

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032318\
 Data File : BF103876.D
 Acq On : 23 Mar 2018 13:27
 Operator : JU/SJ
 Sample : J2002-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-4-MH-B

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-BF031518.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	2.316	33	39	47	rVB	138893	179037	1.01%	0.297%
2	4.163	339	353	356	rBV	7011203	17653883	100.00%	29.296%
3	5.222	530	533	542	rBV	150179	199830	1.13%	0.332%
4	5.604	592	598	601	rBV	3564596	3725581	21.10%	6.182%
5	6.604	761	768	770	rBV	3219895	3743081	21.20%	6.211%
6	6.745	787	792	795	rBV	3791280	3845239	21.78%	6.381%
7	6.975	827	831	834	rBV	1151391	909507	5.15%	1.509%
8	7.133	853	858	861	rBV	3020281	2759431	15.63%	4.579%
9	7.539	922	927	930	rBV	2518929	2454030	13.90%	4.072%
10	8.257	1045	1049	1051	rBV	1474758	1210552	6.86%	2.009%
11	9.327	1227	1231	1235	rBV	3312830	3123365	17.69%	5.183%
12	9.722	1295	1298	1303	rVB	469167	420331	2.38%	0.698%
13	10.016	1344	1348	1351	rBV2	1521497	1280653	7.25%	2.125%
14	10.810	1478	1483	1486	rBV	2539361	2445446	13.85%	4.058%
15	10.904	1496	1499	1509	rVB2	139981	197149	1.12%	0.327%
16	11.510	1598	1602	1604	rBV	1362007	1181826	6.69%	1.961%
17	11.533	1604	1606	1609	rVV	1269095	1005823	5.70%	1.669%
18	11.586	1609	1615	1617	rVV	327429	312766	1.77%	0.519%
19	12.010	1684	1687	1690	rBV	240396	275469	1.56%	0.457%
20	12.039	1690	1692	1695	rVB	305922	233454	1.32%	0.387%
21	12.127	1703	1707	1709	rBV2	337694	401308	2.27%	0.666%
22	12.568	1779	1782	1786	rBV2	155517	237010	1.34%	0.393%
23	12.739	1807	1811	1813	rVV	1277856	1162322	6.58%	1.929%
24	12.763	1813	1815	1818	rVB2	395367	310744	1.76%	0.516%
25	12.963	1842	1849	1854	rVV	1296174	1429828	8.10%	2.373%
26	13.098	1868	1872	1875	rBV	2313723	2087021	11.82%	3.463%
27	13.286	1901	1904	1906	rVB	251709	216835	1.23%	0.360%
28	13.686	1968	1972	1974	rBV2	379350	354243	2.01%	0.588%
29	13.974	2018	2021	2025	rBV3	563654	529780	3.00%	0.879%
30	14.062	2033	2036	2040	rBV3	249020	256445	1.45%	0.426%
31	14.104	2040	2043	2046	rVV	886359	727309	4.12%	1.207%
32	14.162	2046	2053	2056	rVV2	1133075	1657110	9.39%	2.750%
33	14.192	2056	2058	2062	rVV	571197	470270	2.66%	0.780%
34	14.621	2128	2131	2135	rBV4	170201	210710	1.19%	0.350%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.251	2233	2238	2241	rBV	402829	656923	3.72%	1.090%
36	15.574	2289	2293	2300	rVB	283979	372027	2.11%	0.617%
37	15.639	2300	2304	2310	rVB	385233	507390	2.87%	0.842%
38	15.709	2311	2316	2320	rBV	717709	975239	5.52%	1.618%
39	17.286	2580	2584	2591	rVB2	125734	222070	1.26%	0.369%
40	17.762	2661	2665	2676	rVB	141803	319491	1.81%	0.530%

