

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051716\
 Data File : BF087125.D
 Acq On : 18 May 2016 2:19
 Operator : UM/SJ
 Sample : H3085-01
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VITALE-FILL-CSPH3-10K

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-BF051716.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	1.724	10	12	57	rVB	5022684	10844438	100.00%	14.786%
2	3.233	141	144	155	rVB	149907	304946	2.81%	0.416%
3	3.964	206	208	220	rBV	69117	147097	1.36%	0.201%
4	4.719	271	274	290	rBV	1036037	1864293	17.19%	2.542%
5	5.187	312	315	317	rBV	3652350	4680075	43.16%	6.381%
6	5.587	348	350	358	rVB	90303	148760	1.37%	0.203%
7	6.262	406	409	411	rBV	4358493	4890130	45.09%	6.667%
8	6.365	416	418	420	rBV	4771058	4686187	43.21%	6.389%
9	6.593	435	438	444	rVB	770498	905904	8.35%	1.235%
10	6.742	449	451	454	rBV	3057180	3308017	30.50%	4.510%
11	7.165	485	488	491	rBV	2822514	2958651	27.28%	4.034%
12	7.873	548	550	563	rBV2	896783	1323694	12.21%	1.805%
13	8.959	642	645	648	rBV	4452759	4707318	43.41%	6.418%
14	9.359	678	680	684	rBV	509947	485909	4.48%	0.663%
15	9.576	697	699	701	rBV	144985	134455	1.24%	0.183%
16	9.622	701	703	705	rVV2	1091500	1291326	11.91%	1.761%
17	9.656	705	706	710	rVB	463338	406421	3.75%	0.554%
18	9.839	720	722	723	rVV	87486	117453	1.08%	0.160%
19	10.068	740	742	744	rVB	156318	149002	1.37%	0.203%
20	10.171	749	751	754	rBV	244628	234479	2.16%	0.320%
21	10.422	770	773	775	rBV	3011064	3470154	32.00%	4.731%
22	10.536	782	783	787	rVB	131700	193405	1.78%	0.264%
23	10.959	818	820	822	rBV	156915	225763	2.08%	0.308%
24	11.005	822	824	828	rVB	100100	164667	1.52%	0.225%
25	11.108	830	833	834	rBV	1314131	1689433	15.58%	2.303%
26	11.131	834	835	837	rVV	2526344	1848708	17.05%	2.521%
27	11.176	837	839	842	rVV	363272	433897	4.00%	0.592%
28	11.348	850	854	858	rBV	333292	363590	3.35%	0.496%
29	11.656	878	881	884	rVV2	224137	403262	3.72%	0.550%
30	11.714	884	886	892	rVB2	326027	382605	3.53%	0.522%
31	12.308	936	938	941	rBV	1972272	2342685	21.60%	3.194%
32	12.457	948	951	955	rBV2	91286	145805	1.34%	0.199%
33	12.537	955	958	961	rVV	2166417	2332467	21.51%	3.180%
34	12.697	969	972	974	rVV	4481225	5220768	48.14%	7.118%

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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 >
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35	12.765	974	978	980	rVV2	71526	138077	1.27%	0.188%
36	12.868	983	987	991	rBV	274917	365848	3.37%	0.499%
37	12.937	991	993	994	rBV	236303	249631	2.30%	0.340%
38	12.971	994	996	1000	rVV2	110018	204048	1.88%	0.278%
39	13.485	1038	1041	1044	rBV2	127323	279315	2.58%	0.381%
40	13.531	1044	1045	1049	rVV4	121491	256897	2.37%	0.350%
41	13.657	1052	1056	1059	rBV3	96320	285150	2.63%	0.389%
42	13.725	1059	1062	1067	rVV2	1247239	2972880	27.41%	4.053%
43	13.828	1069	1071	1074	rVB	122604	180771	1.67%	0.246%
44	14.217	1101	1105	1106	rBV3	68525	120645	1.11%	0.164%
45	14.571	1134	1136	1141	rBV3	155952	250195	2.31%	0.341%
46	14.720	1147	1149	1154	rBV	592470	1148469	10.59%	1.566%
47	14.948	1166	1169	1170	rBV2	86081	126317	1.16%	0.172%
48	14.983	1170	1172	1175	rVV	384914	424914	3.92%	0.579%
49	15.040	1175	1177	1179	rVV	381856	534311	4.93%	0.729%
50	15.085	1179	1181	1187	rVB2	683265	993717	9.16%	1.355%
51	15.360	1201	1205	1207	rVB	96371	167431	1.54%	0.228%
52	15.828	1243	1246	1250	rBV	103946	211330	1.95%	0.288%
53	16.320	1287	1289	1293	rBV2	181611	370126	3.41%	0.505%
54	16.400	1293	1296	1300	rVB2	99648	232476	2.14%	0.317%
55	16.697	1318	1322	1330	rBV	170835	370195	3.41%	0.505%
56	16.903	1337	1340	1347	rVB	47326	137117	1.26%	0.187%
57	17.097	1352	1357	1362	rVB	59368	164997	1.52%	0.225%
58	17.943	1426	1431	1436	rBV3	30014	109665	1.01%	0.150%
59	18.571	1481	1486	1492	rBV	33366	115447	1.06%	0.157%
60	19.006	1517	1524	1534	rVB8	23204	127624	1.18%	0.174%

