

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052715\
 Data File : BF079527.D
 Acq On : 27 May 2015 19:29
 Operator : TP/IZ
 Sample : G2387-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-15-SS

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-BF052615.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	1.415	18	20	26	rVB	72731	146440	2.02%	0.346%
2	1.575	33	34	42	rBV	698374	1107788	15.25%	2.615%
3	1.690	42	44	79	rVB	2962563	7262278	100.00%	17.146%
4	2.981	153	157	165	rBV	1529404	2910438	40.08%	6.871%
5	4.719	307	309	318	rVB	174159	268483	3.70%	0.634%
6	5.279	356	358	369	rBV	1698669	1947022	26.81%	4.597%
7	6.593	471	473	481	rBV	1999135	2000367	27.54%	4.723%
8	6.719	481	484	486	rBV	1941344	2144162	29.52%	5.062%
9	7.005	506	509	511	rBV	295082	403021	5.55%	0.952%
10	7.187	522	525	527	rBV	1560277	1610984	22.18%	3.803%
11	7.702	567	570	572	rBV	1342122	1330533	18.32%	3.141%
12	8.570	644	646	648	rBV	406812	570043	7.85%	1.346%
13	9.931	762	765	767	rVB	2031626	2602917	35.84%	6.145%
14	10.433	807	809	812	rBV	335161	334515	4.61%	0.790%
15	10.571	819	821	824	rVB	104768	109500	1.51%	0.259%
16	10.742	834	836	839	rVB	695883	695462	9.58%	1.642%
17	11.725	919	922	924	rBV	1703838	1753411	24.14%	4.140%
18	11.759	924	925	928	rVB	254266	207762	2.86%	0.491%
19	11.931	938	940	944	rBV	74359	134607	1.85%	0.318%
20	12.491	987	989	990	rBV	71384	73779	1.02%	0.174%
21	12.582	994	997	998	rBV	595642	679986	9.36%	1.605%
22	12.605	998	999	1001	rVB	304511	281628	3.88%	0.665%
23	12.674	1001	1005	1007	rBV	139781	160057	2.20%	0.378%
24	13.017	1034	1035	1040	rVB2	59718	78373	1.08%	0.185%
25	13.245	1053	1055	1056	rVV	70613	82711	1.14%	0.195%
26	13.337	1059	1063	1069	rVB2	119413	303481	4.18%	0.717%
27	14.011	1119	1122	1126	rBV4	66289	125956	1.73%	0.297%
28	14.079	1126	1128	1130	rBV	696243	674308	9.29%	1.592%
29	14.137	1130	1133	1137	rVV3	178070	266123	3.66%	0.628%
30	14.354	1150	1152	1155	rBV	547125	677446	9.33%	1.599%
31	14.571	1169	1171	1174	rVV	2058456	2893435	39.84%	6.831%
32	14.651	1174	1178	1180	rVV2	54608	91824	1.26%	0.217%
33	14.754	1184	1187	1188	rVV2	43650	87218	1.20%	0.206%
34	14.777	1188	1189	1192	rVB	85315	88560	1.22%	0.209%

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Integration Parameters: rteint.p
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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	14.902	1198	1200	1205	rVB3	91741	201082	2.77%	0.475%
36	15.417	1243	1245	1250	rVB	50825	77359	1.07%	0.183%
37	15.543	1254	1256	1258	rBV	76337	86960	1.20%	0.205%
38	15.600	1258	1261	1263	rBV2	51941	105712	1.46%	0.250%
39	15.748	1271	1274	1278	rBV3	332240	493664	6.80%	1.166%
40	15.851	1280	1283	1285	rBV2	735169	1256629	17.30%	2.967%
41	16.160	1308	1310	1314	rBV3	77177	124139	1.71%	0.293%
42	16.560	1343	1345	1348	rVB	152414	218605	3.01%	0.516%
43	16.937	1376	1378	1381	rBV3	62485	136406	1.88%	0.322%
44	17.120	1390	1394	1400	rBV	440784	1110424	15.29%	2.622%
45	17.234	1403	1404	1406	rVB	85226	100099	1.38%	0.236%
46	17.337	1409	1413	1416	rBV3	162695	469261	6.46%	1.108%
47	17.428	1419	1421	1425	rVV	260808	394805	5.44%	0.932%
48	17.497	1425	1427	1430	rVV	355395	443082	6.10%	1.046%
49	17.554	1430	1432	1441	rVV2	630049	994449	13.69%	2.348%
50	18.080	1476	1478	1480	rVB	110389	134432	1.85%	0.317%
51	18.869	1544	1547	1552	rVB2	209403	488230	6.72%	1.153%
52	19.029	1558	1561	1563	rBV2	126979	252736	3.48%	0.597%
53	19.246	1577	1580	1588	rBV	168554	346350	4.77%	0.818%
54	19.383	1589	1592	1595	rVV3	120274	295098	4.06%	0.697%
55	19.463	1596	1599	1606	rVB2	223049	521605	7.18%	1.231%

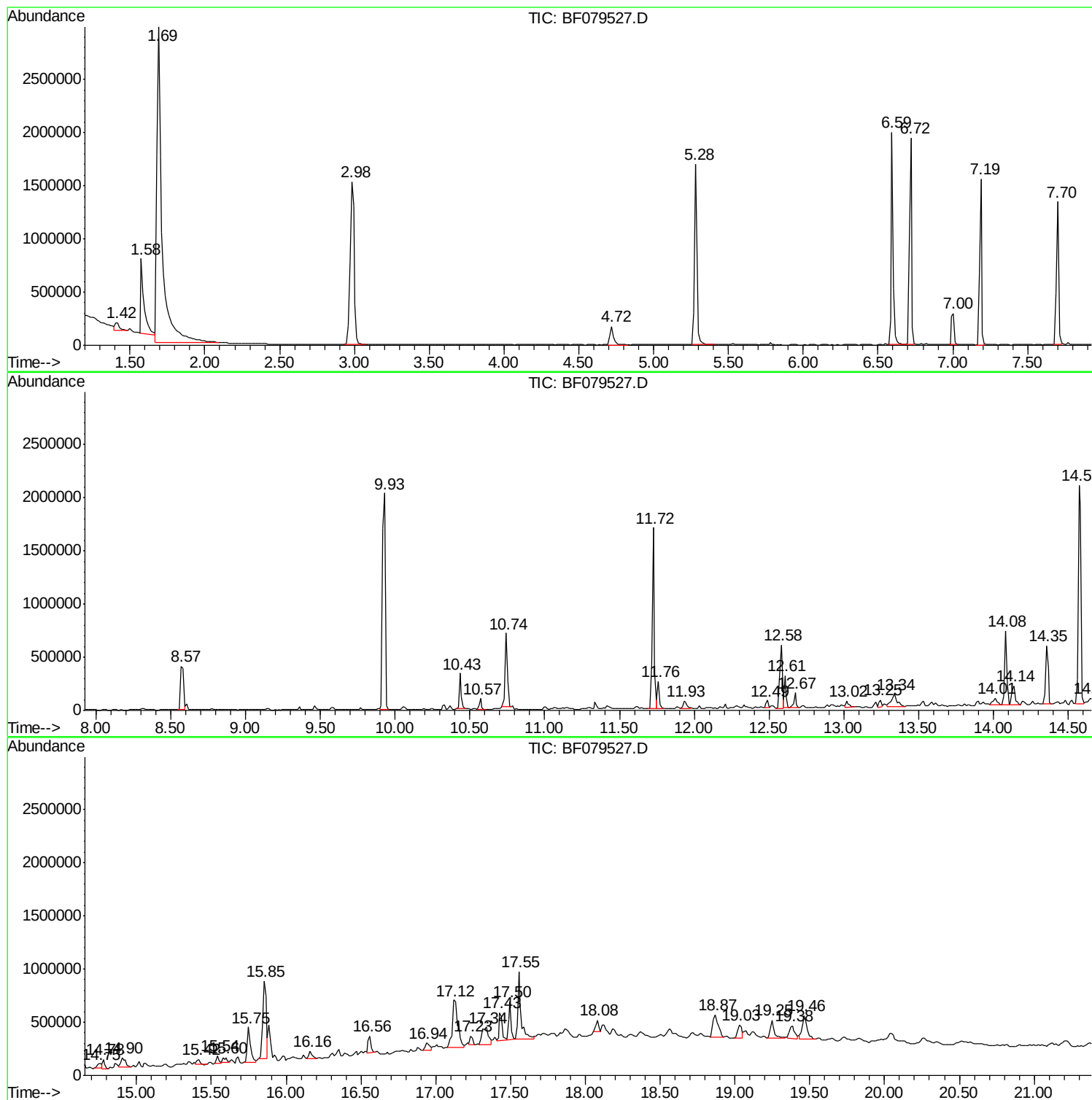
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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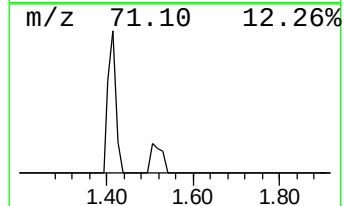
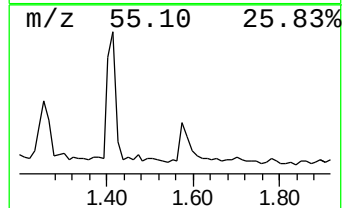
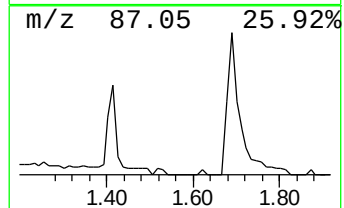
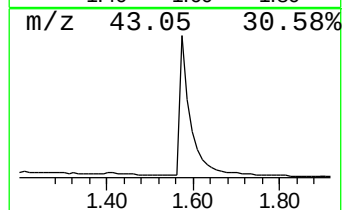
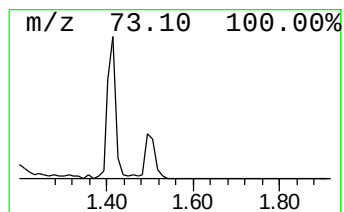
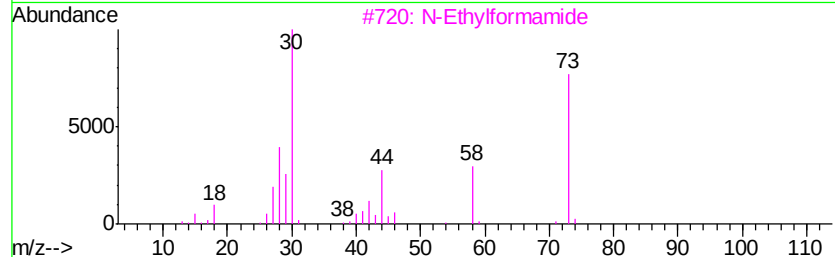
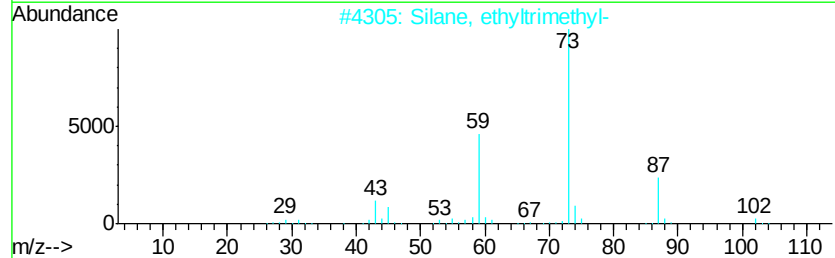
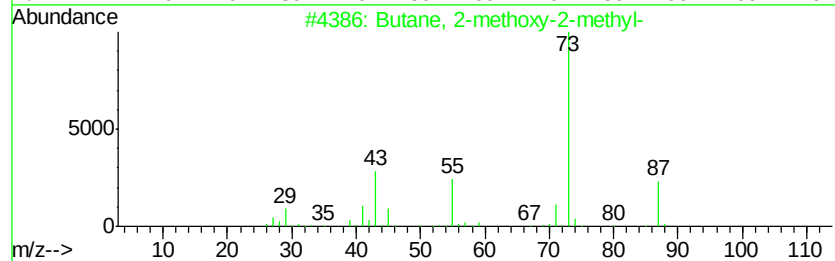
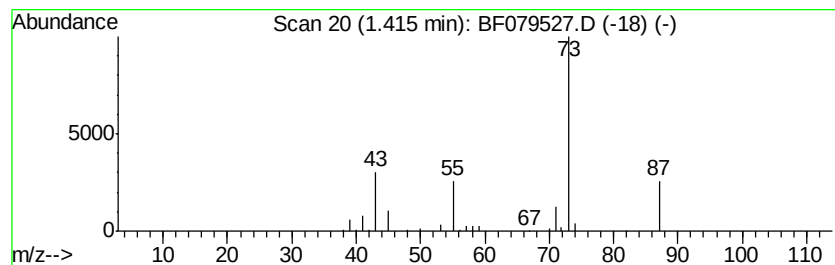
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.42	7.27 ng	146440	1,4-Dichlorobenzene-d4	7.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	33
3			N-Ethylformamide	73	C3H7NO	000627-45-2	9
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9
5			1-Propene-1-thiol	74	C3H6S	000925-89-3	9



