

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF091117\
 Data File : BF098431.D
 Acq On : 12 Sep 2017 4:43
 Operator : SJ/JU
 Sample : I5138-17
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EME-WM-TS003-01

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-BF091117.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	4.869	469	473	487	rBV	443998	741347	9.97%	1.188%
2	5.287	539	544	547	rBV	6558441	7118504	95.73%	11.406%
3	6.328	715	721	731	rBV	5826345	7435918	100.00%	11.915%
4	6.445	736	741	744	rBV	6708180	6763359	90.96%	10.837%
5	6.669	775	779	789	rBV	1695875	1508591	20.29%	2.417%
6	6.822	801	805	809	rBV	4720434	4687837	63.04%	7.511%
7	7.239	870	876	879	rBV	4329062	4532830	60.96%	7.263%
8	7.269	879	881	887	rVB2	63752	74393	1.00%	0.119%
9	7.951	991	997	1000	rBV	2337121	2012422	27.06%	3.225%
10	8.545	1094	1098	1101	rVB	86631	79689	1.07%	0.128%
11	8.975	1168	1171	1174	rVB	97399	76061	1.02%	0.122%
12	9.027	1174	1180	1183	rBV	5708052	5131244	69.01%	8.222%
13	9.110	1189	1194	1198	rBV2	99358	111890	1.50%	0.179%
14	9.422	1244	1247	1251	rVB	1678230	1518126	20.42%	2.433%
15	9.698	1287	1294	1298	rBV2	2144094	2039829	27.43%	3.268%
16	10.416	1412	1416	1422	rBV3	85614	111322	1.50%	0.178%
17	10.492	1424	1429	1432	rBV	3407144	3193952	42.95%	5.118%
18	10.604	1444	1448	1450	rBV2	91041	102641	1.38%	0.164%
19	10.763	1472	1475	1477	rBV2	172570	150295	2.02%	0.241%
20	10.839	1486	1488	1493	rVB	109323	94968	1.28%	0.152%
21	10.886	1493	1496	1501	rVV	138742	130256	1.75%	0.209%
22	10.957	1501	1508	1518	rVB3	123434	218466	2.94%	0.350%
23	11.069	1522	1527	1535	rVB2	117270	160021	2.15%	0.256%
24	11.180	1542	1546	1554	rVV	1981950	1680450	22.60%	2.693%
25	11.251	1554	1558	1561	rVB2	117498	142711	1.92%	0.229%
26	11.592	1611	1616	1618	rVB3	65898	93647	1.26%	0.150%
27	11.868	1658	1663	1670	rBV4	97934	135837	1.83%	0.218%
28	12.263	1725	1730	1732	rBV	163076	167941	2.26%	0.269%
29	12.280	1732	1733	1737	rVB2	98624	83448	1.12%	0.134%
30	12.386	1746	1751	1755	rBV	99765	127950	1.72%	0.205%
31	12.427	1755	1758	1762	rVB	207461	181492	2.44%	0.291%
32	12.616	1788	1790	1793	rVV	102689	85961	1.16%	0.138%
33	12.663	1793	1798	1803	rVB3	52904	96308	1.30%	0.154%
34	12.763	1811	1815	1819	rBV	3211847	2923310	39.31%	4.684%

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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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.657	1964	1967	1972	rBV	442700	501029	6.74%	0.803%
36	13.810	1988	1993	1997	rBV	1241050	1247238	16.77%	1.998%
37	14.292	2071	2075	2079	rBV2	233807	306849	4.13%	0.492%
38	14.821	2162	2165	2167	rBV	73250	88824	1.19%	0.142%
39	14.892	2173	2177	2180	rBV	792324	885276	11.91%	1.418%
40	15.156	2219	2222	2225	rVB	76735	92479	1.24%	0.148%
41	15.215	2227	2232	2239	rBV	930801	1191789	16.03%	1.910%
42	15.633	2298	2303	2307	rBV	530469	760420	10.23%	1.218%
43	16.621	2467	2471	2477	rVB	138261	253880	3.41%	0.407%
44	17.003	2531	2536	2541	rBV3	181414	355842	4.79%	0.570%
45	17.056	2541	2545	2546	rBV3	176980	211305	2.84%	0.339%
46	17.468	2610	2615	2624	rBV3	178205	532179	7.16%	0.853%
47	17.562	2627	2631	2640	rVB8	235791	581262	7.82%	0.931%
48	17.780	2661	2668	2685	rBV7	293847	1245785	16.75%	1.996%
49	18.921	2856	2862	2874	rVB5	154213	442905	5.96%	0.710%

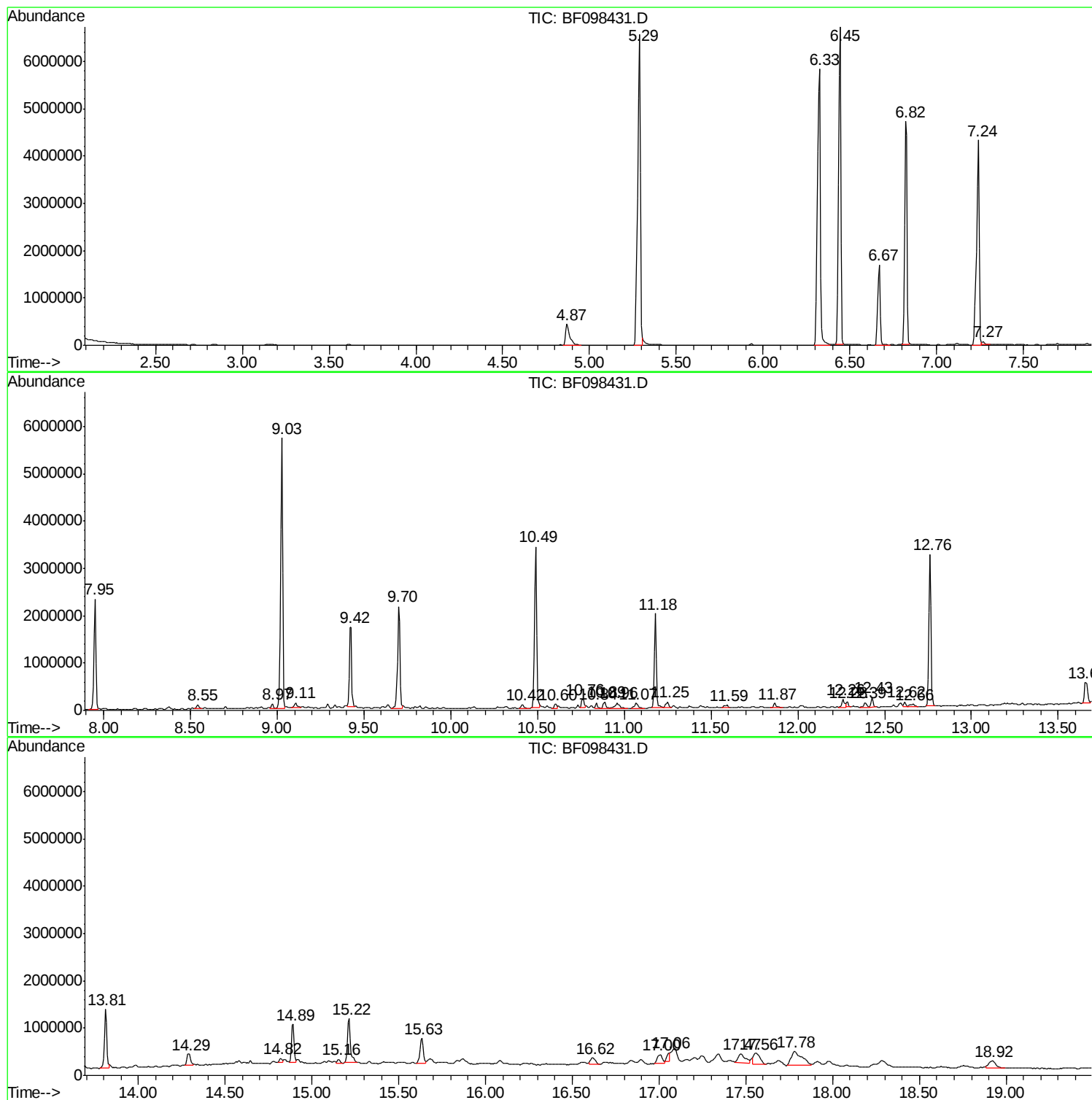
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF091117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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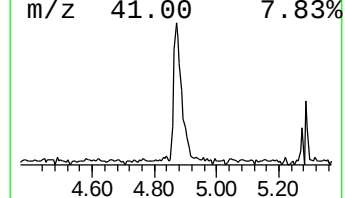
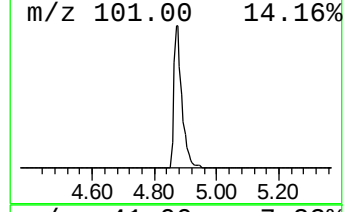
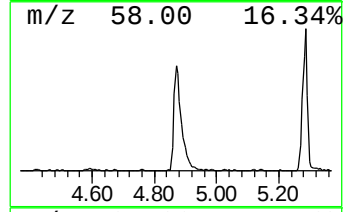
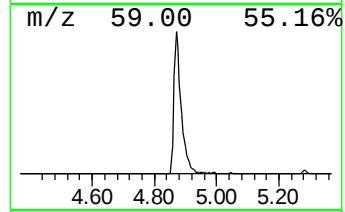
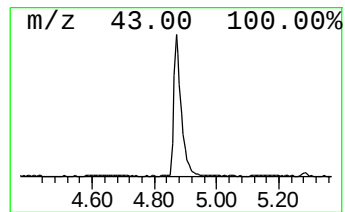
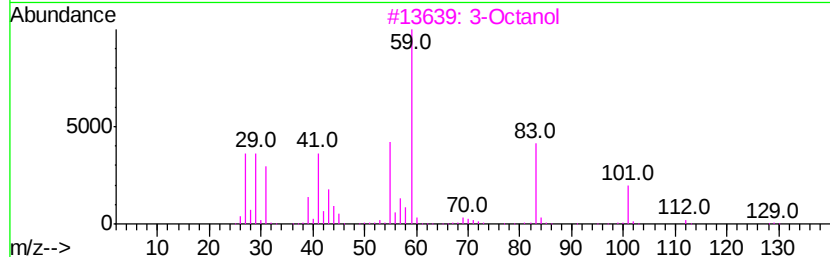
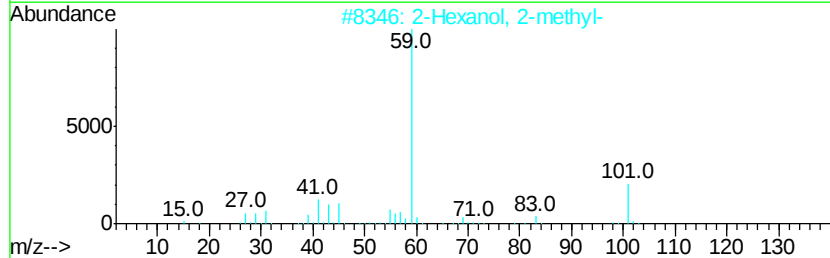
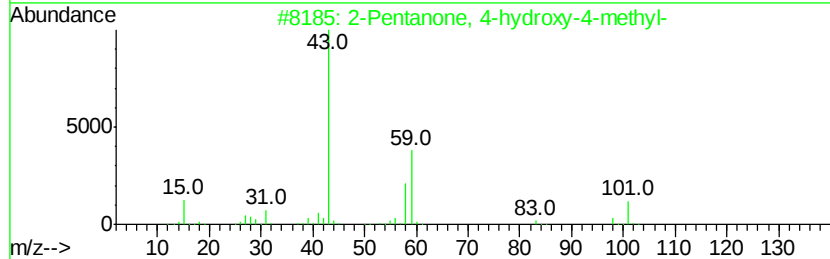
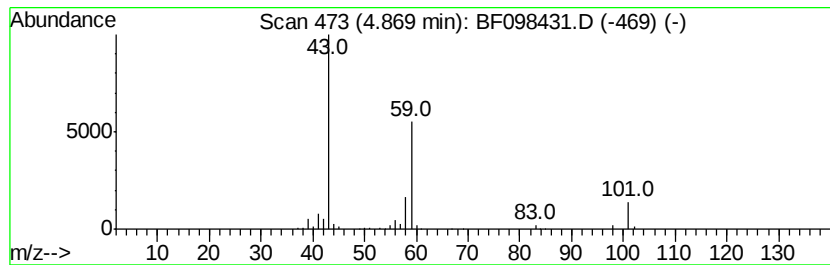
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF091117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.87	9.83 ng	741347	1,4-Dichlorobenzene-d4	6.67

