-93-6	74
5		Tridecane	184	C13H28	000629-50-5	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF092518\
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Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-D

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.92	3.1 ng		114293	1	6.71	737758	20.0
unknown6.48	6.48	80.8 ng		2980140	1	6.71	737758	20.0
Decahydro-4,4,8,9...	9.62	4.2 ng		249406	3	9.75	1198540	20.0
unknown9.66	9.66	6.4 ng		383067	3	9.75	1198540	20.0
Decahydro-8a-ethy...	10.09	6.3 ng		377919	3	9.75	1198540	20.0
unknown10.12	10.12	4.6 ng		273502	3	9.75	1198540	20.0
unknown10.16	10.16	8.2 ng		493555	3	9.75	1198540	20.0
unknown10.21	10.21	2.3 ng		137725	3	9.75	1198540	20.0
Cyclohexadecane	13.72	5.4 ng		186518	5	13.84	692808	20.0
Supraene	14.69	5.9 ng		147176	6	15.25	499207	20.0
Heptadecane	14.94	2.7 ng		67467	6	15.25	499207	20.0