

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042718\
 Data File : BF104886.D
 Acq On : 27 Apr 2018 17:30
 Operator : JU/SJ
 Sample : J2579-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

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-BF040618.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	5.016	493	498	507	rBV	89884	125628	4.44%	0.456%
2	5.416	561	566	580	rBV	2443577	2708763	95.80%	9.842%
3	6.063	672	676	689	rBV	352966	317047	11.21%	1.152%
4	6.440	735	740	743	rBV	2325377	2732079	96.62%	9.927%
5	6.569	757	762	765	rBV	2880682	2827619	100.00%	10.274%
6	6.798	797	801	804	rBV	667755	610426	21.59%	2.218%
7	6.951	823	827	830	rBV	2403437	2136029	75.54%	7.761%
8	7.363	892	897	900	rBV	1595364	1744856	61.71%	6.340%
9	7.457	909	913	923	rVB3	21019	38554	1.36%	0.140%
10	7.534	923	926	930	rVB	33098	30647	1.08%	0.111%
11	8.081	1015	1019	1035	rBV	905310	853864	30.20%	3.103%
12	9.157	1197	1202	1205	rBV	3071370	2732296	96.63%	9.928%
13	9.551	1264	1269	1273	rBV	259001	236485	8.36%	0.859%
14	9.757	1301	1304	1308	rVB	61600	50203	1.78%	0.182%
15	9.834	1314	1317	1327	rVB	968812	947108	33.49%	3.441%
16	10.257	1387	1389	1392	rVB	82722	59817	2.12%	0.217%
17	10.633	1448	1453	1462	rBV	1849738	1894583	67.00%	6.884%
18	10.733	1467	1470	1478	rVB	70170	84303	2.98%	0.306%
19	11.180	1543	1546	1549	rBV	51351	43062	1.52%	0.156%
20	11.328	1567	1571	1577	rBV	1018435	923364	32.66%	3.355%
21	11.610	1615	1619	1622	rBV	33047	30446	1.08%	0.111%
22	12.016	1685	1688	1696	rVB	36220	34684	1.23%	0.126%
23	12.169	1711	1714	1717	rBV2	35781	29282	1.04%	0.106%
24	12.416	1749	1756	1759	rBV3	29461	42186	1.49%	0.153%
25	12.916	1834	1841	1844	rBV	2373126	2200523	77.82%	7.996%
26	13.216	1887	1892	1896	rBV6	20870	31820	1.13%	0.116%
27	13.339	1910	1913	1918	rVB3	68951	83312	2.95%	0.303%
28	13.410	1922	1925	1928	rVB	46635	38202	1.35%	0.139%
29	13.763	1980	1985	1988	rBV2	23225	33937	1.20%	0.123%
30	13.798	1988	1991	1995	rBV	293249	316398	11.19%	1.150%
31	13.839	1995	1998	2001	rVB	164217	153470	5.43%	0.558%
32	13.974	2017	2021	2029	rVB	753637	725794	25.67%	2.637%
33	14.439	2095	2100	2107	rBV2	73662	129106	4.57%	0.469%
34	14.810	2160	2163	2166	rVB	36900	35854	1.27%	0.130%

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

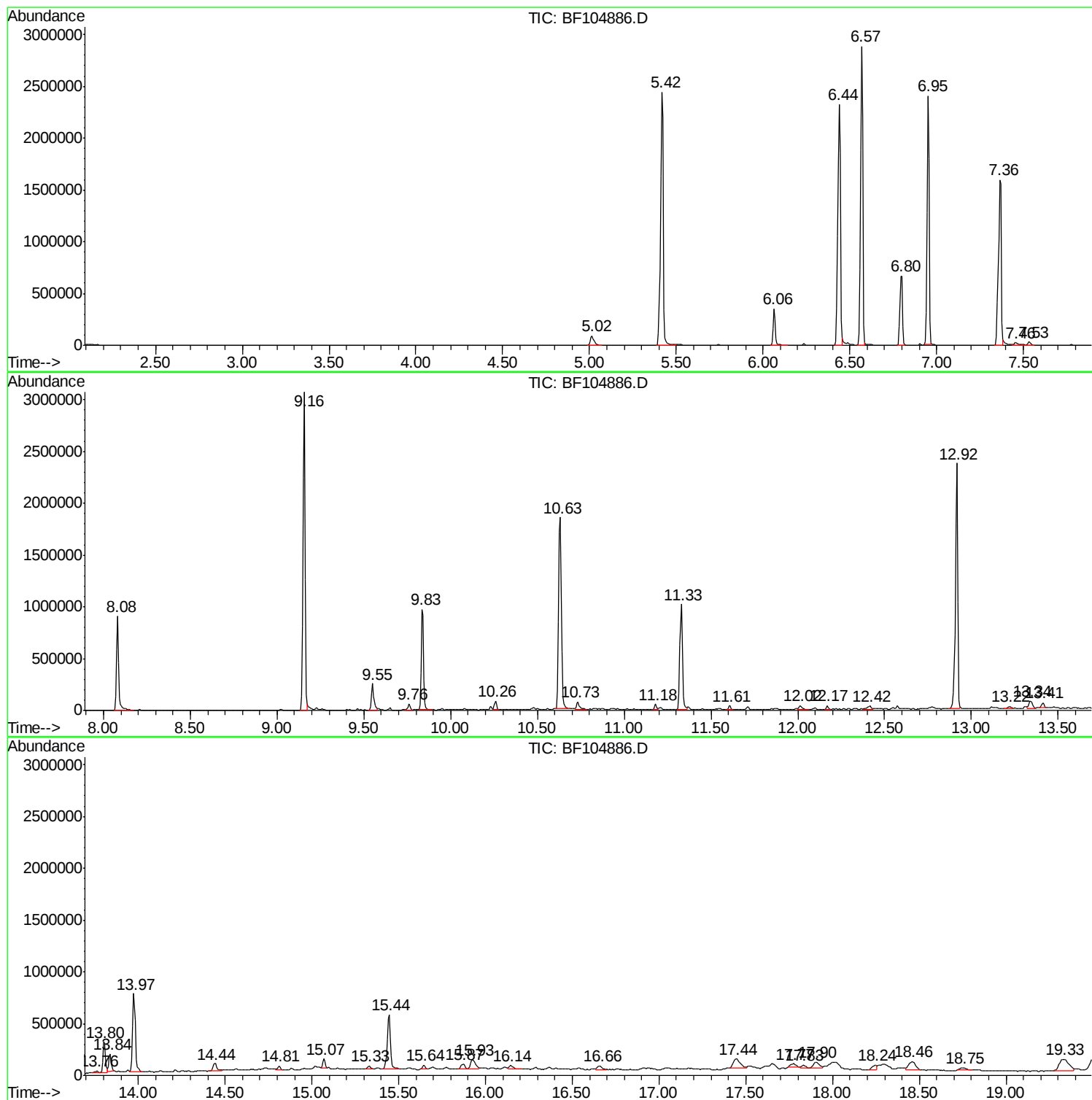
35	15.068	2204	2207	2210	rVB	91636	92213	3.26%	0.335%
36	15.327	2248	2251	2255	rVB3	28067	35604	1.26%	0.129%
37	15.445	2265	2271	2280	rBV	525328	699883	24.75%	2.543%
38	15.645	2302	2305	2308	rBV2	36063	36504	1.29%	0.133%
39	15.868	2339	2343	2348	rBV3	48565	68708	2.43%	0.250%
40	15.927	2348	2353	2359	rBV2	89722	162625	5.75%	0.591%
41	16.145	2387	2390	2400	rVB	33402	64813	2.29%	0.236%
42	16.662	2473	2478	2483	rVB5	35244	66812	2.36%	0.243%
43	17.445	2605	2611	2621	rBV9	89446	240323	8.50%	0.873%
44	17.774	2662	2667	2673	rBV9	37206	75980	2.69%	0.276%
45	17.827	2674	2676	2683	rVB7	32088	53807	1.90%	0.196%
46	17.904	2683	2689	2696	rBV5	61750	163792	5.79%	0.595%
47	18.245	2741	2747	2749	rBV6	40886	81488	2.88%	0.296%
48	18.456	2777	2783	2791	rVB5	81069	207077	7.32%	0.752%
49	18.751	2828	2833	2842	rVB5	23555	57191	2.02%	0.208%
50	19.327	2922	2931	2942	rBV6	112631	432624	15.30%	1.572%