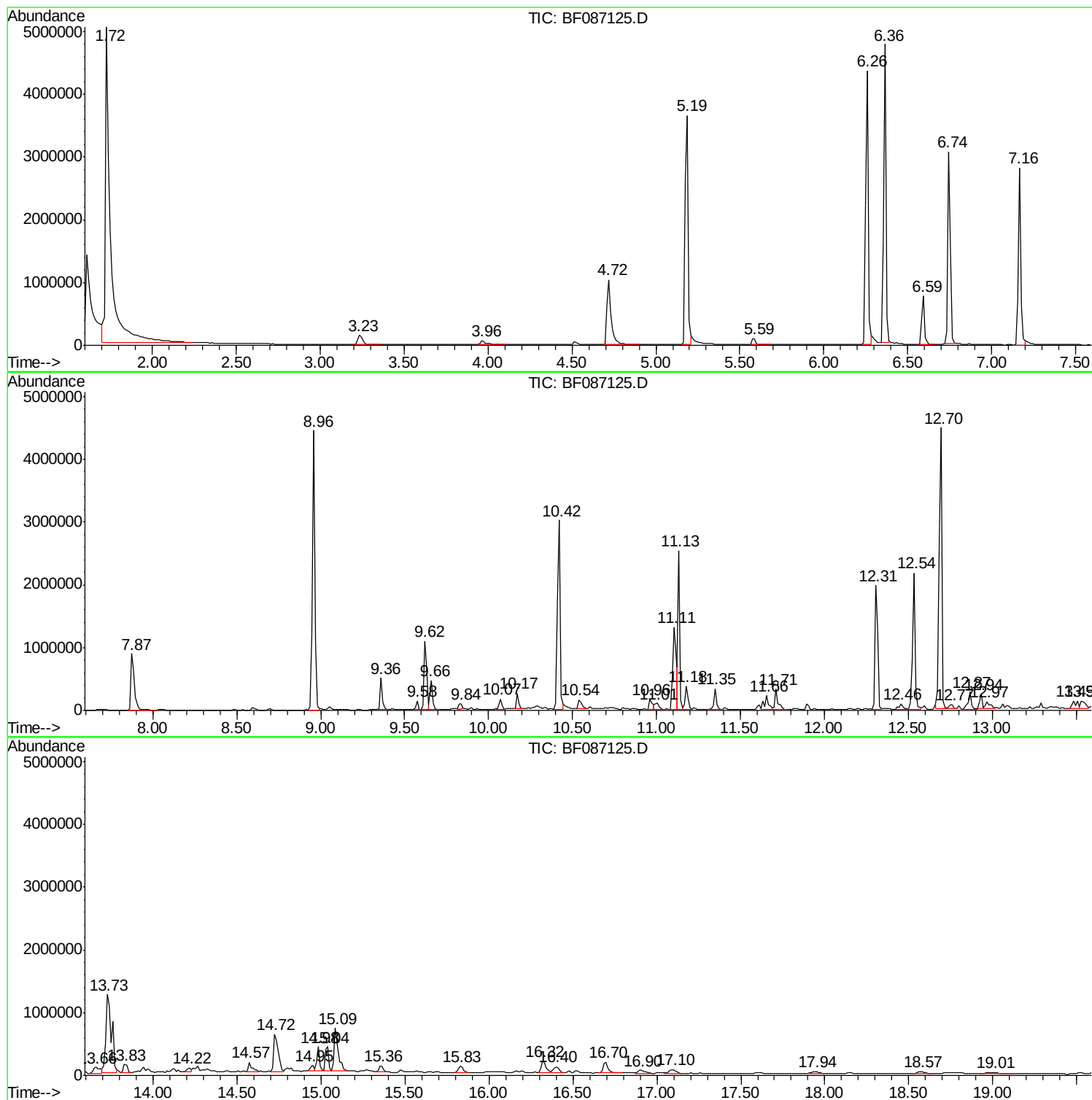
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF051716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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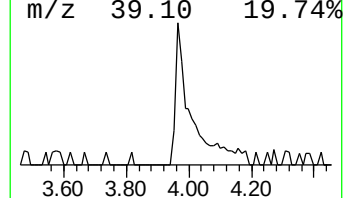
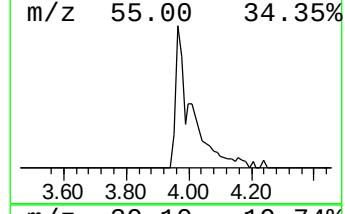
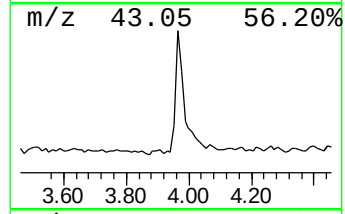
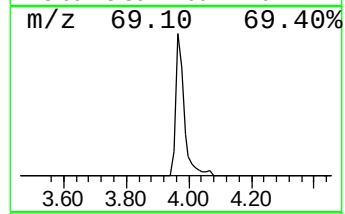
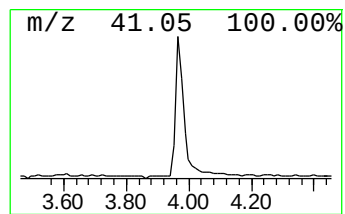
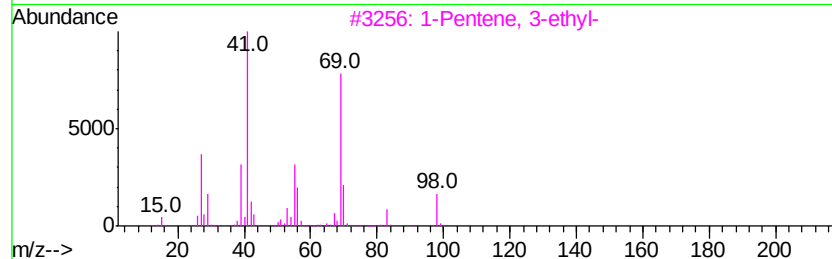
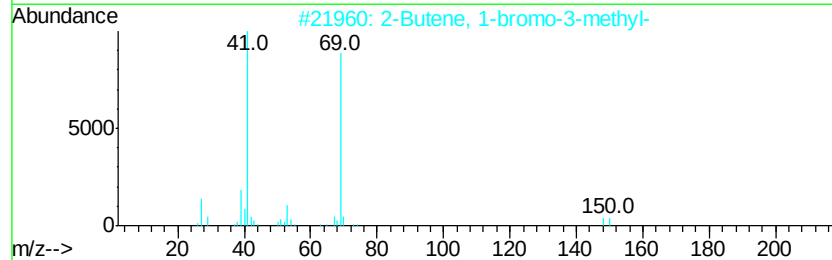
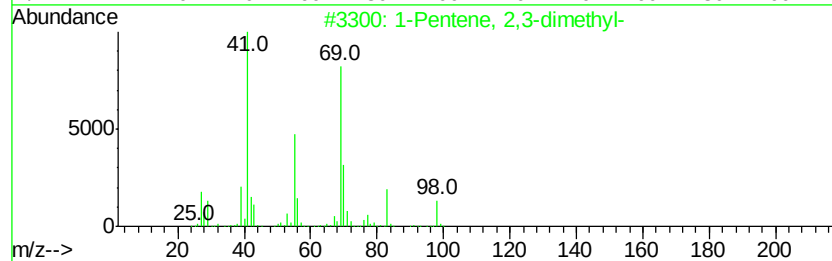
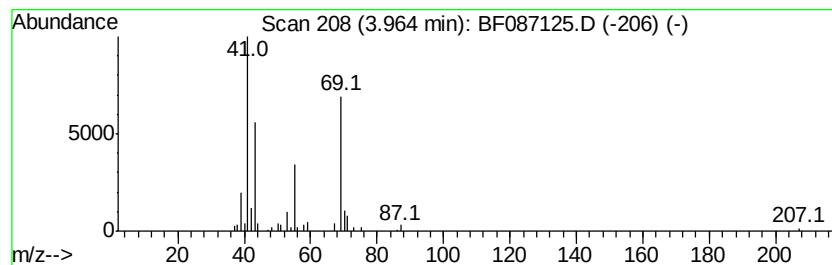
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF051716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 unknown3.96 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.96	3.25 ng	147097	1,4-Dichlorobenzene-d4	6.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Pentene, 2,3-dimethyl-	98	C7H14	003404-72-6	40
2		2-Butene, 1-bromo-3-methyl-	148	C5H9Br	000870-63-3	40
3		1-Pentene, 3-ethyl-	98	C7H14	004038-04-4	40
4		3-Penten-1-ol, 2-methyl-	100	C6H12O	062238-37-3	39
5		Butane, 2,3-dimethyl-2-nitro-	131	C6H13NO2	034075-28-0	39



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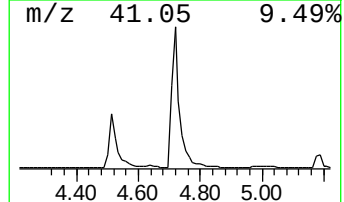
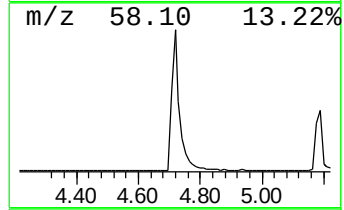
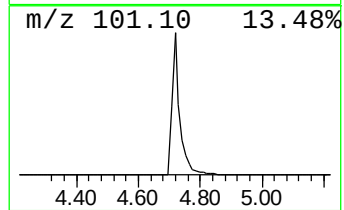
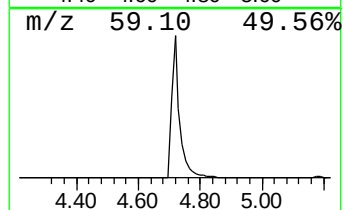
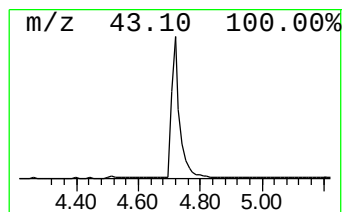
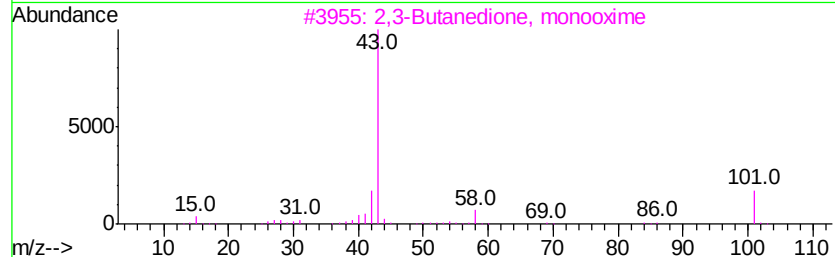
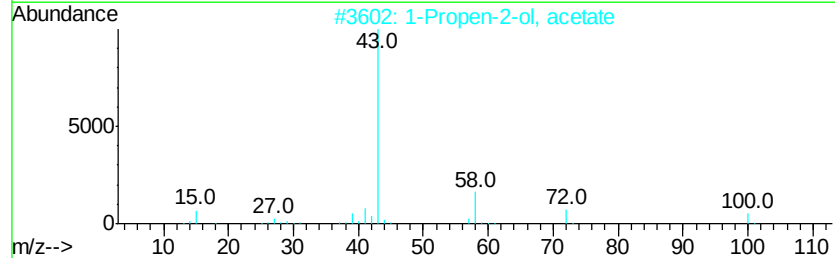
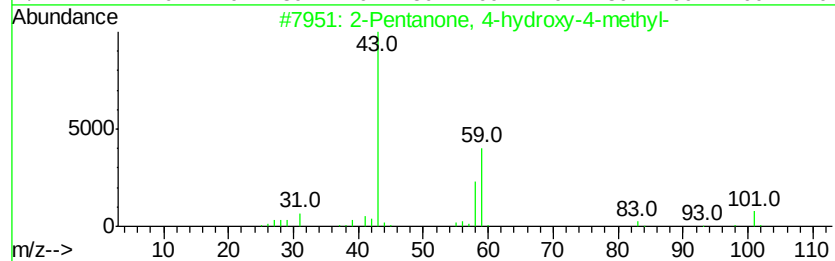
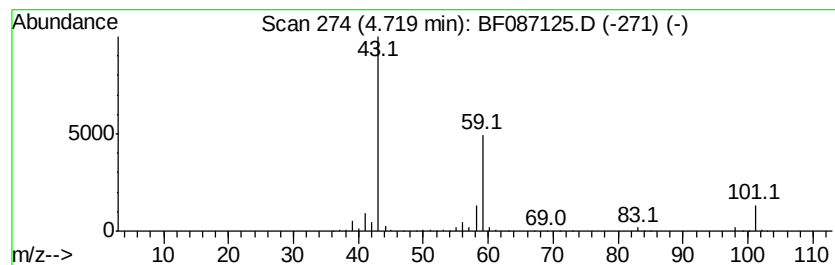
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.72	41.16 ng	1864290	1,4-Dichlorobenzene-d4	6.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		Acetic acid, 2-propenyl ester	100	C5H8O2	000591-87-7	9
5		2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9



