

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051915\
 Data File : BF079362.D
 Acq On : 20 May 2015 3:03
 Operator : TP/IZ
 Sample : G2281-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-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-BF043015.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.255	4	6	9	rVB	1672174	2328149	31.76%	5.792%
2	1.392	16	18	29	rVB3	91047	267349	3.65%	0.665%
3	1.552	29	32	39	rVV	120261	220025	3.00%	0.547%
4	1.644	39	40	48	rVV	381243	607893	8.29%	1.512%
5	1.758	48	50	93	rVB	3288006	7331353	100.00%	18.239%
6	4.878	320	323	331	rBV	222428	289463	3.95%	0.720%
7	5.416	367	370	382	rVB	732505	904762	12.34%	2.251%
8	6.696	480	482	485	rBV	771848	918185	12.52%	2.284%
9	6.833	492	494	497	rBV	769398	979196	13.36%	2.436%
10	7.130	517	520	522	rBV	223278	266402	3.63%	0.663%
11	7.313	533	536	538	rBV	816087	772052	10.53%	1.921%
12	7.816	578	580	583	rBV	496439	570889	7.79%	1.420%
13	8.707	655	658	662	rBV2	270710	486632	6.64%	1.211%
14	10.045	773	775	778	rBV	1116610	1218789	16.62%	3.032%
15	10.559	817	820	822	rBV	72598	87262	1.19%	0.217%
16	10.696	830	832	835	rVB	175338	162625	2.22%	0.405%
17	10.868	846	847	849	rVV	372720	409431	5.58%	1.019%
18	11.851	930	933	936	rBV	810747	840446	11.46%	2.091%
19	12.708	1006	1008	1010	rBV	368240	548961	7.49%	1.366%
20	12.742	1010	1011	1014	rVV	688139	541868	7.39%	1.348%
21	12.811	1014	1017	1019	rVV	246146	320522	4.37%	0.797%
22	13.348	1061	1064	1065	rBV	69825	104840	1.43%	0.261%
23	13.371	1065	1066	1070	rVV	104818	157004	2.14%	0.391%
24	13.439	1070	1072	1073	rVV	99721	135900	1.85%	0.338%
25	13.474	1073	1075	1077	rVV	211478	268105	3.66%	0.667%
26	13.714	1095	1096	1101	rVB2	69020	135665	1.85%	0.338%
27	14.022	1120	1123	1125	rBV2	75616	128740	1.76%	0.320%
28	14.068	1125	1127	1129	rVV	55207	95960	1.31%	0.239%
29	14.125	1131	1132	1135	rVB2	61268	94750	1.29%	0.236%
30	14.217	1138	1140	1142	rBV	1230716	1256753	17.14%	3.126%
31	14.297	1145	1147	1148	rBV	73633	81804	1.12%	0.204%
32	14.331	1148	1150	1153	rVB	130922	171112	2.33%	0.426%
33	14.502	1161	1165	1167	rBV	944894	1354208	18.47%	3.369%
34	14.662	1177	1179	1180	rBV	60555	81485	1.11%	0.203%

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 Sampling : 1
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If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.708	1180	1183	1186	rBV	1517671	1544318	21.06%	3.842%
36	14.777	1186	1189	1192	rBV2	67642	128725	1.76%	0.320%
37	14.891	1195	1199	1200	rBV3	85721	178763	2.44%	0.445%
38	14.914	1200	1201	1203	rVB	180849	169145	2.31%	0.421%
39	15.005	1206	1209	1210	rVV	97852	142696	1.95%	0.355%
40	15.039	1210	1212	1215	rVV	144703	189626	2.59%	0.472%
41	15.154	1220	1222	1224	rVV2	86678	94712	1.29%	0.236%
42	15.188	1224	1225	1228	rVV	67270	89380	1.22%	0.222%
43	15.451	1240	1248	1250	rBV4	45146	182734	2.49%	0.455%
44	15.542	1252	1256	1260	rVB2	70623	144814	1.98%	0.360%
45	15.680	1265	1268	1270	rBV	113277	154179	2.10%	0.384%
46	15.737	1270	1273	1275	rBV2	81988	179768	2.45%	0.447%
47	15.805	1278	1279	1282	rVB	74731	86710	1.18%	0.216%
48	15.885	1283	1286	1289	rBV3	188169	333405	4.55%	0.829%
49	15.988	1292	1295	1297	rBV2	890672	1549851	21.14%	3.856%
50	16.034	1297	1299	1301	rVV	520580	707298	9.65%	1.760%
51	16.125	1304	1307	1308	rBV	94104	111041	1.51%	0.276%
52	16.171	1310	1311	1313	rVB	65436	78968	1.08%	0.196%
53	16.445	1333	1335	1336	rVV	65996	79218	1.08%	0.197%
54	16.491	1336	1339	1341	rVV	152241	233359	3.18%	0.581%
55	16.525	1341	1342	1345	rVV	64580	111729	1.52%	0.278%
56	16.605	1347	1349	1350	rVV	86659	97777	1.33%	0.243%
57	16.640	1350	1352	1353	rVV2	70711	96311	1.31%	0.240%
58	16.685	1355	1356	1359	rVB2	132175	165188	2.25%	0.411%
59	16.880	1371	1373	1377	rBV5	39745	102697	1.40%	0.255%
60	17.074	1387	1390	1393	rBV3	50859	143483	1.96%	0.357%
61	17.234	1401	1404	1405	rBV2	205722	267043	3.64%	0.664%
62	17.268	1405	1407	1412	rVB	696465	1410872	19.24%	3.510%
63	17.383	1414	1417	1419	rVB	216038	273567	3.73%	0.681%
64	17.474	1420	1425	1429	rBV3	333475	898182	12.25%	2.234%
65	17.577	1432	1434	1437	rBV	466581	533593	7.28%	1.327%
66	17.634	1437	1439	1442	rVB	611029	839844	11.46%	2.089%
67	17.703	1442	1445	1446	rBV	489024	572477	7.81%	1.424%
68	18.011	1470	1472	1479	rVB	149718	316088	4.31%	0.786%
69	18.240	1489	1492	1495	rBV2	424076	648784	8.85%	1.614%
70	18.286	1495	1496	1500	rVB3	72544	121915	1.66%	0.303%
71	18.571	1518	1521	1524	rVB4	92426	218254	2.98%	0.543%

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Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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72	18.926	1549	1552	1554	rVB2	64752	132059	1.80%	0.329%
73	18.983	1554	1557	1561	rVB2	83528	150655	2.05%	0.375%
74	19.074	1562	1565	1575	rVB2	359733	929889	12.68%	2.313%
75	19.257	1577	1581	1584	rBV2	199106	557635	7.61%	1.387%
76	19.474	1597	1600	1606	rVB	267939	564340	7.70%	1.404%
77	19.646	1611	1615	1617	rBV3	126341	308912	4.21%	0.768%
78	20.571	1693	1696	1702	rVB5	75636	218241	2.98%	0.543%