Sum of corrected areas: 27521191

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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.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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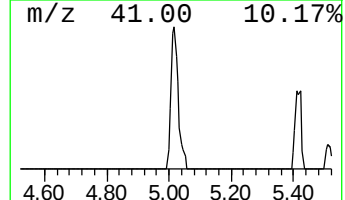
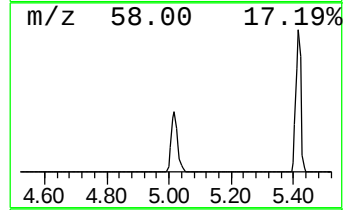
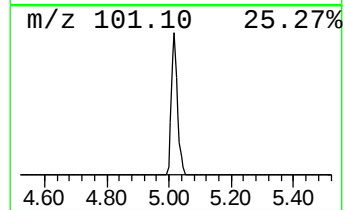
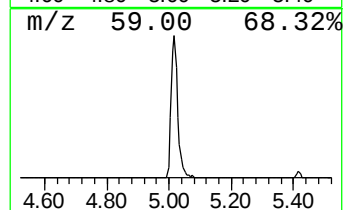
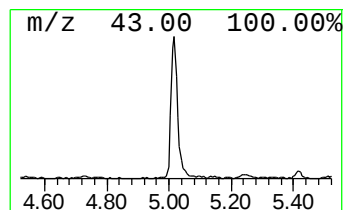
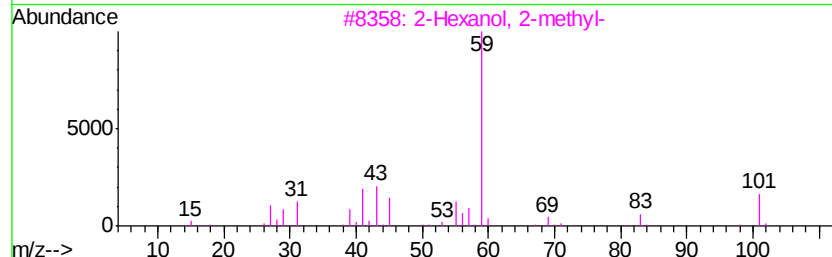
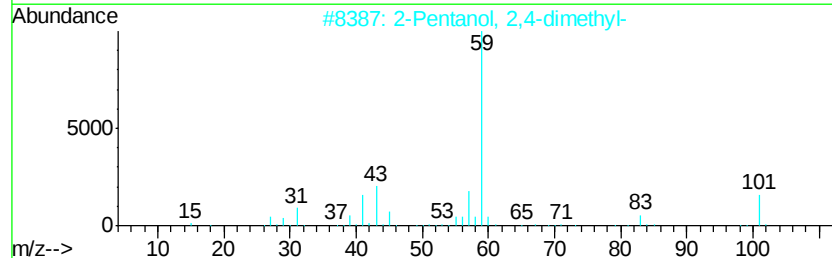
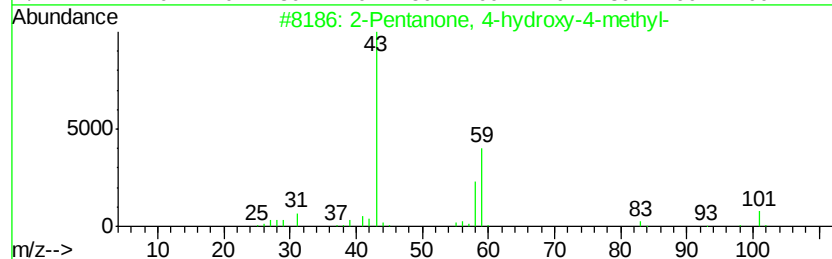
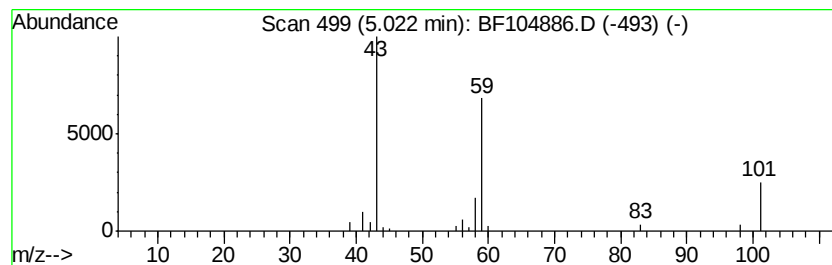
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.02	4.12 ng	125628	1,4-Dichlorobenzene-d4	6.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	42
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	38
4			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
5			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	25



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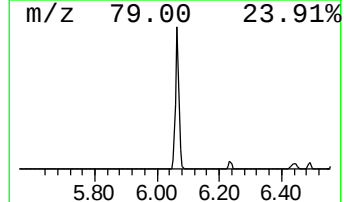
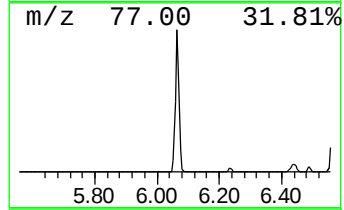
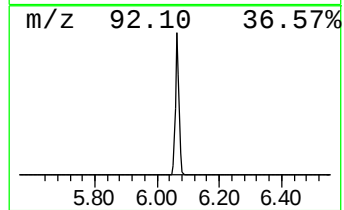
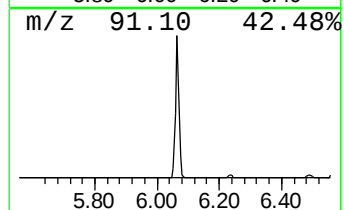
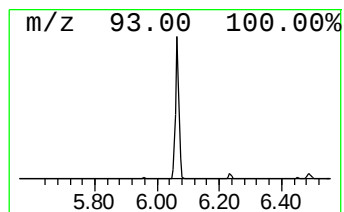
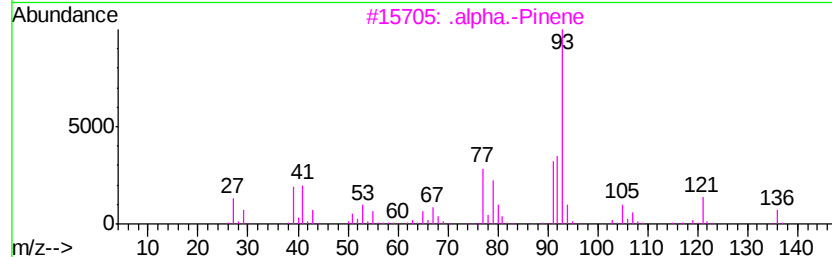
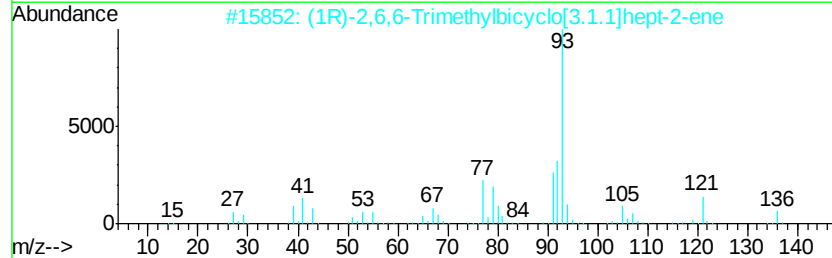
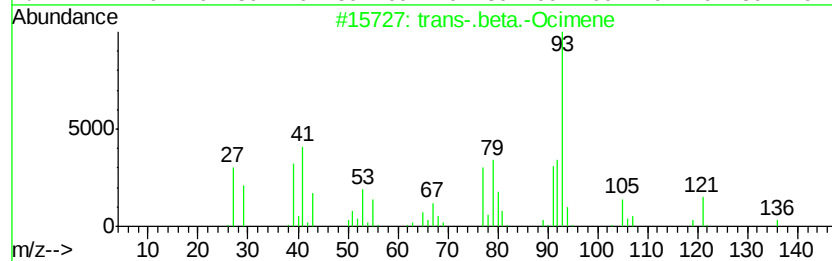
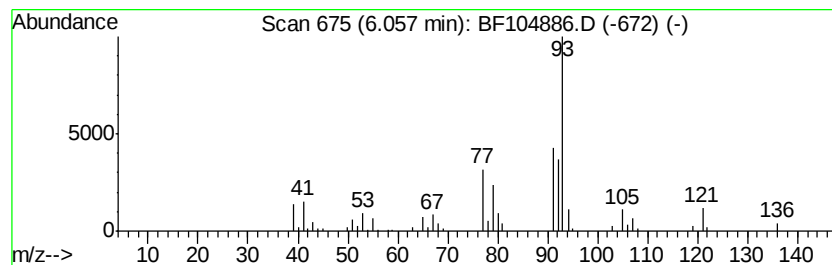
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Peak Number 2 trans-.beta.-Ocimene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.06	10.39 ng	317047	1,4-Dichlorobenzene-d4	6.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-.beta.-Ocimene	136	C10H16	003779-61-1	94
2			(1R)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-70-8	94
3			.alpha.-Pinene	136	C10H16	000080-56-8	94
4			Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	91
5			Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000508-32-7	91



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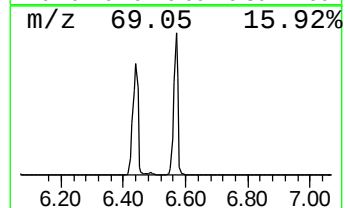
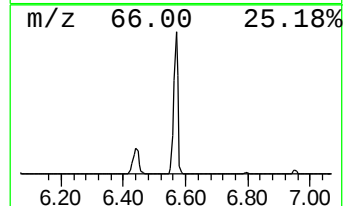
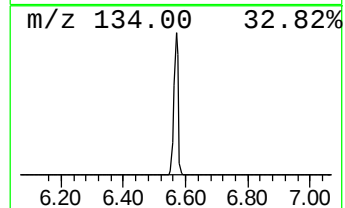
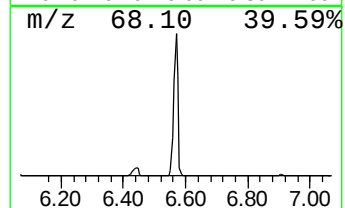
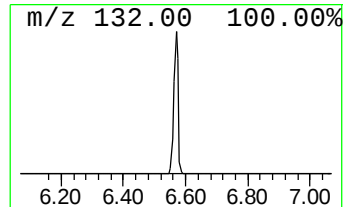
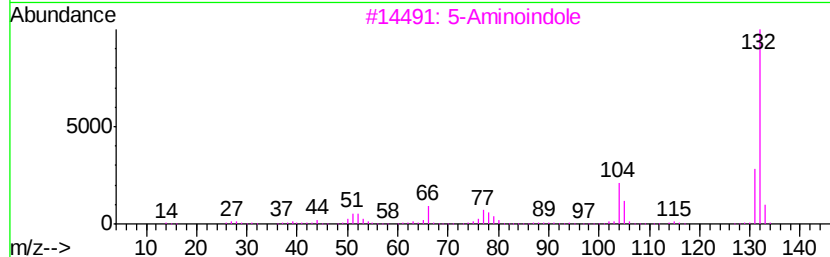
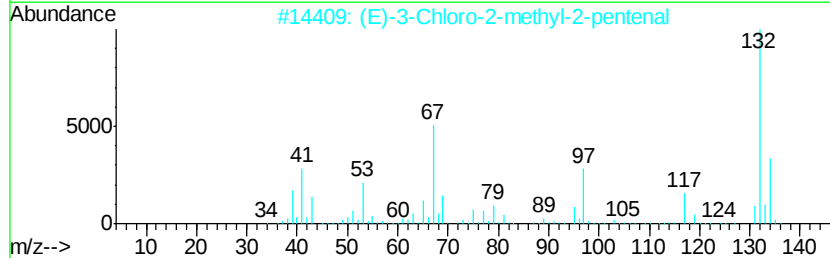
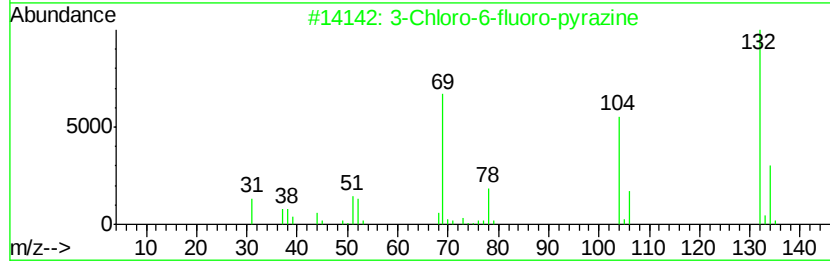
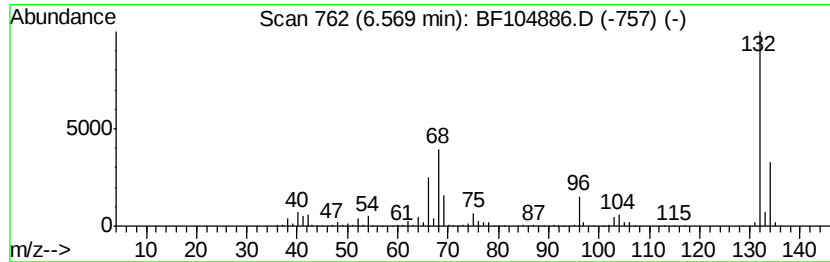
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 unknown6.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	92.64 ng	2827620	1,4-Dichlorobenzene-d4	6.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	27
2	(E)-3-Chloro-2-methyl-2-pentenal			132	C6H9ClO	031357-76-3	16
3	5-Aminoindole			132	C8H8N2	005192-03-0	12
4	5-Fluoro-2-chloropyrimidine			132	C4H2ClFN2	062802-42-0	9
5	1H-Benzimidazole, 2-methyl-			132	C8H8N2	000615-15-6	9