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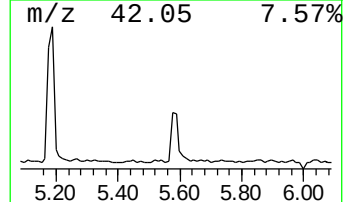
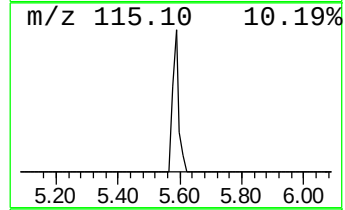
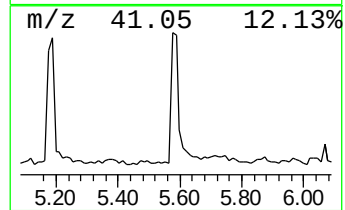
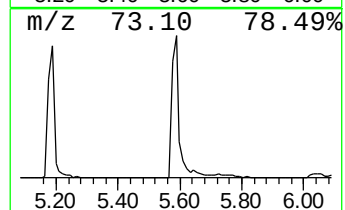
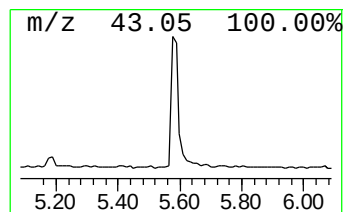
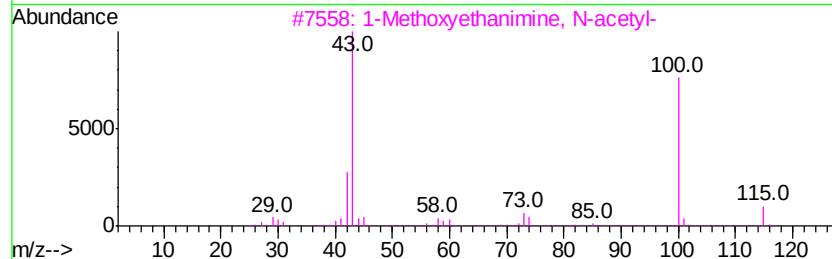
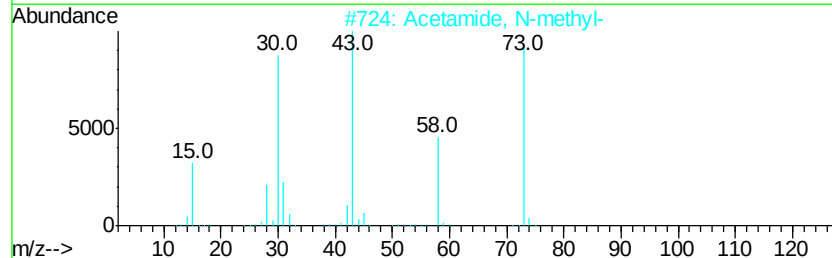
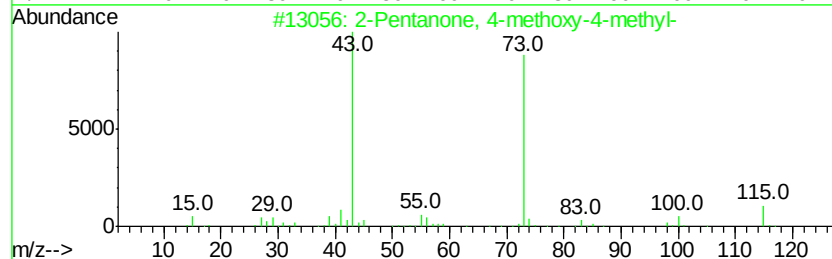
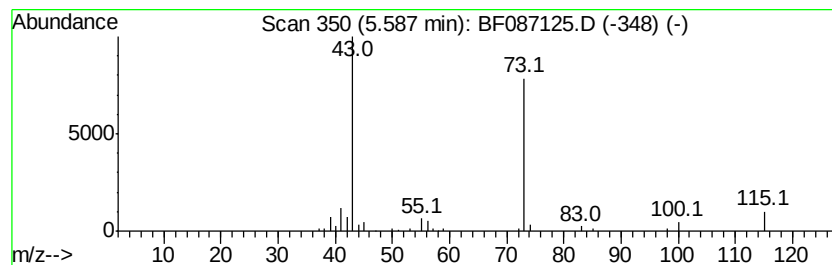
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2-Pentanone, 4-methoxy-4-me... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.59	3.28 ng	148760	1,4-Dichlorobenzene-d4	6.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	83
2		Acetamide, N-methyl-	73	C3H7NO	000079-16-3	43
3		1-Methoxyethanimine, N-acetyl-	115	C5H9NO2	099028-43-0	23
4		3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	12
5		2-Pentanone, 3-methyl-	100	C6H12O	000565-61-7	9