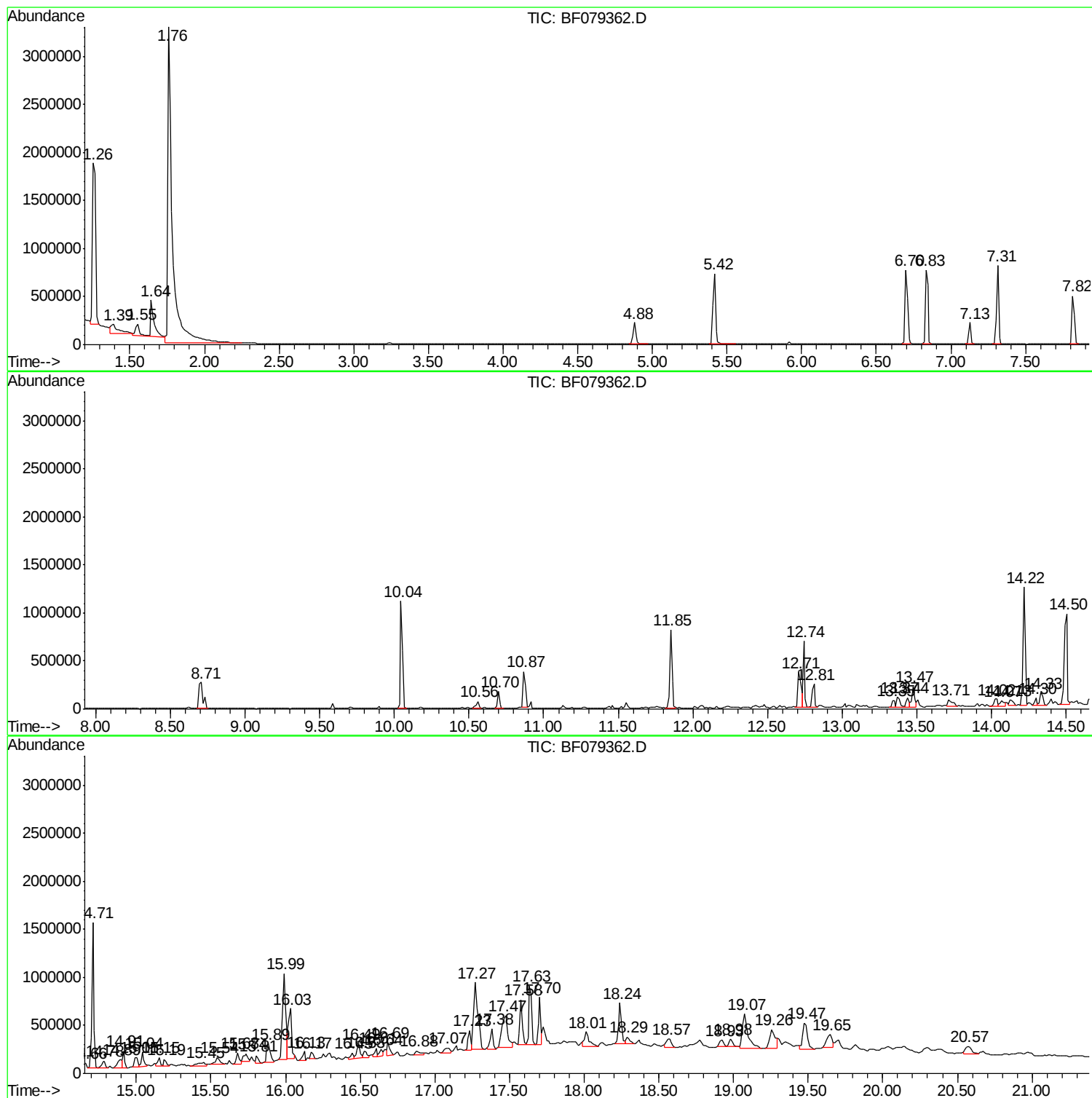
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Instrument :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Operator : TP/IZ
Sample : G2281-01
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Instrument :
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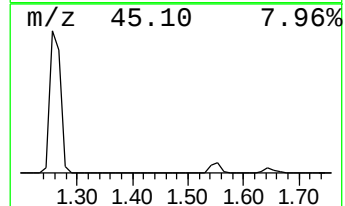
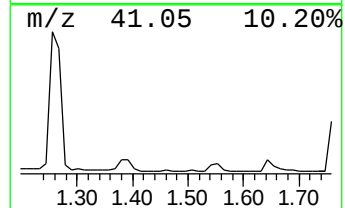
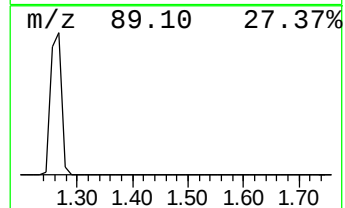
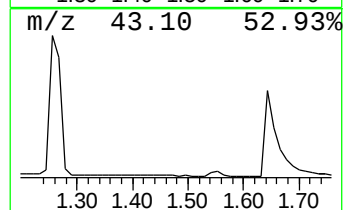
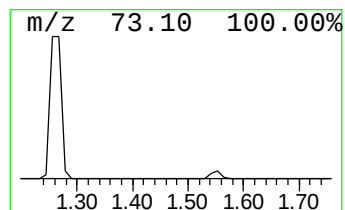
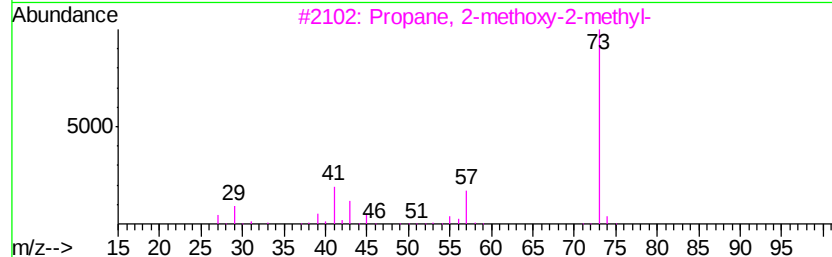
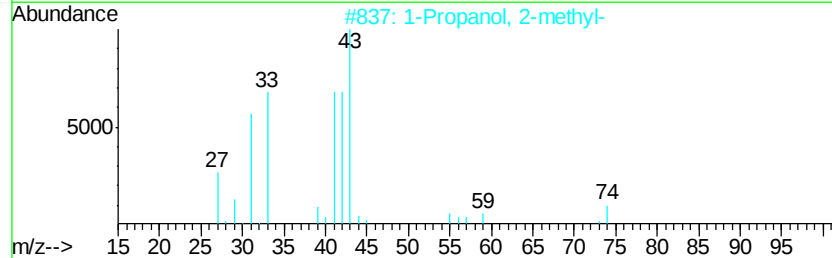
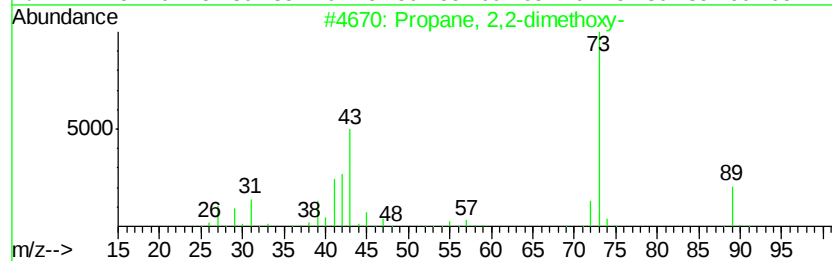
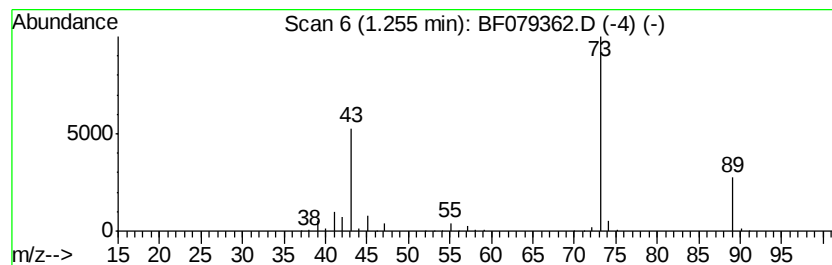
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.26	174.79 ng	2328150	1,4-Dichlorobenzene-d4	7.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	72
2		1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	9
3		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
4		2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9
5		1-Hexen-4-ol, 1-chloro-3,5-dimet...	162	C8H15ClO	100244-09-5	9