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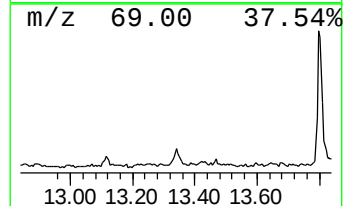
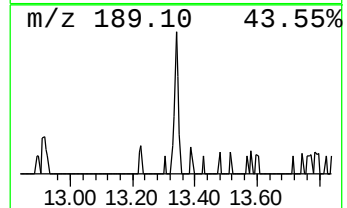
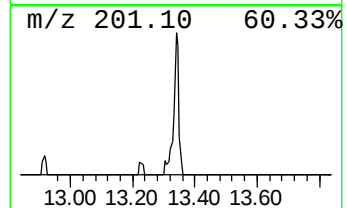
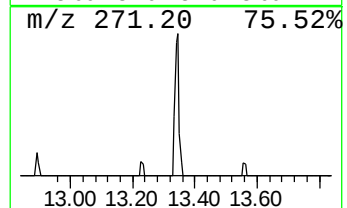
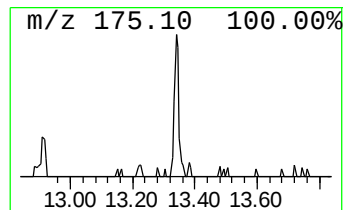
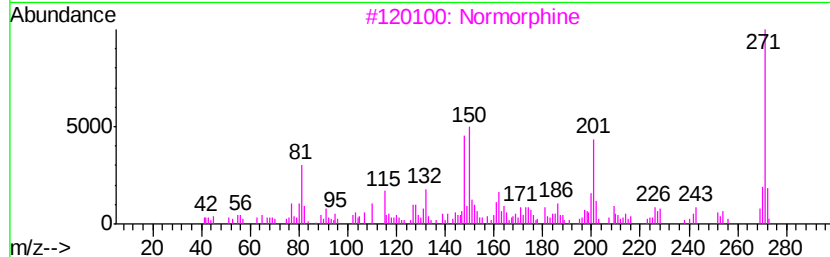
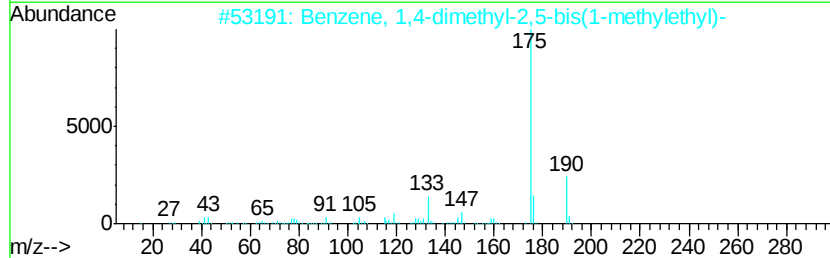
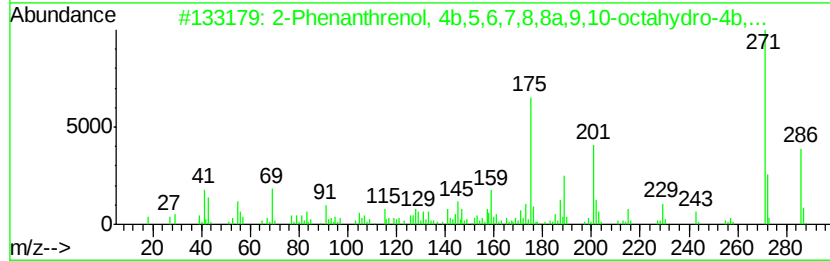
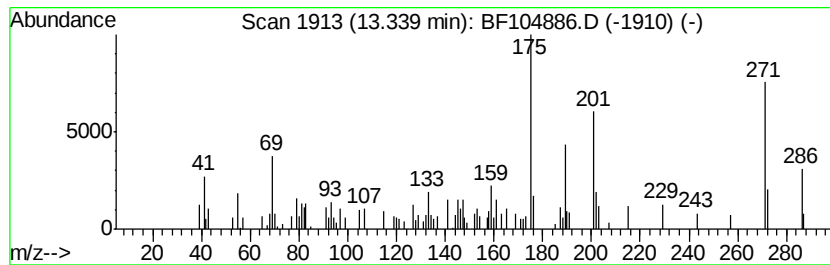
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Peak Number 5 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	2.30 ng	83312	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	60
2		Benzene, 1,4-dimethyl-2,5-bis(1-...	190	C14H22	010375-96-9	35
3		Normorphine	271	C16H17N03	000466-97-7	35
4		2,4-Diamino-6-methylpvrido[3,2-d...	175	C8H9N5	001955-57-3	14
5		6-Amino-5,8-dimethoxy-4-methyl-2...	286	C13H13F3N2O2	085868-17-3	14



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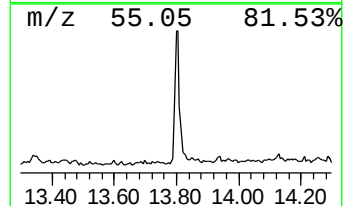
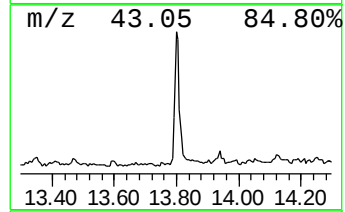
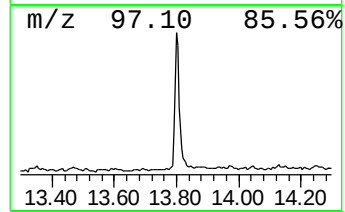
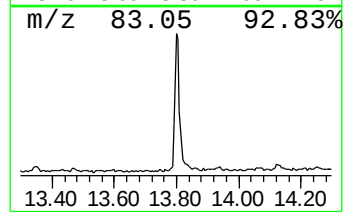
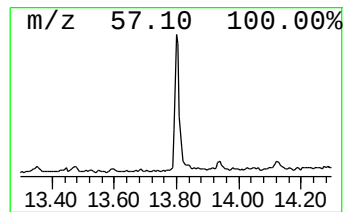
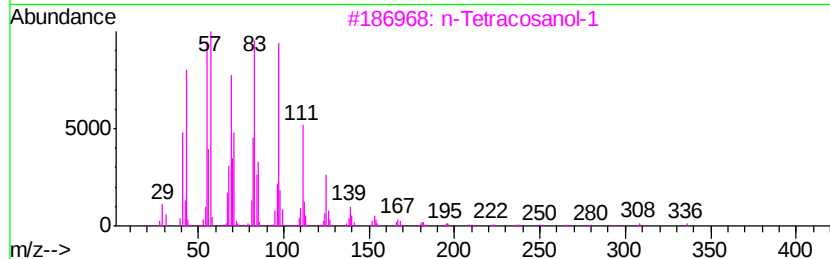
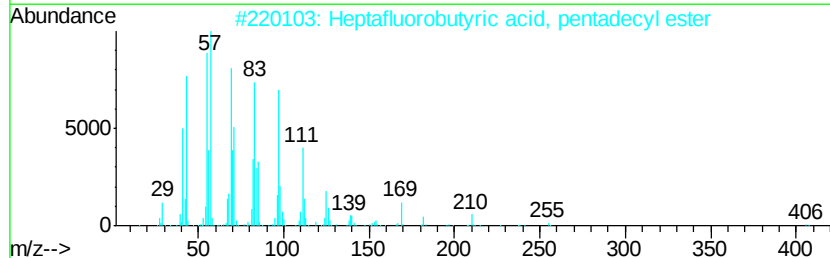
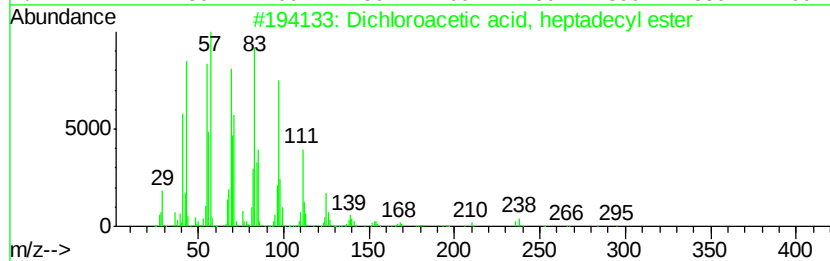
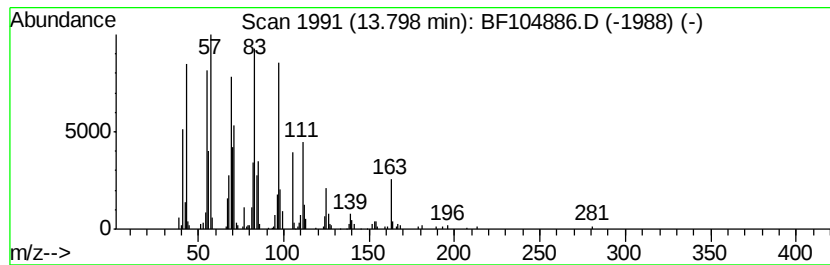
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Dichloroacetic acid, heptad... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.80	8.72 ng	316398	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	95
2		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	95
3		n-Tetracosanol-1	354	C24H50O	000506-51-4	95
4		Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	95
5		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042718\
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Acq On : 27 Apr 2018 17:30
Operator : JU/SJ
Sample : J2579-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