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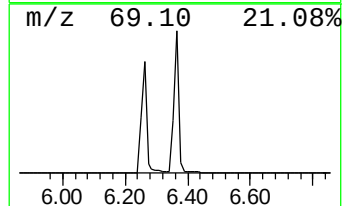
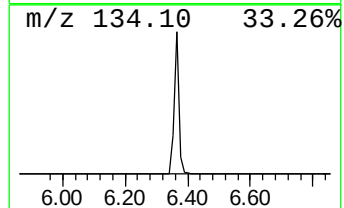
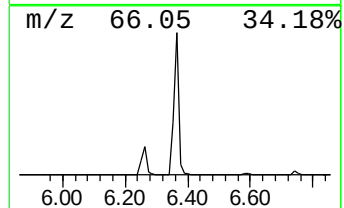
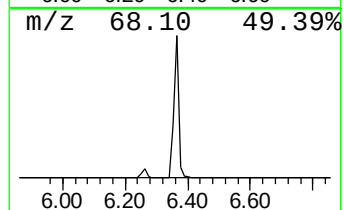
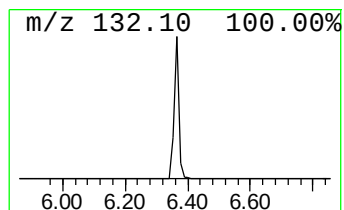
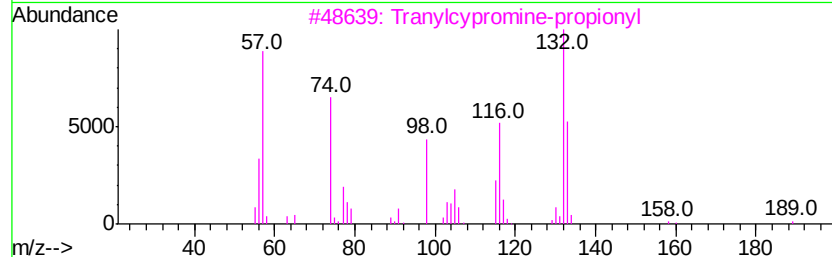
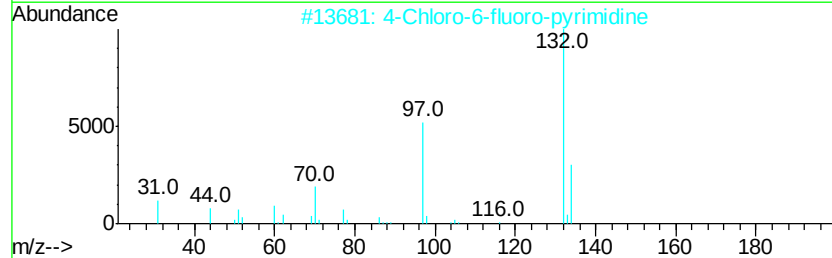
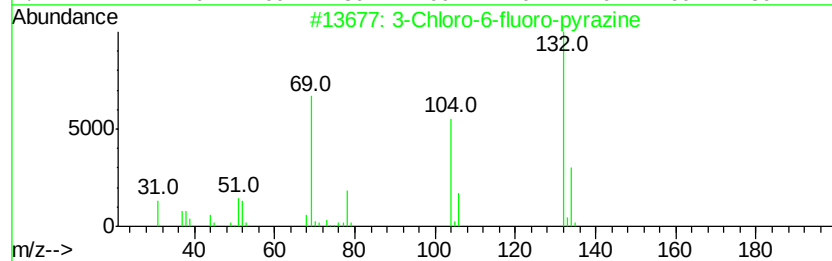
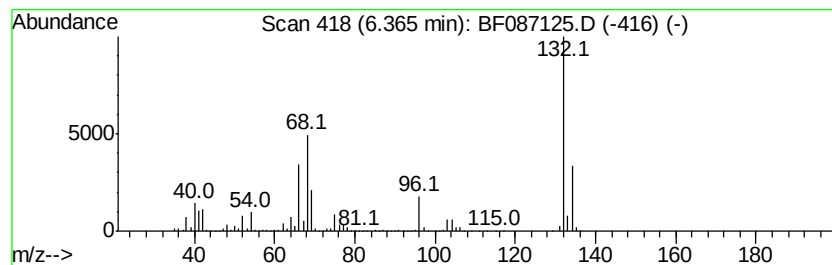
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown6.36 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.36	103.46 ng	4686190	1,4-Dichlorobenzene-d4	6.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	22
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16



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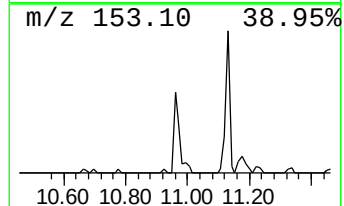
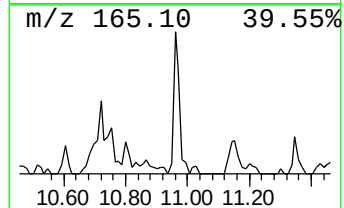
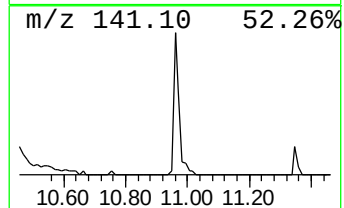
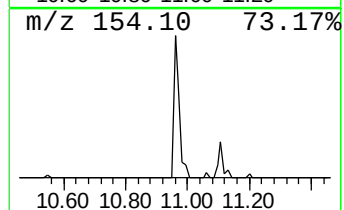
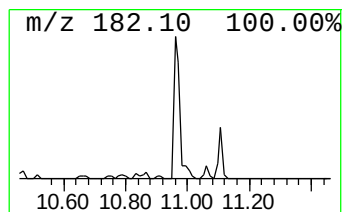
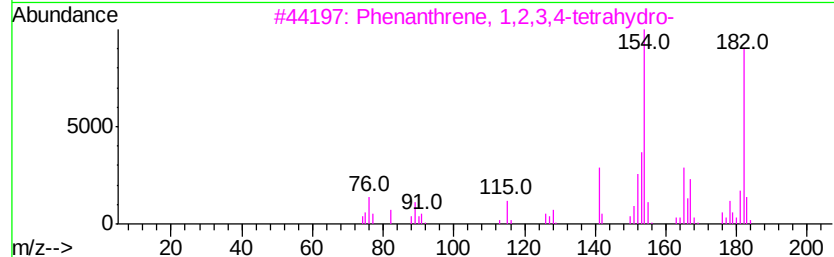
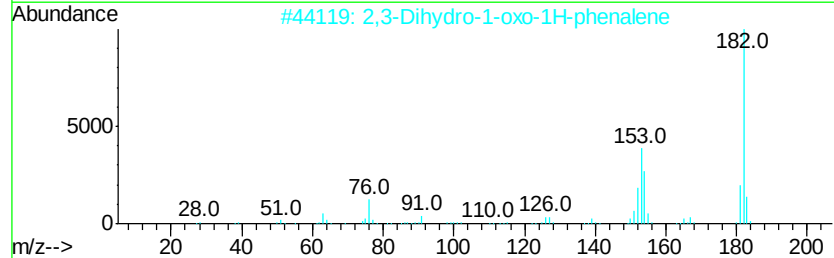
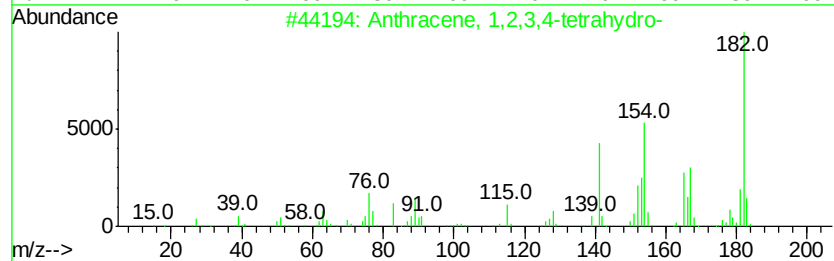
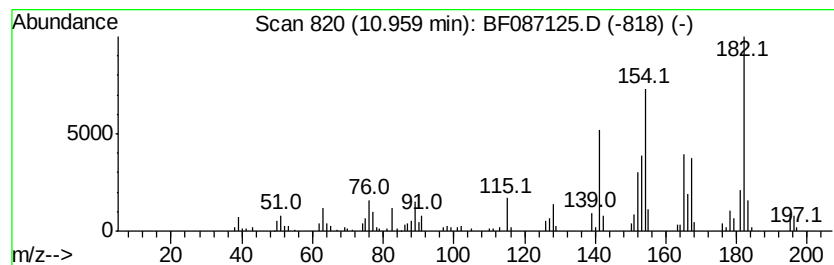
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Anthracene, 1,2,3,4-tetrahy... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.96	2.67 ng	225763	Phenanthrene-d10	11.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 1,2,3,4-tetrahydro-	182	C14H14	002141-42-6	98
2		2,3-Dihydro-1-oxo-1H-phenalene	182	C13H10O	000518-85-4	83
3		Phenanthrene, 1,2,3,4-tetrahydro-	182	C14H14	001013-08-7	83
4		3,3'-Dimethylbiphenyl	182	C14H14	000612-75-9	55
5		7-Phenylbicyclo[3.2.1]octa-2,6-d...	182	C14H14	1000201-01-7	52



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051716\
Data File : BF087125.D
Acq On : 18 May 2016 2:19
Operator : UM/SJ
Sample : H3085-01
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VITALE-FILL-CSPH3-10K