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3		3-Octanol	130	C8H18O	000589-98-0	33
4		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
5		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23



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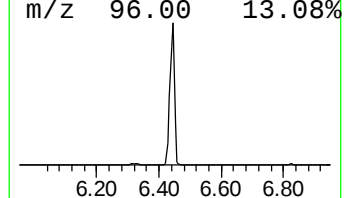
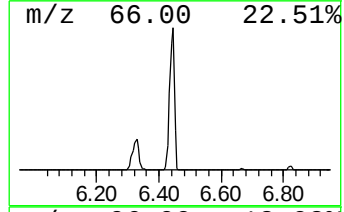
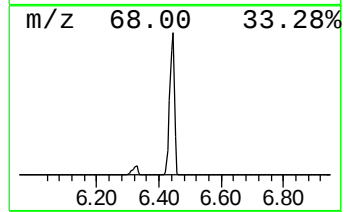
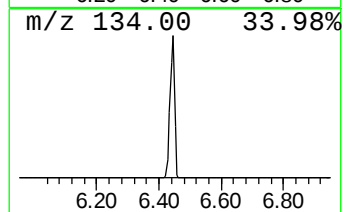
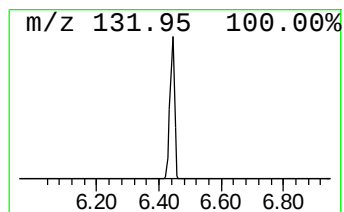
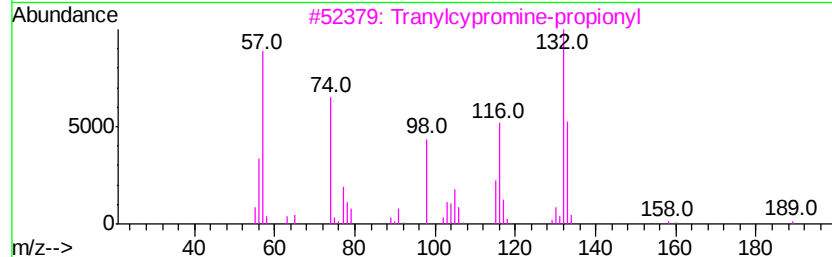
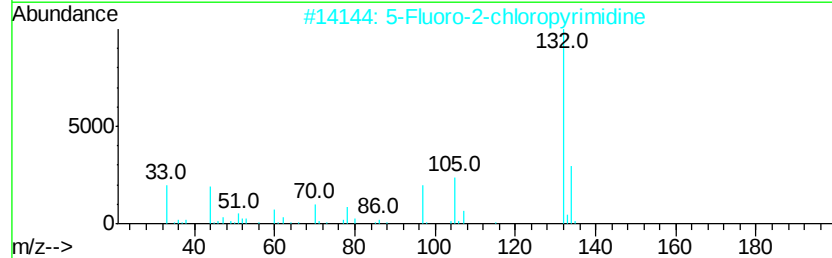
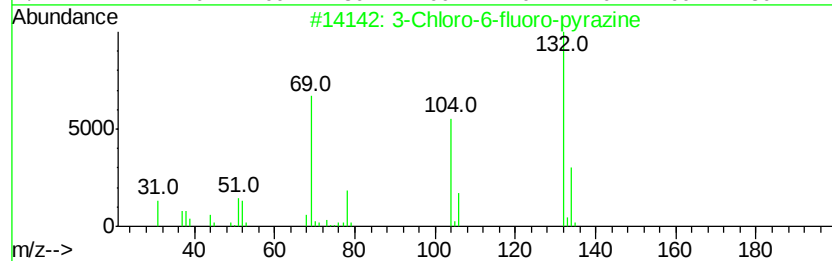
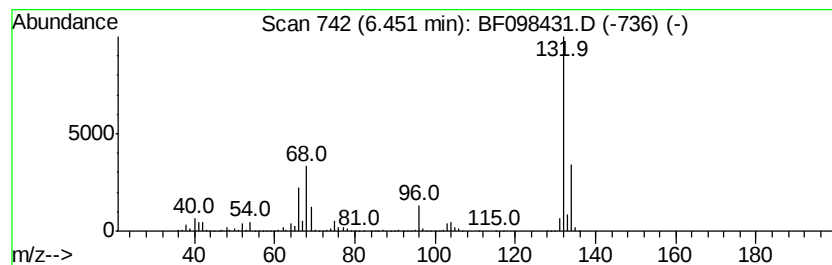
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown6.45 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.45	89.66 ng	6763360	1,4-Dichlorobenzene-d4	6.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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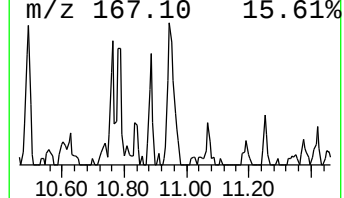
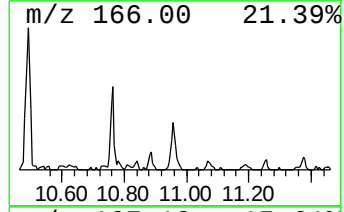
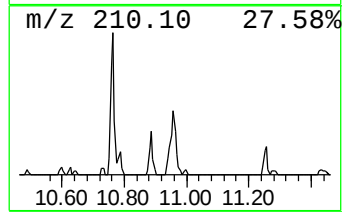
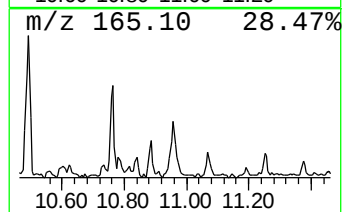
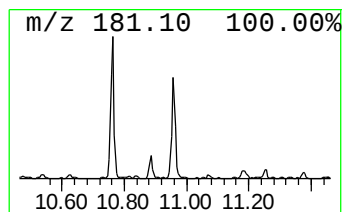
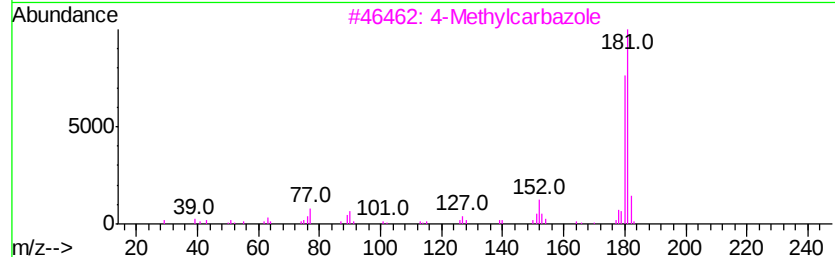
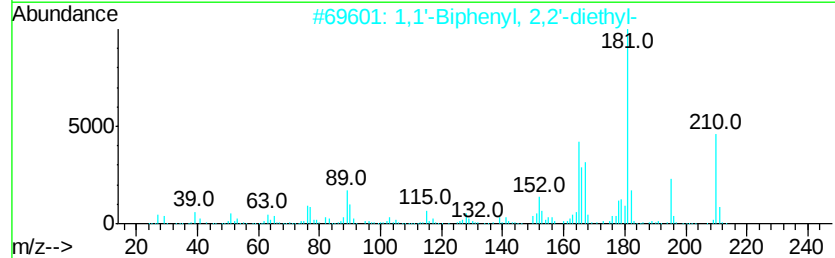
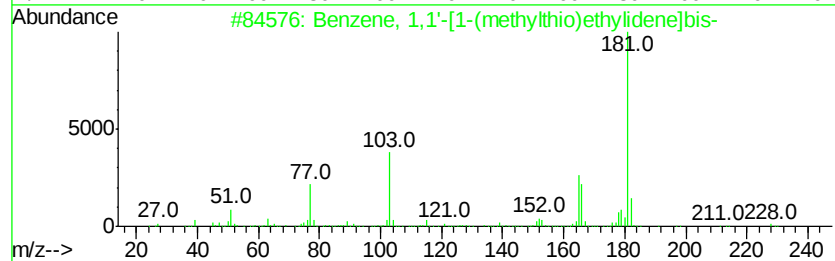
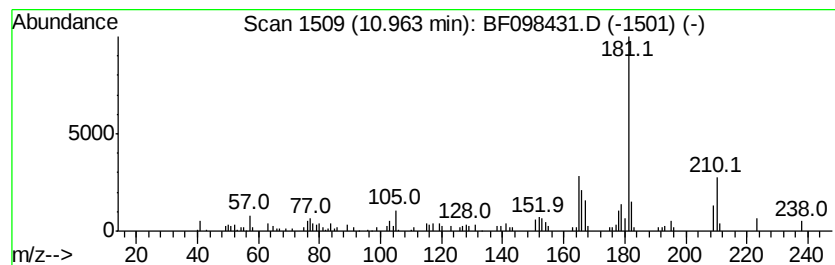
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1,1'-[1-(methylthio)eth... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.96	2.60 ng	218466	Phenanthrene-d10	11.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,1'-[1-(methylthio)eth...	228	C15H16S	054851-63-7	72
2		1,1'-Biphenyl, 2,2'-diethyl-	210	C16H18	013049-35-9	68
3		4-Methylcarbazole	181	C13H11N	003770-48-7	60
4		1-Methylcarbazole	181	C13H11N	006510-65-2	60
5		9H-Carbazole, 2-methyl-	181	C13H11N	003652-91-3	60