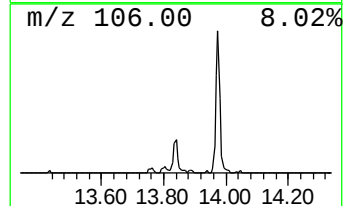
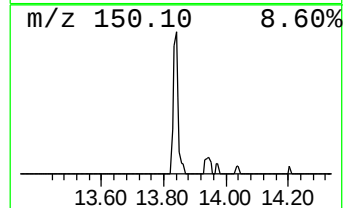
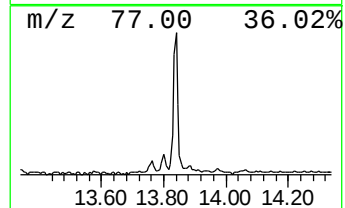
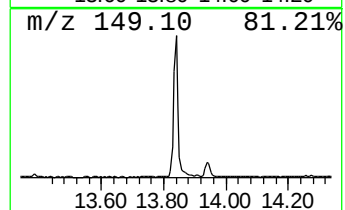
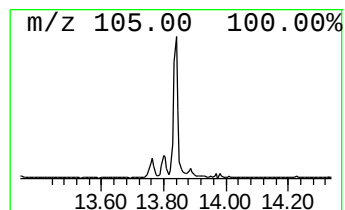
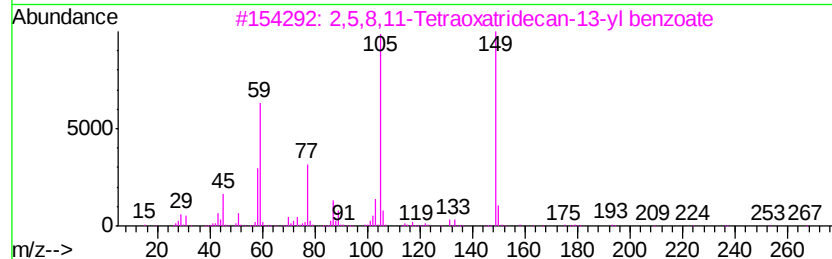
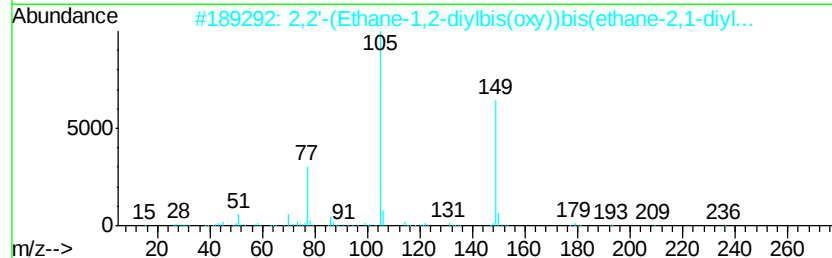
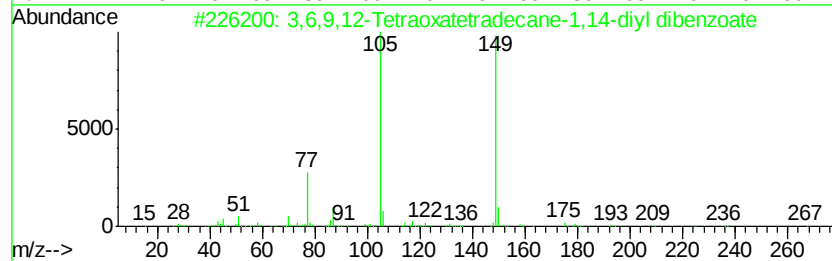
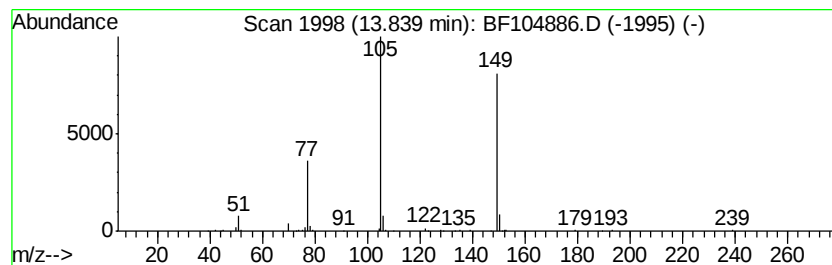
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 3,6,9,12-Tetraoxatetradecan... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.84	4.23 ng	153470	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3,6,9,12-Tetraoxatetradecane-1,1...	446	C24H30O8	1000366-97-0	83
2		2,2'-(Ethane-1,2-diylbis(oxy))bi...	358	C20H22O6	1000366-99-4	83
3		2,5,8,11-Tetraoxatridecan-13-yl ...	312	C16H24O6	1000366-96-7	78
4		2-(2-(2-Methoxyethoxy)ethoxy)eth...	268	C14H20O5	1000366-99-1	74
5		Benzoic acid, 2-(3-nitrophenyl)e...	271	C15H13NO4	1000368-22-4	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042718\
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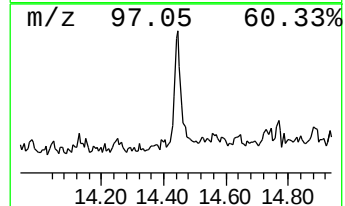
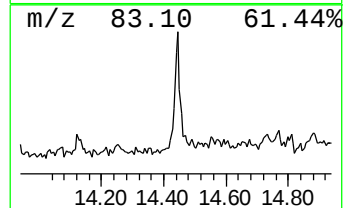
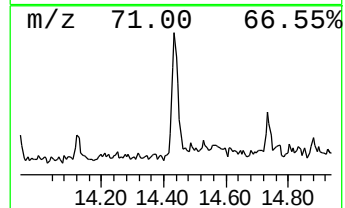
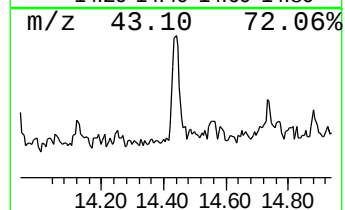
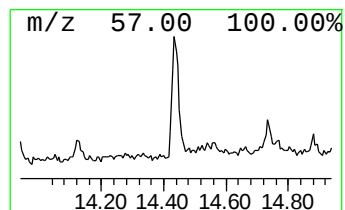
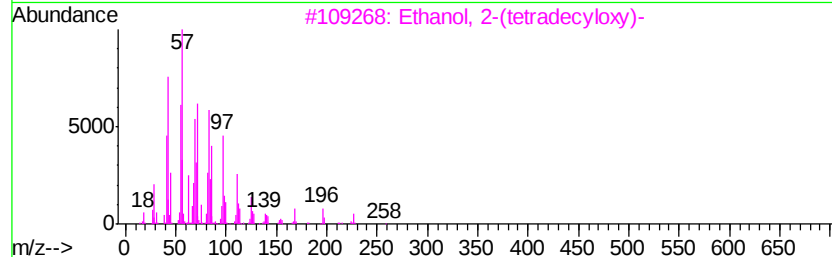
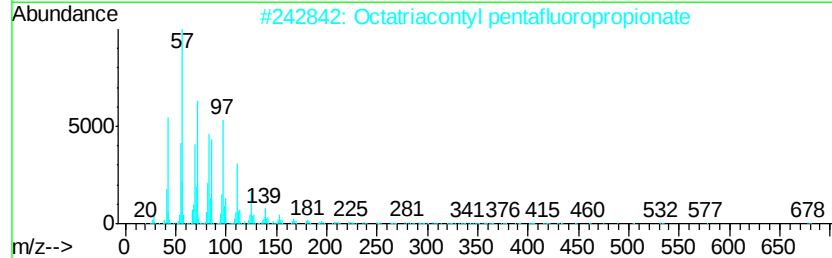
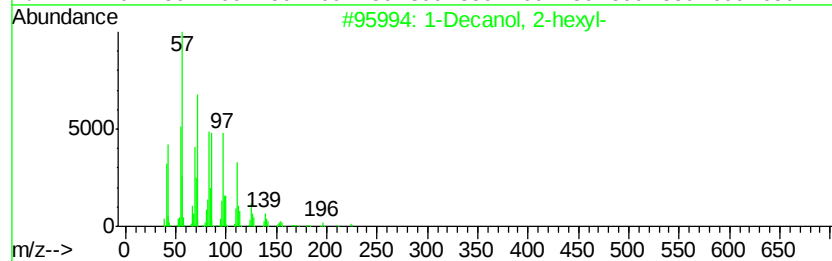
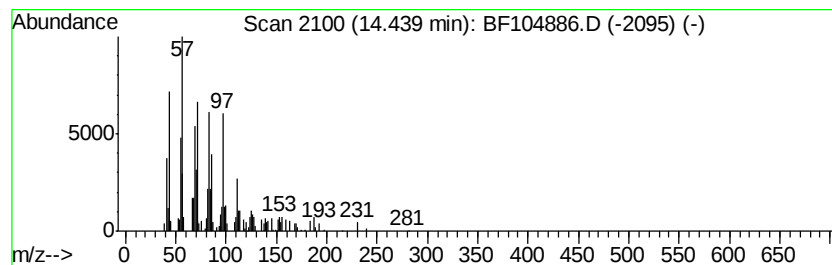
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1-Decanol, 2-hexyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.44	3.56 ng	129106	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	91
2		Octatriacontyl pentafluoropropio...	697	C41H77F5O2	1000351-89-1	90
3		Ethanol, 2-(tetradecyl)-	258	C16H34O2	002136-70-1	87
4		Heptadecyl heptafluorobutyrates	452	C21H35F7O2	959085-66-6	87
5		Heptacosyl pentafluoropropionate	542	C30H55F5O2	1000351-81-6	81