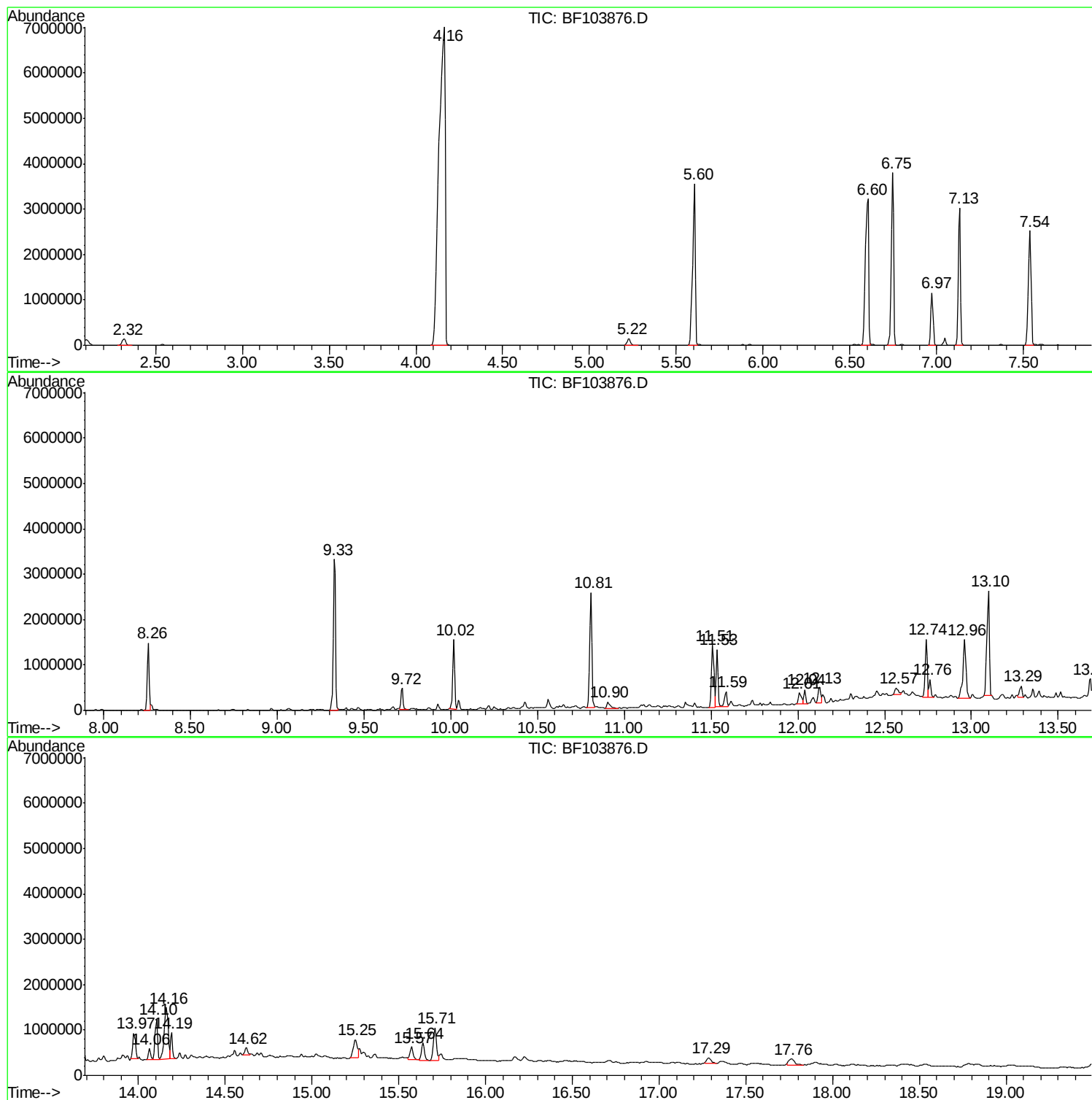
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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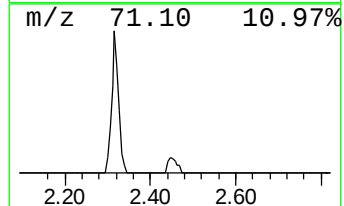
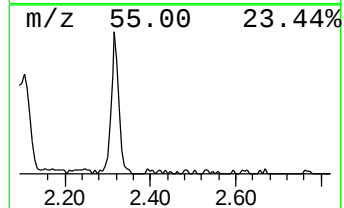
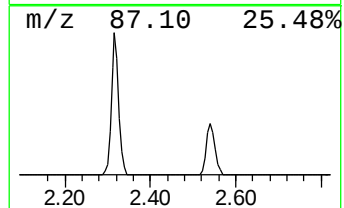
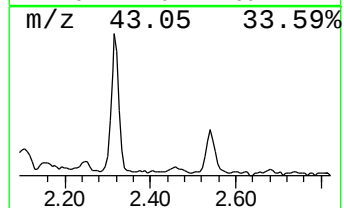
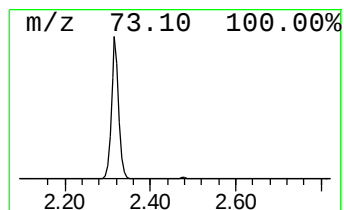
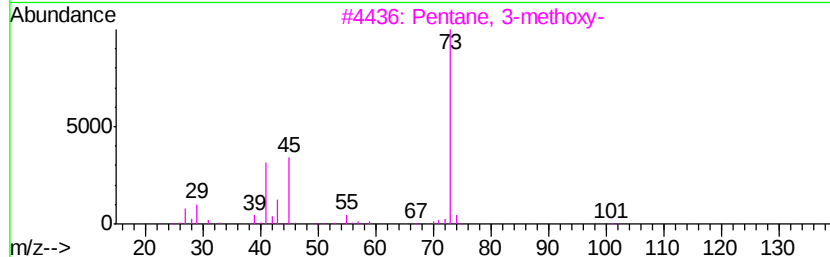
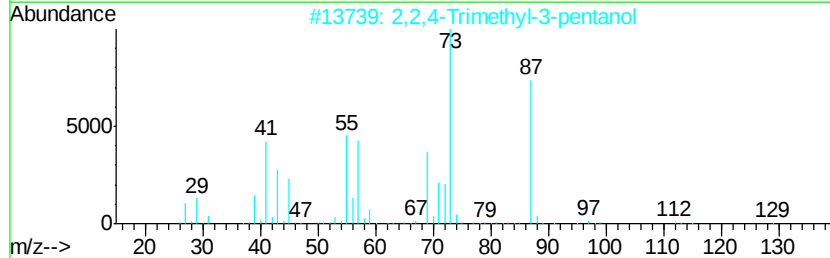
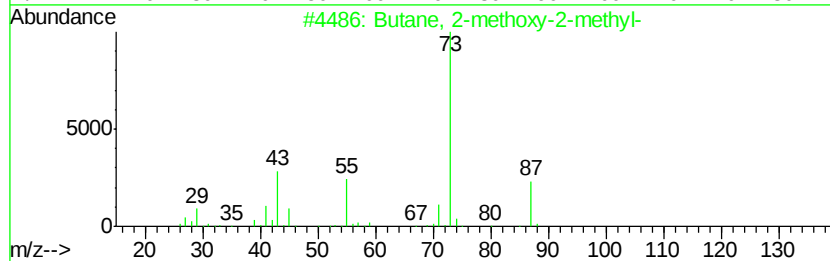
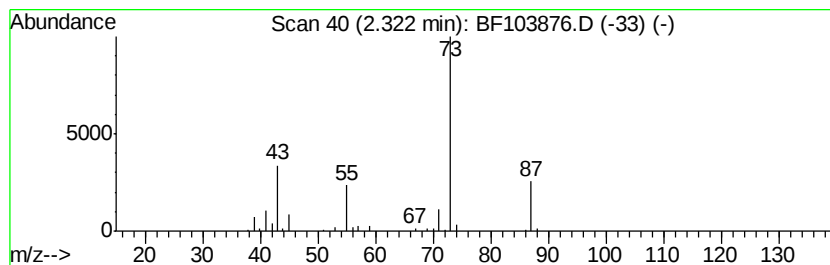
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.32	3.94 ng	179037	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	17
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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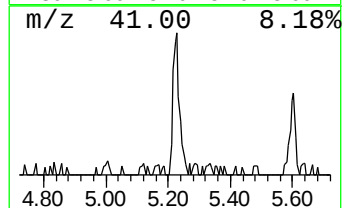
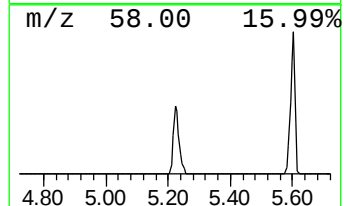
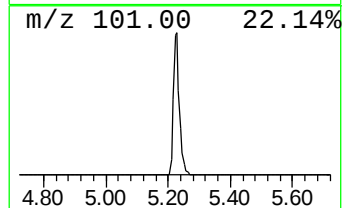
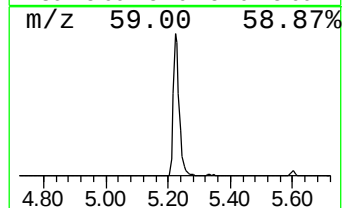
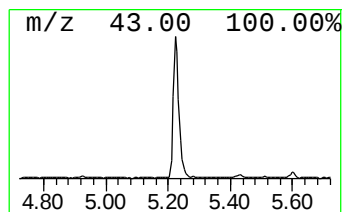
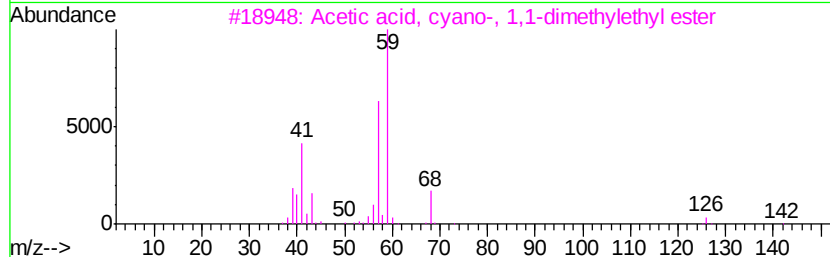
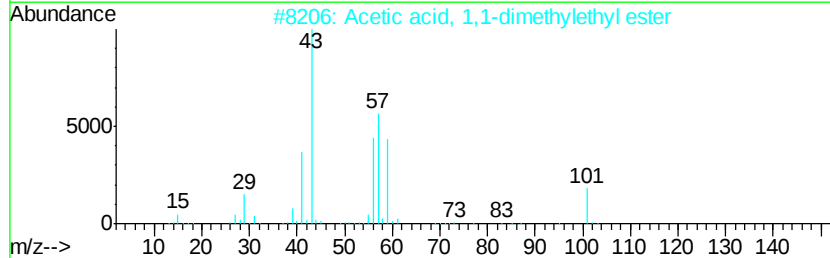
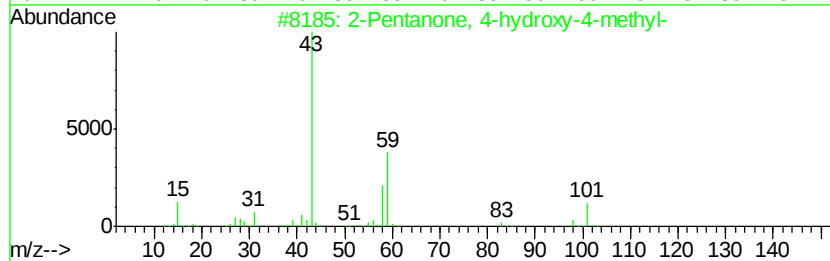
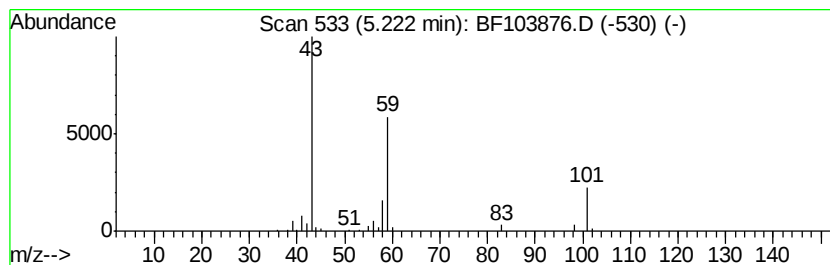
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	4.39 ng	199830	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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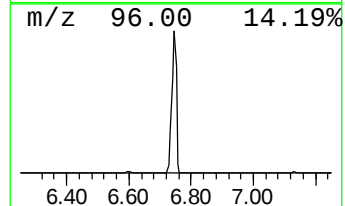
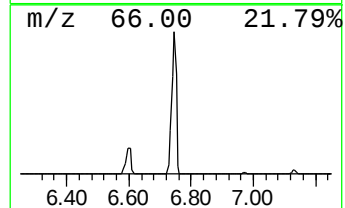
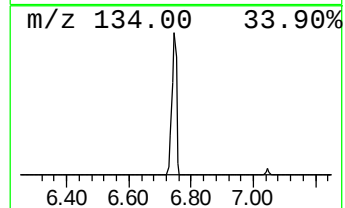
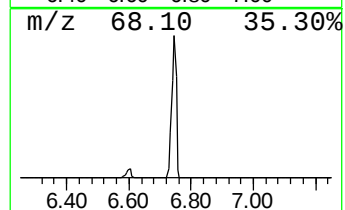
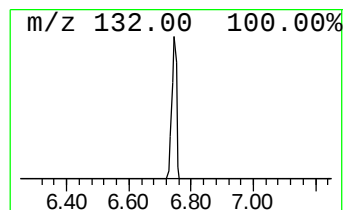
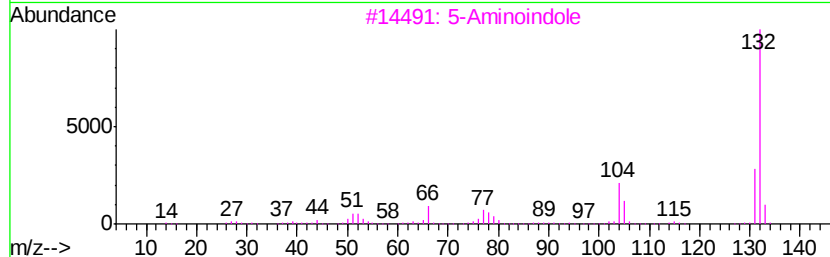
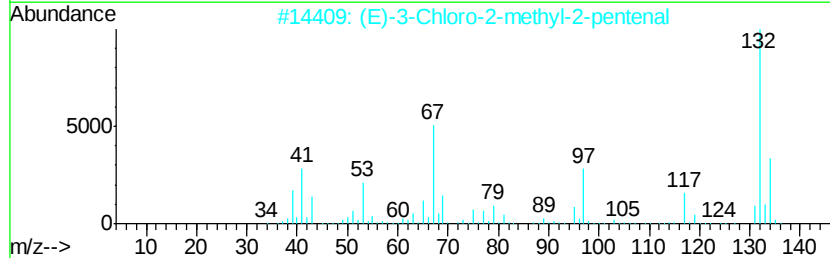
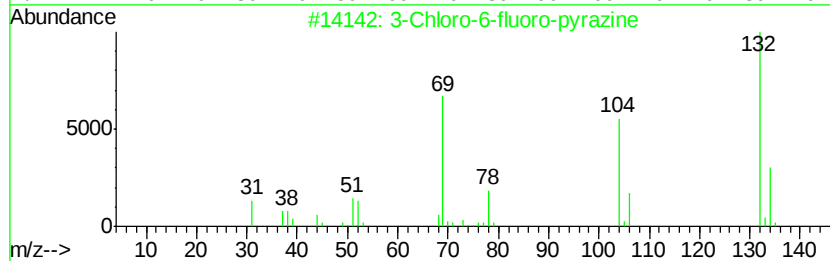
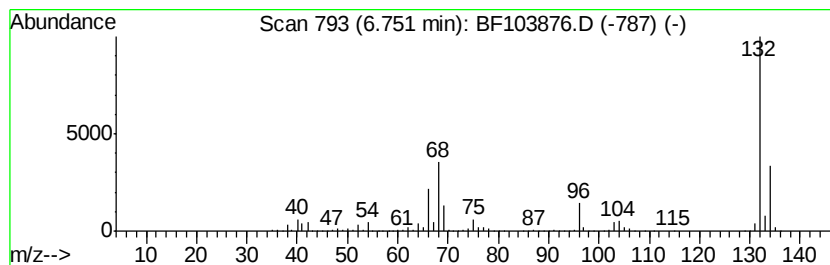
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown6.75 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	84.56 ng	3845240	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	9