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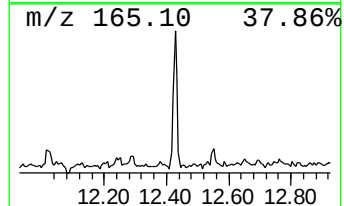
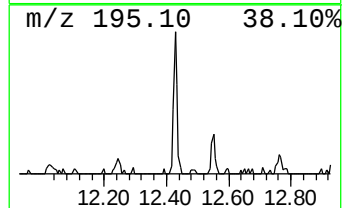
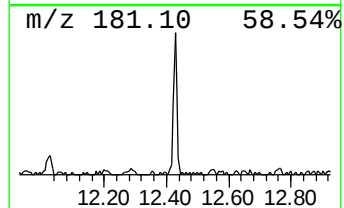
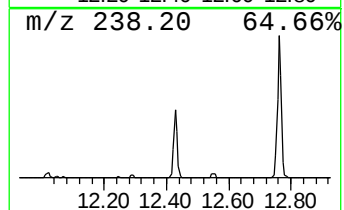
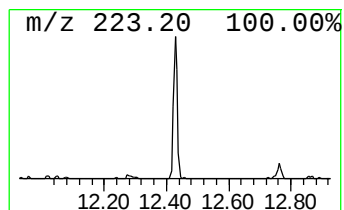
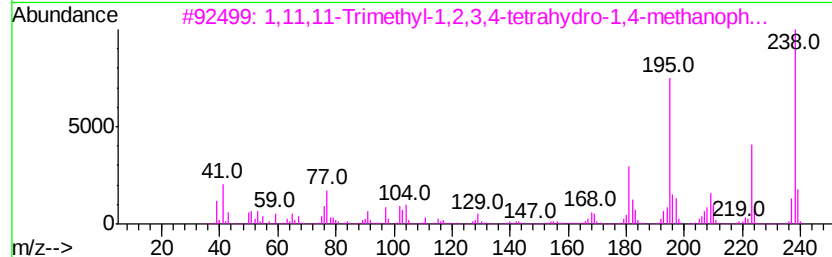
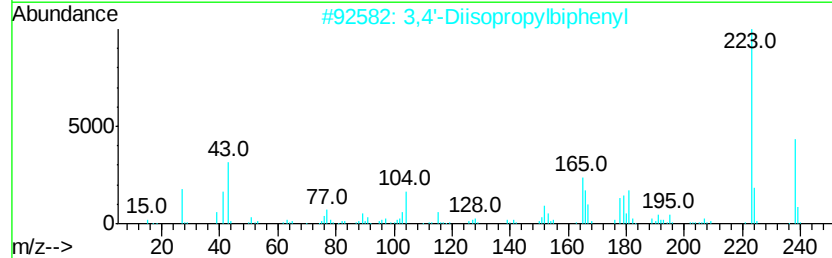
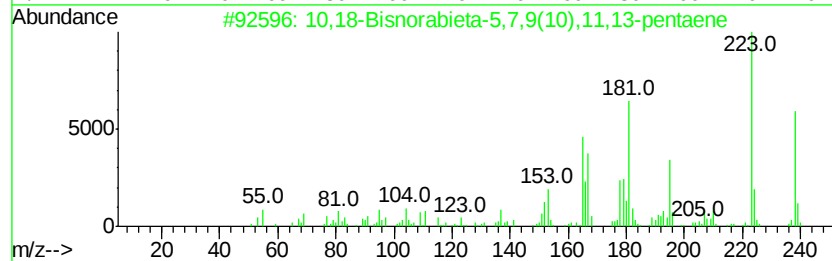
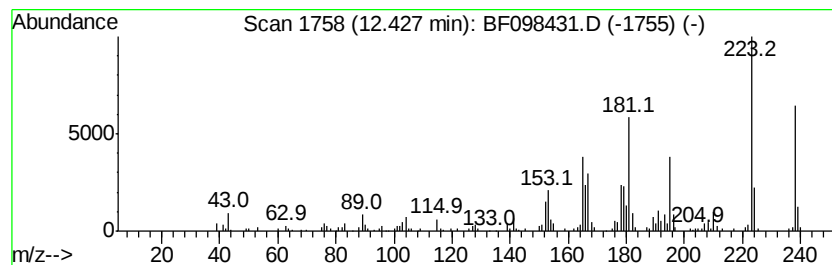
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.43	2.16 ng	181492	Phenanthrene-d10	11.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99
2		3,4'-Diisopropylbiphenyl	238	C18H22	061434-46-6	64
3		1,11,11-Trimethyl-1,2,3,4-tetra...	238	C16H18N2	1000210-51-9	55
4		3-(N,N-Diethylamino)carbazole	238	C16H18N2	054994-31-9	53
5		2-Morpholin-4-yl-5,6-dihydro-4H-...	238	C11H14N2O2S	313251-21-7	45