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Sample : J2579-01
Misc :
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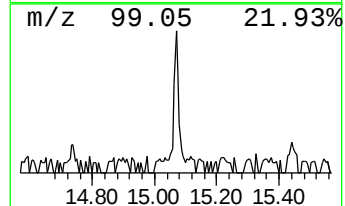
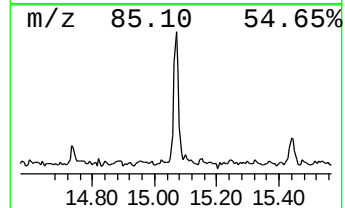
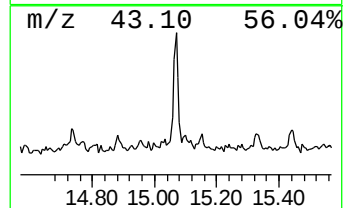
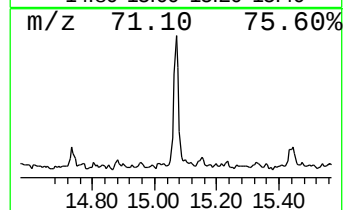
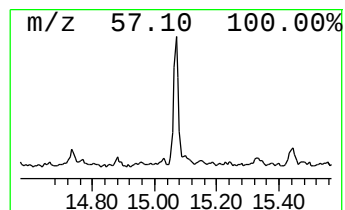
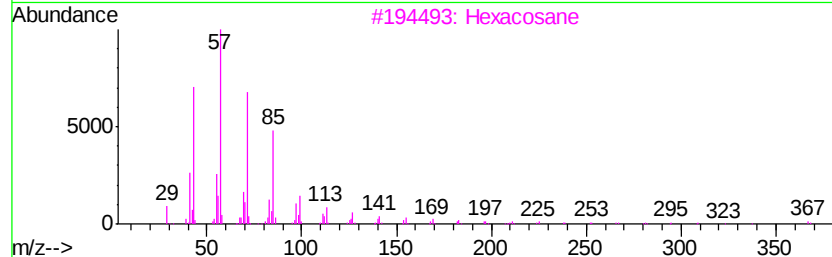
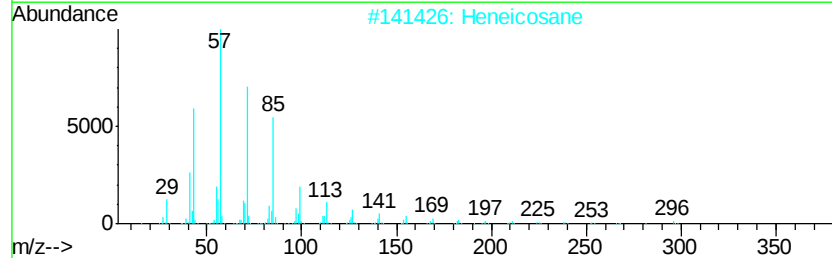
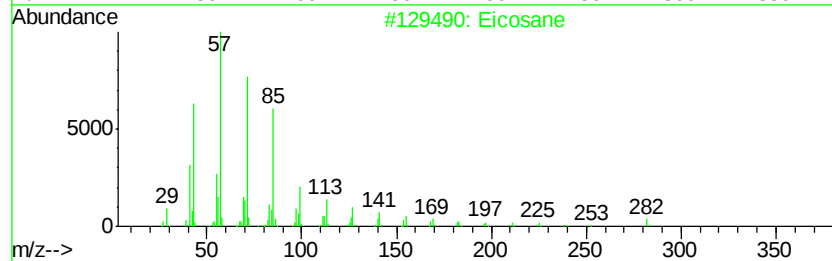
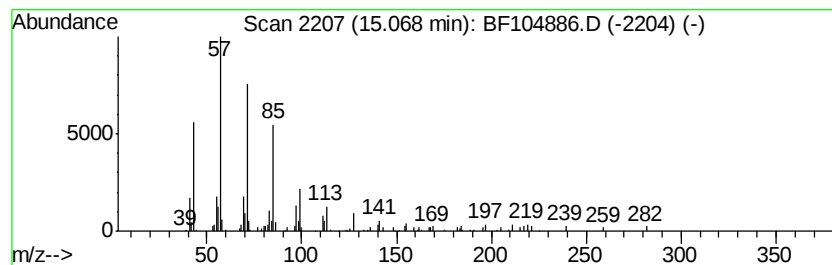
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Eicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	2.64 ng	92213	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	94
2		Heneicosane	296	C21H44	000629-94-7	91
3		Hexacosane	366	C26H54	000630-01-3	90
4		2-methyloctacosane	408	C29H60	1000376-72-8	90
5		Dodecane, 2-methyl-	184	C13H28	001560-97-0	86



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042718\
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Sample : J2579-01
Misc :
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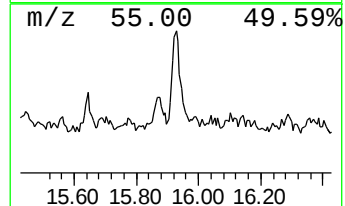
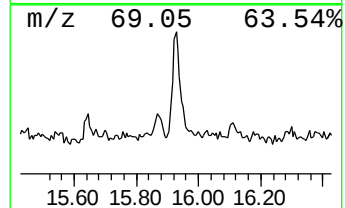
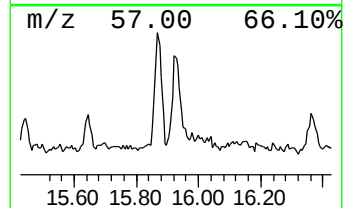
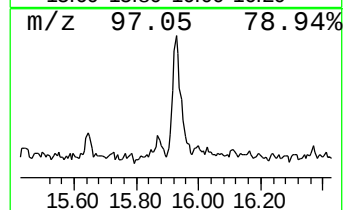
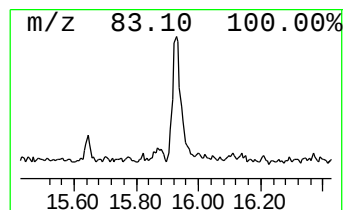
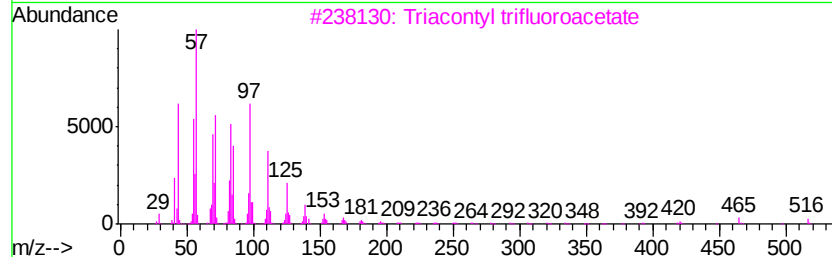
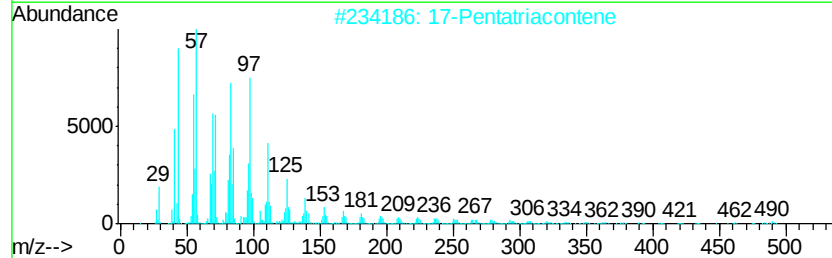
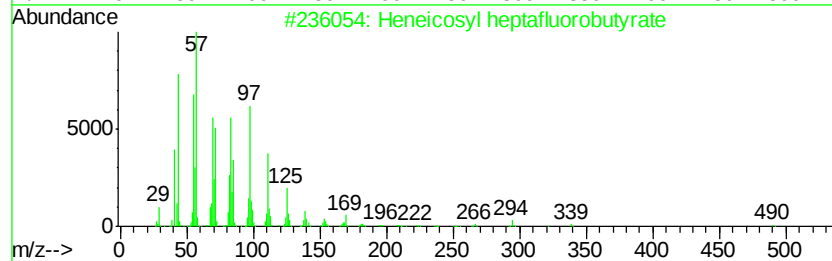
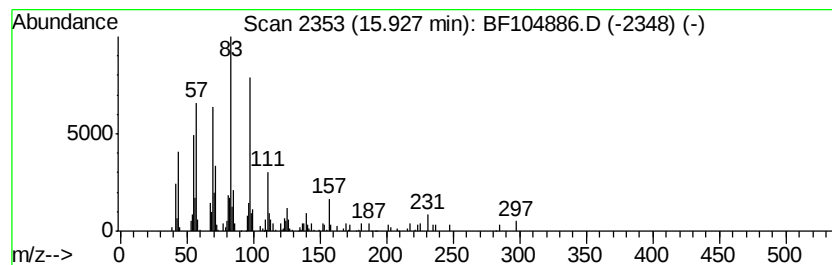
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 unknown15.93 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.93	4.65 ng	162625	Perylene-d12	15.44

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosyl heptafluorobutyrte	508	C25H43F7O2	1000351-83-8	42
2			17-Pentatriacontene	491	C35H70	006971-40-0	42
3			Triacetyl trifluoroacetate	534	C32H61F3O2	1000351-75-7	42
4			Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	42
5			Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	42