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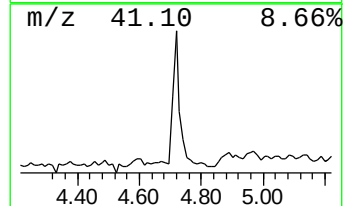
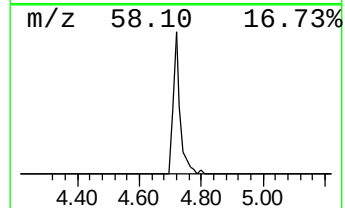
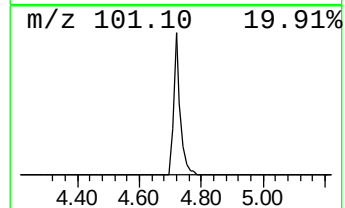
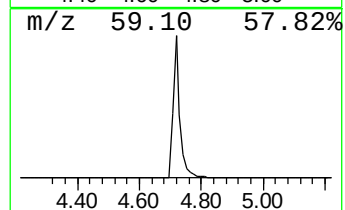
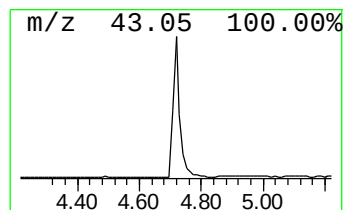
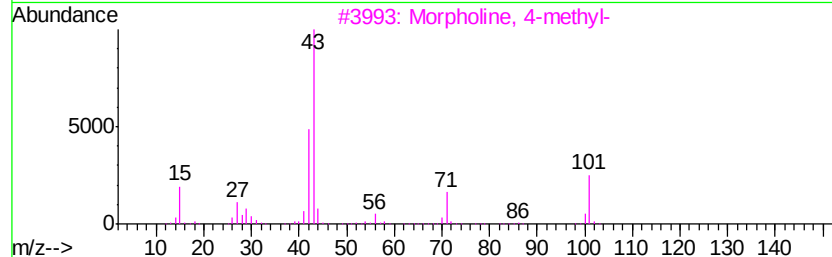
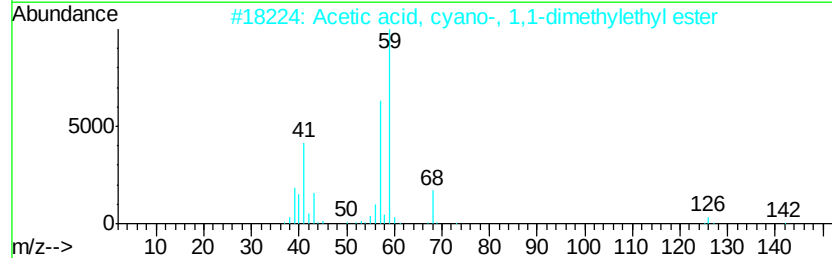
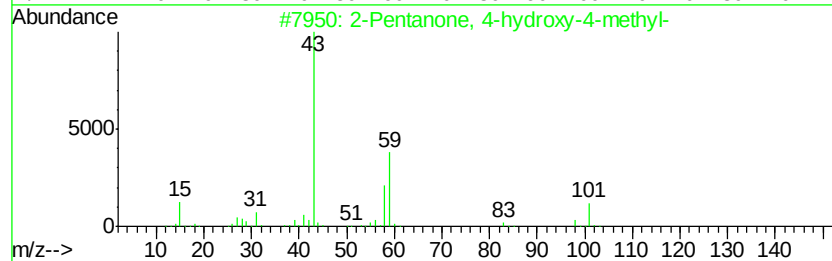
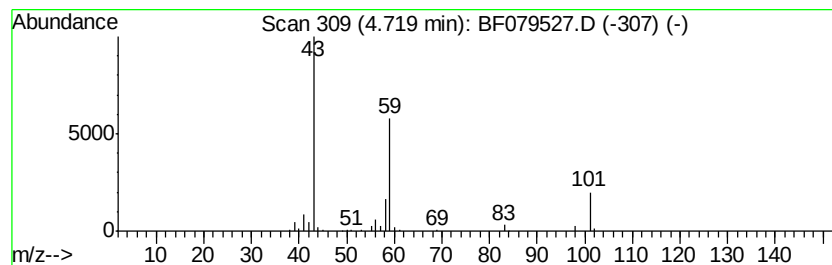
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.72	13.32 ng	268483	1,4-Dichlorobenzene-d4	7.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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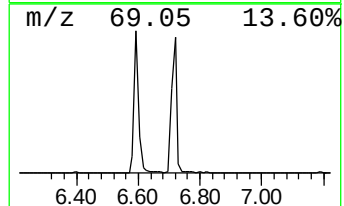
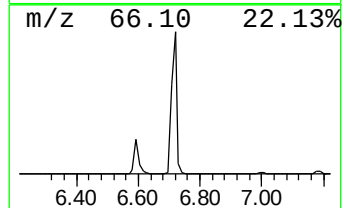
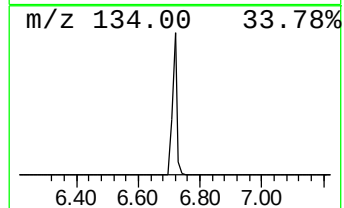
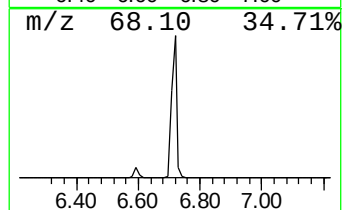
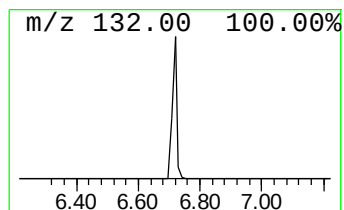
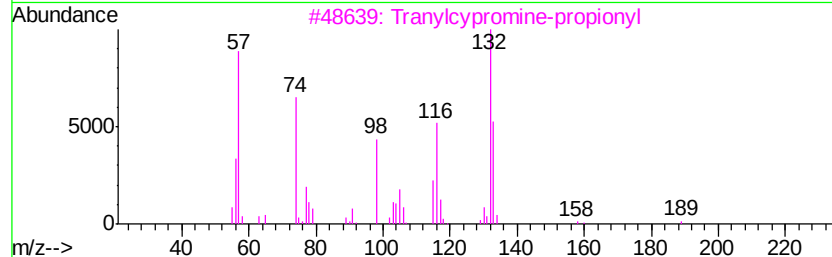
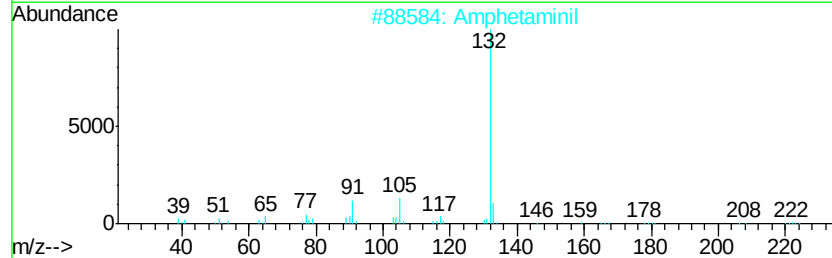
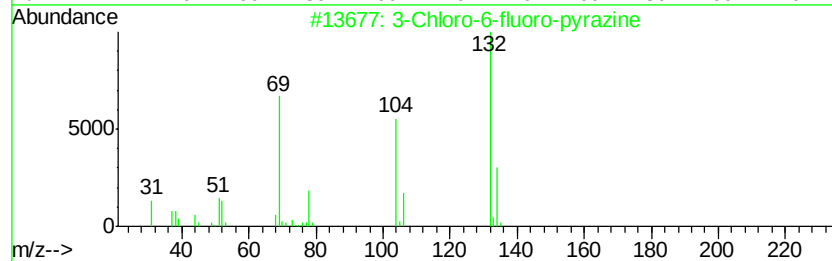
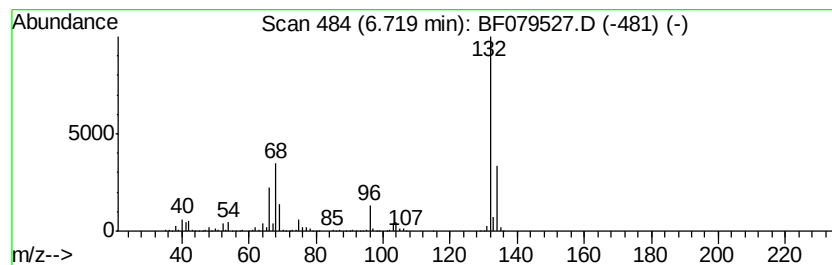
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown6.72 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.72	106.40 ng	2144160	1,4-Dichlorobenzene-d4	7.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		Amphetaminil	250	C17H18N2	017590-01-1	12
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12