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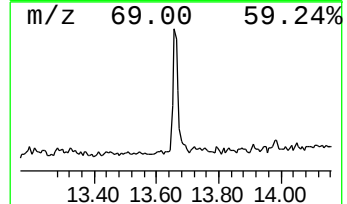
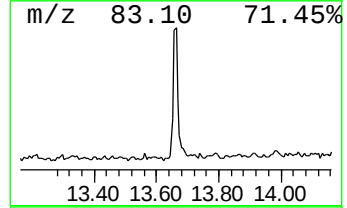
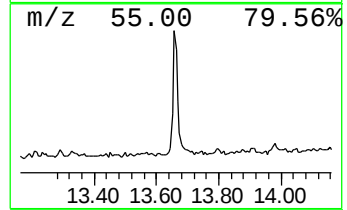
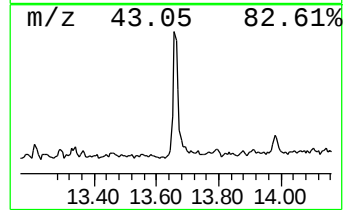
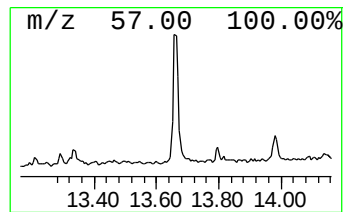
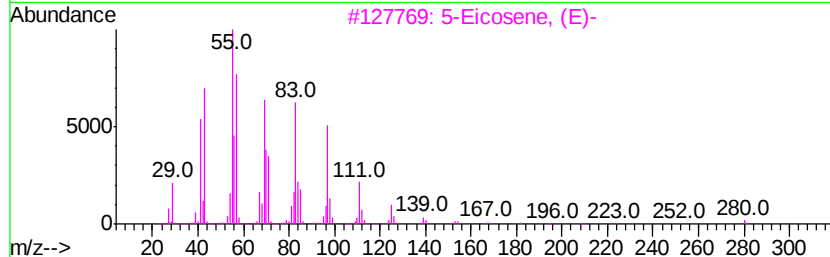
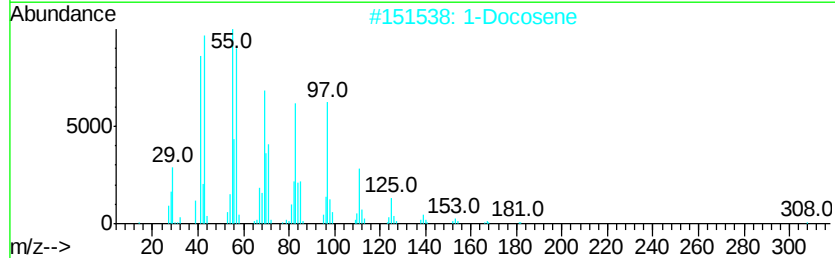
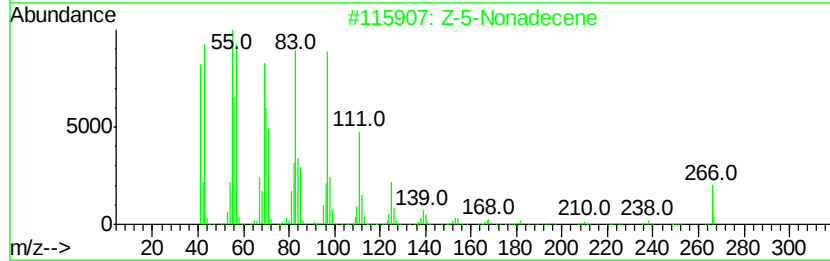
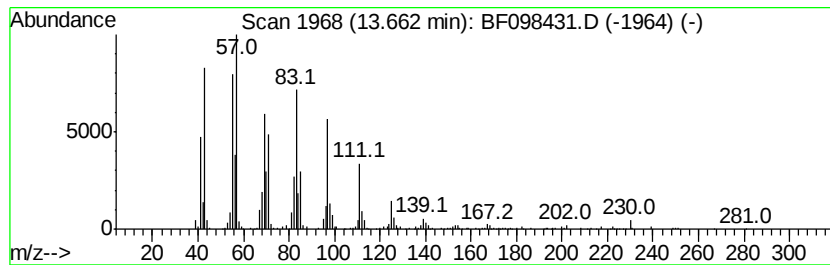
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Z-5-Nonadecene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.66	8.03 ng	501029	Chrysene-d12	13.81	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Z-5-Nonadecene	266	C19H38	1000131-11-8	95
2	1-Docosene	308	C22H44	001599-67-3	94
3	5-Eicosene, (E)-	280	C20H40	074685-30-6	91
4	Behenic alcohol	326	C22H46O	000661-19-8	91
5	n-Tetracosanol-1	354	C24H50O	000506-51-4	91



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091117\
Data File : BF098431.D
Acq On : 12 Sep 2017 4:43
Operator : SJ/JU
Sample : I5138-17
Misc :
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Instrument :
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ClientSampleId :
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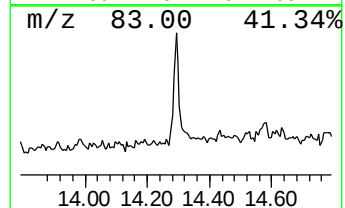
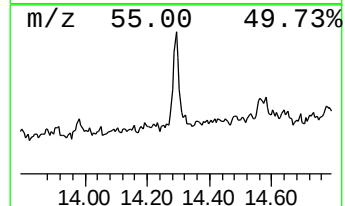
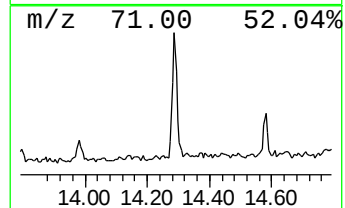
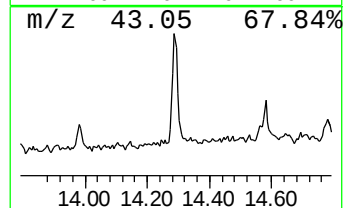
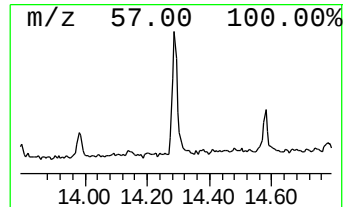
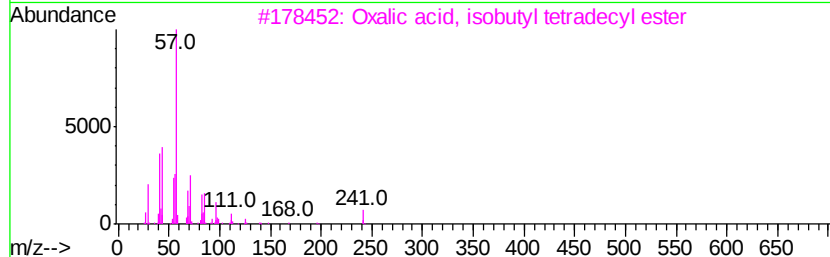
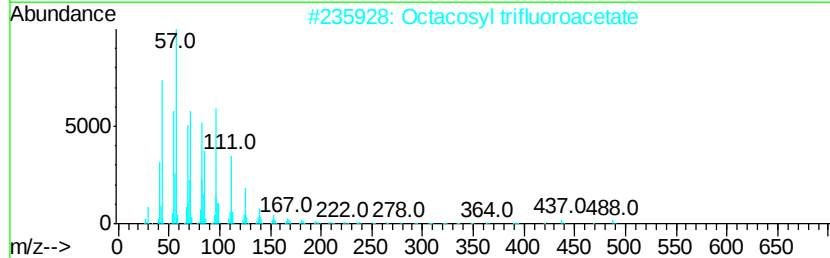
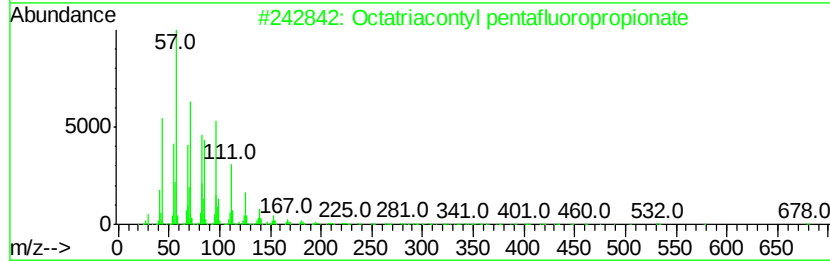
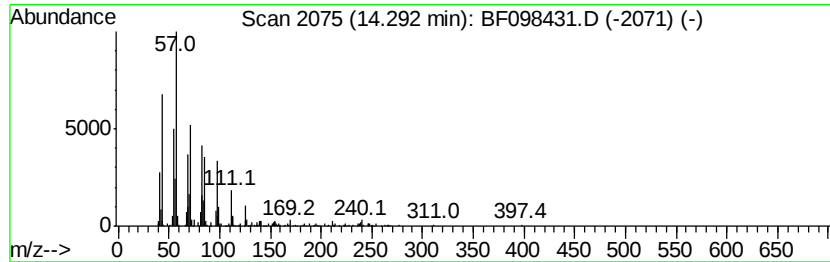
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octatriacontyl pentafluorop... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.29	4.92 ng	306849	Chrysene-d12	13.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octatriacontyl pentafluoropropio...	697	C41H77F5O2	1000351-89-1	91
2		Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	91
3		Oxalic acid, isobutyl tetradecyl...	342	C20H38O4	1000309-37-9	91
4		Tetratriacontyl pentafluoropropi...	641	C37H69F5O2	1000351-81-5	91
5		Tetratriacontyl heptafluorobutyrate	691	C38H69F7O2	1000351-84-1	91