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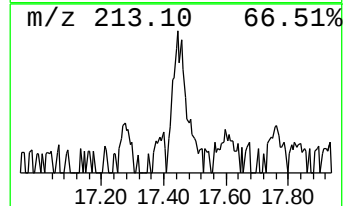
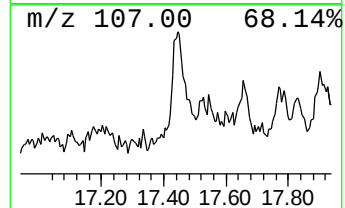
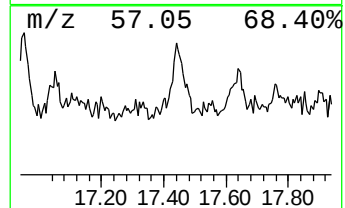
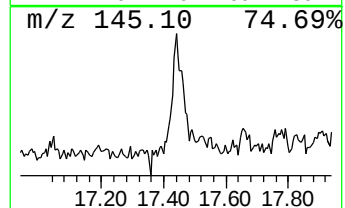
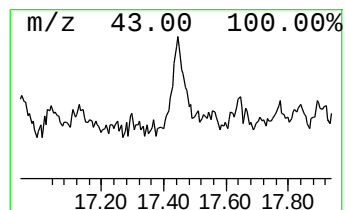
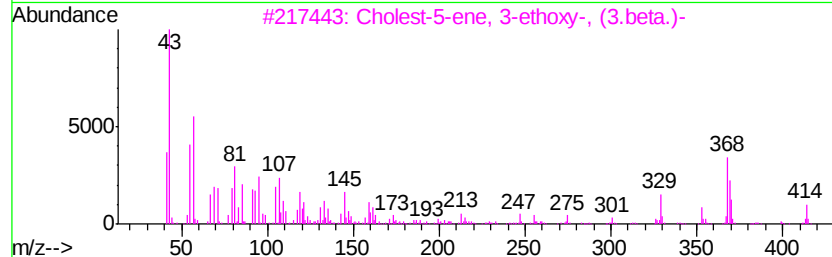
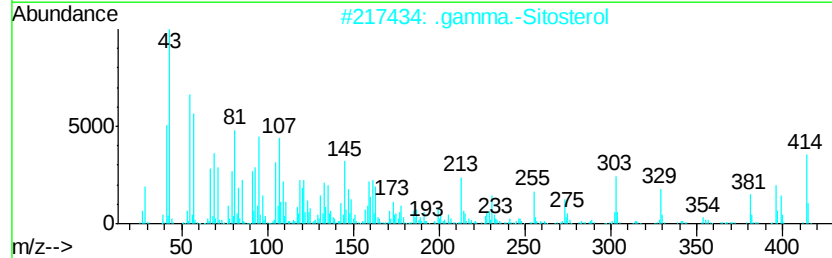
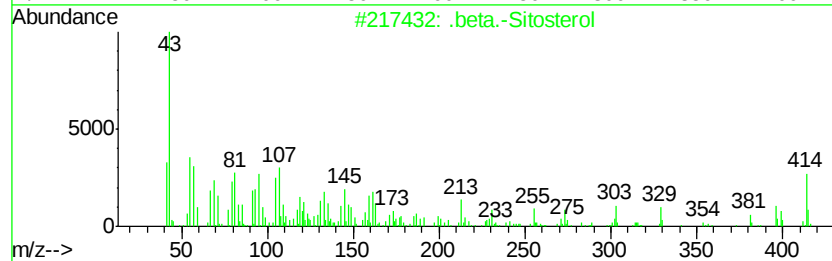
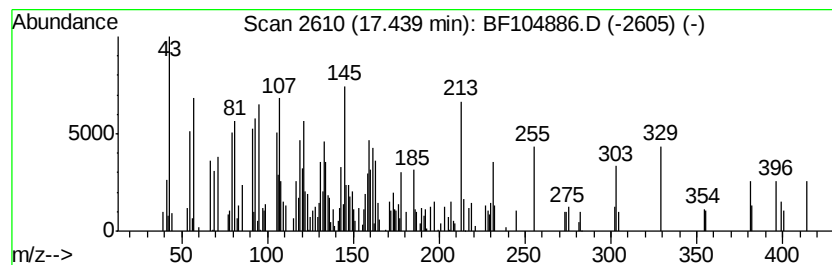
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 .beta.-Sitosterol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.44	6.87 ng	240323	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.beta.-Sitosterol	414	C29H50O	000083-46-5	93
2		.gamma.-Sitosterol	414	C29H50O	000083-47-6	45
3		Cholest-5-ene, 3-ethoxy-, (3.beta...	414	C29H50O	000986-19-6	11
4		5-(1-Benzoxazolyl)-2(1H)-pyrimid...	213	C11H7N3O2	349488-95-5	11
5		Norflurazon	303	C12H9ClF3N3O	027314-13-2	10



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Sample : J2579-01
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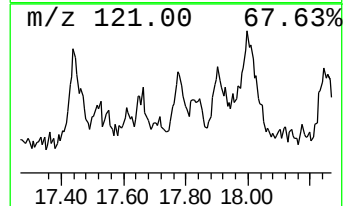
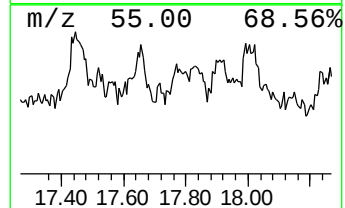
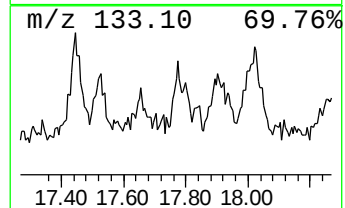
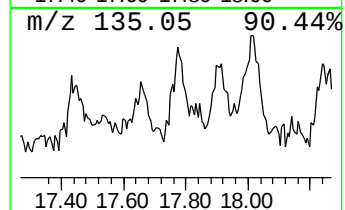
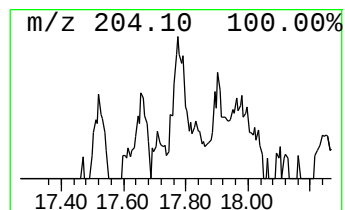
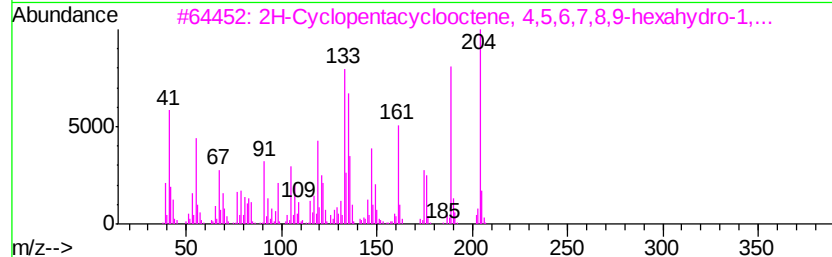
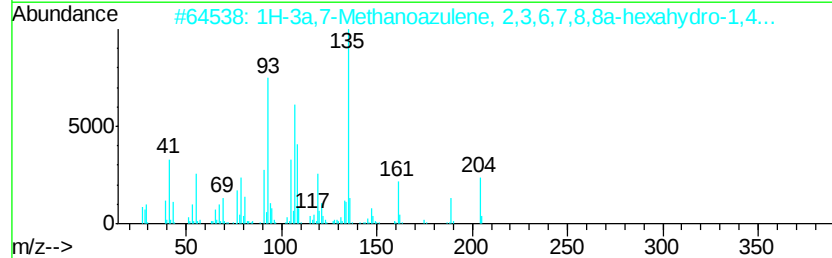
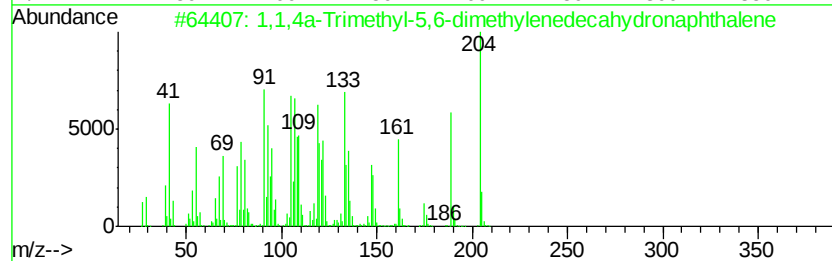
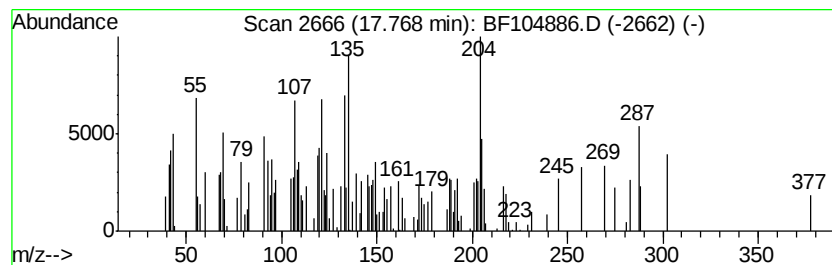
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown17.77 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.77	2.17 ng	75980	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1,4a-Trimethyl-5,6-dimethylene...	204	C15H24	1000193-60-8	38
2		1H-3a,7-Methanoazulene, 2,3,6,7,...	204	C15H24	000560-32-7	38
3		2H-Cyclopentacyclooctene, 4,5,6,...	204	C15H24	1000221-85-8	35
4		Guaia-3,9-diene	204	C15H24	000489-83-8	35
5		Isocaryophyllene	204	C15H24	1000140-07-2	30



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Misc :
ALS Vial : 8 Sample Multiplier: 1