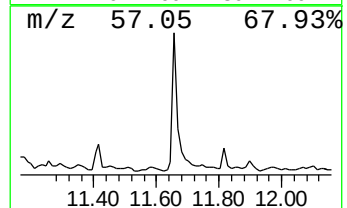
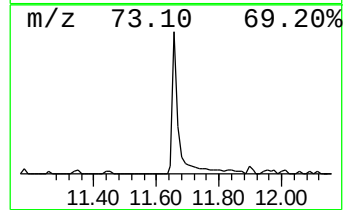
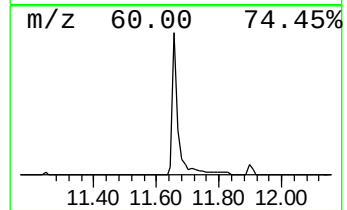
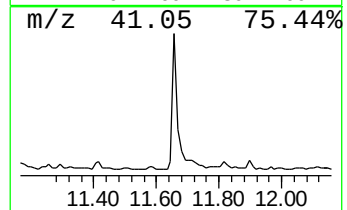
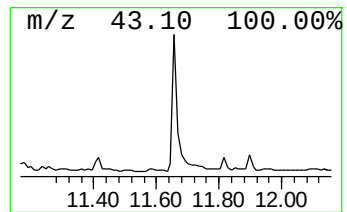
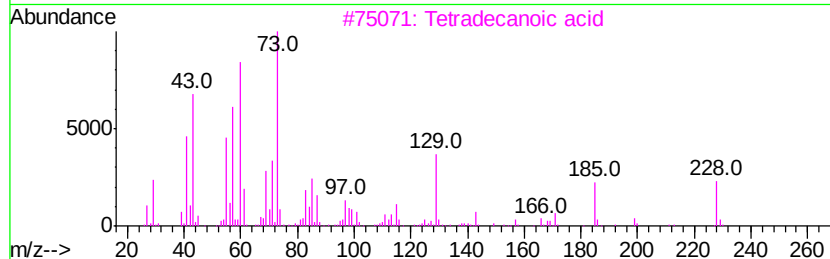
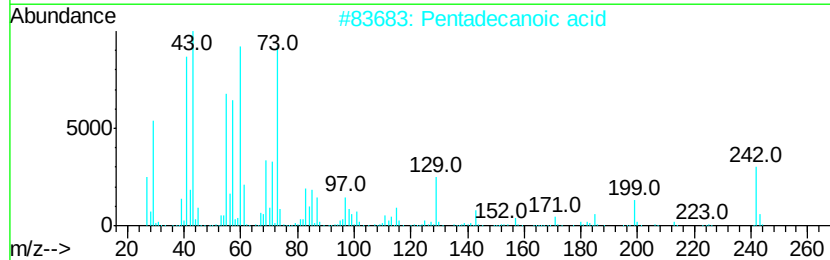
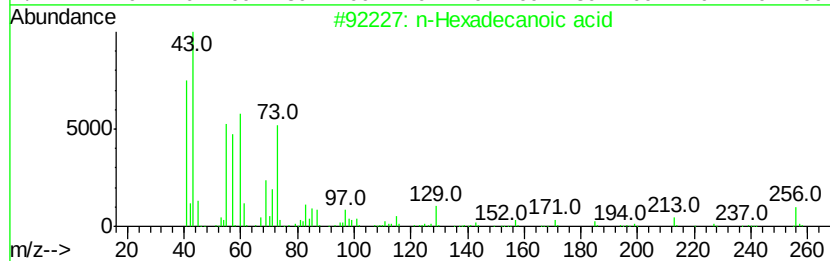
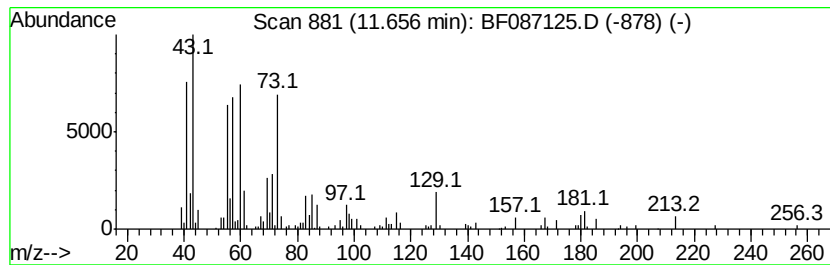
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	4.77 ng	403262	Phenanthrene-d10	11.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
4		Undecanoic acid	186	C11H22O2	000112-37-8	80
5		Tridecanoic acid	214	C13H26O2	000638-53-9	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051716\
Data File : BF087125.D
Acq On : 18 May 2016 2:19
Operator : UM/SJ
Sample : H3085-01
Misc :
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ClientSampleId :
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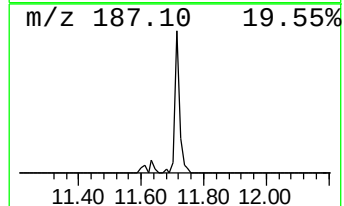
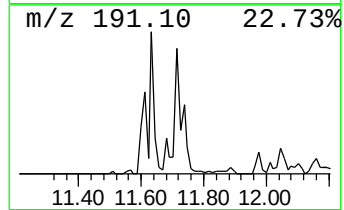
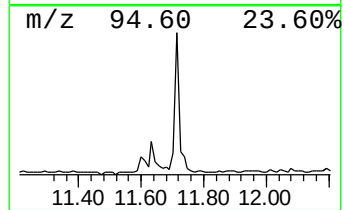
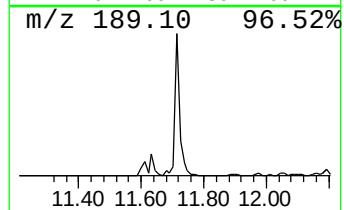
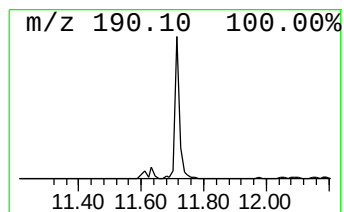
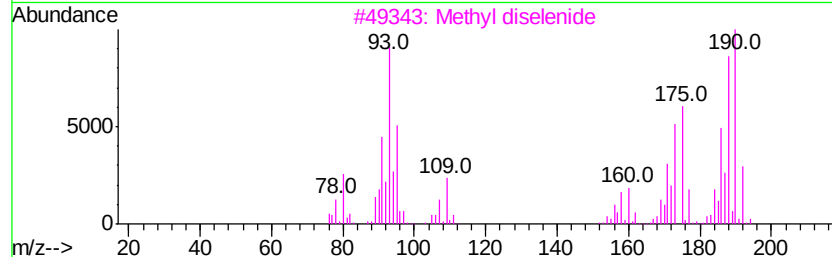
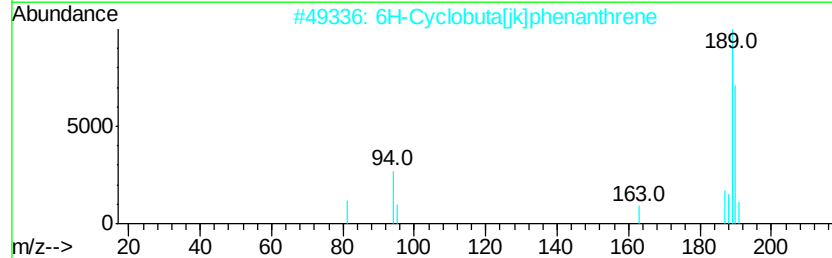
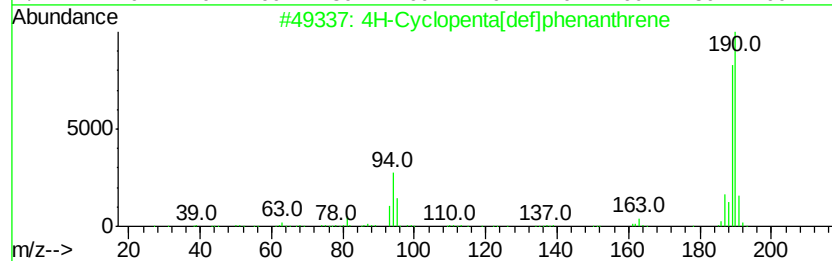
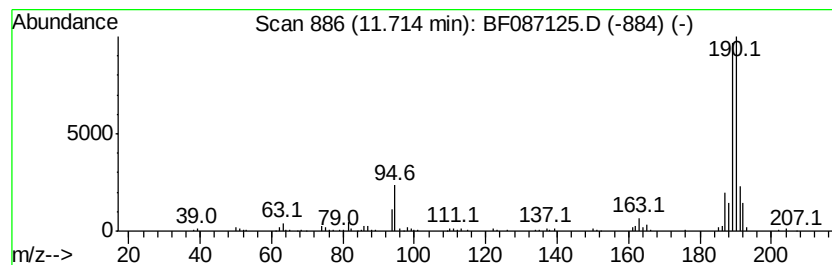
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.71	4.53 ng	382605	Phenanthrene-d10	11.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	94
2			6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	72
3			Methyl diselenide	190	C2H6Se2	007101-31-7	49
4			1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	25
5			2-Naphthaldehyde, 1,2,3,4-tetra...	190	C12H14O2	002472-02-8	14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051716\
Data File : BF087125.D
Acq On : 18 May 2016 2:19
Operator : UM/SJ
Sample : H3085-01
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
VITALE-FILL-CSPH3-10K