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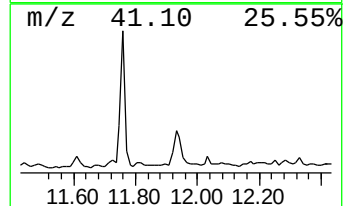
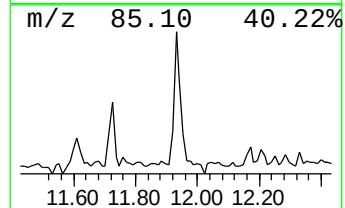
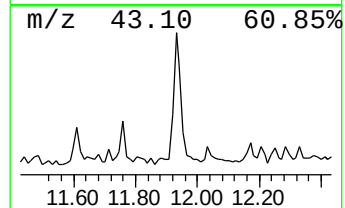
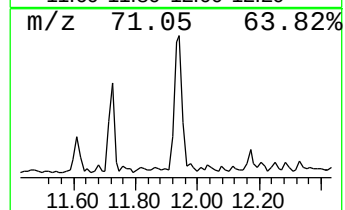
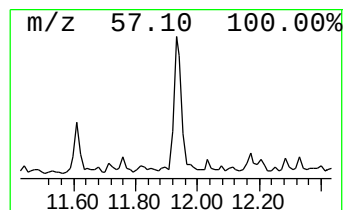
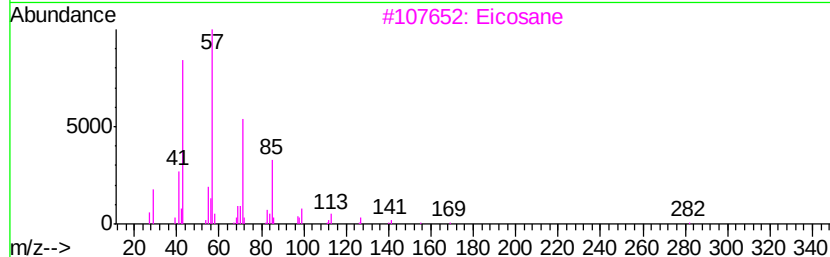
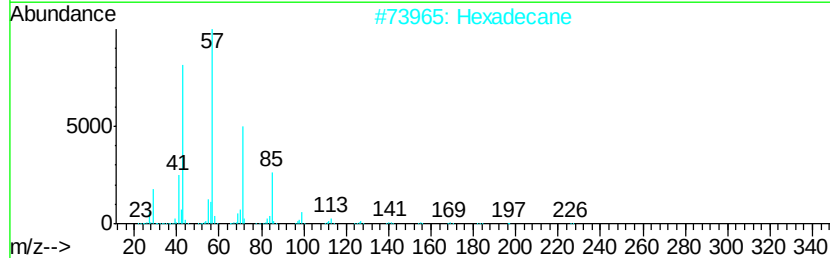
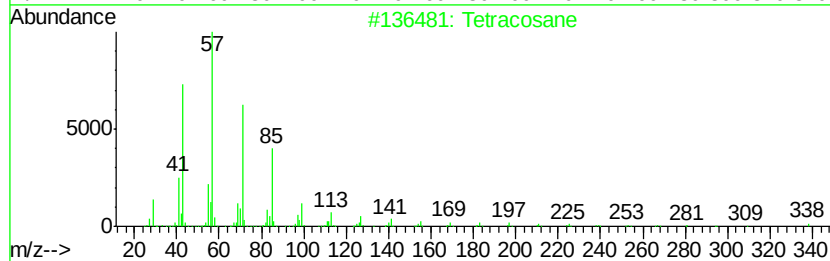
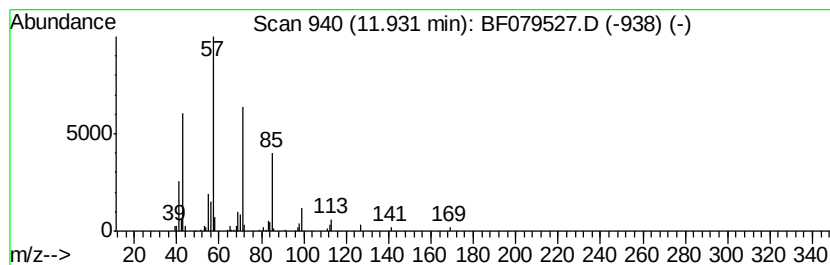
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tetracosane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	3.96 ng	134607	Phenanthrene-d10	12.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	90
2		Hexadecane	226	C16H34	000544-76-3	83
3		Eicosane	282	C20H42	000112-95-8	80
4		Heptacosane	380	C27H56	000593-49-7	78
5		Undecane, 5-ethyl-	184	C13H28	017453-94-0	72