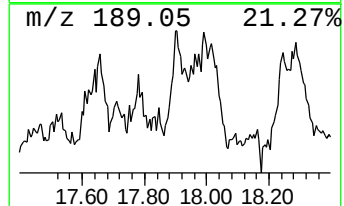
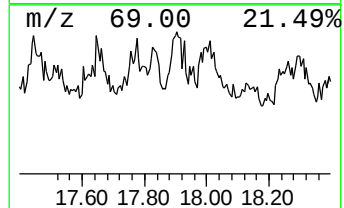
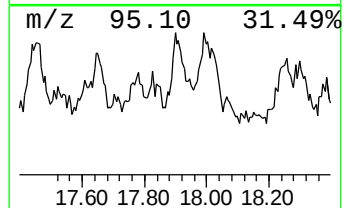
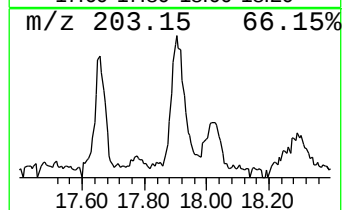
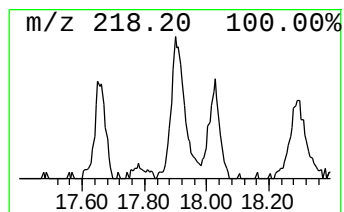
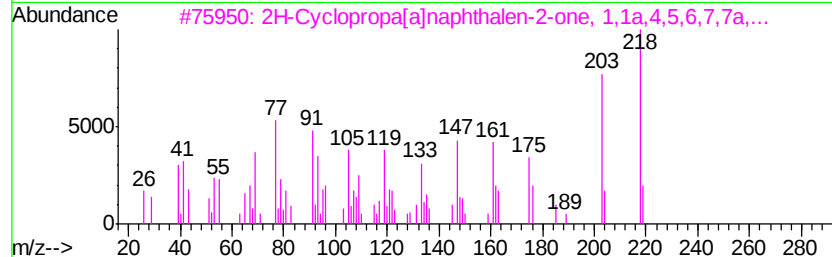
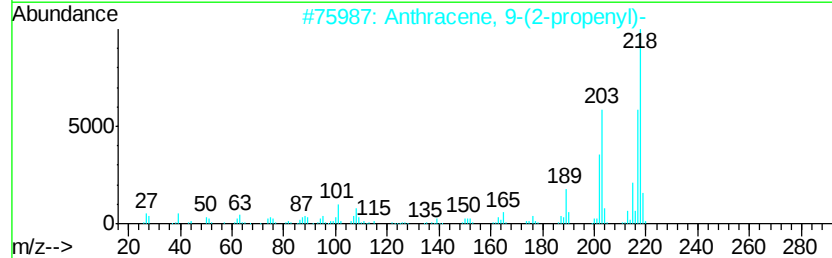
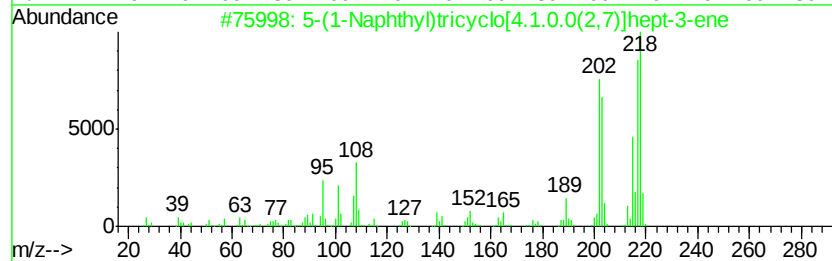
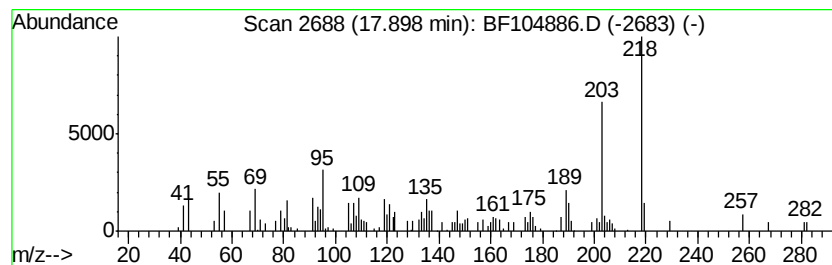
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 5-(1-Naphthyl)tricyclo[4.1.... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.90	4.68 ng	163792	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-(1-Naphthyl)tricyclo[4.1.0.0(2...]	218	C17H14	107729-05-5	80
2		Anthracene, 9-(2-propenyl)-	218	C17H14	023707-65-5	76
3		2H-Cyclopropa[alnaphthalen-2-one...	218	C15H22O	006831-17-0	50
4		Pvrene, hexadecahydro-	218	C16H26	002435-85-0	46
5		Benzo[b]naphtho[2,3-d]furan	218	C16H10O	000243-42-5	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042718\
Data File : BF104886.D
Acq On : 27 Apr 2018 17:30
Operator : JU/SJ
Sample : J2579-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

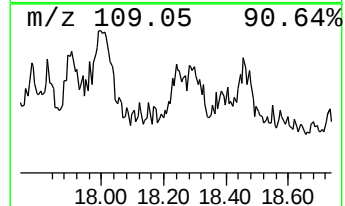
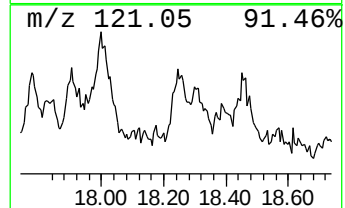
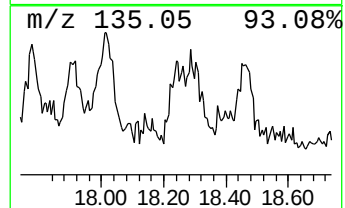
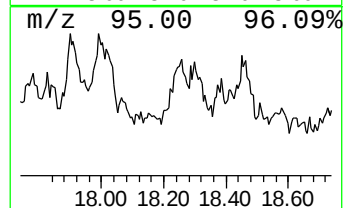
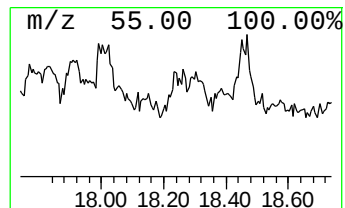
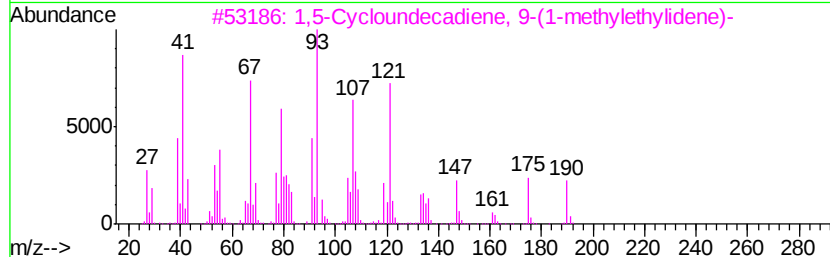
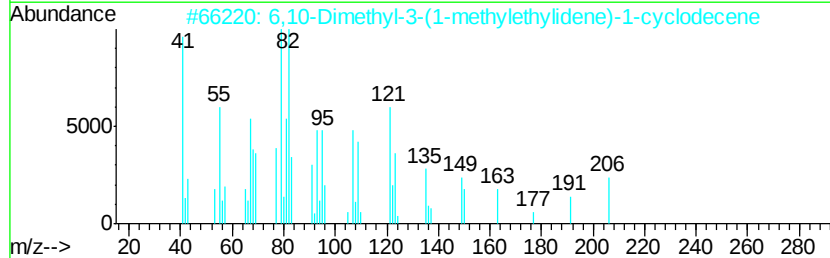
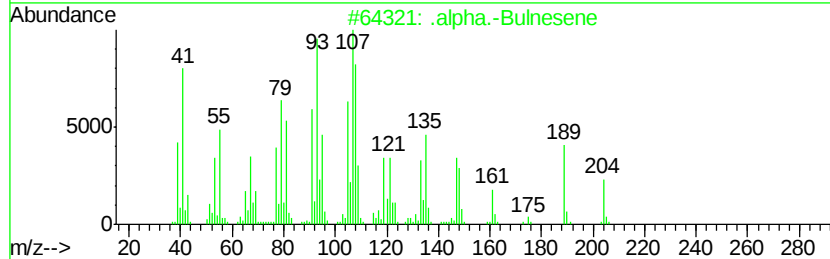
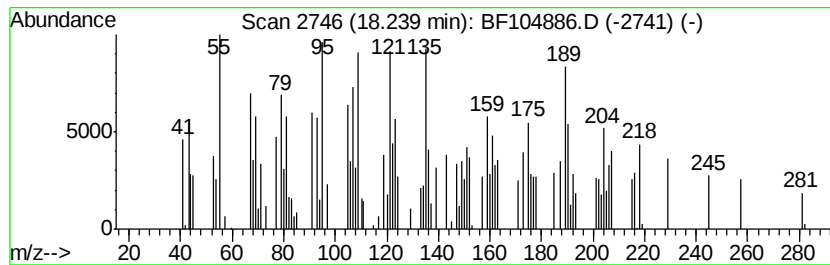
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 .alpha.-Bulnesene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.24	2.33 ng	81488	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.alpha.-Bulnesene	204	C15H24	1000374-19-9	64
2		6,10-Dimethyl-3-(1-methylethylidene)-1-cyclodecene	206	C15H26	069239-71-0	64
3		1,5-Cycloundecadiene, 9-(1-methylethylidene)-	190	C14H22	062338-55-0	62
4		Naphthalene, decahydro-4a-methyl-	204	C15H24	017066-67-0	60
5		Guaia-9,11-diene	204	C15H24	1000374-19-8	56