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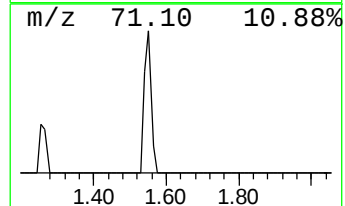
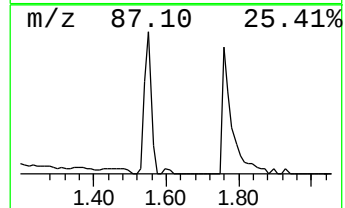
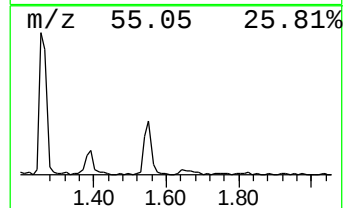
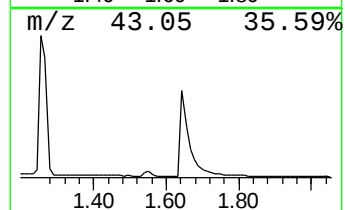
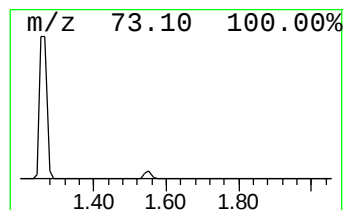
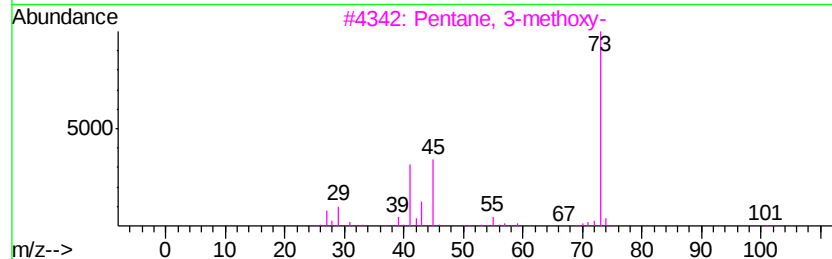
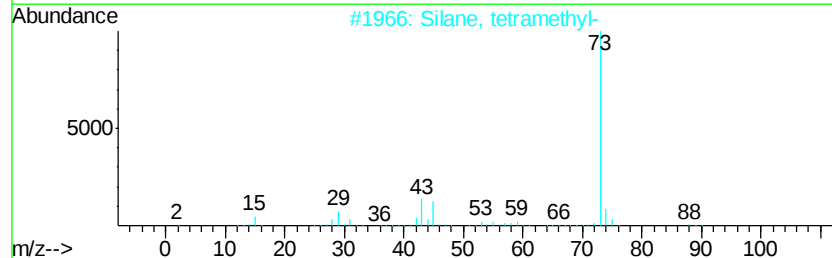
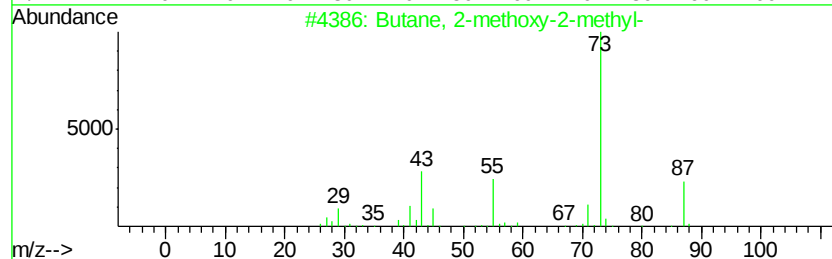
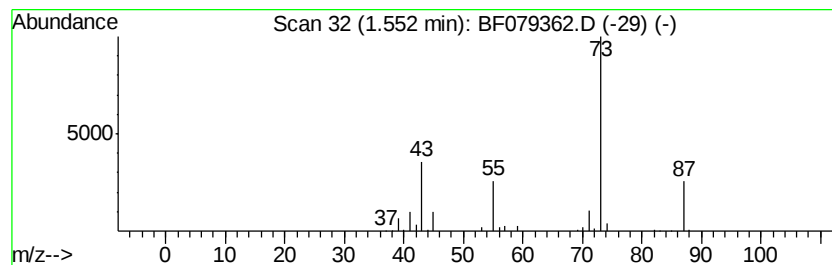
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.55	16.52 ng	220025	1,4-Dichlorobenzene-d4	7.13

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		Silane, tetramethyl-	88	C4H12Si	000075-76-3	43
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4		3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	16
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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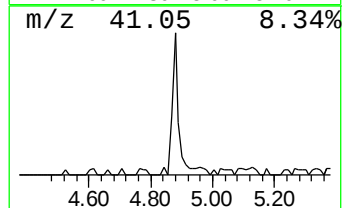
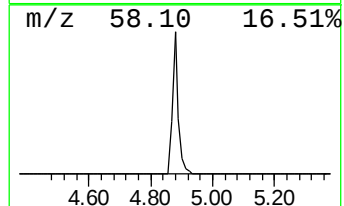
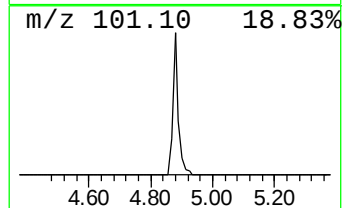
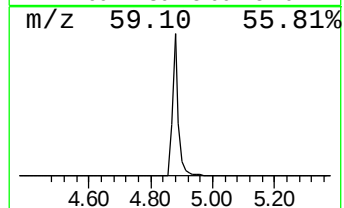
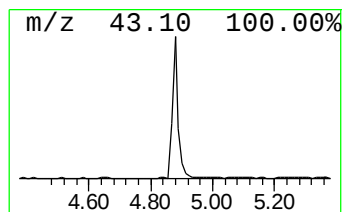
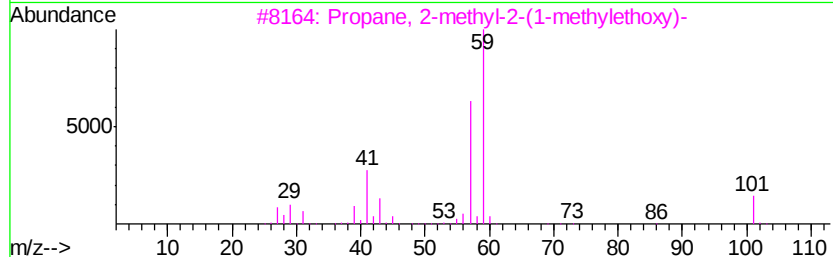
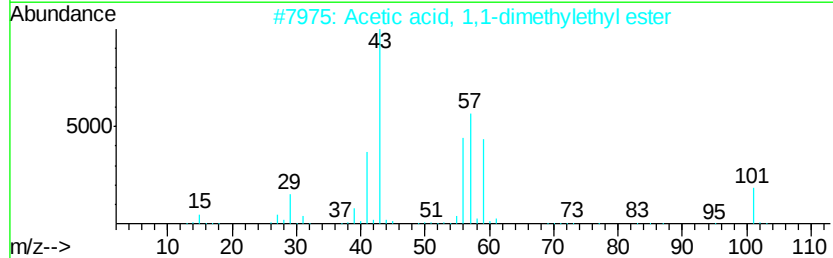
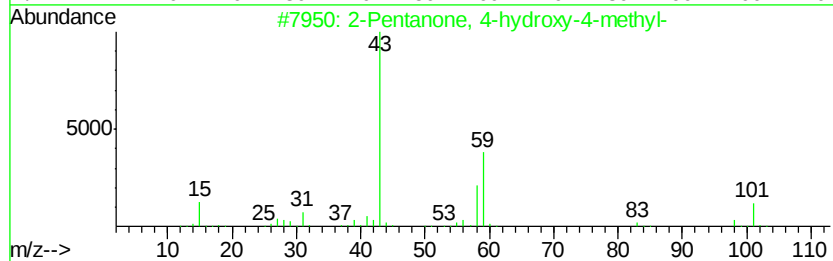
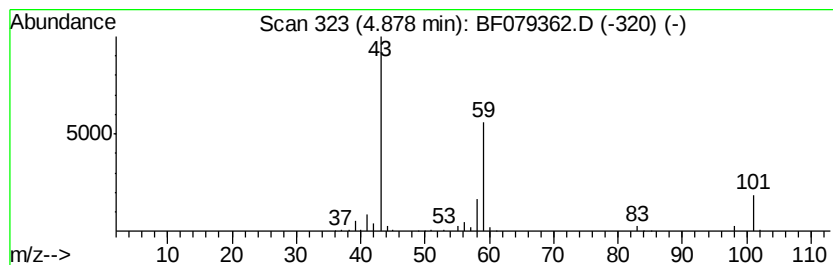
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.88	21.73 ng	289463	1,4-Dichlorobenzene-d4	7.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16