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091117\
Data File : BF098431.D
Acq On : 12 Sep 2017 4:43
Operator : SJ/JU
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ClientSampleId :
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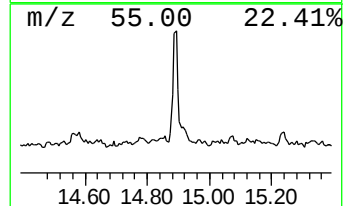
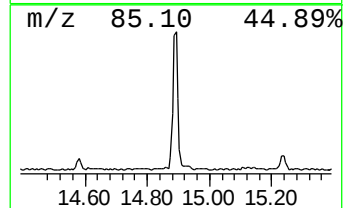
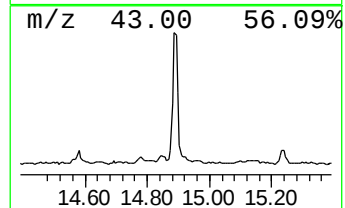
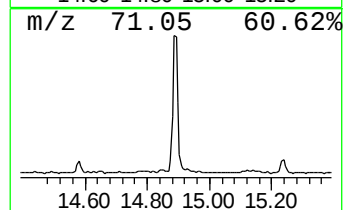
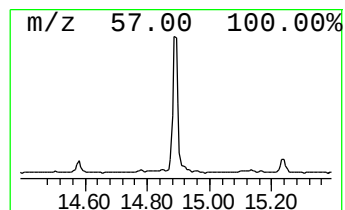
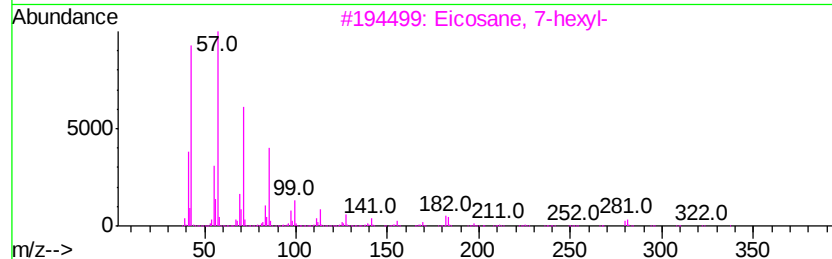
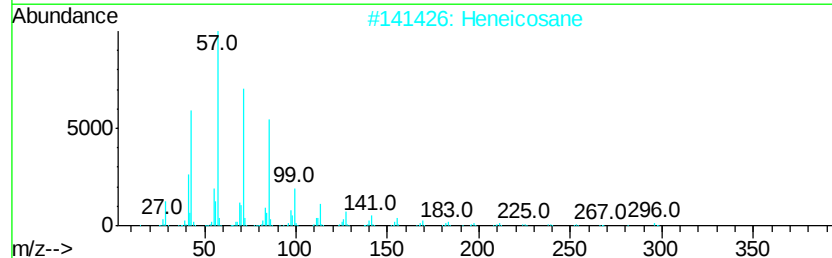
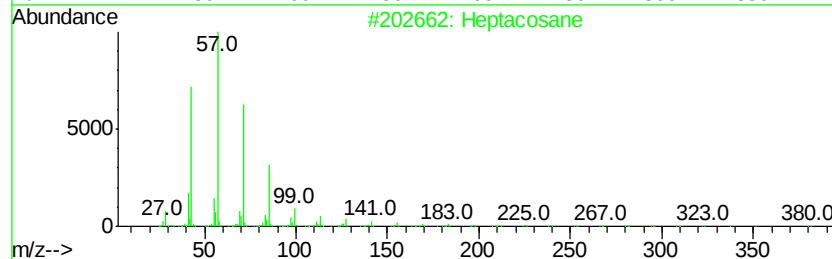
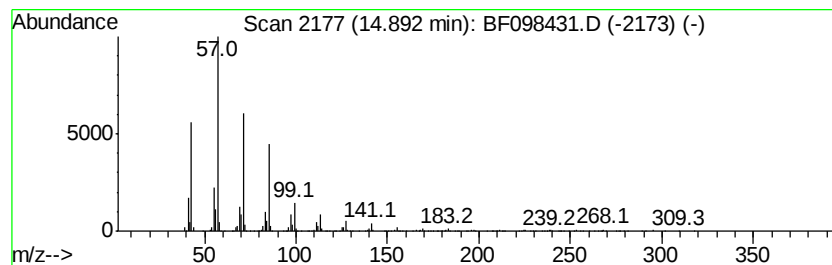
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.89	14.86 ng	885276	Perylene-d12	15.22

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane		380	C27H56	000593-49-7	91
2	Heneicosane		296	C21H44	000629-94-7	90
3	Eicosane, 7-hexyl-		366	C26H54	055333-99-8	90
4	triacontane		422	C30H62	000638-68-6	90
5	Nonadecane		268	C19H40	000629-92-5	87



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF091117\
Data File : BF098431.D
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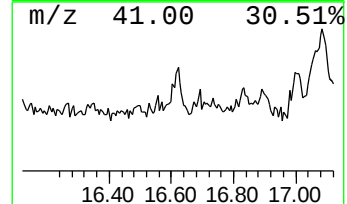
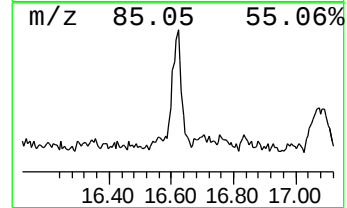
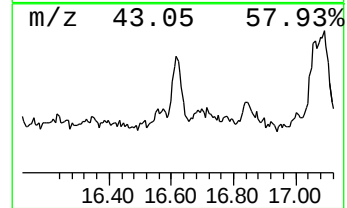
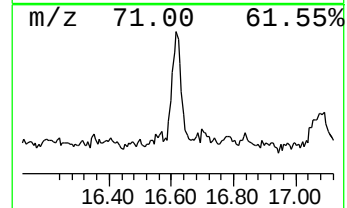
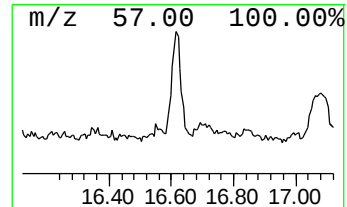
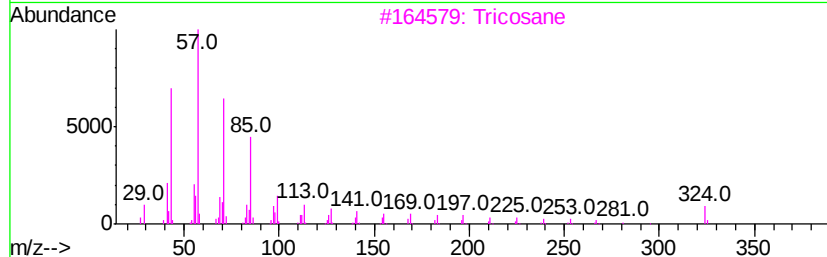
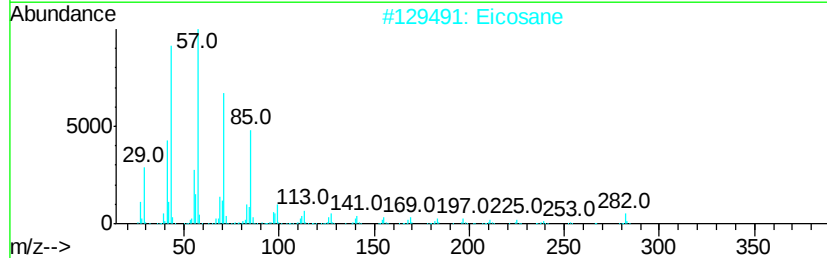
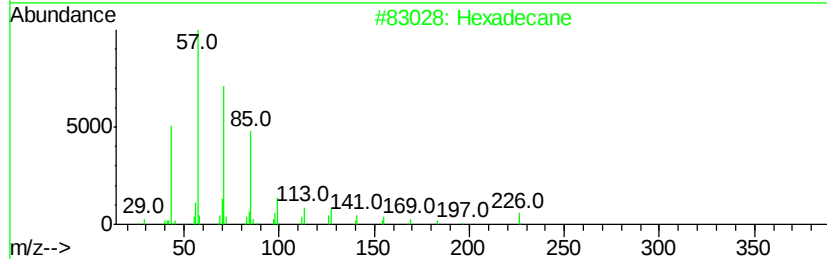
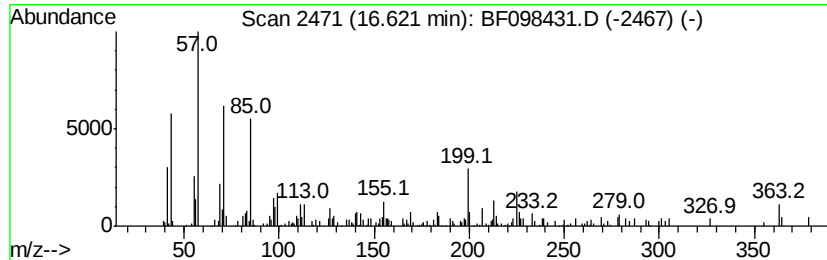
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.62	4.26 ng	253880	Perylene-d12	15.22

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	70
2			Eicosane	282	C20H42	000112-95-8	55
3			Tricosane	324	C23H48	000638-67-5	49
4			Heneicosane	296	C21H44	000629-94-7	49
5			Hentriacontane	437	C31H64	000630-04-6	49