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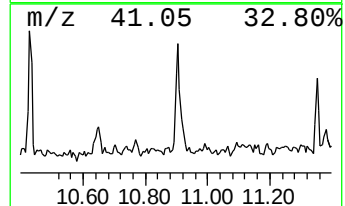
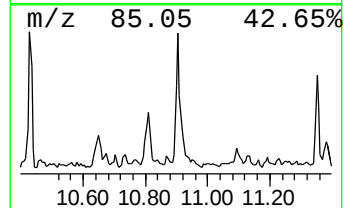
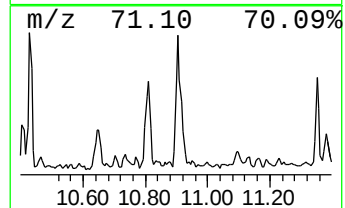
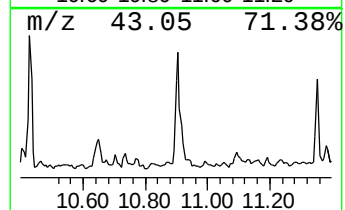
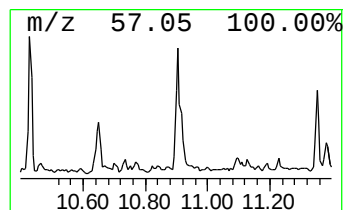
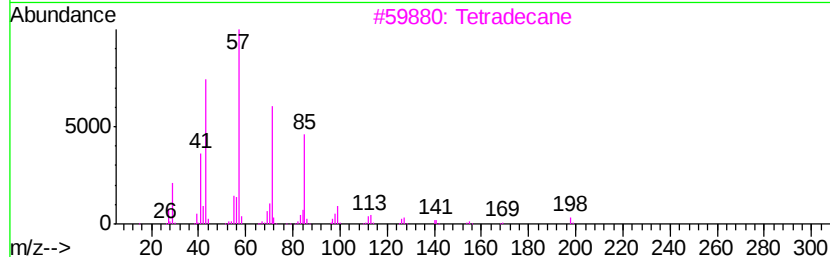
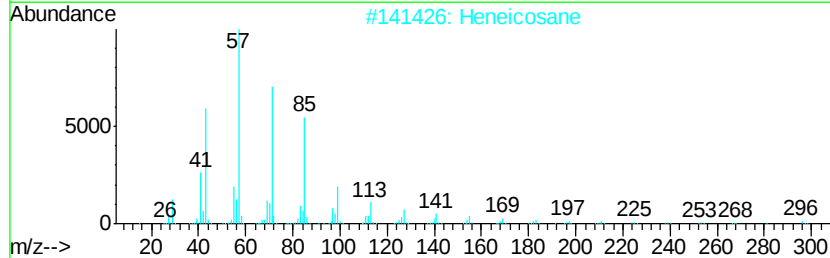
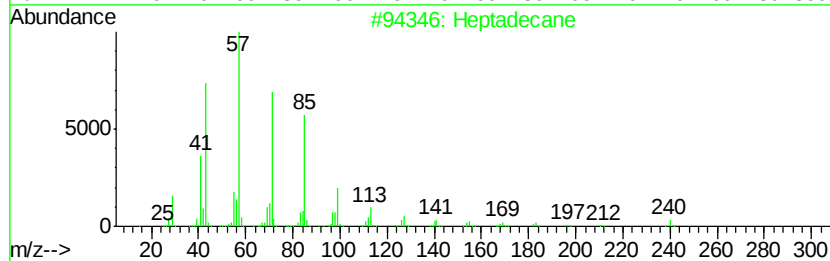
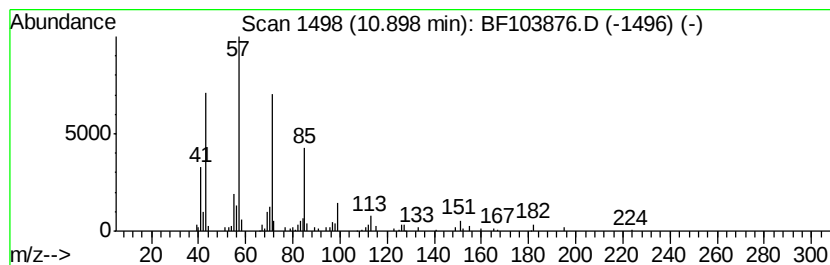
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Heptadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.90	3.34 ng	197149	Phenanthrene-d10		11.51
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	83
2	Heneicosane	296	C21H44	000629-94-7	83
3	Tetradecane	198	C14H30	000629-59-4	83
4	Hexadecane	226	C16H34	000544-76-3	80
5	Sulfurous acid, 2-propyl tetrade...	320	C17H36O3S	1000309-12-5	80