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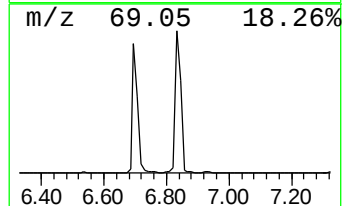
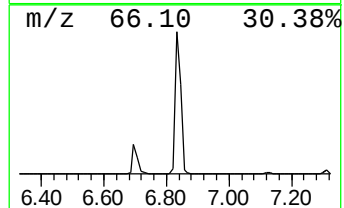
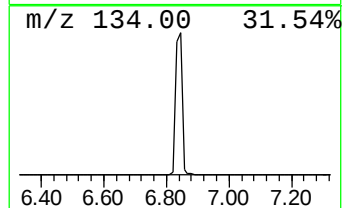
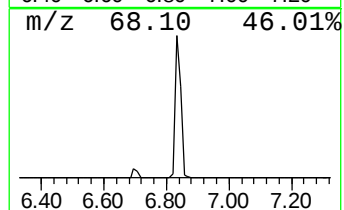
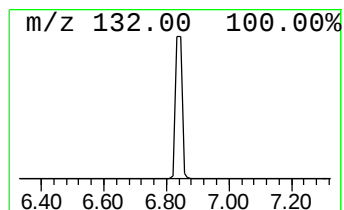
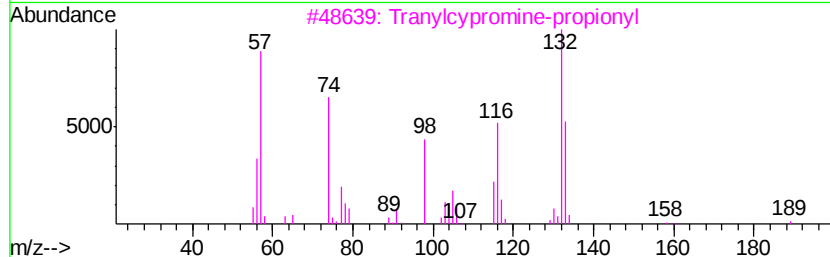
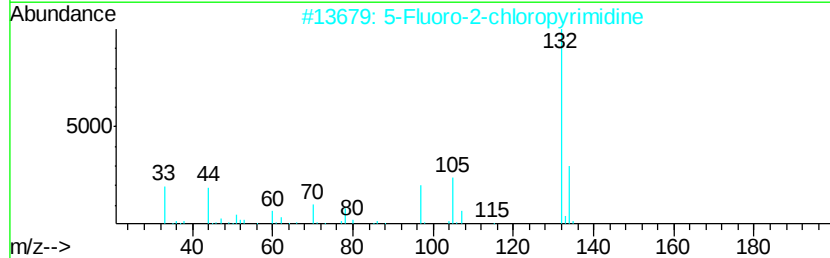
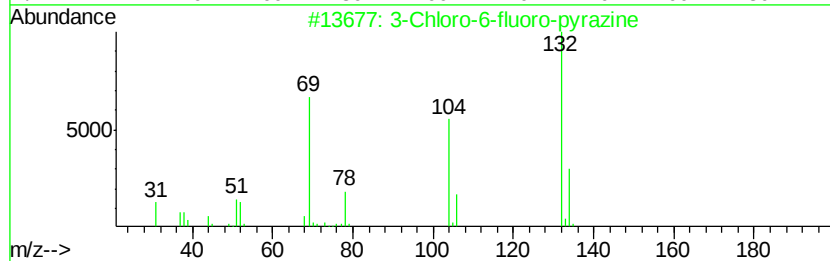
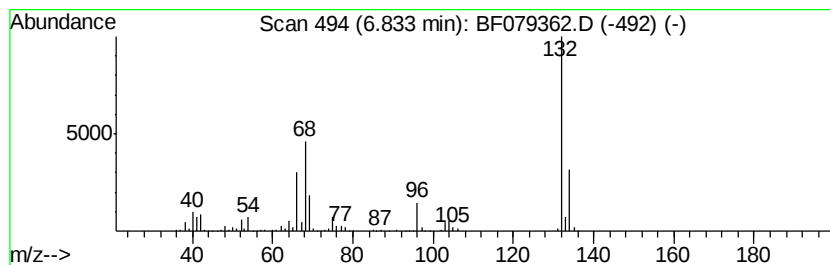
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown6.83 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.83	73.51 ng	979196	1,4-Dichlorobenzene-d4	7.13

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	12
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		5-Aminoindole	132	C8H8N2	005192-03-0	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051915\
Data File : BF079362.D
Acq On : 20 May 2015 3:03
Operator : TP/IZ
Sample : G2281-01
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-01

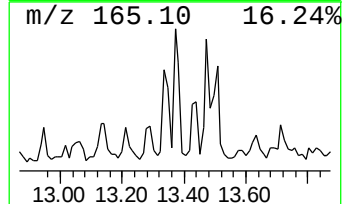
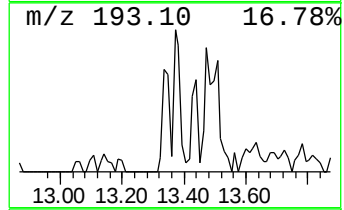
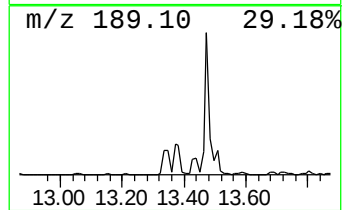
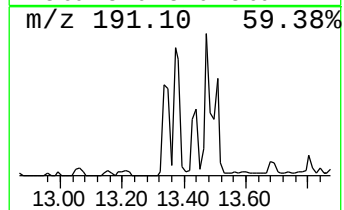
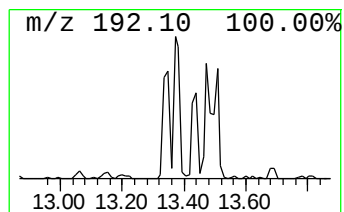
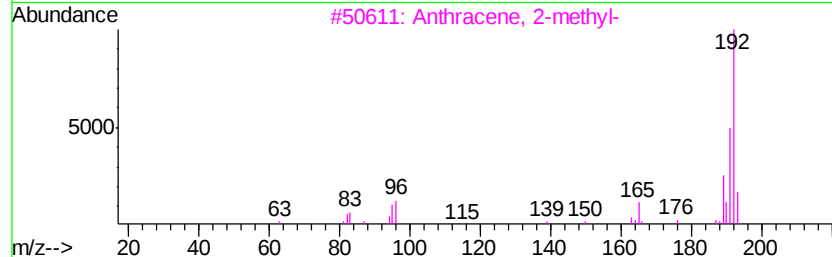
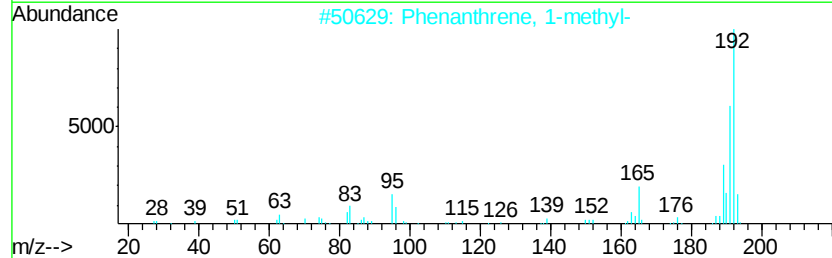
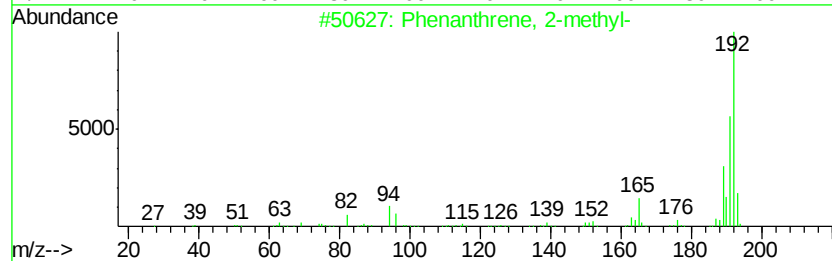
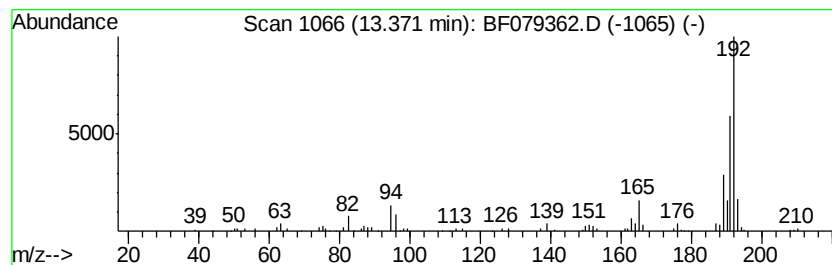
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenanthrene, 2-methyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.37	5.72 ng	157004	Phenanthrene-d10	12.71

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
4		1H-Cyclopropa[1,1]phenanthrene, 1a,...	192	C15H12	000949-41-7	95
5		Anthracene, 1-methyl-	192	C15H12	000610-48-0	94