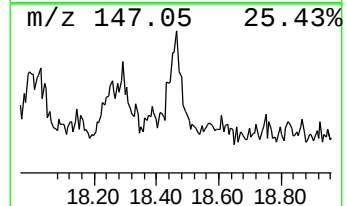
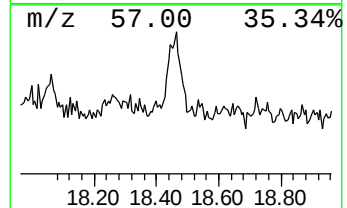
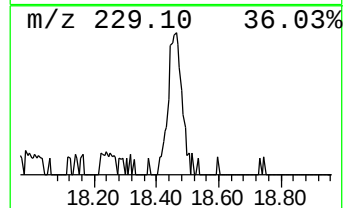
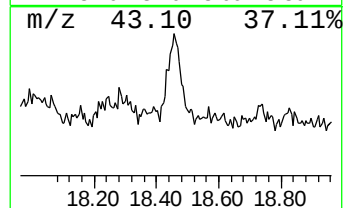
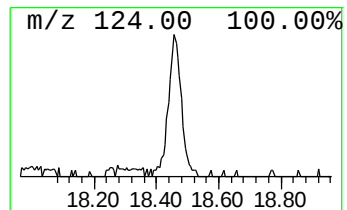
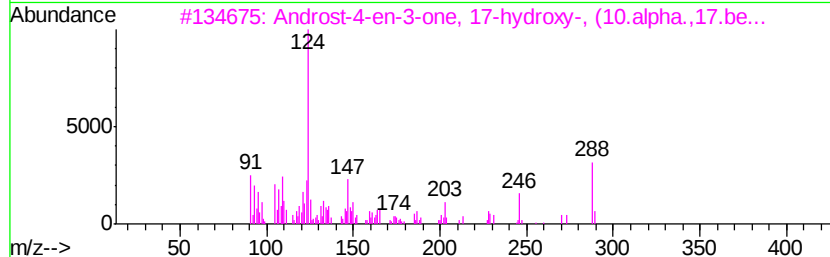
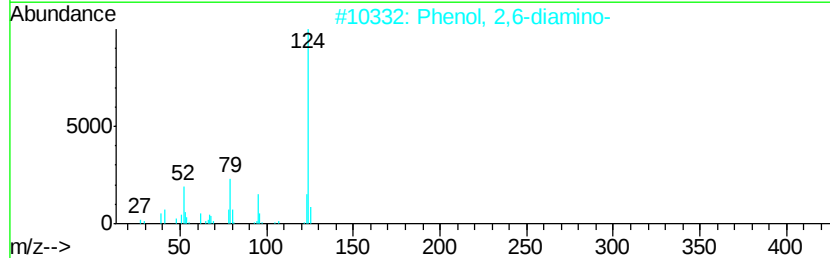
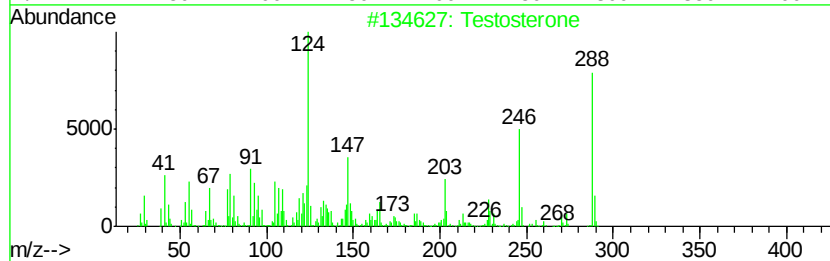
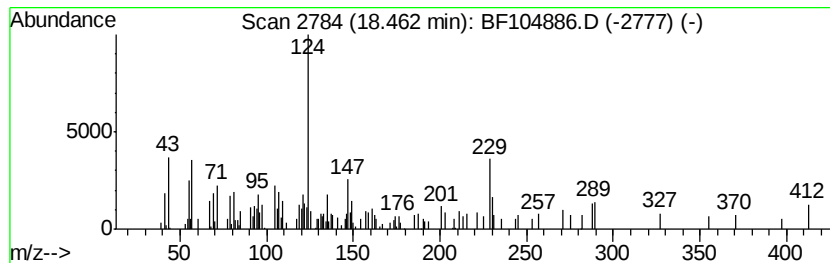
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Acq On : 27 Apr 2018 17:30
Operator : JU/SJ
Sample : J2579-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Testosterone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.46	5.92 ng	207077	Perylene-d12	15.44
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Testosterone	288 C19H28O2	000058-22-0	50
2	Phenol, 2,6-diamino-	124 C6H8N2O	022440-82-0	38
3	Androst-4-en-3-one, 17-hydroxy-,...	288 C19H28O2	000604-39-7	38
4	8-Azabicyclo[3.2.1]octane, 3-chl...	159 C8H14ClN	051275-31-1	38
5	Succinic acid, decyl 4-methoxyph...	364 C21H32O5	1000325-80-0	35



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042718\
Data File : BF104886.D
Acq On : 27 Apr 2018 17:30
Operator : JU/SJ
Sample : J2579-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

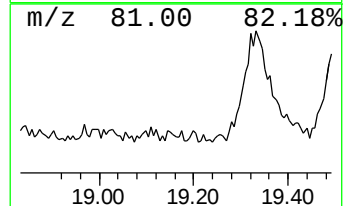
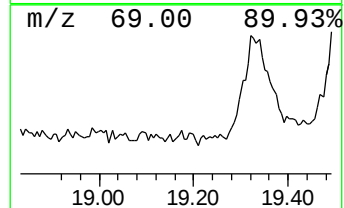
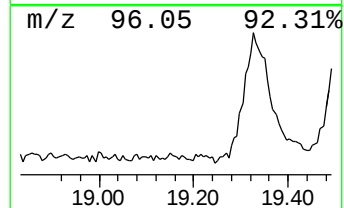
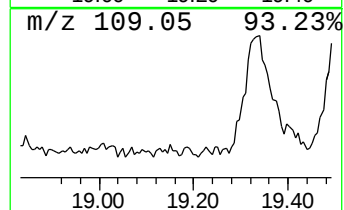
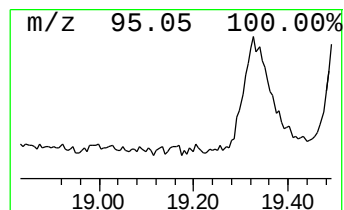
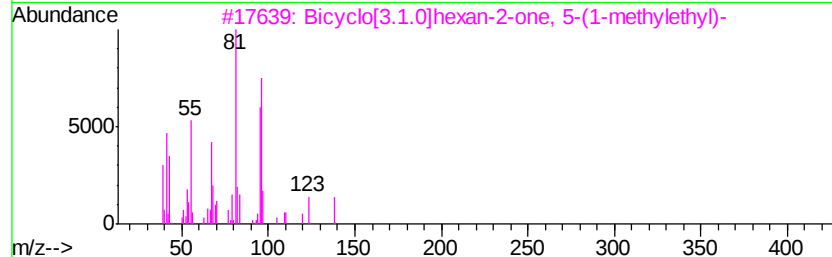
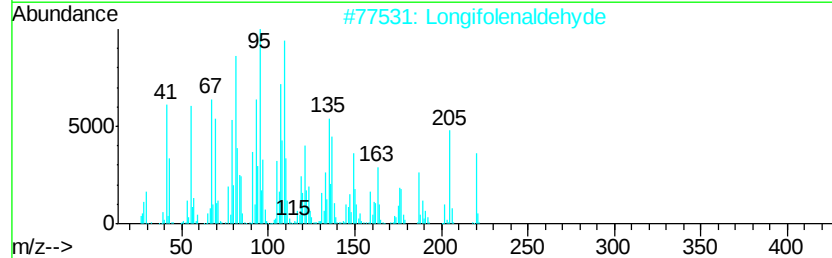
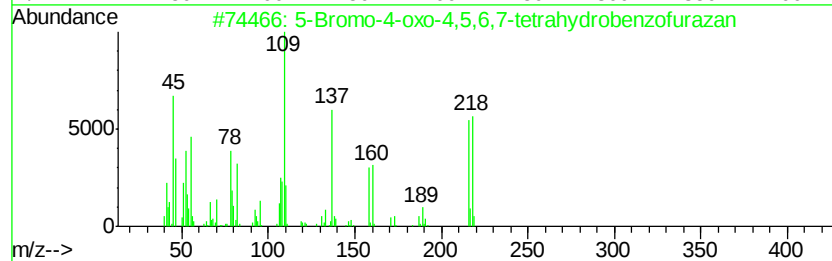
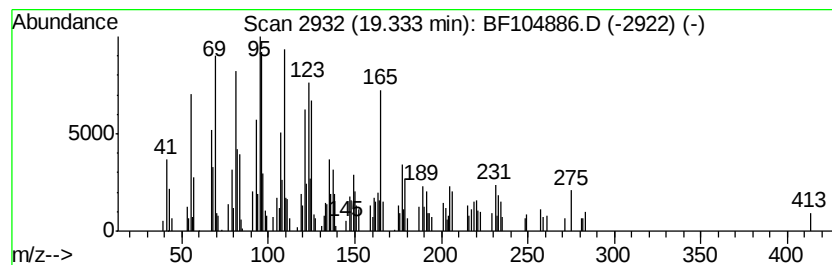
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 5-Bromo-4-oxo-4,5,6,7-tetra... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.33	12.36 ng	432624	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	59
2		Longifolenaldehyde	220	C15H24O	019890-84-7	40
3		Bicyclo[3.1.0]hexan-2-one, 5-(1-...	138	C9H14O	000513-20-2	38
4		1,1,6-trimethyl-3-methylene-2-(3...	452	C33H56	1000373-94-5	38
5		Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	020536-40-7	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042718\
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 Acq On : 27 Apr 2018 17:30
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 Sample : J2579-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.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...	5.02	4.1 ng		125628	1	6.80	610426	20.0
trans-.beta.-Ocimene	6.06	10.4 ng		317047	1	6.80	610426	20.0
unknown6.57	6.57	92.6 ng		2827620	1	6.80	610426	20.0
2-Phenanthrenol, ...	13.34	2.3 ng		83312	5	13.97	725794	20.0
Dichloroacetic ac...	13.80	8.7 ng		316398	5	13.97	725794	20.0
3,6,9,12-Tetraoxa...	13.84	4.2 ng		153470	5	13.97	725794	20.0
1-Decanol, 2-hexyl-	14.44	3.6 ng		129106	5	13.97	725794	20.0
Eicosane	15.07	2.6 ng		92213	6	15.44	699883	20.0
unknown15.93	15.93	4.7 ng		162625	6	15.44	699883	20.0
.beta.-Sitosterol	17.44	6.9 ng		240323	6	15.44	699883	20.0
unknown17.77	17.77	2.2 ng		75980	6	15.44	699883	20.0
5-(1-Naphthyl)tri...	17.90	4.7 ng		163792	6	15.44	699883	20.0
.alpha.-Bulnesene	18.24	2.3 ng		81488	6	15.44	699883	20.0
Testosterone	18.46	5.9 ng		207077	6	15.44	699883	20.0
5-Bromo-4-oxo-4,5...	19.33	12.4 ng		432624	6	15.44	699883	20.0