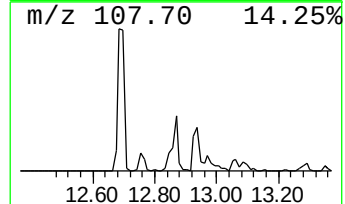
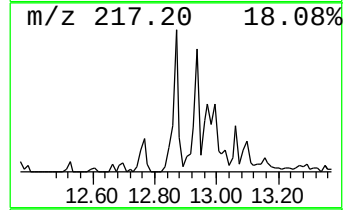
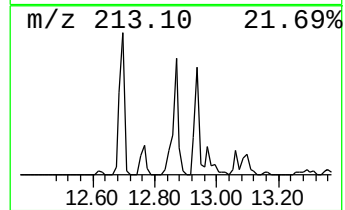
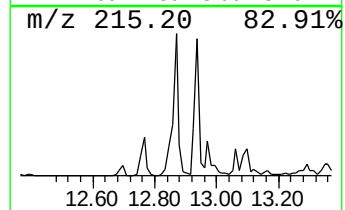
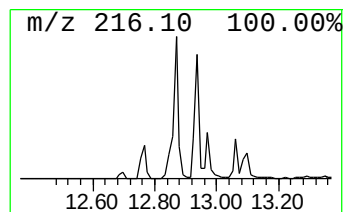
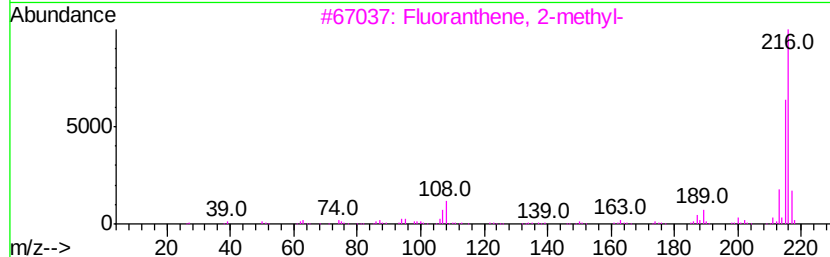
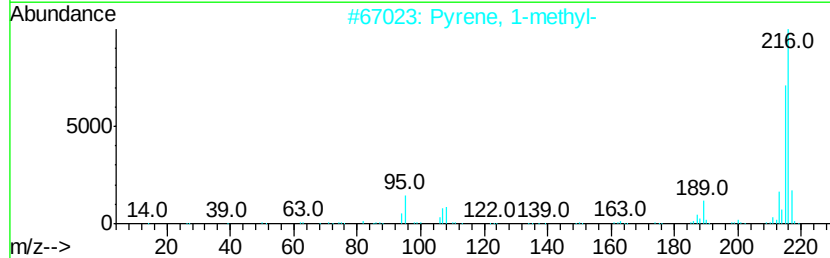
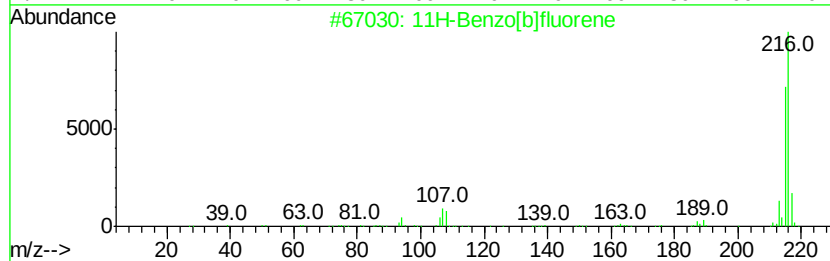
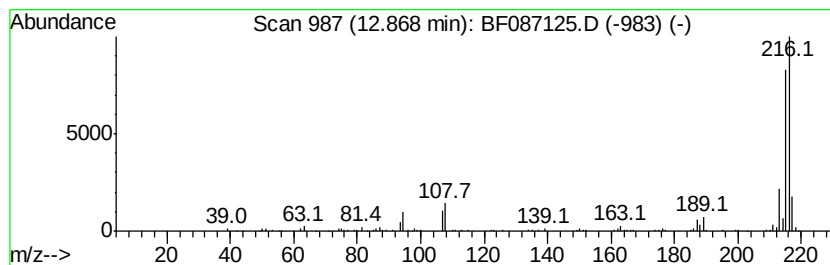
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 11H-Benzo[b]fluorene Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.87	2.46 ng	365848	Chrysene-d12	13.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
3			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	93
4			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	92
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	83



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051716\
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Sample : H3085-01
Misc :
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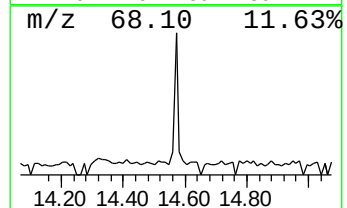
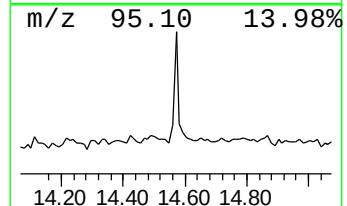
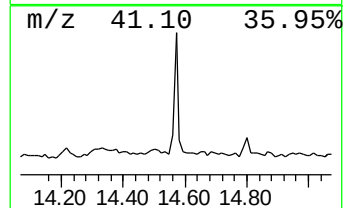
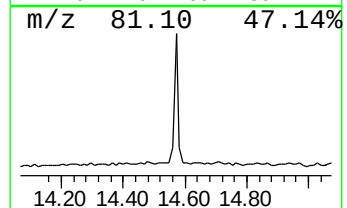
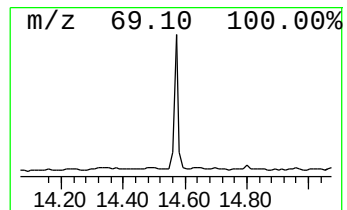
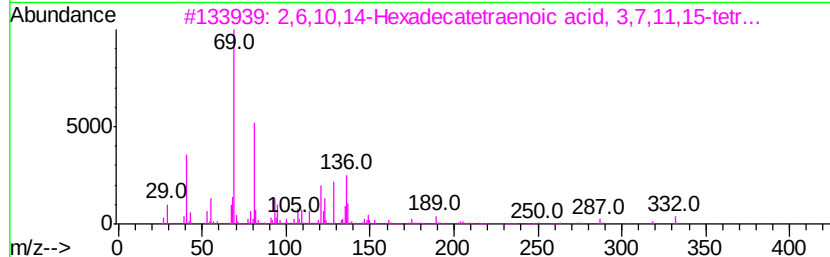
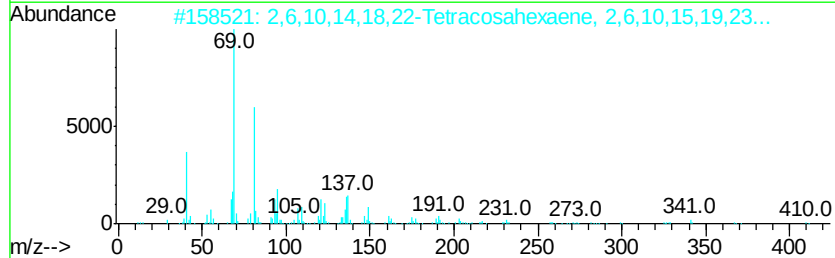
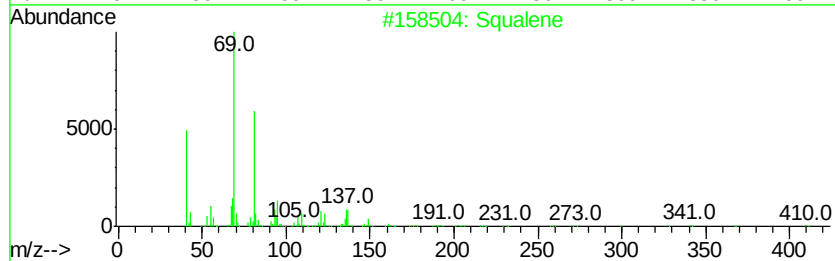
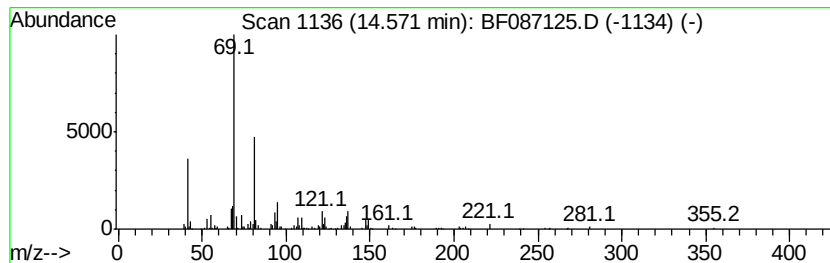
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ClientSampleId :
VITALE-FILL-CSPH3-10K