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051915\
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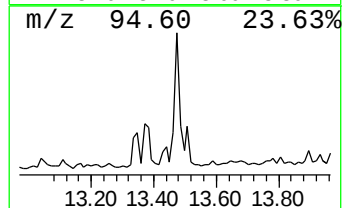
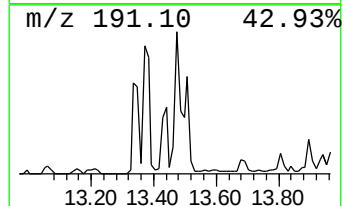
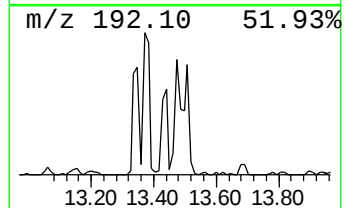
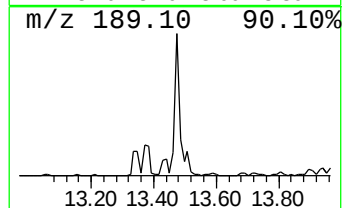
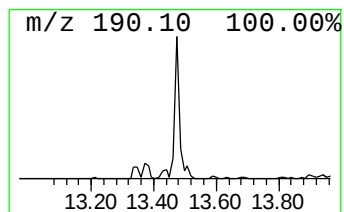
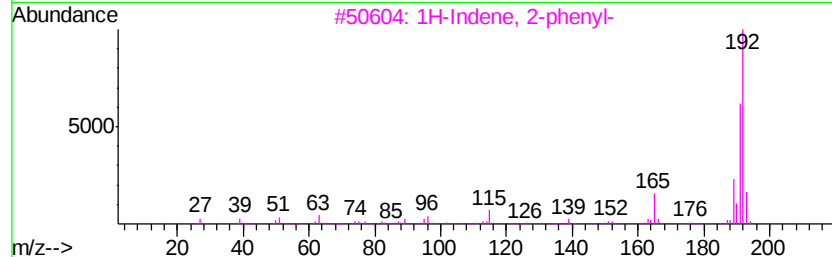
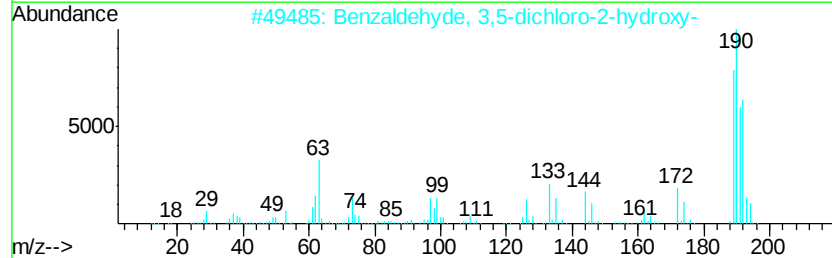
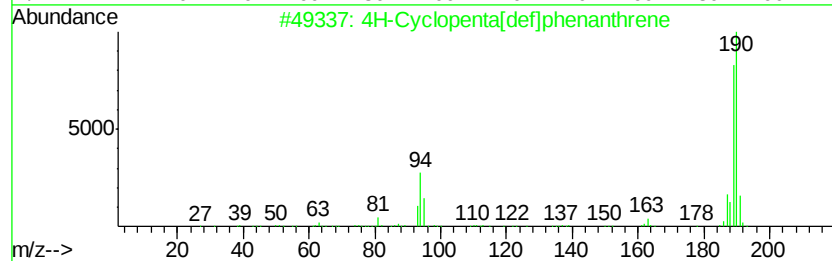
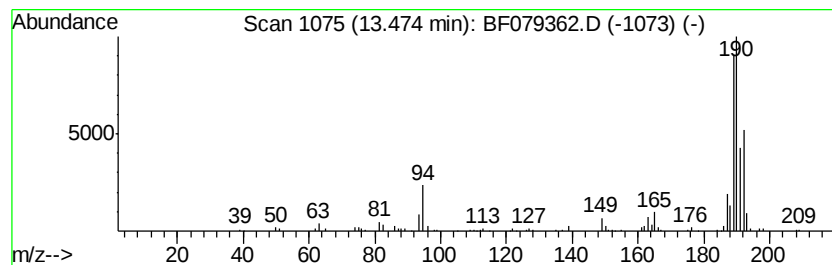
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 4H-Cyclopenta[def]phenanthrene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	9.77 ng	268105	Phenanthrene-d10	12.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2		Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	53
3		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	30
4		Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BCl03S	036155-87-0	27
5		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	25



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Sample : G2281-01
Misc :
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ClientSampleId :
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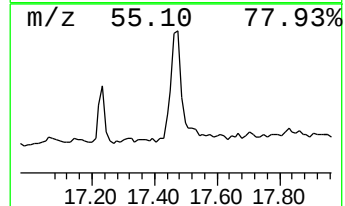
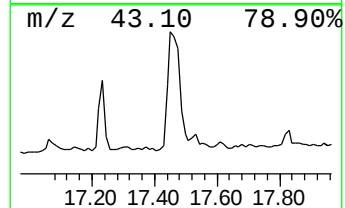
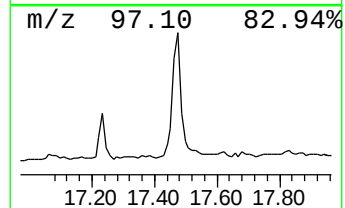
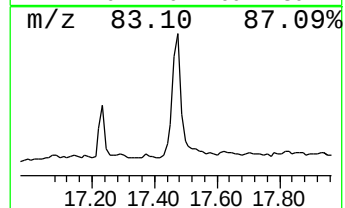
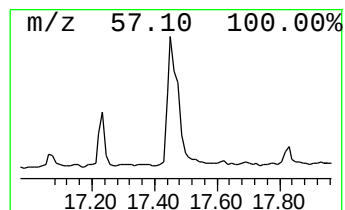
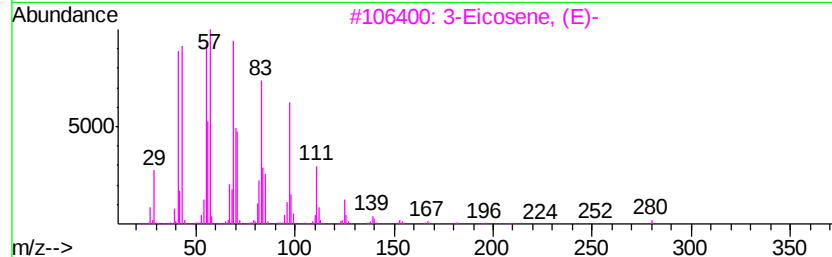
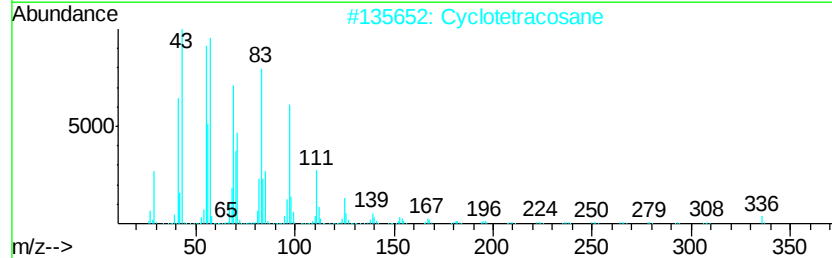
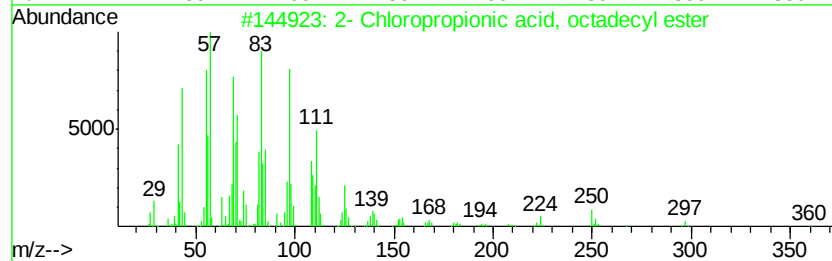
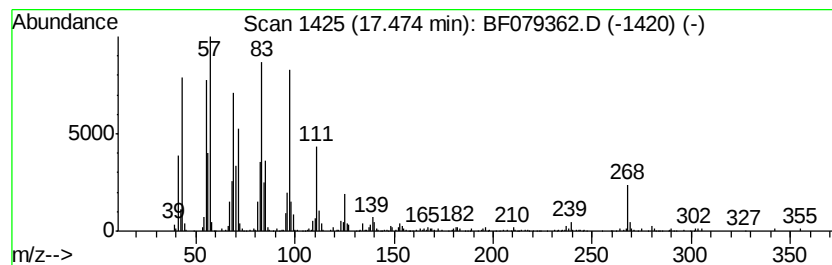
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 2- Chloropropionic acid, oc... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.47	31.38 ng	898182	Perylene-d12	17.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	91
2		Cyclotetracosane	336	C24H48	000297-03-0	91
3		3-Eicosene, (E)-	280	C20H40	074685-33-9	90
4		5-Eicosene, (E)-	280	C20H40	074685-30-6	89
5		10-Heneicosene (c,t)	294	C21H42	095008-11-0	83