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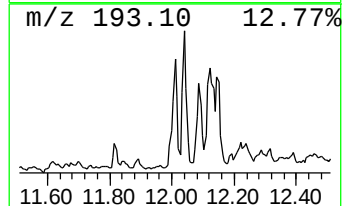
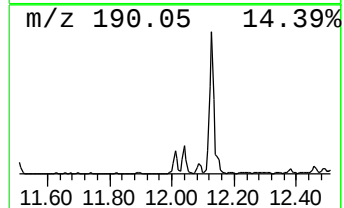
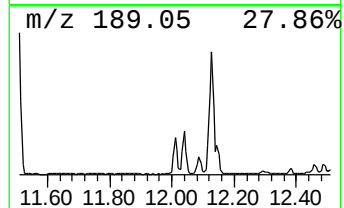
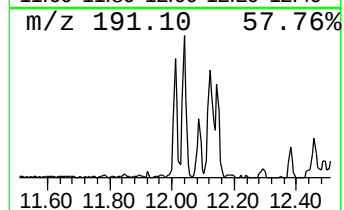
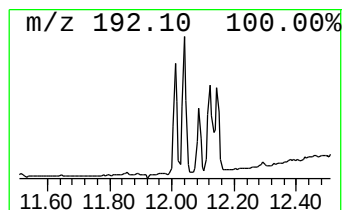
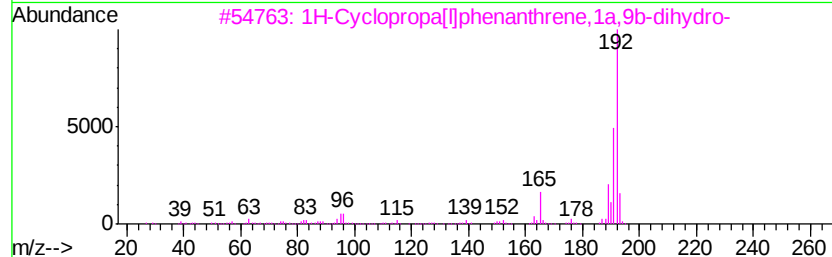
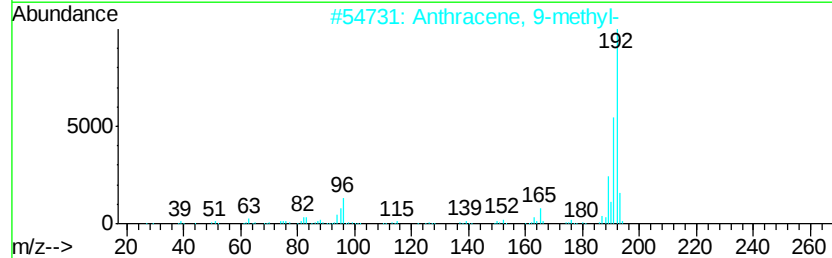
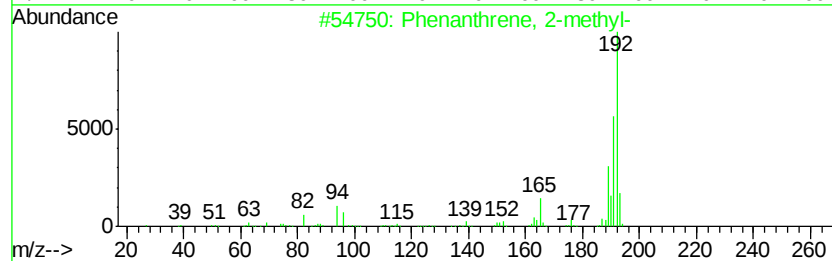
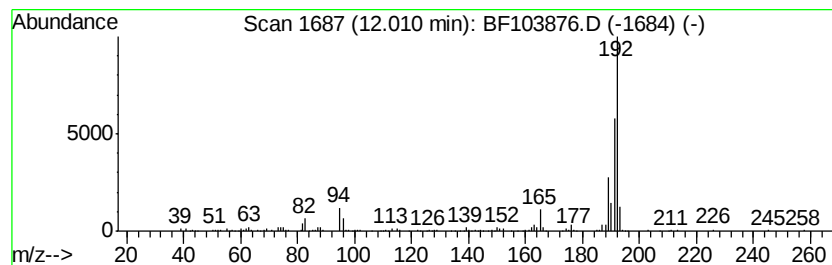
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenanthrene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	4.66 ng	275469	Phenanthrene-d10	11.51

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
2			Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
3			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	89
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	89
5			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	81



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032318\
Data File : BF103876.D
Acq On : 23 Mar 2018 13:27
Operator : JU/SJ
Sample : J2002-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-4-MH-B

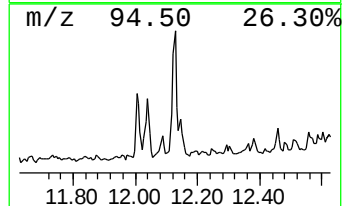
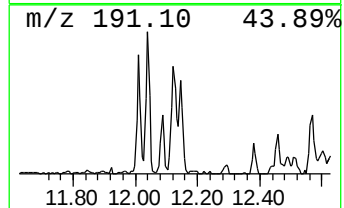
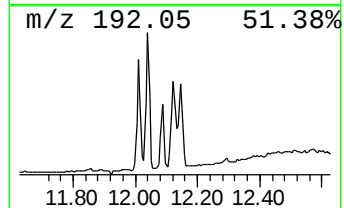
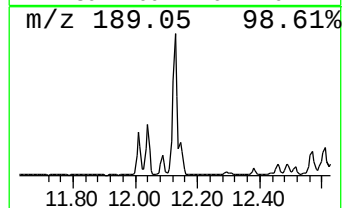
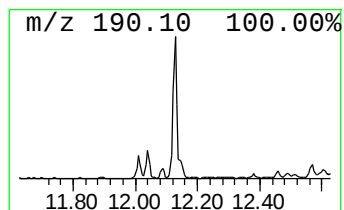
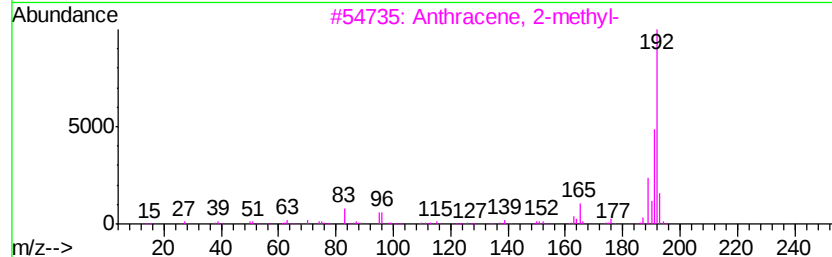
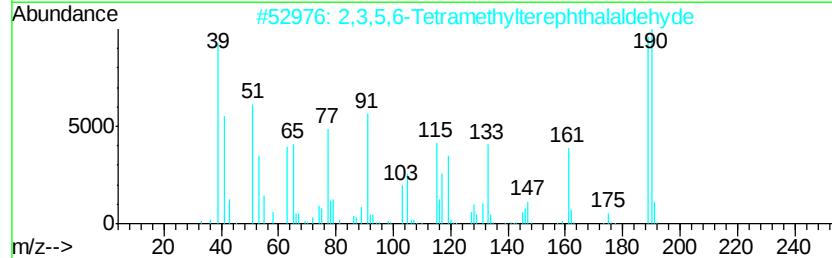
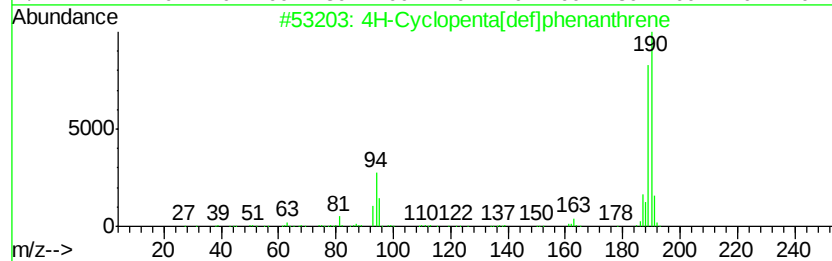
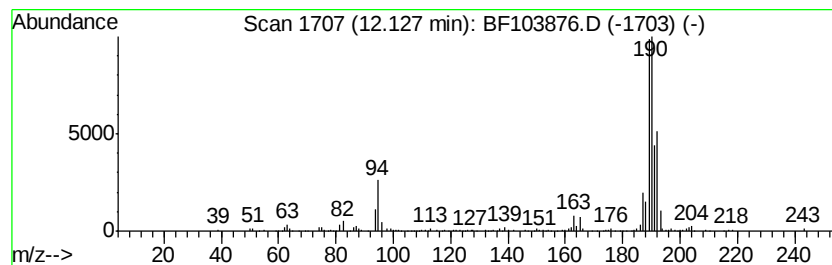
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.13	6.79 ng	401308	Phenanthrene-d10	11.51

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	92
2		2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	43
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	25
4		5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	18
5		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	18



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032318\
Data File : BF103876.D
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Sample : J2002-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