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091117\
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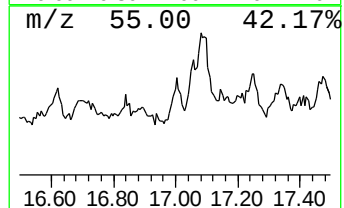
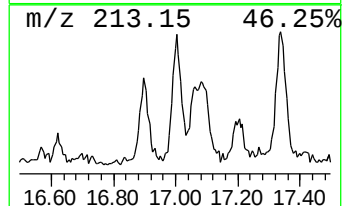
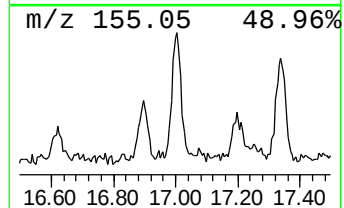
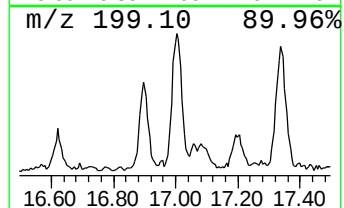
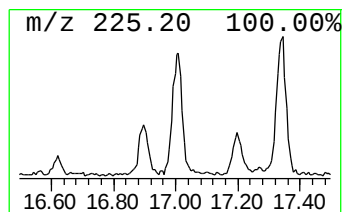
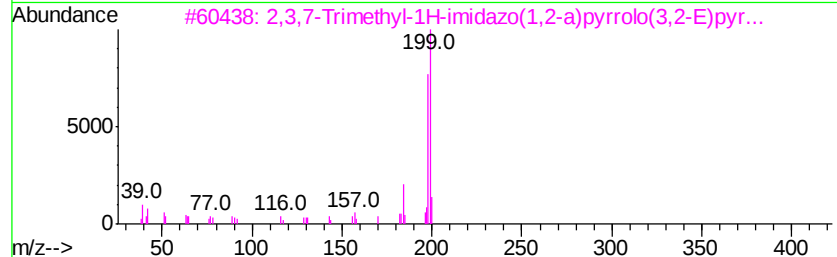
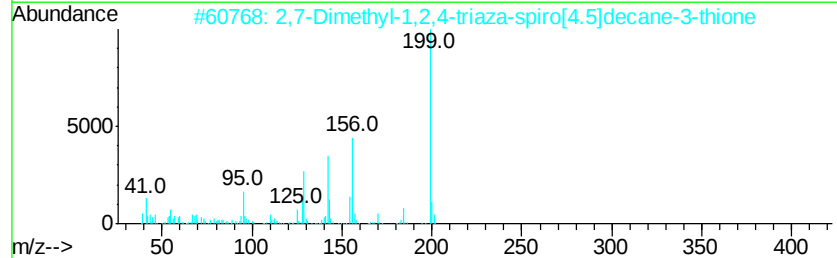
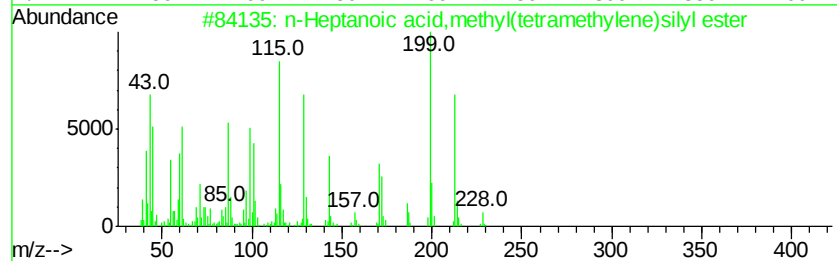
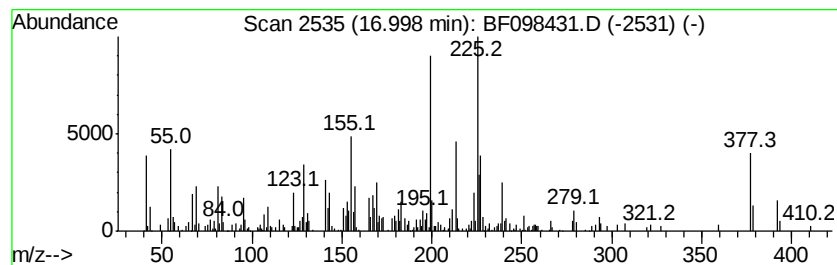
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown17.00 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.00	5.97 ng	355842	Perylene-d12	15.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Heptanoic acid,methyl(tetramet...	228	C12H24O2Si	1000217-03-6	22
2		2,7-Dimethyl-1,2,4-triaza-spiro[...	199	C9H17N3S	1000275-07-0	15
3		2,3,7-Trimethyl-1H-imidazo(1,2-a...	199	C12H13N3	1000147-53-1	15
4		4-Methyl-thieno(2,3-b)quinoline	199	C12H9NS	059280-92-1	15
5		2-Thia-3,5,6-triazaacenanthrylen...	225	C12H7N3S	209617-42-5	11



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF091117\
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ClientSampleId :
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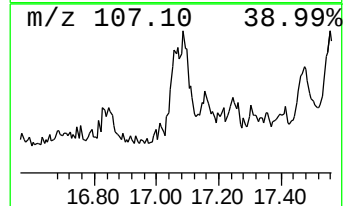
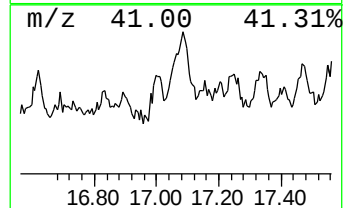
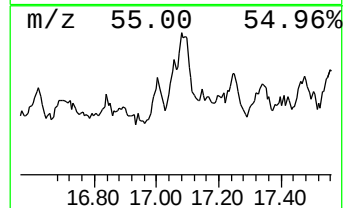
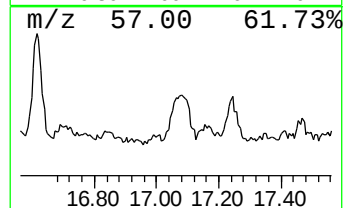
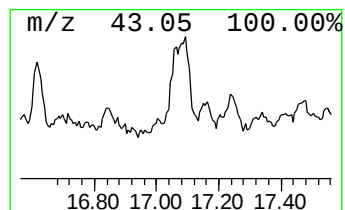
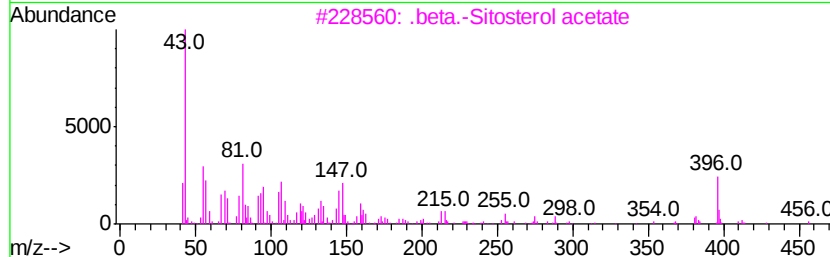
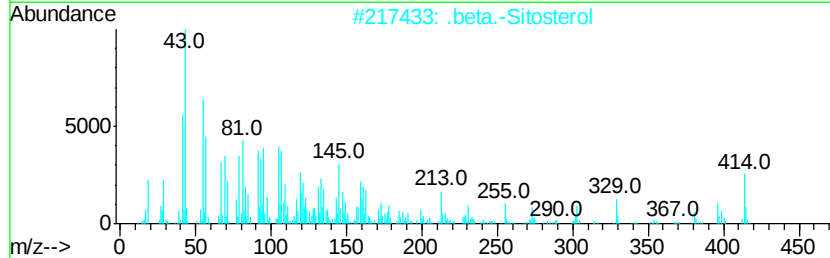
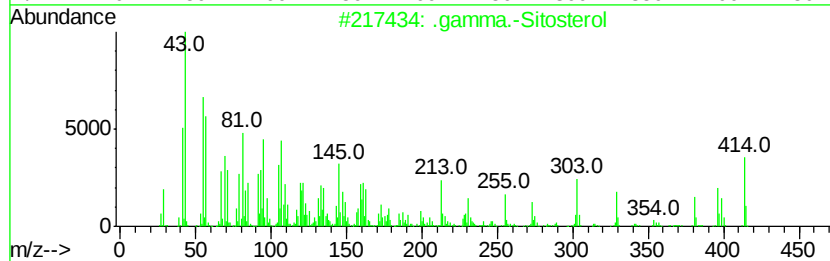
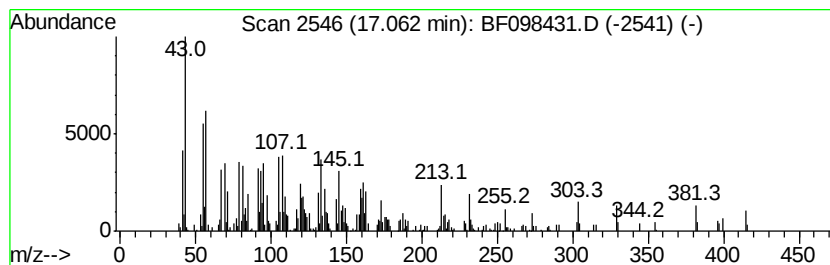
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 .gamma.-Sitosterol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.06	3.55 ng	211305	Perylene-d12	15.22

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.gamma.-Sitosterol	414	C29H50O	000083-47-6	90
2			.beta.-Sitosterol	414	C29H50O	000083-46-5	64
3			.beta.-Sitosterol acetate	456	C31H52O2	000915-05-9	32
4			Stigmastan-3,5-diene	396	C29H48	1000214-16-4	17
5			Methyl 12-acetyl-7-desoxycholate	448	C27H44O5	055547-48-3	17