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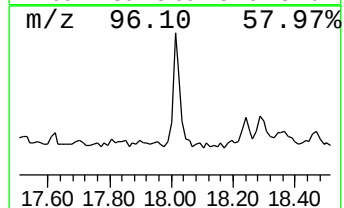
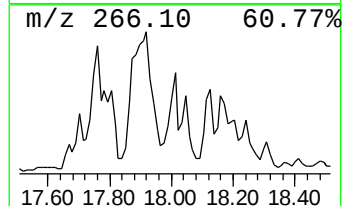
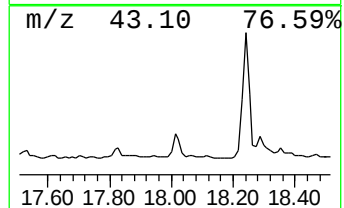
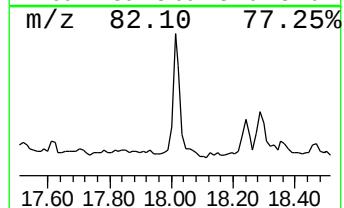
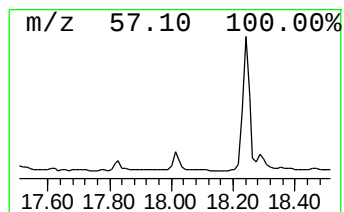
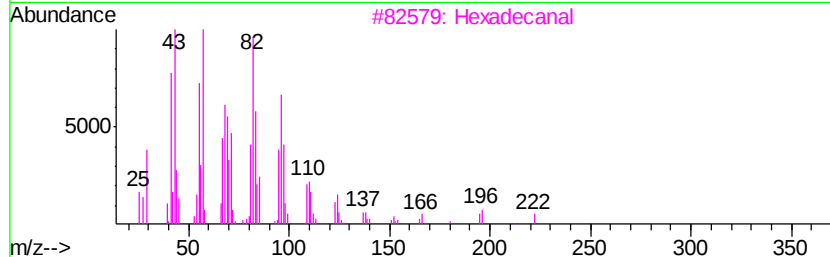
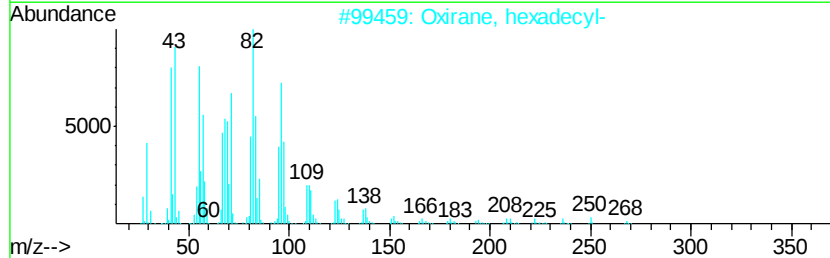
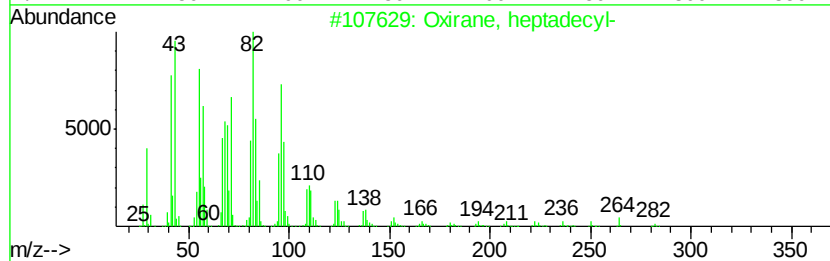
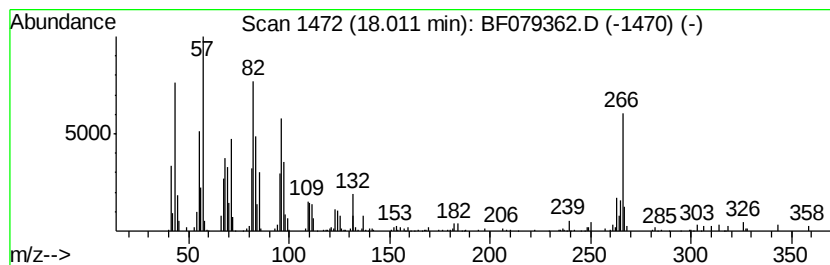
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Oxirane, heptadecyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.01	11.04 ng	316088	Perylene-d12	17.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	94
2		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	70
3		Hexadecanal	240	C16H32O	000629-80-1	70
4		Octadecanal	268	C18H36O	000638-66-4	70
5		1,19-Eicosadiene	278	C20H38	014811-95-1	58



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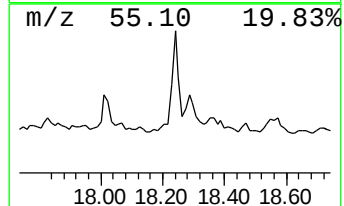
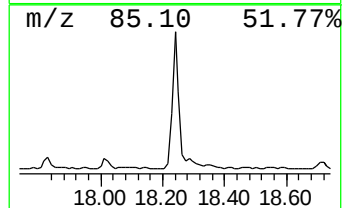
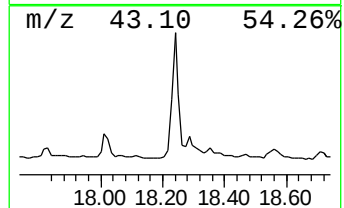
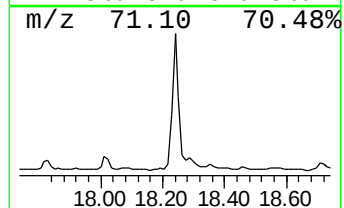
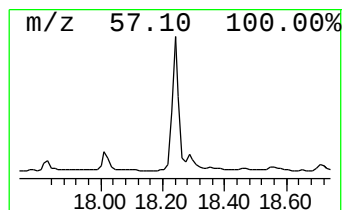
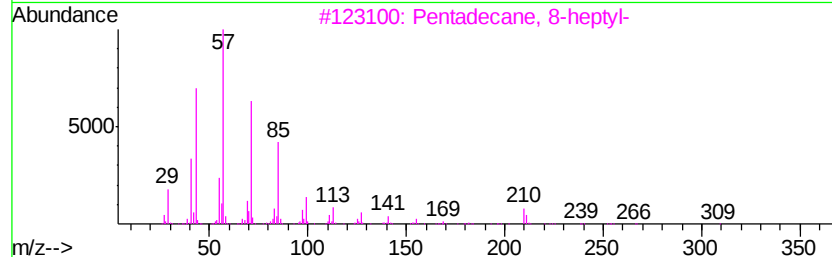
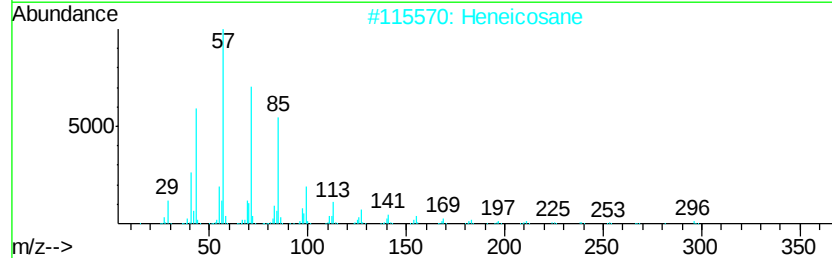
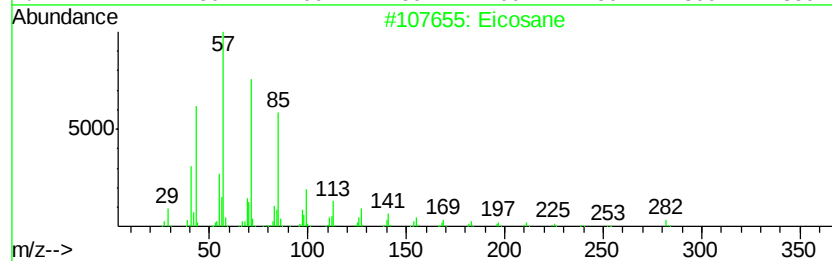
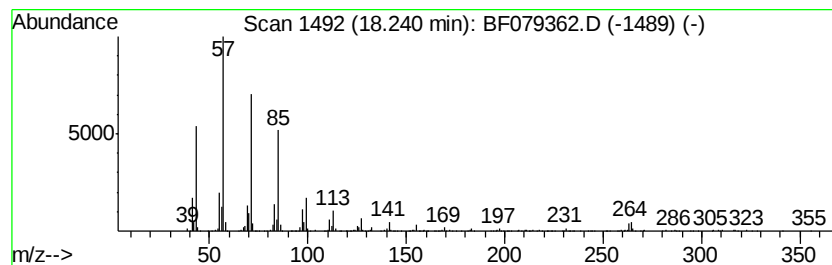
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Eicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.24	22.67 ng	648784	Perylene-d12	17.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	93
2		Heneicosane	296	C21H44	000629-94-7	90
3		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	87
4		Tetratriacontane	479	C34H70	014167-59-0	87
5		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	87