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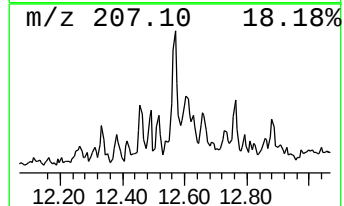
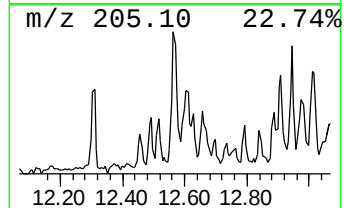
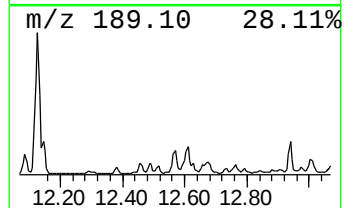
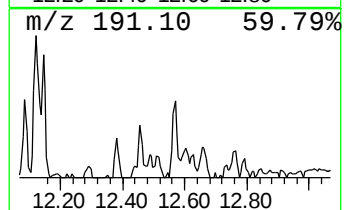
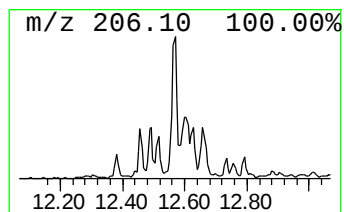
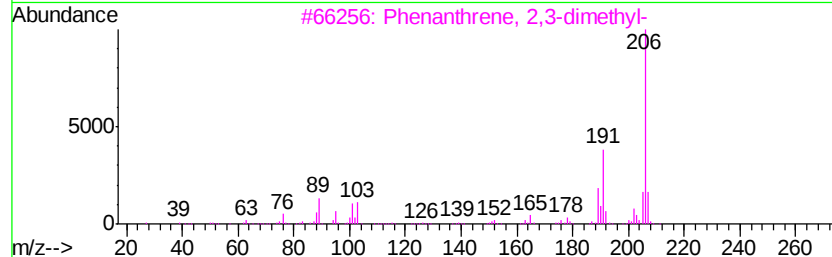
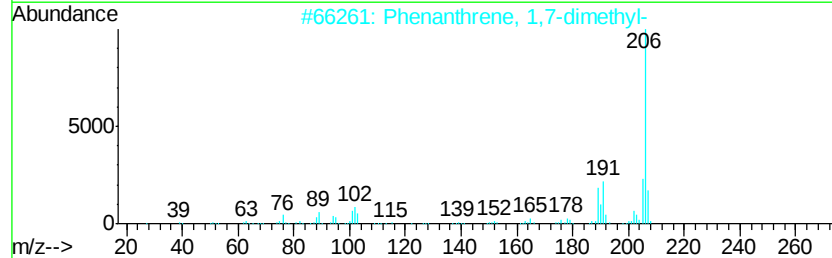
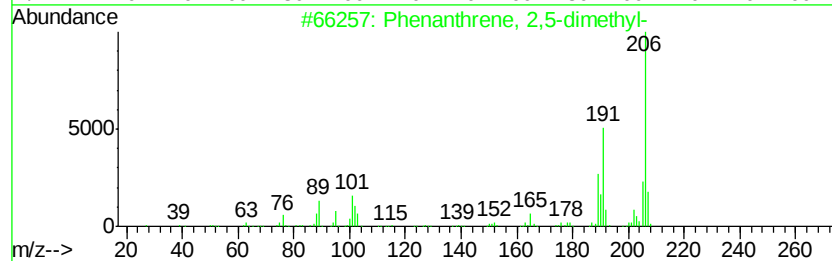
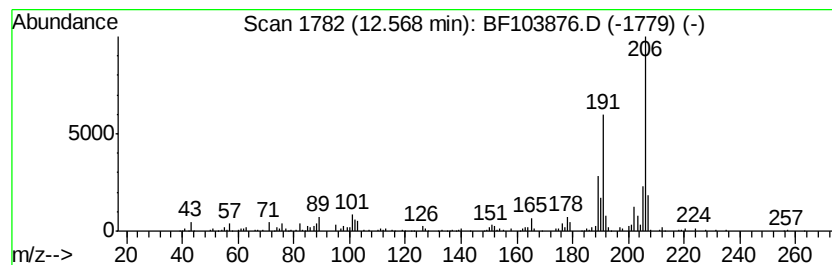
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Phenanthrene, 2,5-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.57	4.01 ng	237010	Phenanthrene-d10	11.51

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	89
3			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	87
4			Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	87
5			Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032318\
Data File : BF103876.D
Acq On : 23 Mar 2018 13:27
Operator : JU/SJ
Sample : J2002-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

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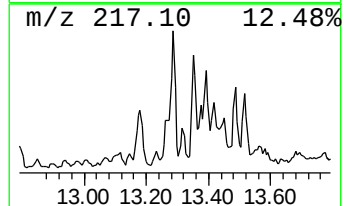
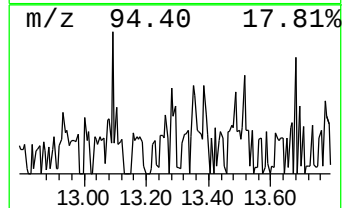
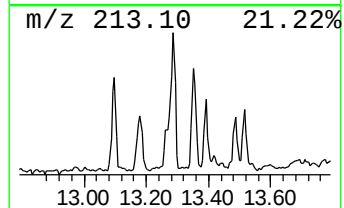
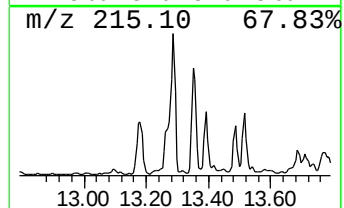
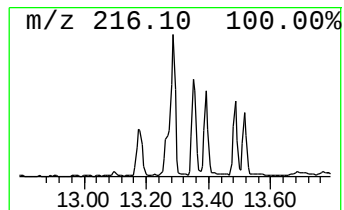
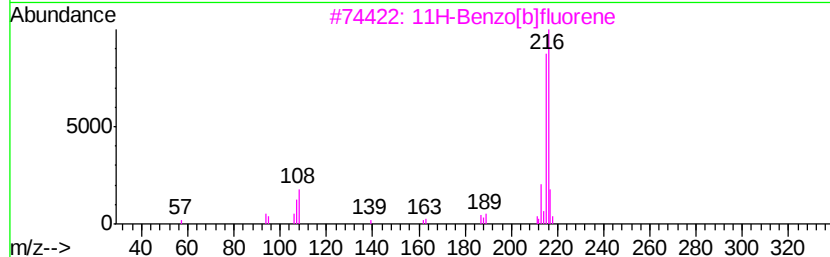
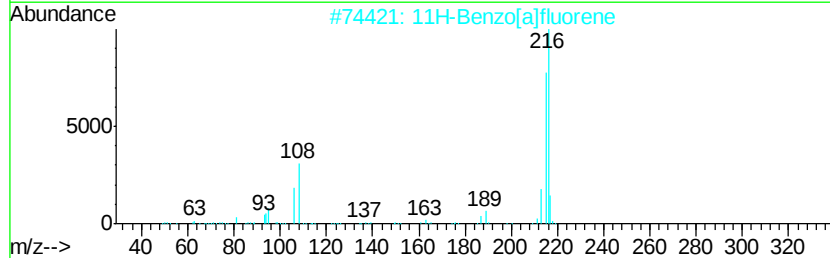
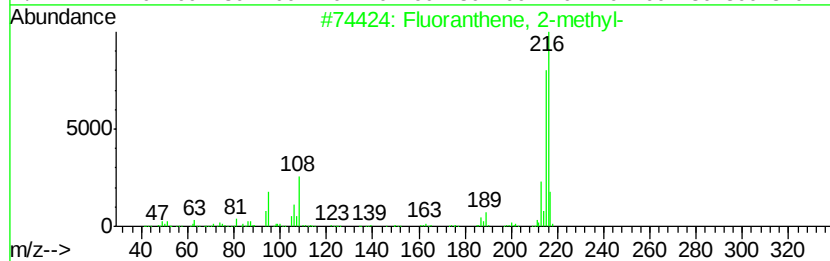
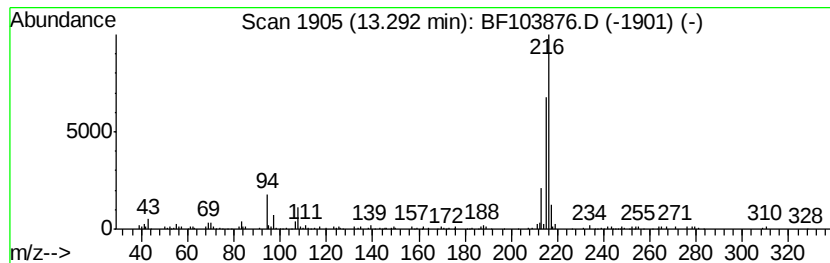
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Fluoranthene, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.29	2.62 ng	216835	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	96
2		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
3		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
4		7H-Benzo[c]fluorene	216	C17H12	000205-12-9	74
5		Pyrene, 1-methyl-	216	C17H12	002381-21-7	74