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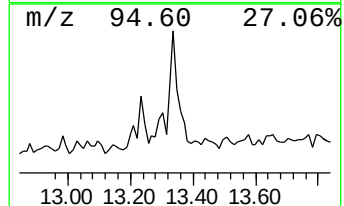
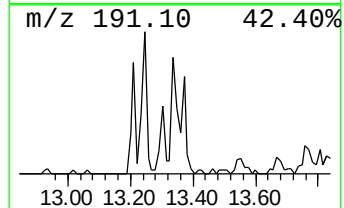
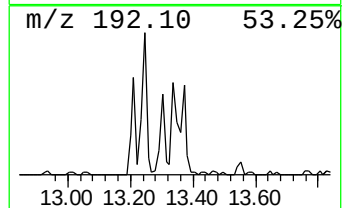
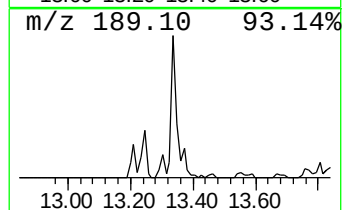
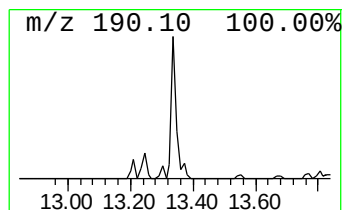
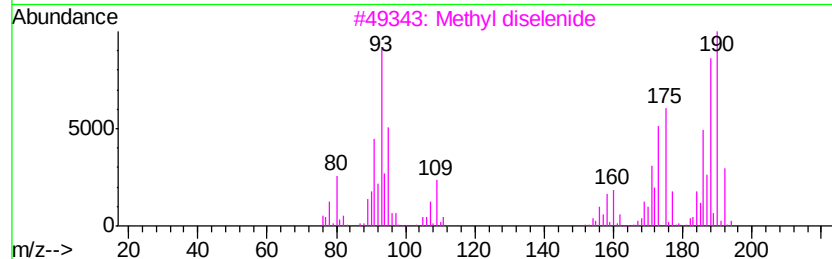
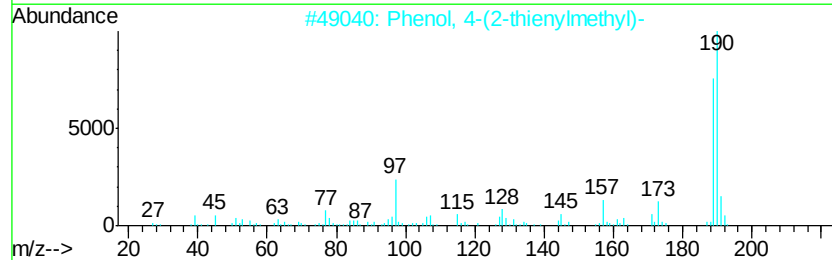
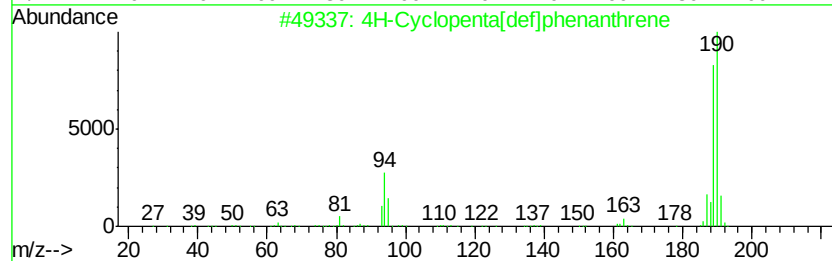
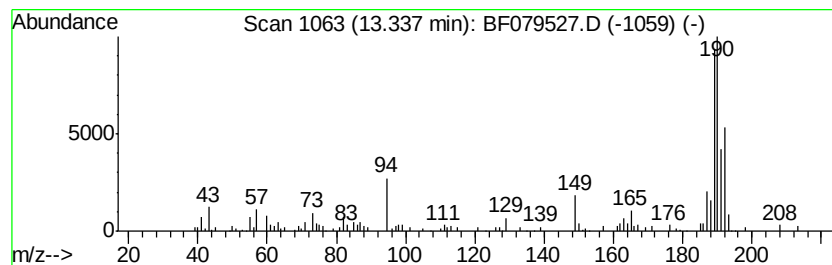
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.34	8.93 ng	303481	Phenanthrene-d10	12.58	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	62
2	Phenol, 4-(2-thienylmethyl)-	190	C11H10OS	091680-55-6	46
3	Methyl diselenide	190	C2H6Se2	007101-31-7	45
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	20
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	18



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052715\
Data File : BF079527.D
Acq On : 27 May 2015 19:29
Operator : TP/IZ
Sample : G2387-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

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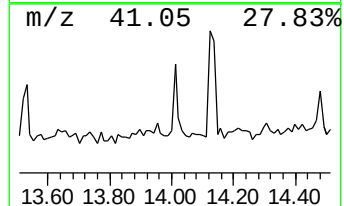
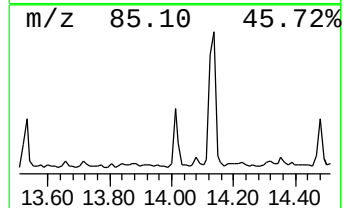
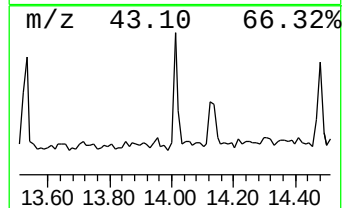
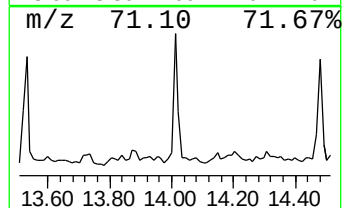
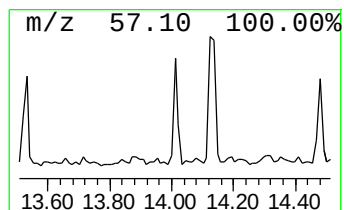
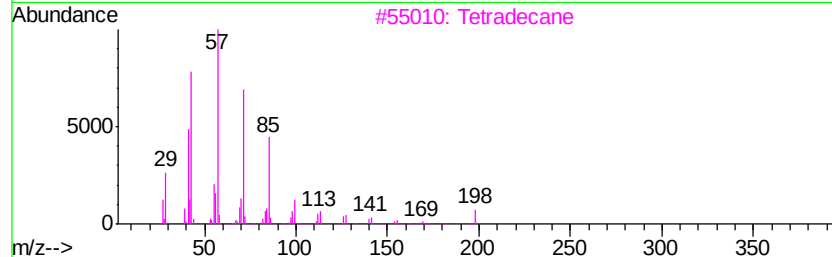
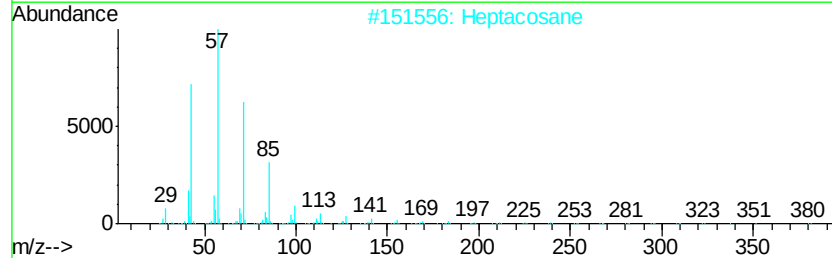
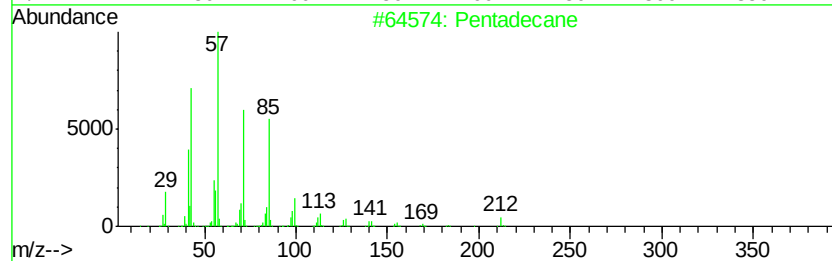
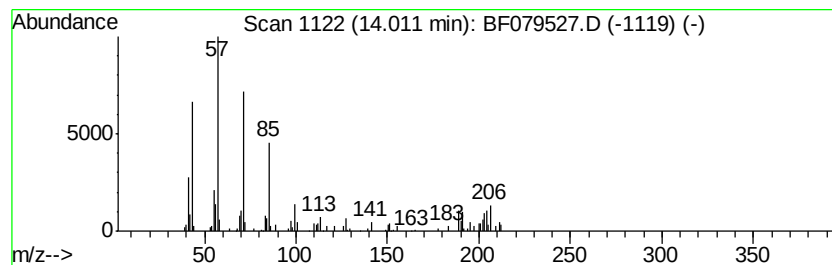
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Pentadecane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.01	3.70 ng	125956	Phenanthrene-d10	12.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane	212	C15H32	000629-62-9	86
2		Heptacosane	380	C27H56	000593-49-7	58
3		Tetradecane	198	C14H30	000629-59-4	58
4		Heptadecane	240	C17H36	000629-78-7	58
5		Heneicosane	296	C21H44	000629-94-7	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052715\
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Acq On : 27 May 2015 19:29
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ALS Vial : 11 Sample Multiplier: 1