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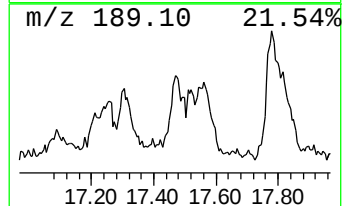
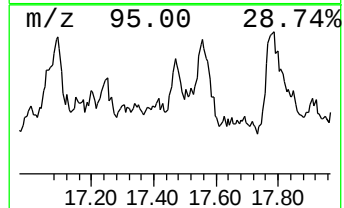
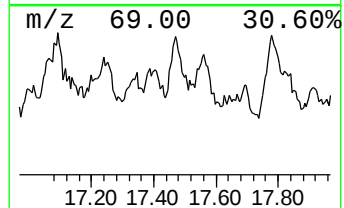
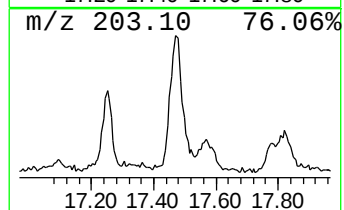
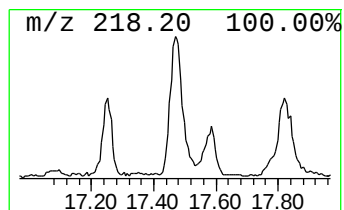
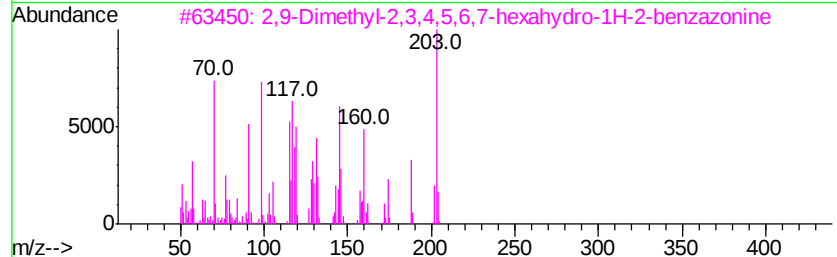
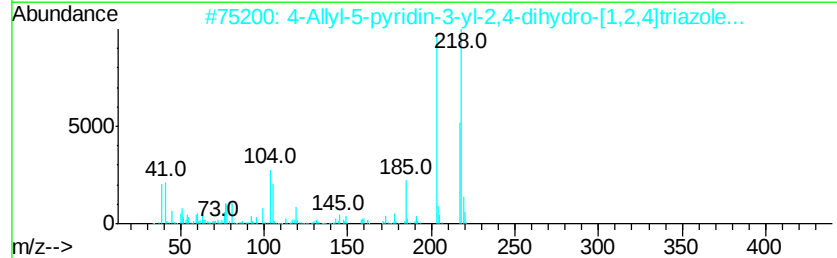
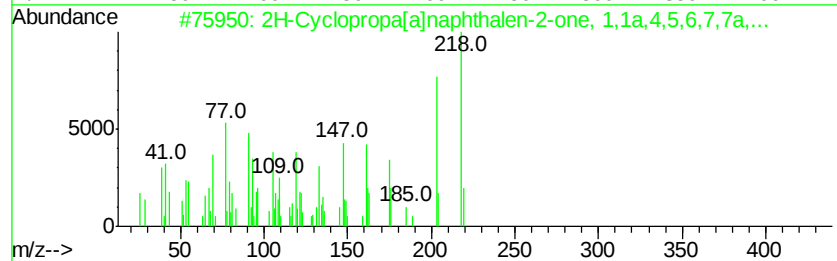
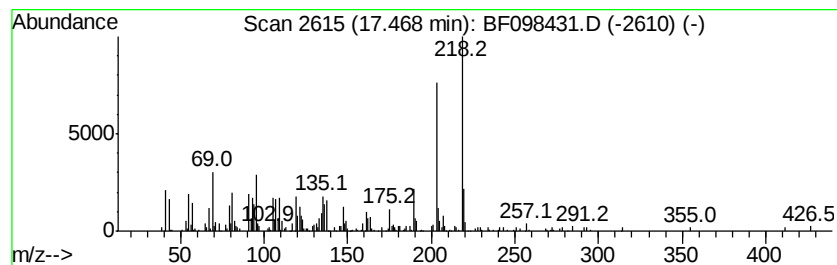
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 2H-Cyclopropa[a]naphthalen-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.47	8.93 ng	532179	Perylene-d12	15.22
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2H-Cyclopropa[a]naphthalen-2-one...	218 C15H22O	006831-17-0	76
2	4-Allyl-5-pyridin-3-yl-2,4-dihyd...	218 C10H10N4S	1000300-40-5	62
3	2,9-Dimethyl-2,3,4,5,6,7-hexahyd...	203 C14H21N	077581-11-4	60
4	.beta.-Amyrin	426 C30H50O	000559-70-6	60
5	8-Amino-6,7-dimethoxy-4-methylqu...	218 C12H14N2O2	1000213-07-2	59



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Data File : BF098431.D
Acq On : 12 Sep 2017 4:43
Operator : SJ/JU
Sample : I5138-17
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EME-WM-TS003-01

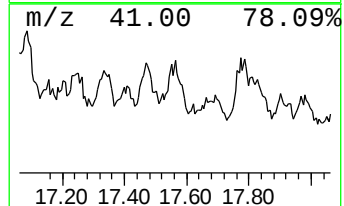
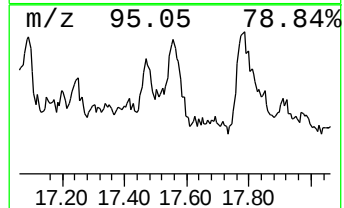
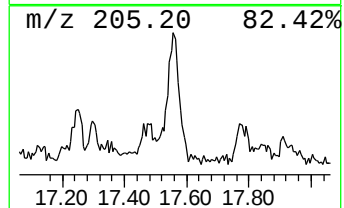
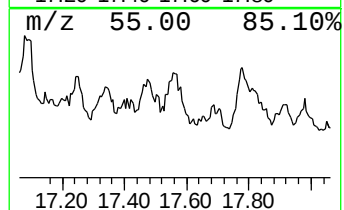
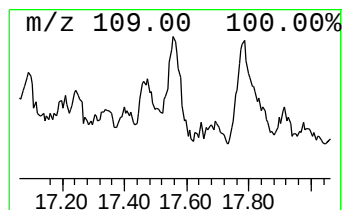
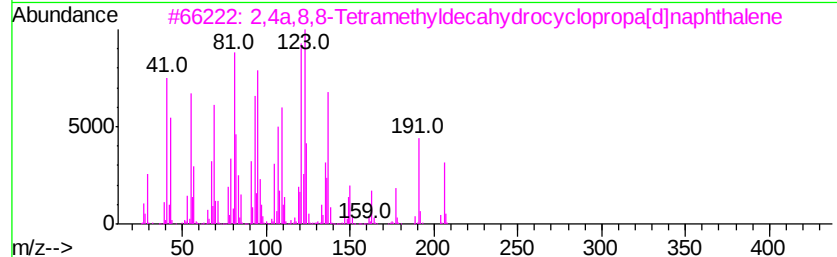
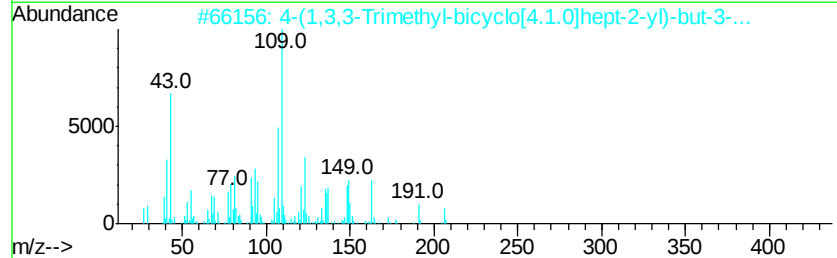
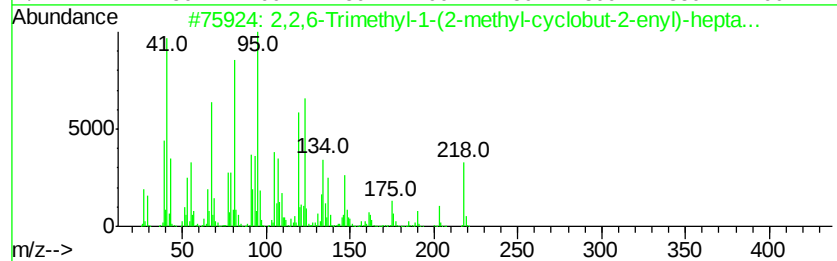
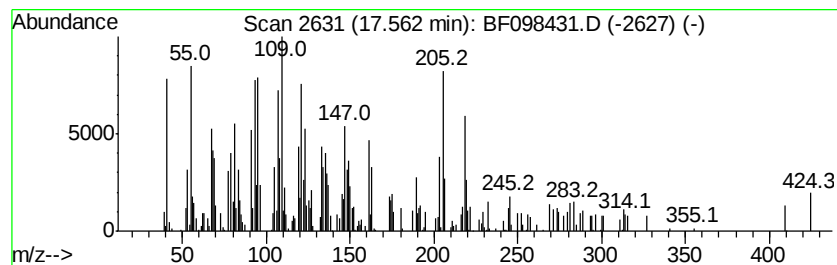
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 2,2,6-Trimethyl-1-(2-methyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.56	9.75 ng	581262	Perylene-d12	15.22

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2,6-Trimethyl-1-(2-methyl-cycl...	218	C15H22O	1000188-72-8	68		
2	4-(1,3,3-Trimethyl-bicyclo[4.1.0...	206	C14H22O	077143-31-8	40		
3	2,4a,8,8-Tetramethyldecahydrocyc...	206	C15H26	074022-04-1	38		
4	2(1H)Naphthalenone, 3,5,6,7,8,8a...	218	C15H22O	1000188-66-5	35		
5	1,4-Methanoazulen-7-ol, decahydr...	222	C15H26O	018319-27-2	30		



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF091117\
Data File : BF098431.D
Acq On : 12 Sep 2017 4:43
Operator : SJ/JU
Sample : I5138-17
Misc :
ALS Vial : 21 Sample Multiplier: 1