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032318\
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Acq On : 23 Mar 2018 13:27
Operator : JU/SJ
Sample : J2002-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

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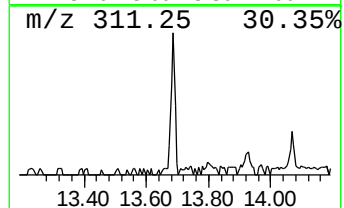
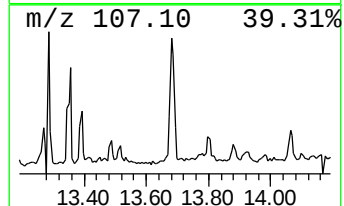
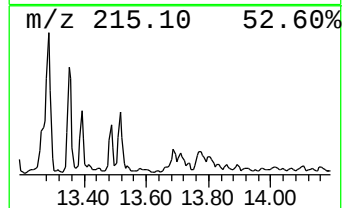
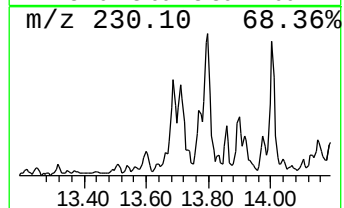
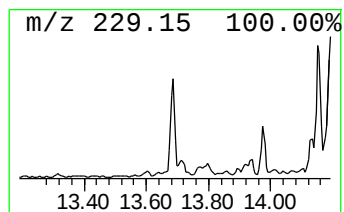
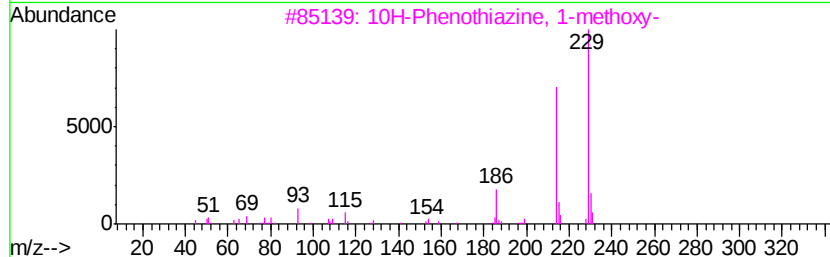
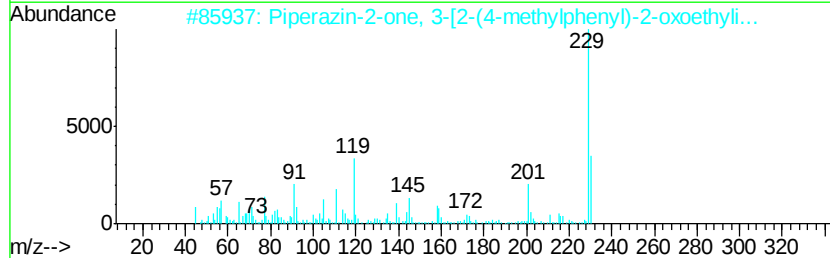
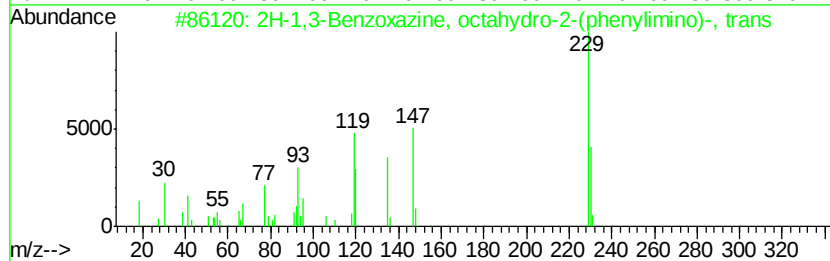
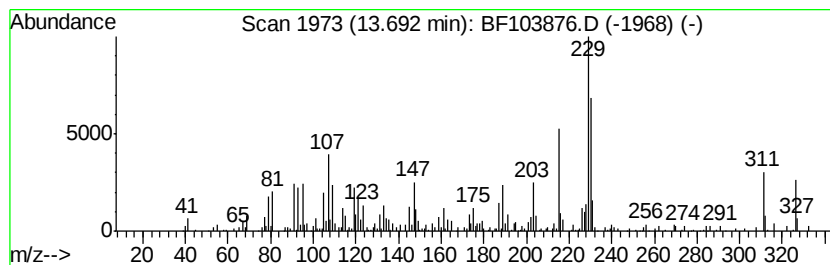
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown13.69 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.69	4.28 ng	354243	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2H-1,3-Benzoxazine, octahydro-2-...	230	C14H18N2O	1000138-91-8	25
2		Piperazin-2-one, 3-[2-(4-methylp...	230	C13H14N2O2	063656-21-3	22
3		10H-Phenothiazine, 1-methoxy-	229	C13H11NOS	001576-70-1	22
4		2H-1,3-Benzoxazine, octahydro-2-...	230	C14H18N2O	1000138-91-9	22
5		1-Piperazinepropanenitrile, .bet...	229	C13H15N3O	014761-40-1	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032318\
Data File : BF103876.D
Acq On : 23 Mar 2018 13:27
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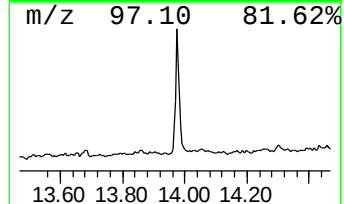
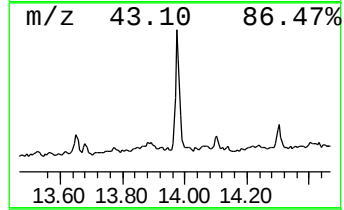
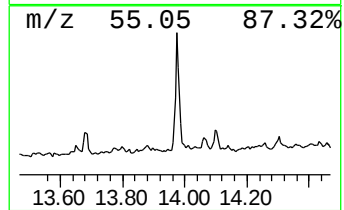
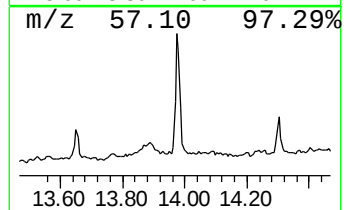
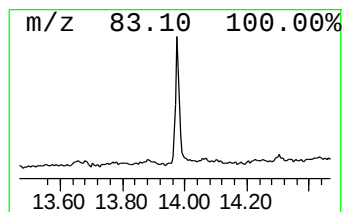
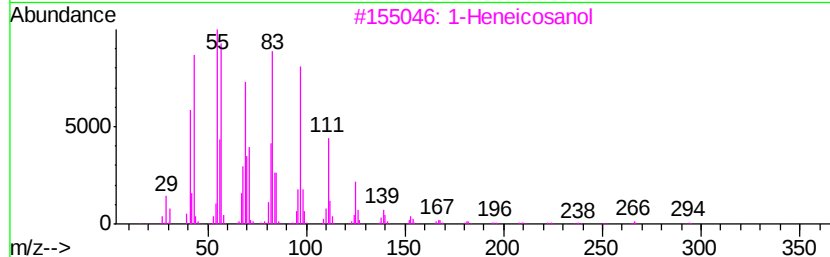
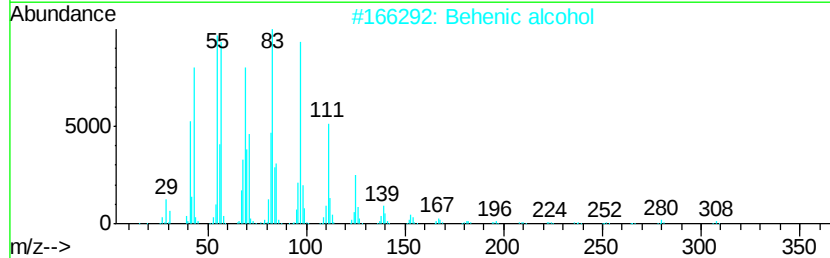
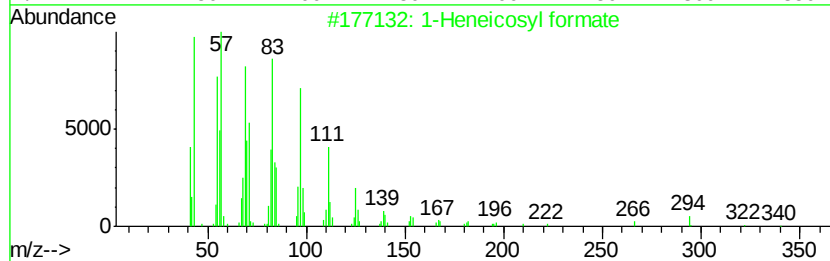
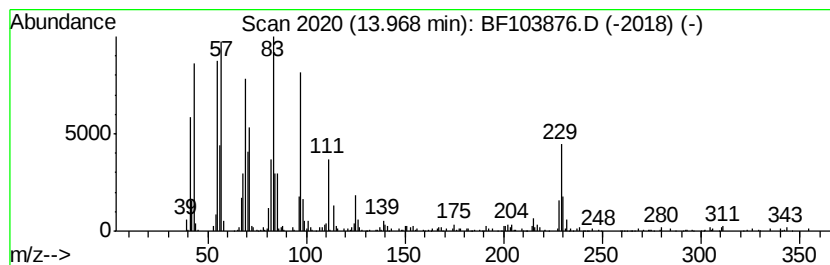
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1-Heneicosyl formate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	6.39 ng	529780	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Heneicosyl formate	340	C22H44O2	077899-03-7	93
2		Behenic alcohol	326	C22H46O	000661-19-8	91
3		1-Heneicosanol	312	C21H44O	015594-90-8	91
4		Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	90
5		n-Nonadecanol-1	284	C19H40O	001454-84-8	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032318\
Data File : BF103876.D