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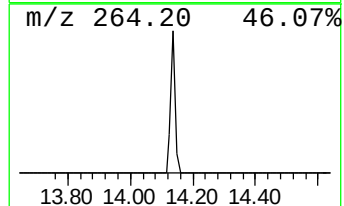
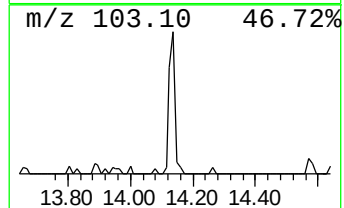
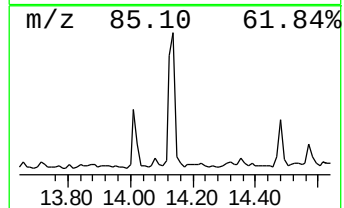
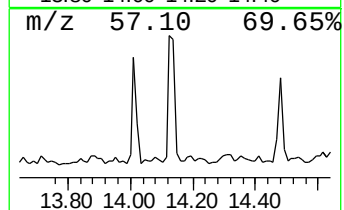
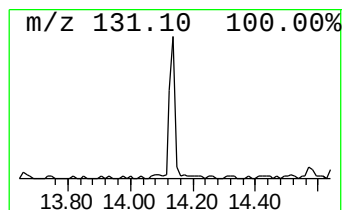
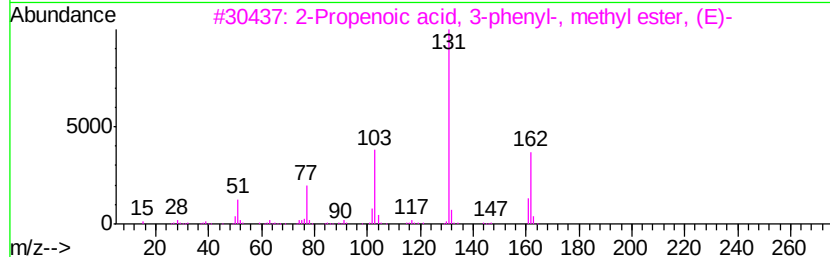
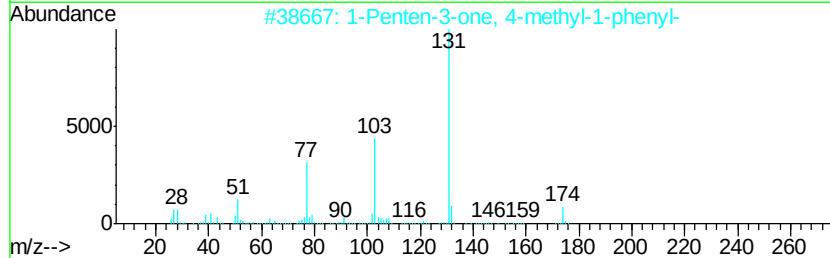
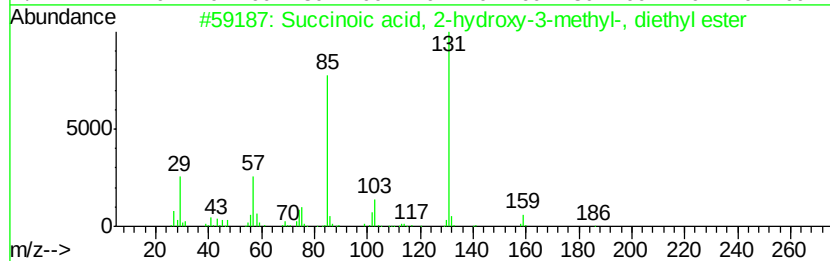
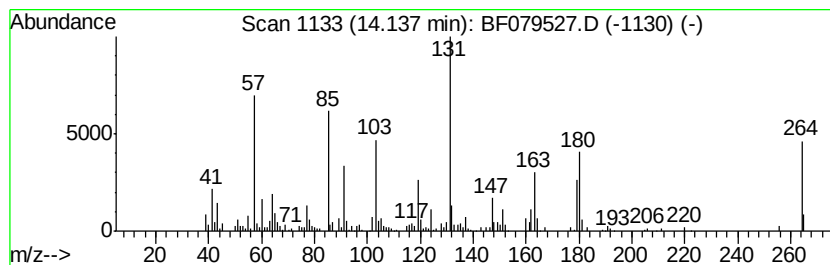
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown14.14 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.14	7.83 ng	266123	Phenanthrene-d10	12.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Succinic acid, 2-hydroxy-3-meth...	204	C9H16O5	1000196-85-6	27
2		1-Penten-3-one, 4-methyl-1-phenyl-	174	C12H14O	003160-32-5	22
3		2-Propenoic acid, 3-phenyl-, met...	162	C10H10O2	001754-62-7	22
4		3-Butenoic acid, 2-oxo-4-phenyl-...	190	C11H10O3	1000142-22-4	22
5		3-Butenoic acid, 2-oxo-4-phenyl-	176	C10H8O3	1000142-21-0	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052715\
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Sample : G2387-02
Misc :
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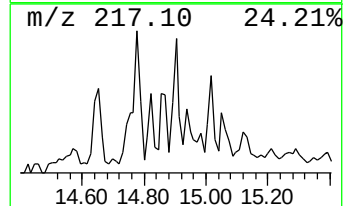
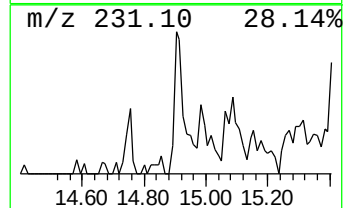
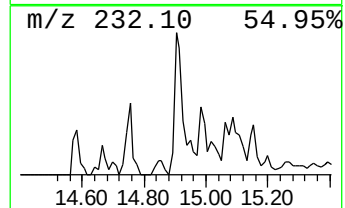
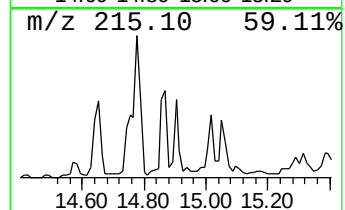
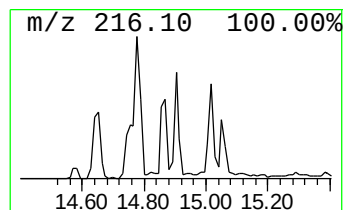
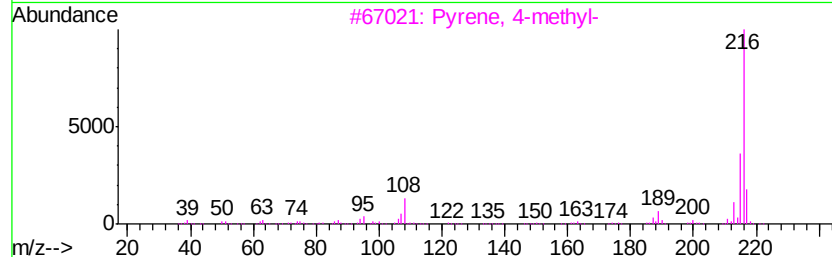
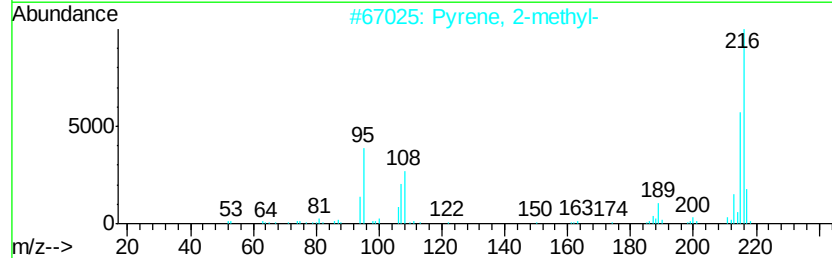
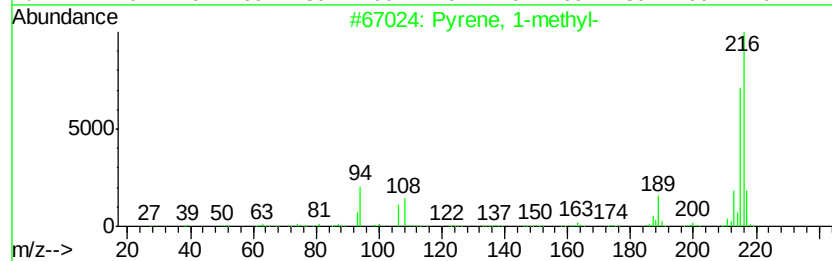
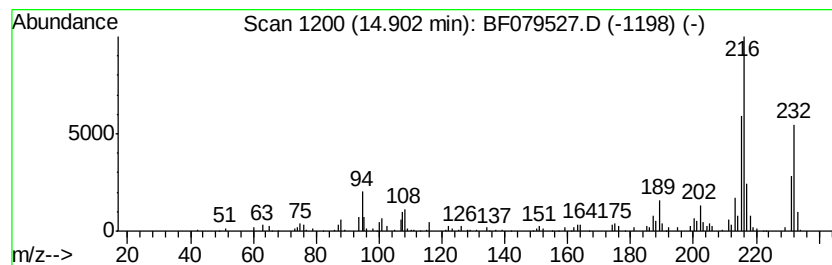
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Pyrene, 1-methyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD		R.T.
14.90	3.20 ng	201082	Chrysene-d12		15.86
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	95
2	Pyrene, 2-methyl-	216	C17H12	003442-78-2	83
3	Pyrene, 4-methyl-	216	C17H12	003353-12-6	50
4	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	50
5	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	49