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF051716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Squalene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	5.04 ng	250195	Perylene-d12	15.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Squalene	410 C30H50	007683-64-9	91
2	2,6,10,14,18,22-Tetracosahexaene...	410 C30H50	000111-02-4	90
3	2,6,10,14-Hexadecatetraenoic aci...	332 C22H36O2	024035-35-6	72
4	4,8,12-Tetradecatrienal, 5,9,13-...	248 C17H28O	066408-55-7	72
5	1,5-Heptadiene, 3,3,6-trimethyl-	138 C10H18	035387-63-4	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051716\
Data File : BF087125.D
Acq On : 18 May 2016 2:19
Operator : UM/SJ
Sample : H3085-01
Misc :
ALS Vial : 27 Sample Multiplier: 1

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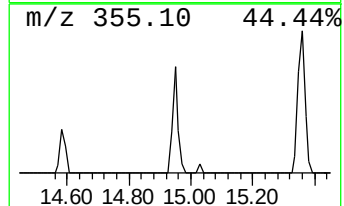
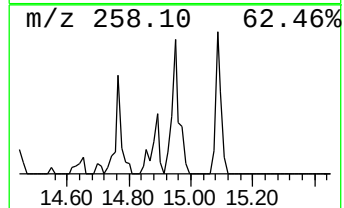
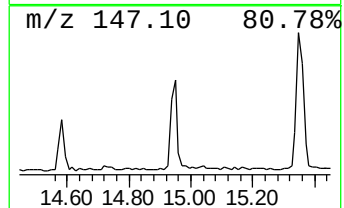
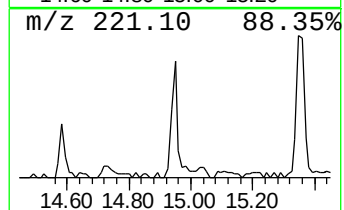
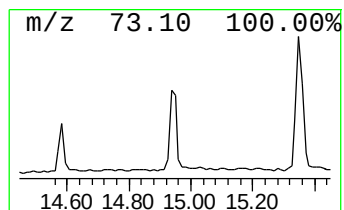
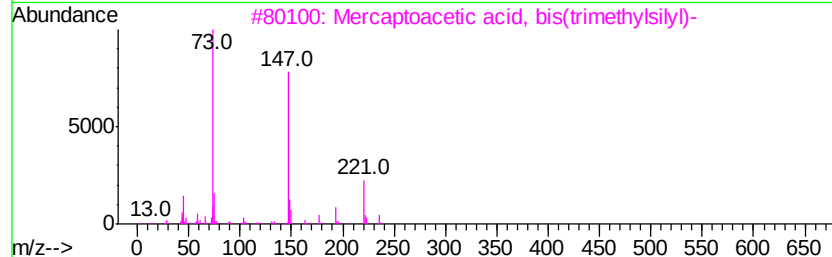
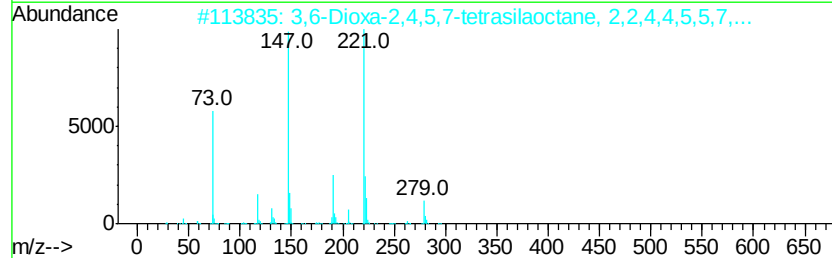
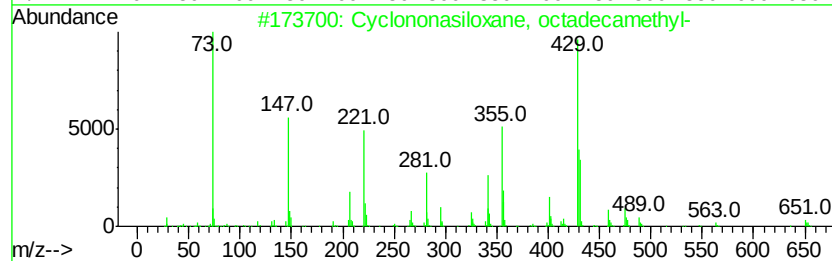
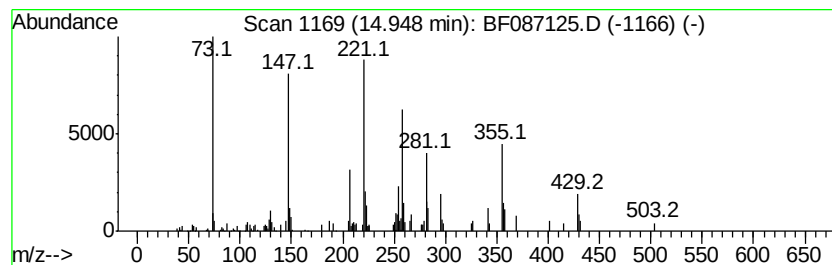
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown14.95 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.95	2.54 ng	126317	Perylene-d12	15.09

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	45
2			3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	32
3			Mercaptoacetic acid, bis(trimeth...	236	C8H20O2SSi2	006398-62-5	32
4			N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	30
5			Pentasiloxane, dodecamethyl-	384	C12H36O4Si5	000141-63-9	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051716\
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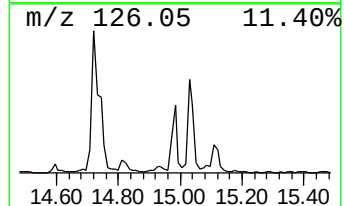
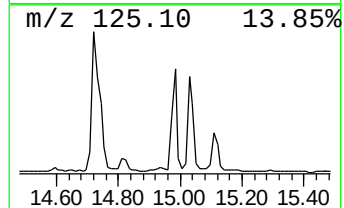
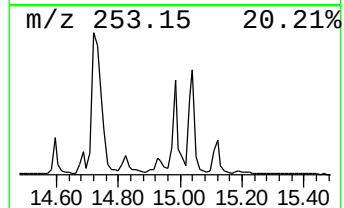
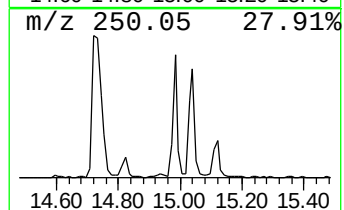
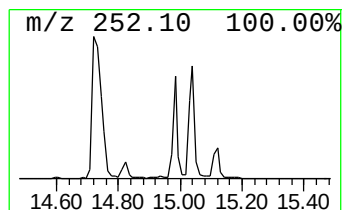
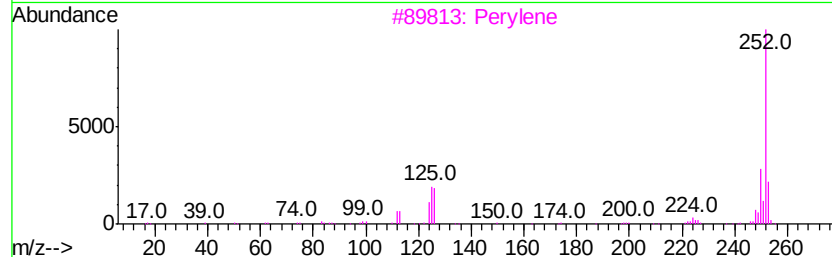
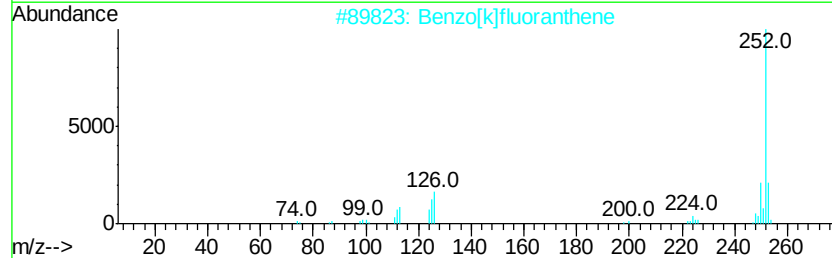
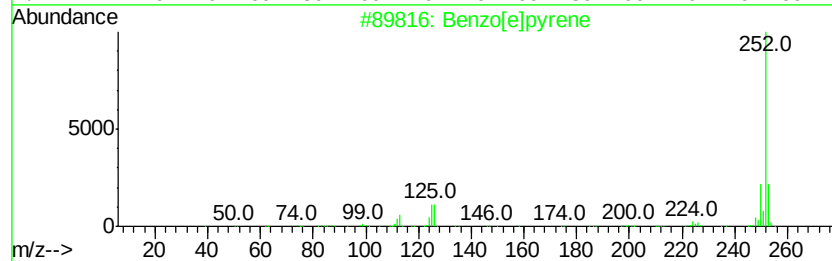
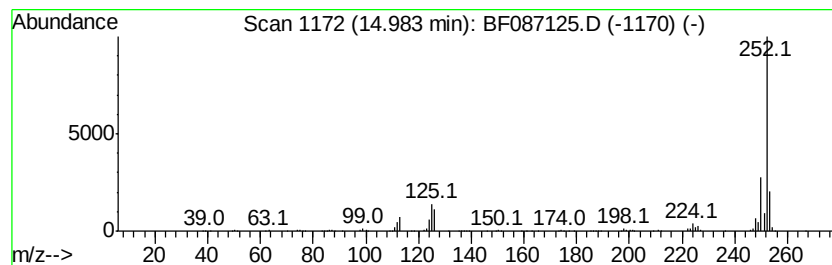
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Benzo[e]pyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.98	8.55 ng	424914	Perylene-d12	15.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	99
2		Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3		Perylene	252	C20H12	000198-55-0	96
4		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	95
5		Benzo[j]fluoranthene	252	C20H12	000205-82-3	94