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ClientSampleId :
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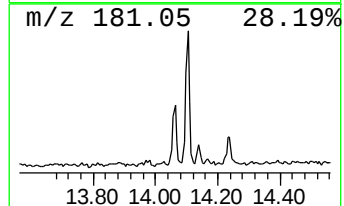
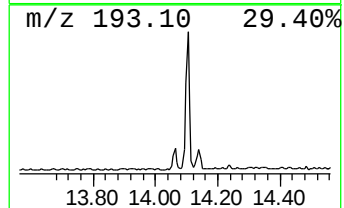
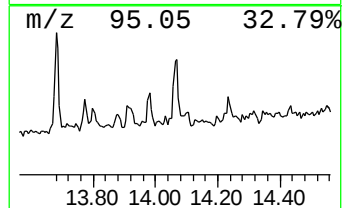
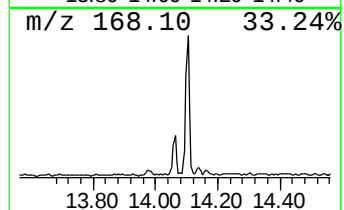
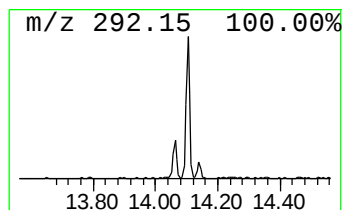
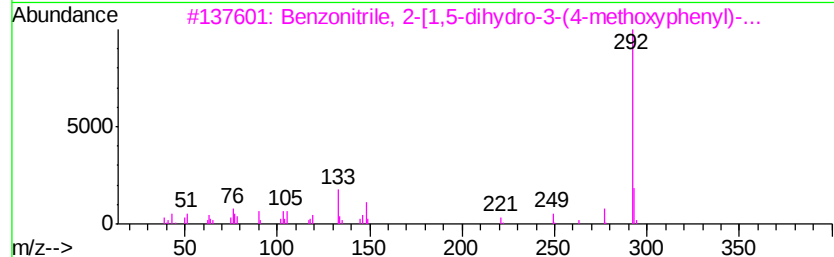
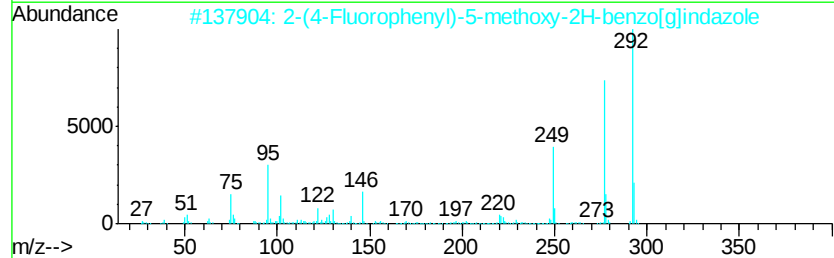
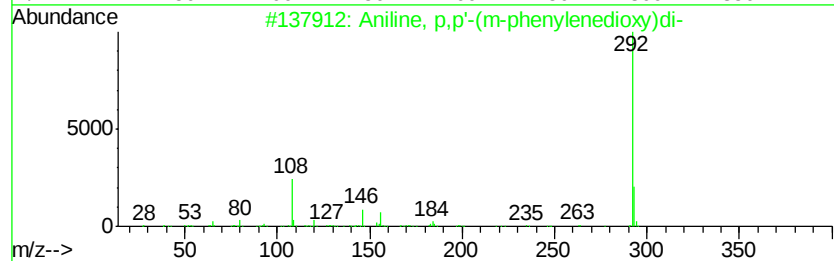
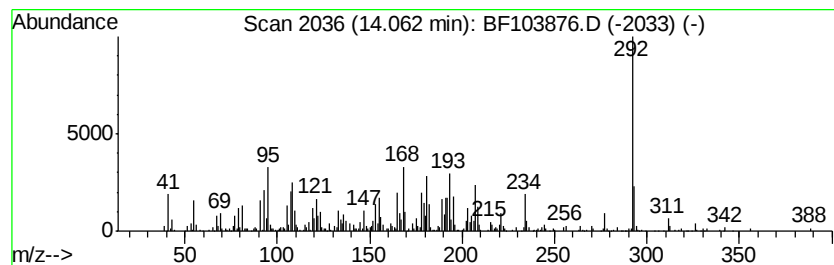
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown14.06 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.06	3.10 ng	256445	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Aniline, p,p'-(m-phenylenedioxy)di-	292	C18H16N2O2	002479-46-1	41
2		2-(4-Fluorophenyl)-5-methoxy-2H-...	292	C18H13FN2O	1000194-51-1	38
3		Benzonitrile, 2-[1,5-dihydro-3-(...	292	C16H12N4O2	1000349-81-5	38
4		4,6-Dihydroxy-5-[3,4,5-trimethox...	292	C14H16N2O5	118507-85-0	38
5		1-Benzylpyrene	292	C23H16	042211-34-7	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032318\
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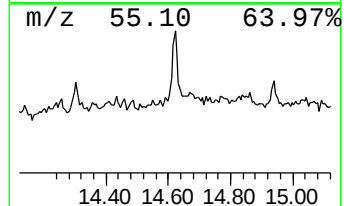
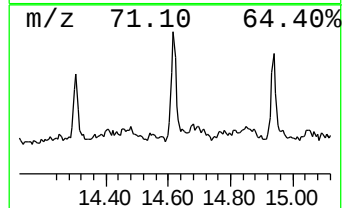
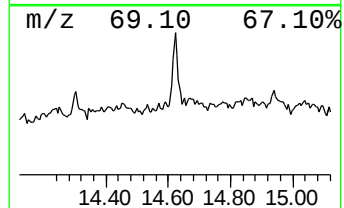
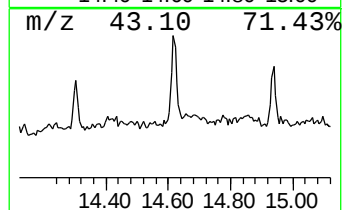
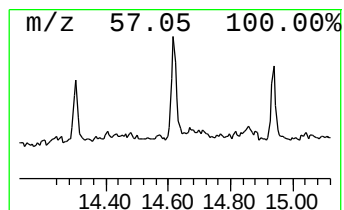
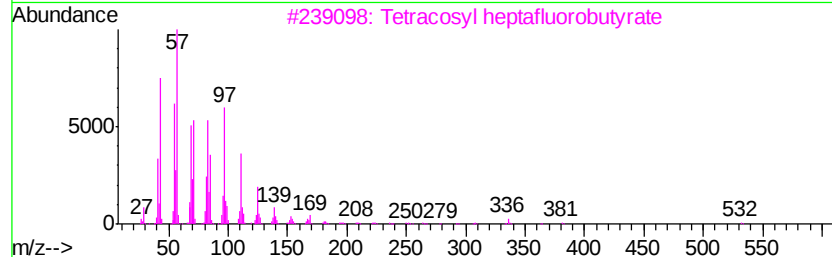
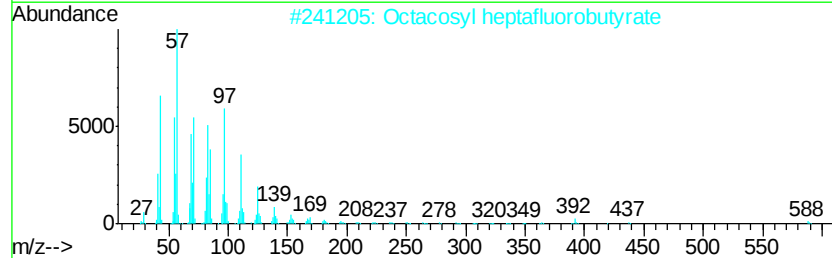
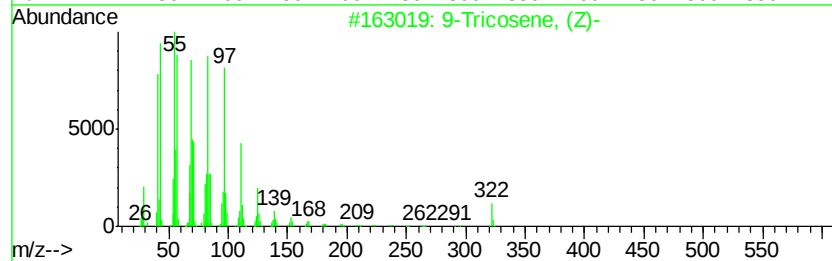
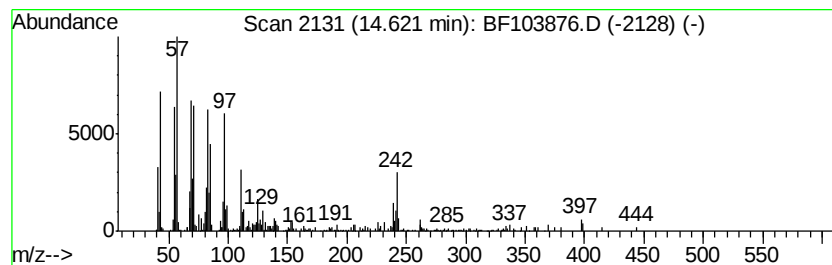
Instrument :
BNA_F
ClientSampleId :
TP-4-MH-B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 9-Tricosene, (Z)- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.62	2.54 ng	210710	Chrysene-d12	14.16
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	9-Tricosene, (Z)-	322 C23H46	027519-02-4	91
2	Octacosyl heptafluorobutyrte	606 C32H57F7O2	1000351-83-6	91
3	Tetracosyl heptafluorobutyrte	550 C28H49F7O2	1000351-83-7	91
4	Octatriacontyl pentafluoropropio...	697 C41H77F5O2	1000351-89-1	90
5	1-Decanol, 2-hexyl-	242 C16H34O	002425-77-6	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032318\
Data File : BF103876.D
Acq On : 23 Mar 2018 13:27
Operator : JU/SJ
Sample : J2002-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-4-MH-B