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052715\
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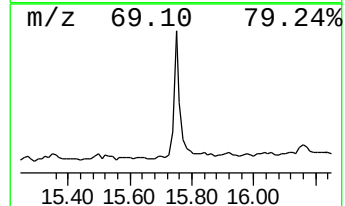
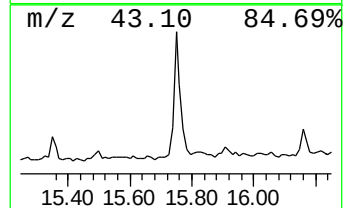
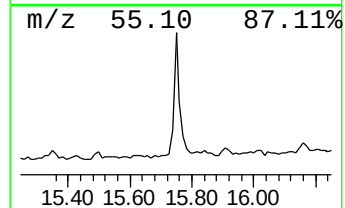
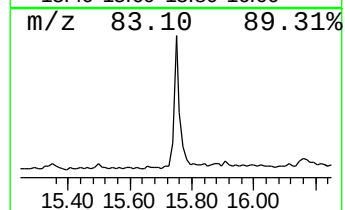
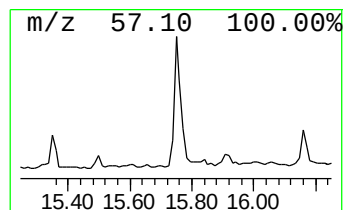
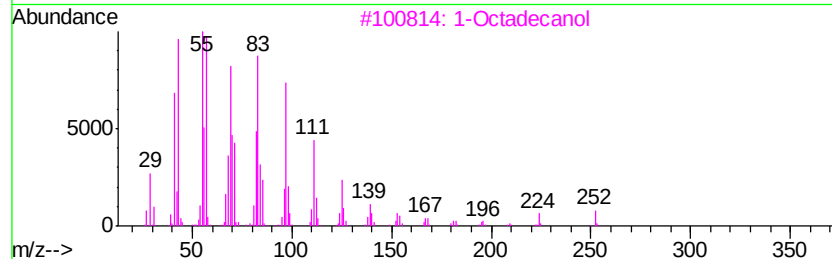
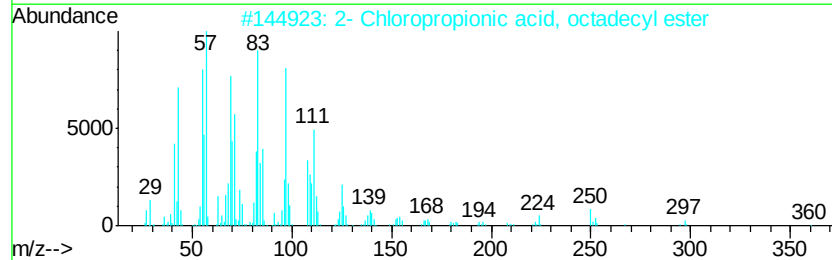
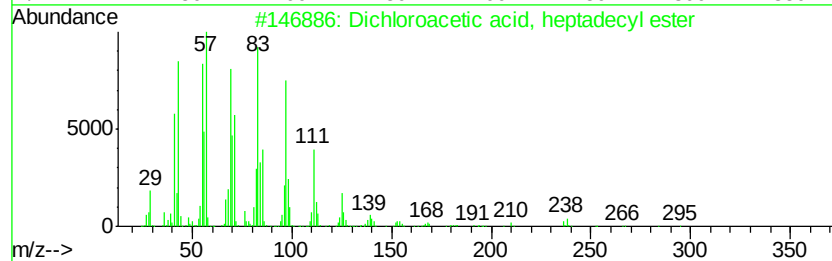
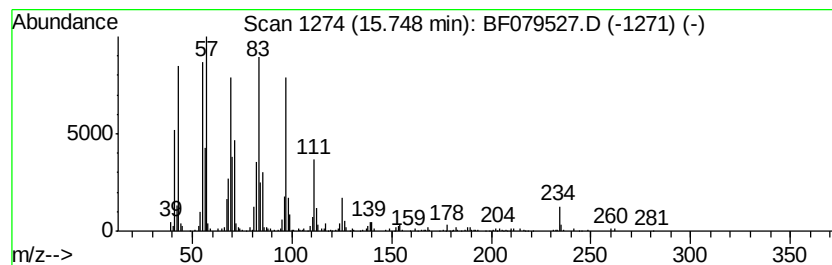
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Dichloroacetic acid, heptad... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.75	7.86 ng	493664	Chrysene-d12	15.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	94
2		2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	91
3		1-Octadecanol	270	C18H38O	000112-92-5	91
4		Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	91
5		3-Eicosene, (E)-	280	C20H40	074685-33-9	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052715\
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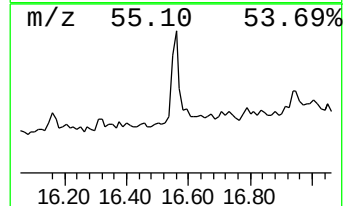
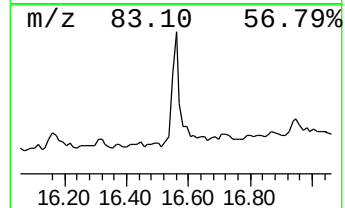
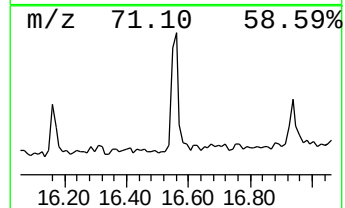
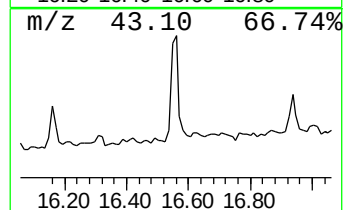
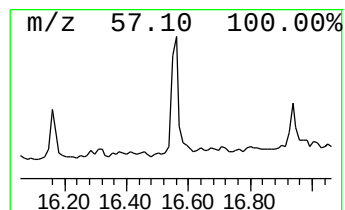
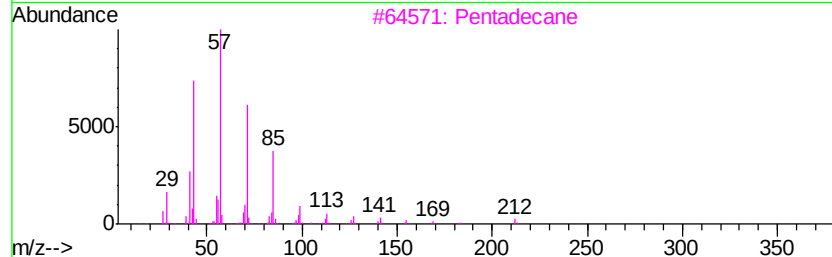
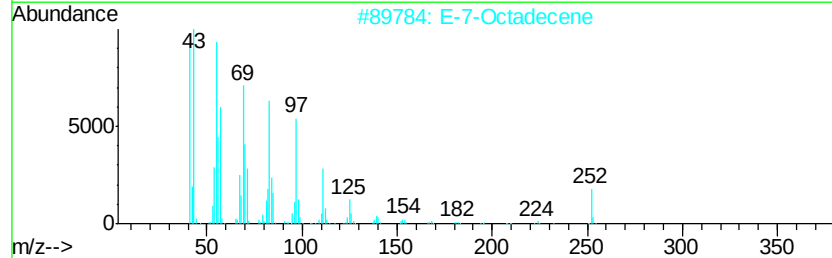
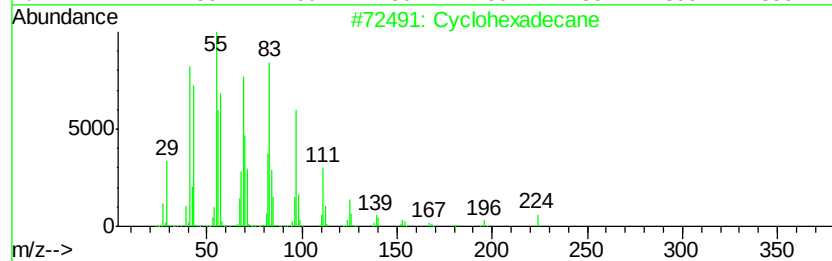
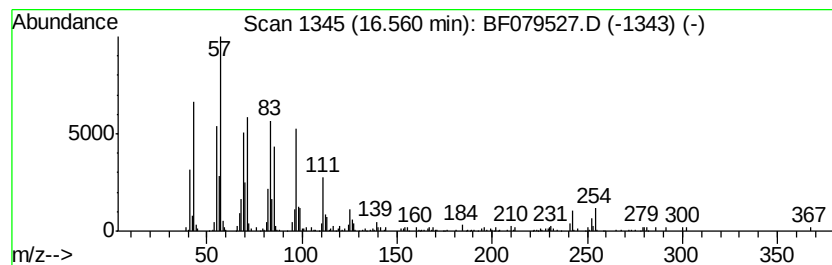
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Cyclohexadecane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.56	3.48 ng	218605	Chrysene-d12	15.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	95
2			E-7-Octadecene	252	C18H36	1000130-92-0	90
3			Pentadecane	212	C15H32	000629-62-9	83
4			1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	81
5			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	70