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051716\
Data File : BF087125.D
Acq On : 18 May 2016 2:19
Operator : UM/SJ
Sample : H3085-01
Misc :
ALS Vial : 27 Sample Multiplier: 1

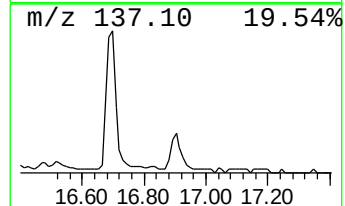
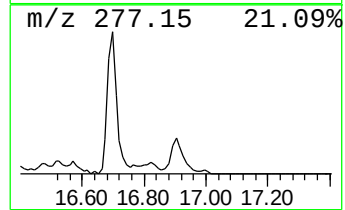
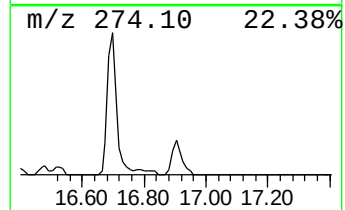
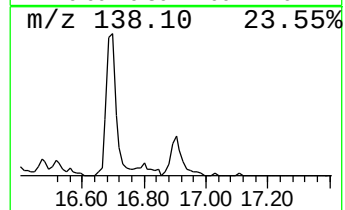
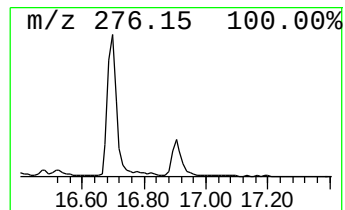
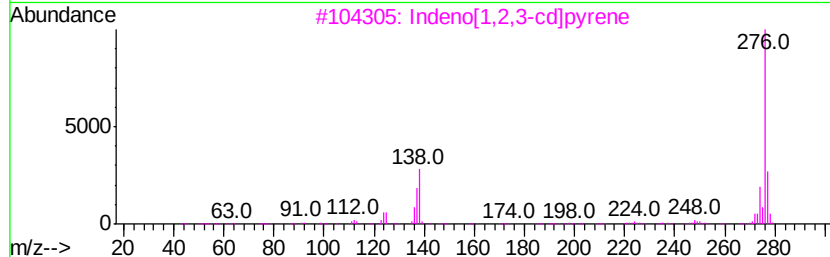
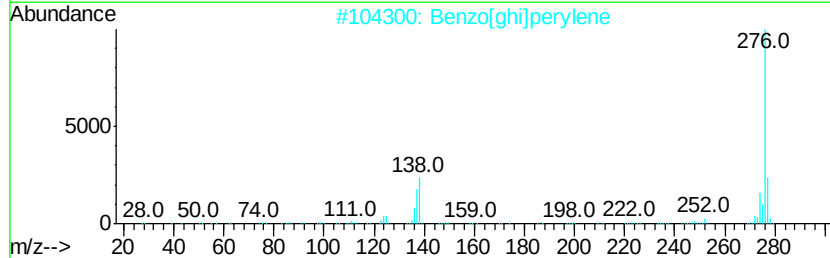
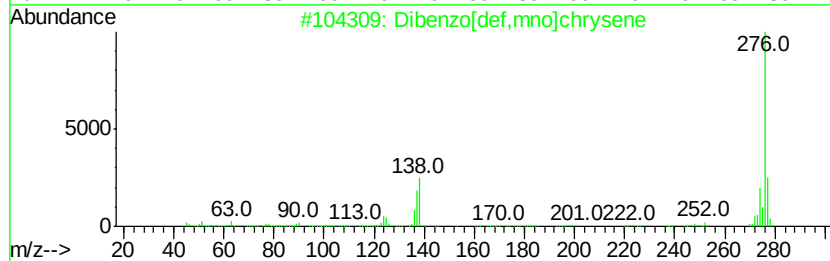
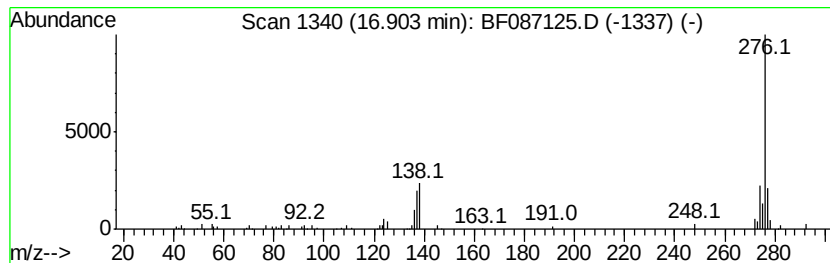
Instrument :
BNA_F
ClientSampleId :
VITALE-FILL-CSPH3-10K

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF051716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Dibenzo[def,mno]chrysene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.90	2.76 ng	137117	Perylene-d12	15.09	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzo[def,mno]chrysene	276	C22H12	000191-26-4	96
2	Benzo[ghi]perylene	276	C22H12	000191-24-2	95
3	Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	91
4	Indeno[1,2,3-cd]fluoranthene	276	C22H12	000193-43-1	87
5	Naphthalen-1-yl-(3-nitro-benzyl)...	276	C17H12N2O2	200421-53-0	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051716\
 Data File : BF087125.D
 Acq On : 18 May 2016 2:19
 Operator : UM/SJ
 Sample : H3085-01
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VITALE-FILL-CSPH3-10K

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF051716.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
unknown3.96	3.96	3.3	ng	147097	1	6.59	905904	20.0
2-Pentanone, 4-hy...	4.72	41.2	ng	1864290	1	6.59	905904	20.0
2-Pentanone, 4-me...	5.59	3.3	ng	148760	1	6.59	905904	20.0
unknown6.36	6.36	103.5	ng	4686190	1	6.59	905904	20.0
Anthracene, 1,2,3...	10.96	2.7	ng	225763	4	11.11	1689430	20.0
n-Hexadecanoic acid	11.66	4.8	ng	403262	4	11.11	1689430	20.0
4H-Cyclopenta[def...	11.71	4.5	ng	382605	4	11.11	1689430	20.0
11H-Benzo[b]fluorene	12.87	2.5	ng	365848	5	13.74	2972880	20.0
Squalene	14.57	5.0	ng	250195	6	15.09	993717	20.0
unknown14.95	14.95	2.5	ng	126317	6	15.09	993717	20.0
Benzo[e]pyrene	14.98	8.6	ng	424914	6	15.09	993717	20.0
Dibenzo[def,mno]c...	16.90	2.8	ng	137117	6	15.09	993717	20.0