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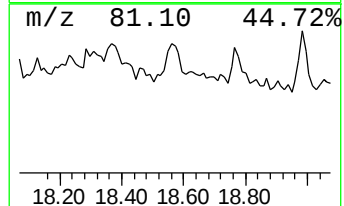
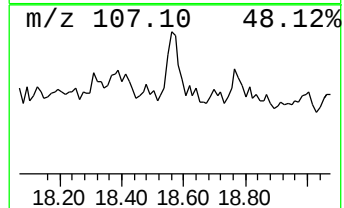
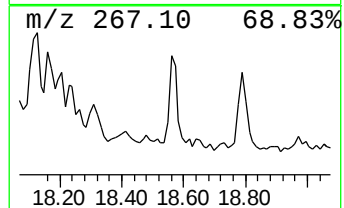
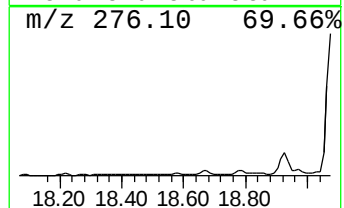
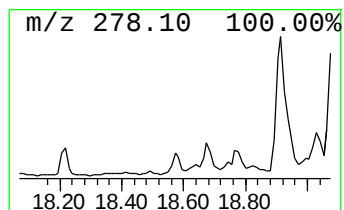
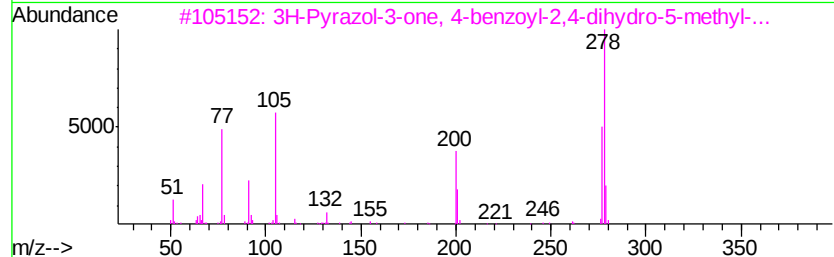
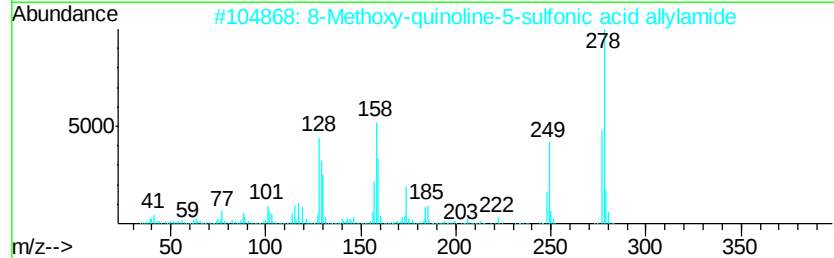
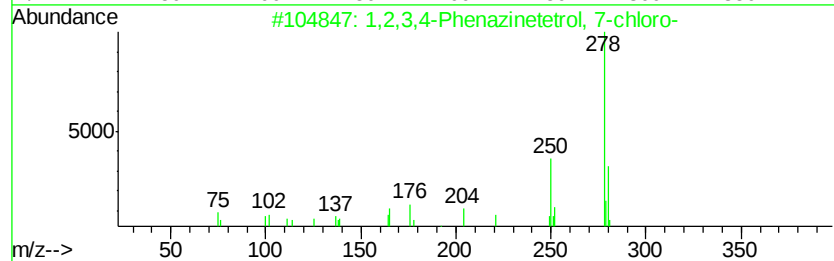
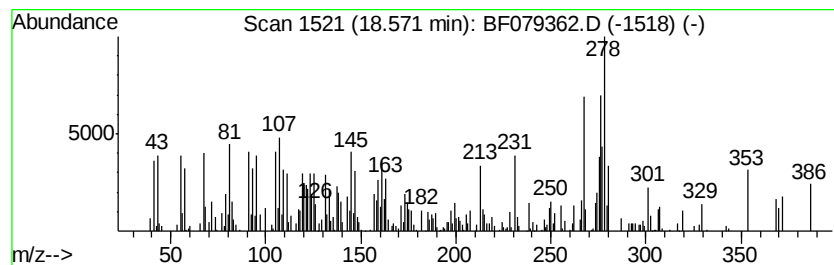
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 unknown18.57 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.57	7.62 ng	218254	Perylene-d12	17.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2,3,4-Phenazinetetrol, 7-chloro-	278	C12H7ClN2O4	023774-10-9	35
2		8-Methoxy-quinoline-5-sulfonic a...	278	C13H14N2O3S	1000274-64-1	25
3		3H-Pyrazol-3-one, 4-benzovl-2,4-...	278	C17H14N2O2	004551-69-3	22
4		2-Methyl-4,6-di(2-hydroxyphenyl)...	278	C17H14N2O2	1000070-55-8	22
5		Indeno[2,1-b]pyran, 2-(4-chlorop...	278	C18H11ClO	062096-34-8	18



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051915\
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ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-01

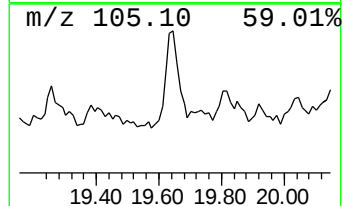
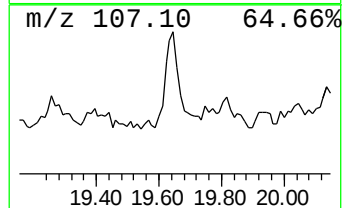
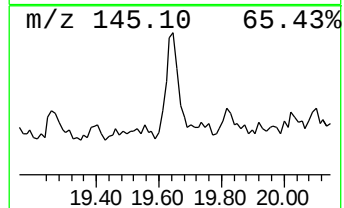
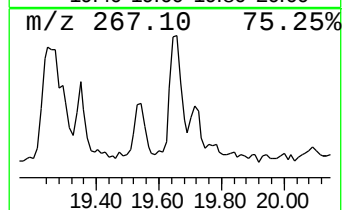
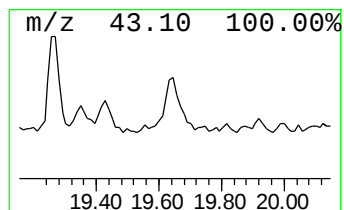
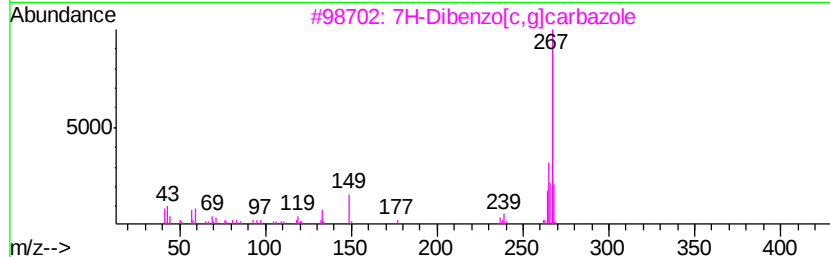
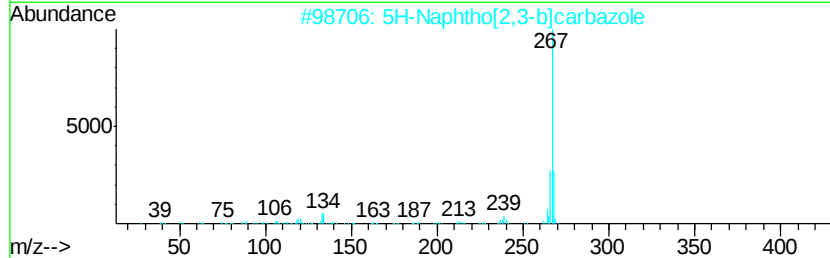
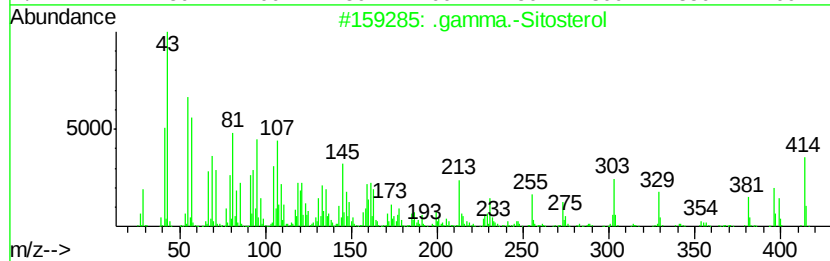
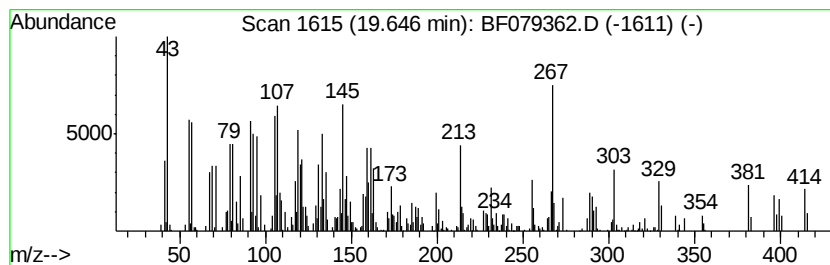
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 .gamma.-Sitosterol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.65	10.79 ng	308912	Perylene-d12	17.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.gamma.-Sitosterol	414	C29H50O	000083-47-6	90
2		5H-Naphtho[2,3-b]carbazole	267	C20H13N	000248-96-4	30
3		7H-Dibenzo[c,g]carbazole	267	C20H13N	000194-59-2	30
4		13H-Dibenzo[a,i]carbazole	267	C20H13N	000239-64-5	25
5		.beta.-Sitosterol	414	C29H50O	000083-46-5	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051915\
Data File : BF079362.D
Acq On : 20 May 2015 3:03
Operator : TP/IZ
Sample : G2281-01
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-01