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052715\
Data File : BF079527.D
Acq On : 27 May 2015 19:29
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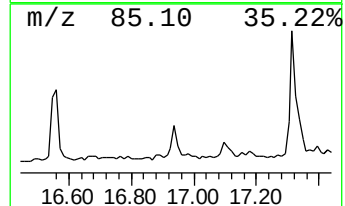
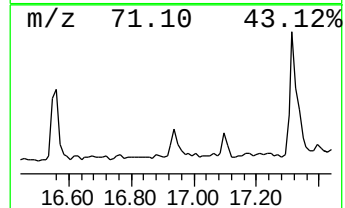
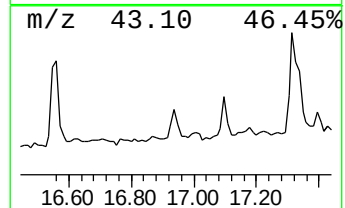
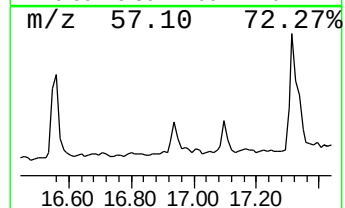
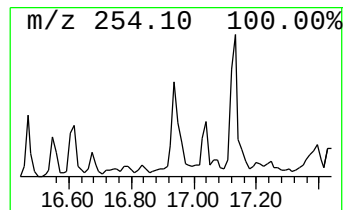
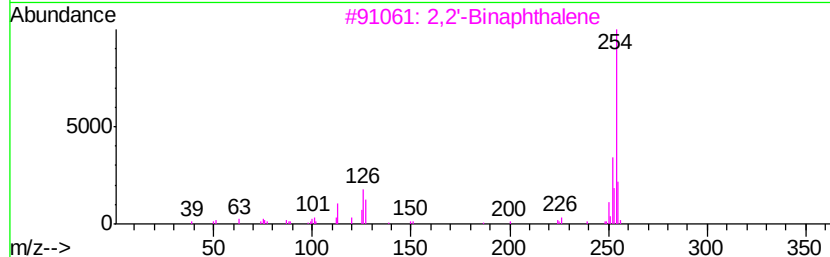
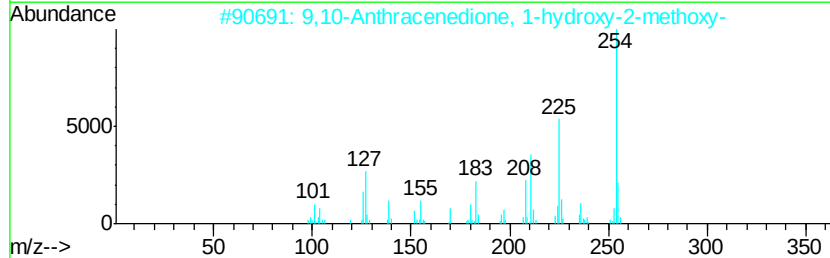
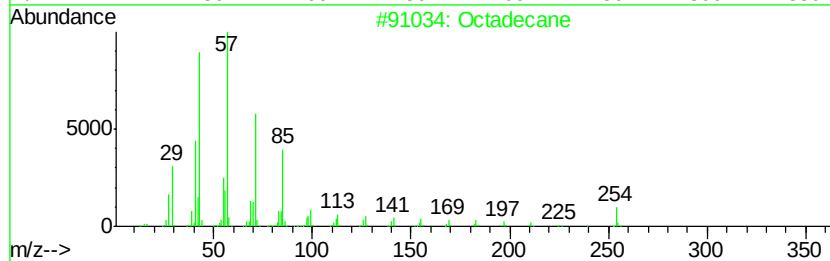
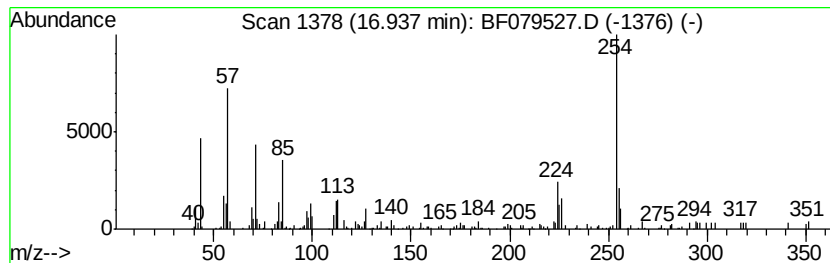
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Octadecane Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.94	2.74 ng	136406	Perylene-d12	17.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	70
2		9,10-Anthracenedione, 1-hydroxy-...	254	C15H10O4	006003-11-8	53
3		2,2'-Binaphthalene	254	C20H14	000612-78-2	43
4		Anthracene, 9-phenyl-	254	C20H14	000602-55-1	38
5		Phenanthrene, 2-phenyl-	254	C20H14	004325-77-3	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052715\
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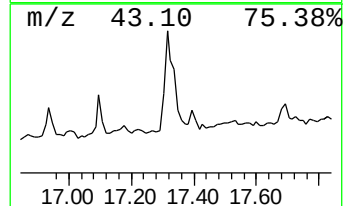
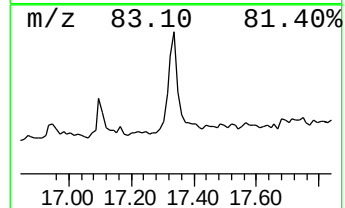
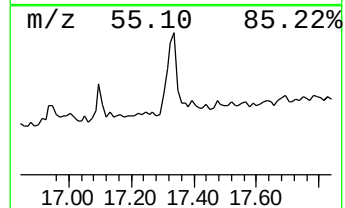
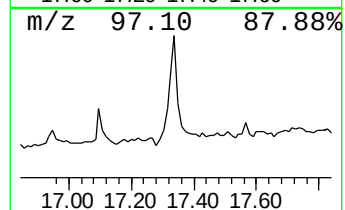
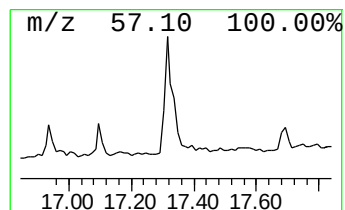
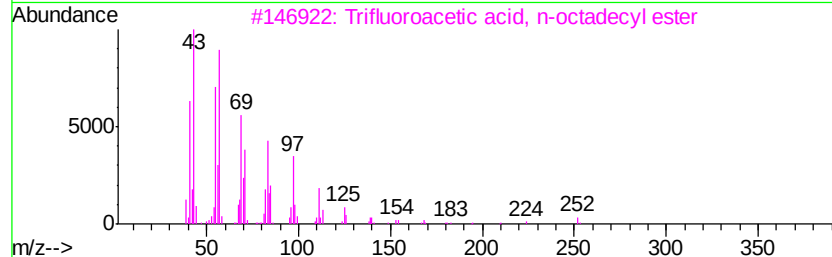
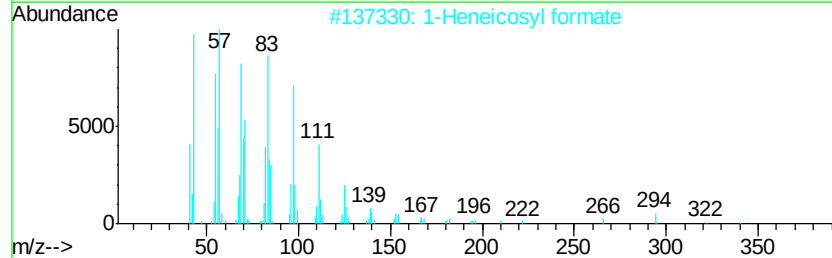
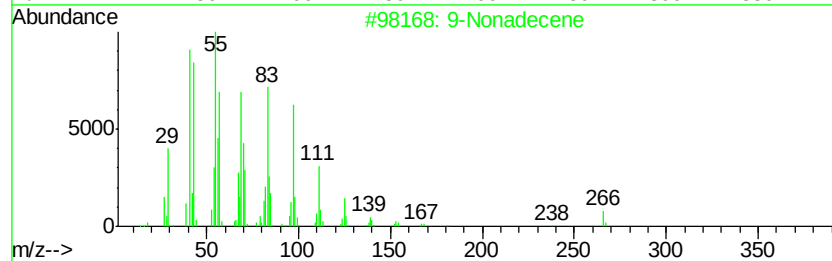
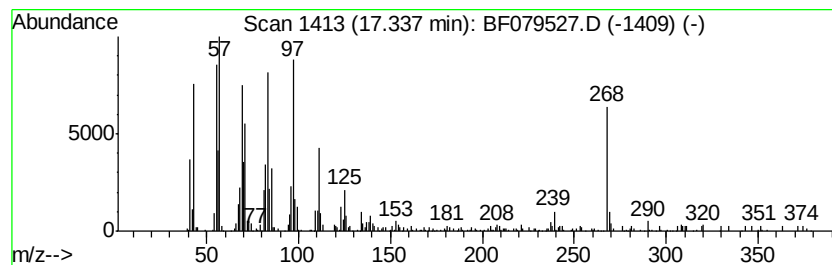
Instrument :
BNA_F
ClientSampleId :
TP-15-SS

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 9-Nonadecene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.34	9.44 ng	469261	Perylene-d12	17.55	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Nonadecene	266	C19H38	031035-07-1	78
2	1-Heneicosyl formate	340	C22H44O2	077899-03-7	76
3	Trifluoroacetic acid, n-octadecy...	366	C20H37F3O2	079392-43-1	74
4	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	70
5	1-Docosene	308	C22H44	001599-67-3	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052715\
Data File : BF079527.D
Acq On : 27 May 2015 19:29
Operator : TP/IZ
Sample : G2387-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-15-SS