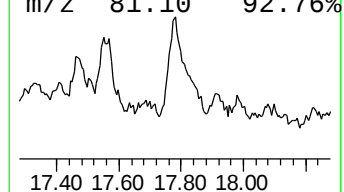
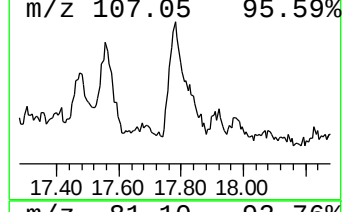
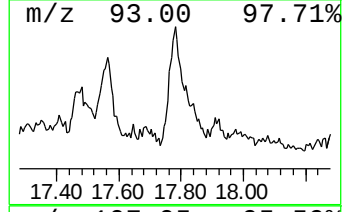
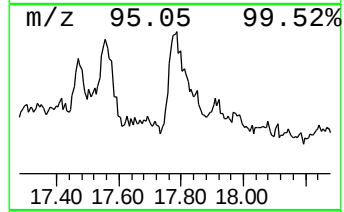
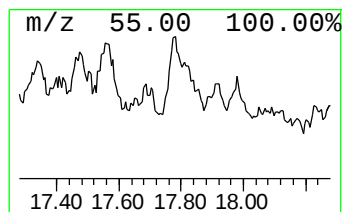
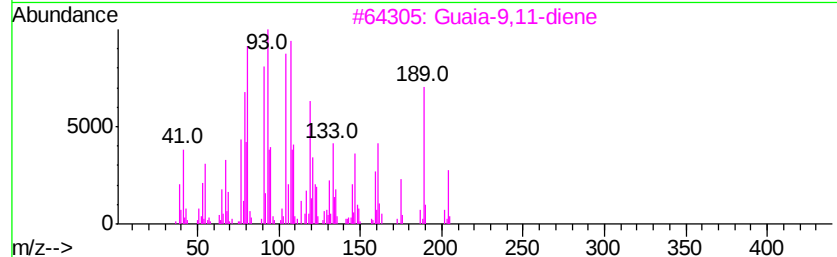
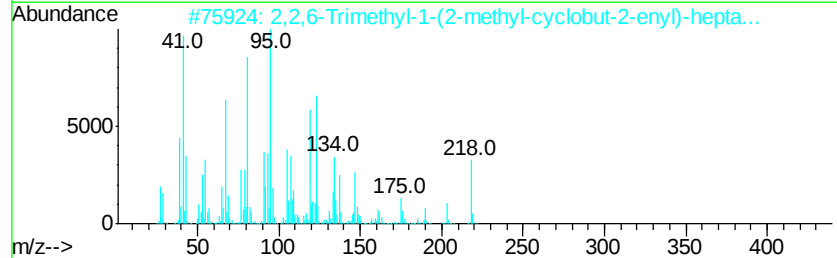
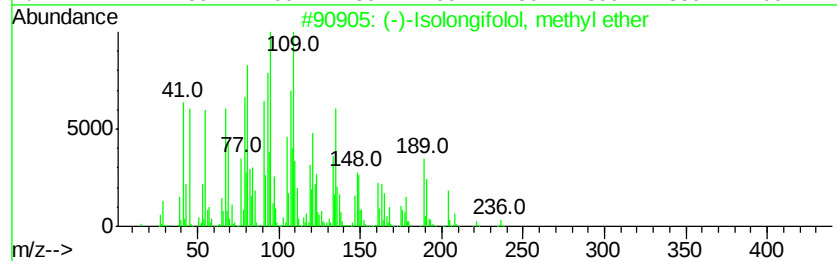
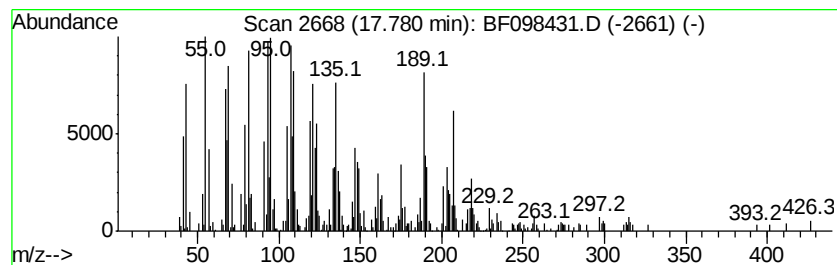
Instrument :
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ClientSampleId :
EME-WM-TS003-01

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF091117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown17.78 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.78	20.91 ng	1245790	Perylene-d12	15.22
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	(-)-Isolongifolol, methyl ether	236 C16H28O	1000333-80-8	48
2	2,2,6-Trimethyl-1-(2-methyl-cycl...	218 C15H22O	1000188-72-8	44
3	Guaia-9,11-diene	204 C15H24	1000374-19-8	42
4	Caparratriene	206 C15H26	1000374-08-7	41
5	1-Naphthalenecarboxylic acid, de...	316 C21H32O2	001235-39-8	38



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF091117\
Data File : BF098431.D
Acq On : 12 Sep 2017 4:43
Operator : SJ/JU
Sample : I5138-17
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EME-WM-TS003-01

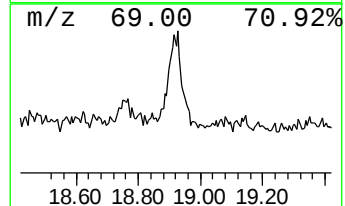
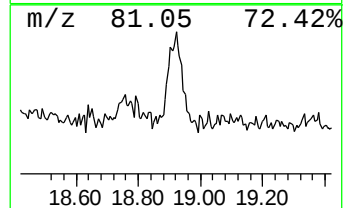
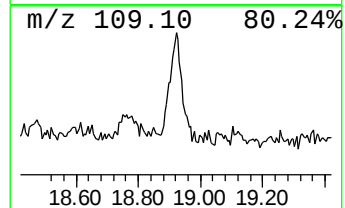
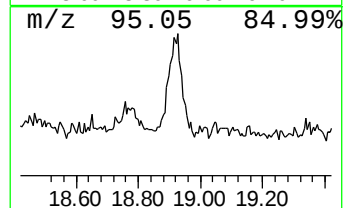
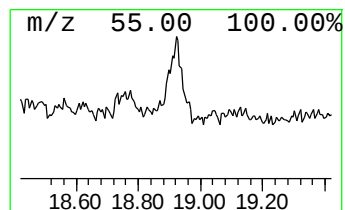
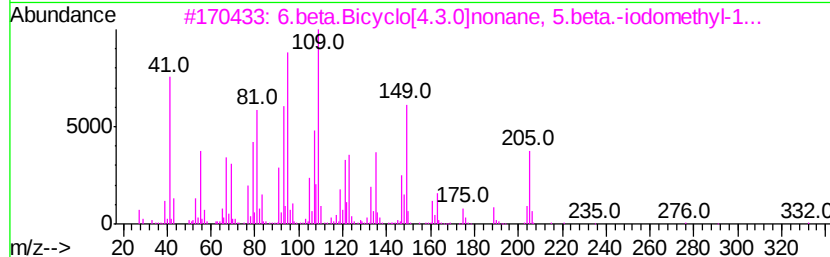
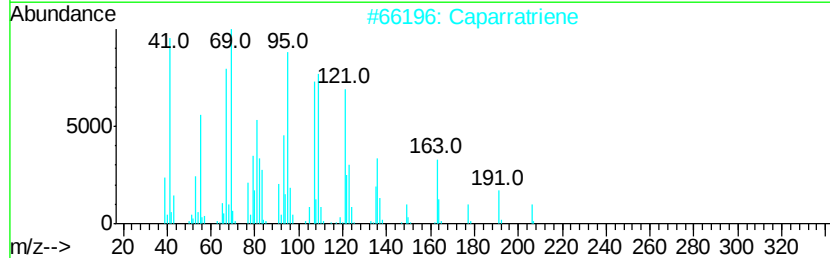
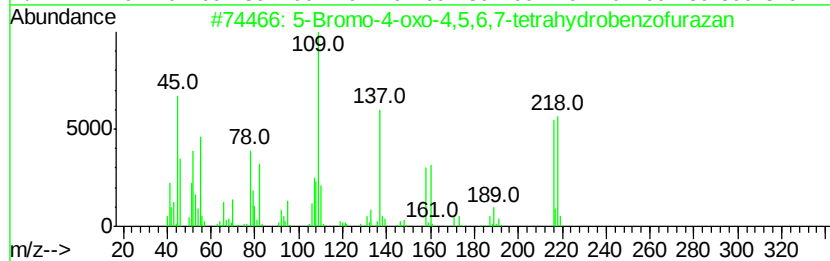
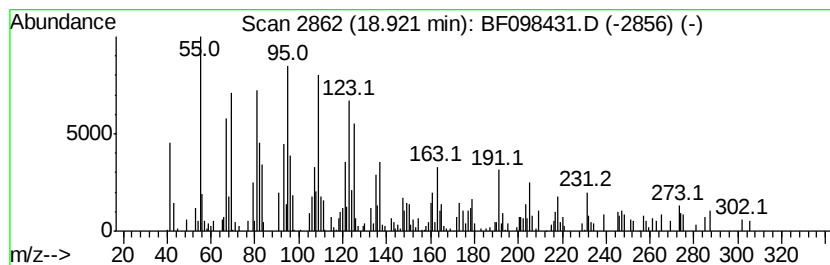
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 5-Bromo-4-oxo-4,5,6,7-tetra... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.92	7.43 ng	442905	Perylene-d12	15.22

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	68
2			Caparratriene	206	C15H26	1000374-08-7	49
3			6.beta.Bicyclo[4.3.0]nonane, 5.b...	332	C15H25I	1000195-85-9	38
4			Methyl 5-ethylcyclopent-1-ene-1-...	154	C9H14O2	182683-18-7	38
5			(2Z,4E)-3,7,11-Trimethyl-2,4,10-...	206	C15H26	172549-29-0	38



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF091117\
 Data File : BF098431.D
 Acq On : 12 Sep 2017 4:43
 Operator : SJ/JU
 Sample : I5138-17
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EME-WM-TS003-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF091117.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...	4.87	9.8	ng	741347	1	6.67	1508590	20.0
unknown6.45	6.45	89.7	ng	6763360	1	6.67	1508590	20.0
Benzene, 1,1'-[1-...	10.96	2.6	ng	218466	4	11.18	1680450	20.0
10,18-Bisnorabiet...	12.43	2.2	ng	181492	4	11.18	1680450	20.0
Z-5-Nonadecene	13.66	8.0	ng	501029	5	13.81	1247240	20.0
Octatriacontyl pe...	14.29	4.9	ng	306849	5	13.81	1247240	20.0
Heptacosane	14.89	14.9	ng	885276	6	15.22	1191790	20.0
Hexadecane	16.62	4.3	ng	253880	6	15.22	1191790	20.0
unknown17.00	17.00	6.0	ng	355842	6	15.22	1191790	20.0
.gamma.-Sitosterol	17.06	3.5	ng	211305	6	15.22	1191790	20.0
2H-Cyclopropa[a]n...	17.47	8.9	ng	532179	6	15.22	1191790	20.0
2,2,6-Trimethyl-1...	17.56	9.8	ng	581262	6	15.22	1191790	20.0
unknown17.78	17.78	20.9	ng	1245790	6	15.22	1191790	20.0
5-Bromo-4-oxo-4,5...	18.92	7.4	ng	442905	6	15.22	1191790	20.0