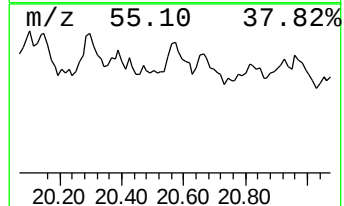
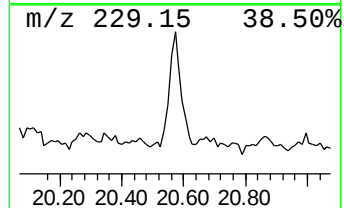
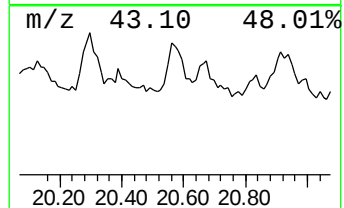
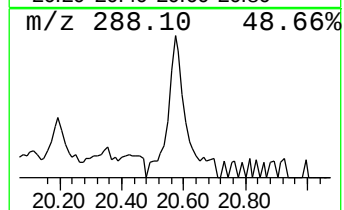
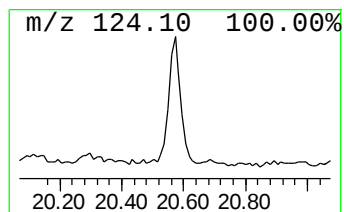
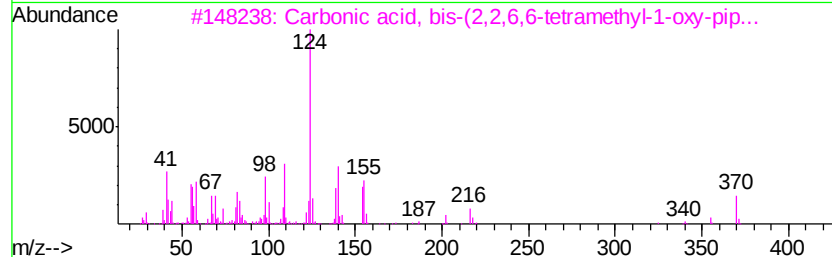
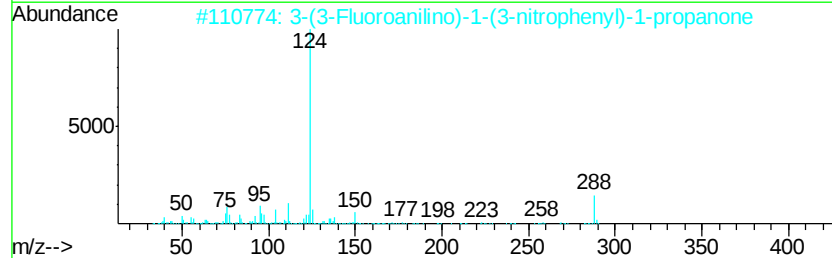
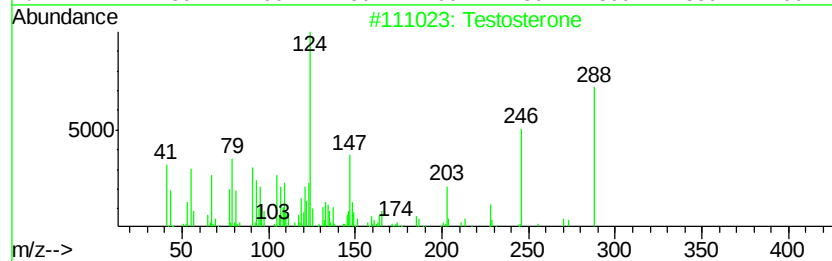
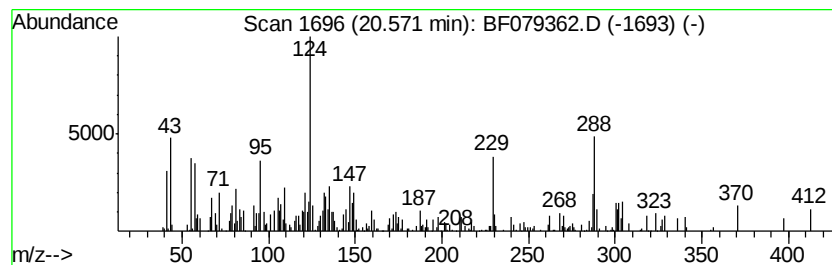
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Testosterone Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.57	7.62 ng	218241	Perylene-d12	17.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Testosterone	288	C19H28O2	000058-22-0	62
2		3-(3-Fluoroanilino)-1-(3-nitroph...	288	C15H13FN2O3	300726-64-1	43
3		Carbonic acid, bis-(2,2,6,6-tetr...	370	C19H34N2O5	1000188-97-0	22
4		Beta-(O-methoxyphenoxy)cinnamic ...	270	C16H14O4	1000243-81-2	20
5		Bicyclo[2.2.1]hept-5-ene-1,2-dic...	210	C11H14O4	1000194-11-4	18



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051915\
 Data File : BF079362.D
 Acq On : 20 May 2015 3:03
 Operator : TP/IZ
 Sample : G2281-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-01

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 2,2-dime...	1.26	174.8	ng	2328150	1	7.13	266402	20.0
Butane, 2-methoxy...	1.55	16.5	ng	220025	1	7.13	266402	20.0
2-Pentanone, 4-hy...	4.88	21.7	ng	289463	1	7.13	266402	20.0
unknown6.83	6.83	73.5	ng	979196	1	7.13	266402	20.0
Phenanthrene, 2-m...	13.37	5.7	ng	157004	4	12.71	548961	20.0
4H-Cyclopenta[def...	13.47	9.8	ng	268105	4	12.71	548961	20.0
2-Chloropropioni...	17.47	31.4	ng	898182	6	17.70	572477	20.0
Oxirane, heptadecyl-	18.01	11.0	ng	316088	6	17.70	572477	20.0
Eicosane	18.24	22.7	ng	648784	6	17.70	572477	20.0
unknown18.57	18.57	7.6	ng	218254	6	17.70	572477	20.0
.gamma.-Sitosterol	19.65	10.8	ng	308912	6	17.70	572477	20.0
Testosterone	20.57	7.6	ng	218241	6	17.70	572477	20.0