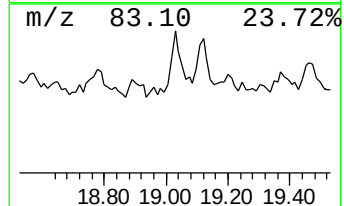
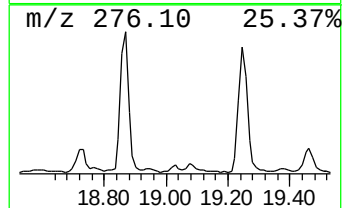
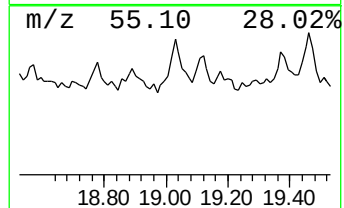
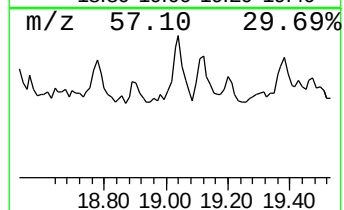
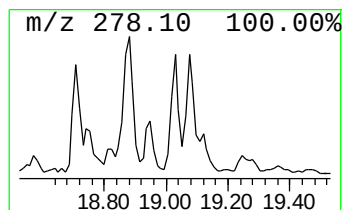
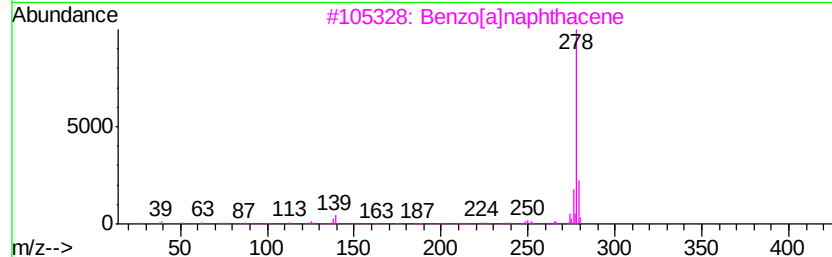
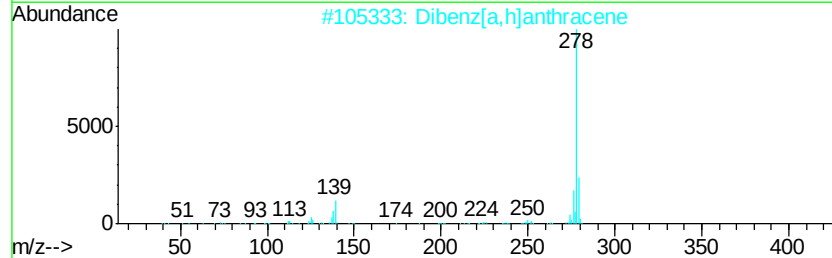
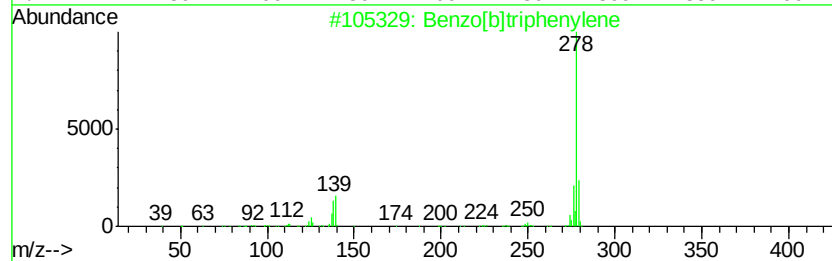
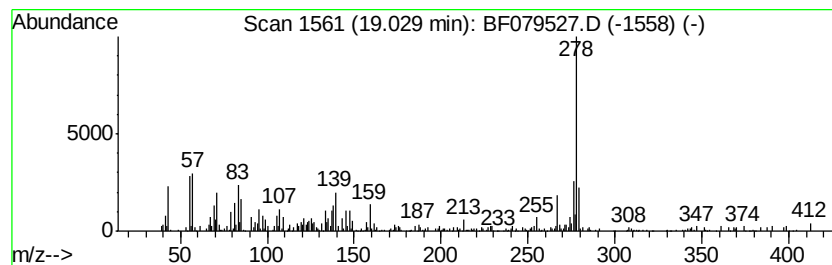
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Benzo[b]triphenylene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.03	5.08 ng	252736	Perylene-d12	17.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]triphenylene	278	C22H14	000215-58-7	93
2		Dibenz[a,h]anthracene	278	C22H14	000053-70-3	89
3		Benzo[a]naphthacene	278	C22H14	000226-88-0	89
4		Pentaphene	278	C22H14	000222-93-5	89
5		1,2:7,8-Dibenzophenanthrene	278	C22H14	000213-46-7	89



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052715\
Data File : BF079527.D
Acq On : 27 May 2015 19:29
Operator : TP/IZ
Sample : G2387-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-15-SS

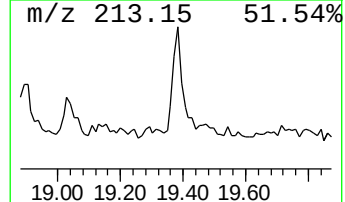
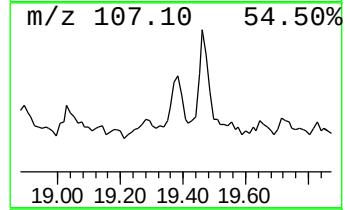
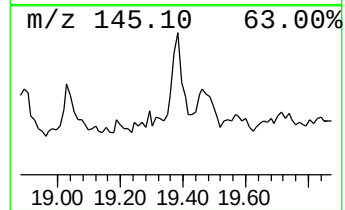
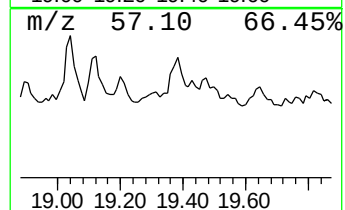
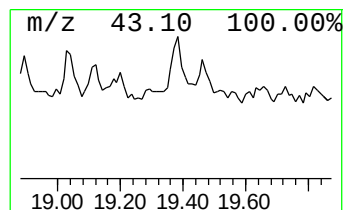
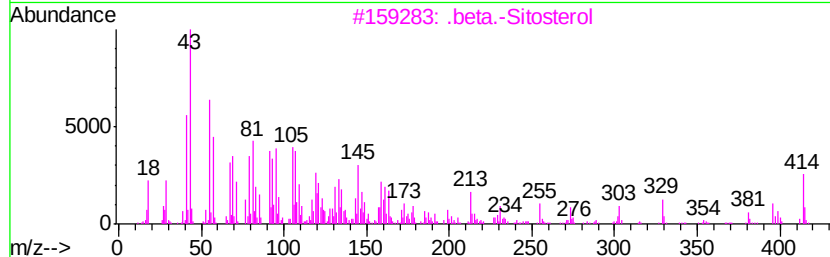
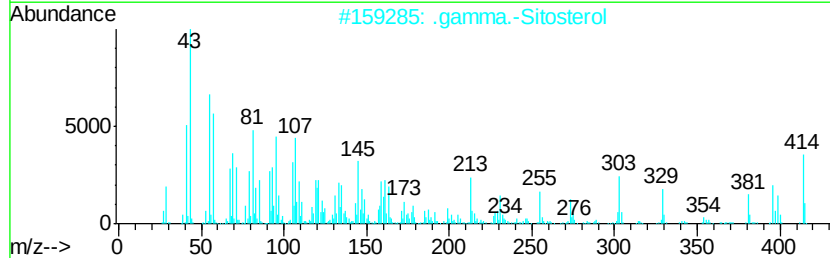
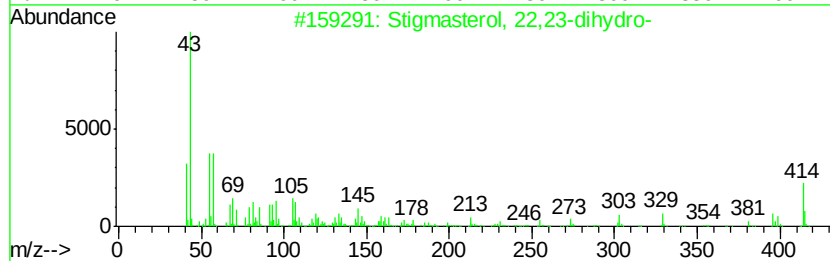
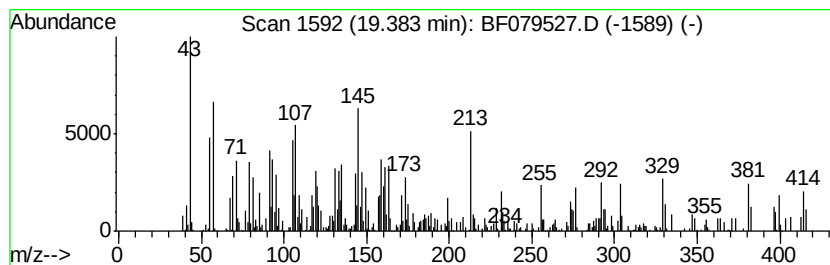
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Stigmasterol, 22,23-dihydro- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.38	5.93 ng	295098	Perylene-d12	17.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Stigmasterol, 22,23-dihydro-	414	C29H50O	1000214-20-7	52
2		.gamma.-Sitosterol	414	C29H50O	000083-47-6	38
3		.beta.-Sitosterol	414	C29H50O	000083-46-5	30
4		Pregn-5-ene-3,12,14,17,20-pentol...	366	C21H34O5	028417-32-5	10
5		3-[(3,7,11,15-Tetramethylhexadec...	374	C25H46N2	067374-05-4	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052715\
 Data File : BF079527.D
 Acq On : 27 May 2015 19:29
 Operator : TP/IZ
 Sample : G2387-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-15-SS

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052615.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
Butane, 2-methoxy...	1.42	7.3	ng	146440	1	7.00	403021	20.0
2-Pentanone, 4-hy...	4.72	13.3	ng	268483	1	7.00	403021	20.0
unknown6.72	6.72	106.4	ng	2144160	1	7.00	403021	20.0
Tetracosane	11.93	4.0	ng	134607	4	12.58	679986	20.0
4H-Cyclopenta[def...	13.34	8.9	ng	303481	4	12.58	679986	20.0
Pentadecane	14.01	3.7	ng	125956	4	12.58	679986	20.0
unknown14.14	14.14	7.8	ng	266123	4	12.58	679986	20.0
Pyrene, 1-methyl-	14.90	3.2	ng	201082	5	15.86	1256630	20.0
Dichloroacetic ac...	15.75	7.9	ng	493664	5	15.86	1256630	20.0
Cyclohexadecane	16.56	3.5	ng	218605	5	15.86	1256630	20.0
Octadecane	16.94	2.7	ng	136406	6	17.55	994449	20.0
9-Nonadecene	17.34	9.4	ng	469261	6	17.55	994449	20.0
Benzo[b]triphenylene	19.03	5.1	ng	252736	6	17.55	994449	20.0
Stigmasterol, 22,...	19.38	5.9	ng	295098	6	17.55	994449	20.0