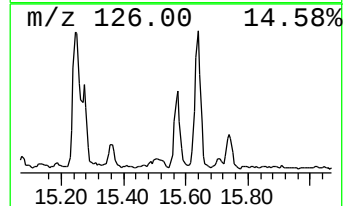
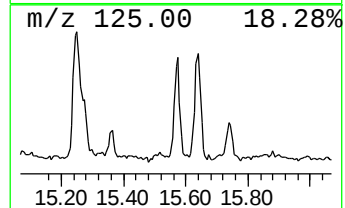
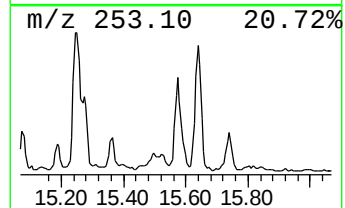
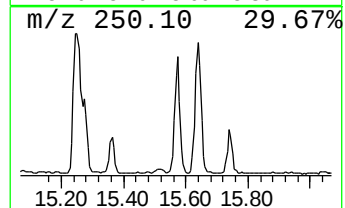
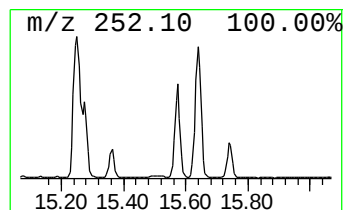
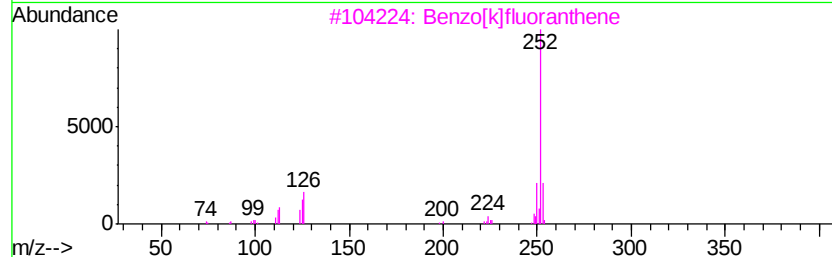
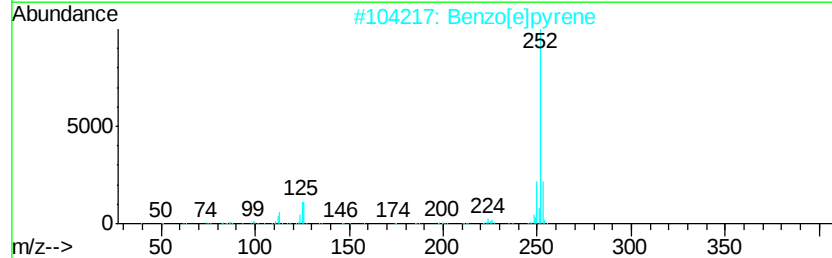
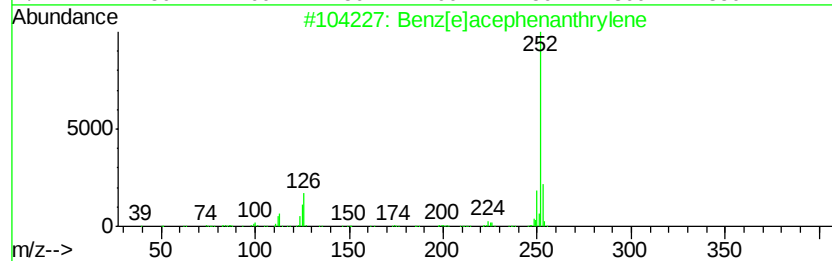
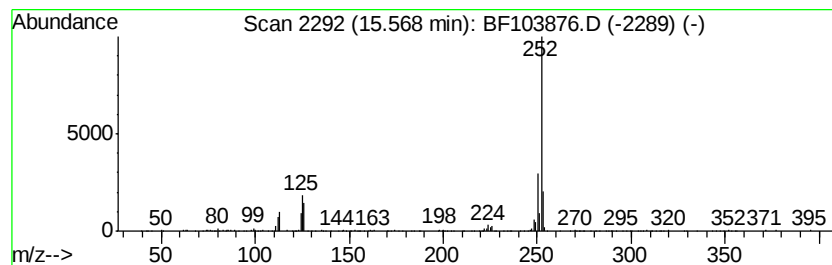
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Benz[e]acephenanthrylene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.57	7.63 ng	372027	Perylene-d12	15.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benz[e]acephenanthrylene	252	C20H12	000205-99-2	98
2			Benzo[e]pyrene	252	C20H12	000192-97-2	98
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
4			Perylene	252	C20H12	000198-55-0	93
5			Benzo[a]pyrene	252	C20H12	000050-32-8	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032318\
 Data File : BF103876.D
 Acq On : 23 Mar 2018 13:27
 Operator : JU/SJ
 Sample : J2002-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-4-MH-B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.32	3.9 ng		179037	1	6.97	909507	20.0
2-Pentanone, 4-hy...	5.22	4.4 ng		199830	1	6.97	909507	20.0
unknown6.75	6.75	84.6 ng		3845240	1	6.97	909507	20.0
Heptadecane	10.90	3.3 ng		197149	4	11.51	1181830	20.0
Phenanthrene, 2-m...	12.01	4.7 ng		275469	4	11.51	1181830	20.0
4H-Cyclopenta[def...	12.13	6.8 ng		401308	4	11.51	1181830	20.0
Phenanthrene, 2,5...	12.57	4.0 ng		237010	4	11.51	1181830	20.0
Fluoranthene, 2-m...	13.29	2.6 ng		216835	5	14.16	1657110	20.0
unknown13.69	13.69	4.3 ng		354243	5	14.16	1657110	20.0
1-Heneicosyl formate	13.97	6.4 ng		529780	5	14.16	1657110	20.0
unknown14.06	14.06	3.1 ng		256445	5	14.16	1657110	20.0
9-Tricosene, (Z)-	14.62	2.5 ng		210710	5	14.16	1657110	20.0
Benz[e]acephenant...	15.57	7.6 ng		372027	6	15.71	975239	20.0