

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF030118\
 Data File : BF103332.D
 Acq On : 1 Mar 2018 13:15
 Operator : SJ/JU
 Sample : J1720-15
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-29

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-BF022218.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	2.134	3	8	15	rVB	251316	335389	8.18%	0.825%
2	5.116	511	515	523	rBV	92954	134184	3.27%	0.330%
3	5.504	576	581	600	rVB	3554427	4101530	100.00%	10.087%
4	6.516	748	753	771	rBV	3303377	4096204	99.87%	10.074%
5	6.645	771	775	779	rVV	3727441	3870624	94.37%	9.519%
6	6.698	781	784	790	rVB2	42695	51993	1.27%	0.128%
7	6.875	810	814	820	rBV	1215899	1093727	26.67%	2.690%
8	7.028	836	840	844	rBV	3021910	3009732	73.38%	7.402%
9	7.439	906	910	913	rBV	2583576	2661006	64.88%	6.544%
10	7.528	923	925	935	rVB2	30922	53256	1.30%	0.131%
11	8.157	1028	1032	1035	rBV	1497524	1332092	32.48%	3.276%
12	8.875	1150	1154	1160	rBV	73706	82045	2.00%	0.202%
13	8.969	1166	1170	1173	rBV	45621	45047	1.10%	0.111%
14	9.233	1210	1215	1224	rBV	3587483	3699743	90.20%	9.099%
15	9.298	1224	1226	1229	rVV	122703	104494	2.55%	0.257%
16	9.333	1229	1232	1236	rVB2	43170	50492	1.23%	0.124%
17	9.498	1256	1260	1264	rBV3	31491	50914	1.24%	0.125%
18	9.575	1268	1273	1275	rVV4	39349	54149	1.32%	0.133%
19	9.622	1275	1281	1288	rVV	378763	418883	10.21%	1.030%
20	9.774	1304	1307	1312	rBV	86309	90625	2.21%	0.223%
21	9.827	1313	1316	1319	rVV	234484	178440	4.35%	0.439%
22	9.916	1327	1331	1334	rBV	1249237	1220612	29.76%	3.002%
23	9.945	1334	1336	1339	rVB	84497	72416	1.77%	0.178%
24	10.057	1352	1355	1358	rBV3	34837	51082	1.25%	0.126%
25	10.116	1362	1365	1368	rVV3	87069	96381	2.35%	0.237%
26	10.151	1368	1371	1375	rVB3	74362	75566	1.84%	0.186%
27	10.327	1398	1401	1404	rVB	290677	231439	5.64%	0.569%
28	10.463	1421	1424	1431	rVB2	55840	74492	1.82%	0.183%
29	10.545	1433	1438	1441	rBV	89980	114615	2.79%	0.282%
30	10.710	1461	1466	1475	rBV	1589847	1796292	43.80%	4.418%
31	10.804	1479	1482	1492	rVB	258533	356843	8.70%	0.878%
32	10.998	1512	1515	1516	rBV3	44364	46908	1.14%	0.115%
33	11.127	1534	1537	1541	rBV4	27459	50897	1.24%	0.125%
34	11.251	1555	1558	1561	rVV	217372	184165	4.49%	0.453%

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35	11.280	1561	1563	1565	rVV	73382	65778	1.60%	0.162%
36	11.404	1580	1584	1586	rBV	1059403	971177	23.68%	2.388%
37	11.427	1586	1588	1593	rVV	653628	577947	14.09%	1.421%
38	11.480	1594	1597	1602	rVB	156793	152464	3.72%	0.375%
39	11.639	1620	1624	1628	rBV2	55272	81516	1.99%	0.200%
40	11.674	1628	1630	1633	rVV	80433	67434	1.64%	0.166%
41	11.916	1667	1671	1673	rBV2	203562	239867	5.85%	0.590%
42	12.021	1686	1689	1691	rBV2	147468	159360	3.89%	0.392%
43	12.086	1697	1700	1703	rBV2	75274	61445	1.50%	0.151%
44	12.204	1717	1720	1723	rBV	64465	62585	1.53%	0.154%
45	12.239	1723	1726	1731	rVB2	72509	71657	1.75%	0.176%
46	12.457	1760	1763	1765	rBV	83450	78425	1.91%	0.193%
47	12.621	1788	1791	1795	rBV	1010938	920688	22.45%	2.264%
48	12.857	1824	1831	1836	rBV	969946	1103715	26.91%	2.714%
49	12.992	1849	1854	1857	rBV	2309076	2213266	53.96%	5.443%
50	13.074	1864	1868	1872	rBV3	70887	126495	3.08%	0.311%
51	13.251	1895	1898	1900	rBV	68977	65609	1.60%	0.161%
52	13.410	1923	1925	1931	rBV3	73556	119799	2.92%	0.295%
53	13.698	1972	1974	1977	rVB	96117	78050	1.90%	0.192%
54	13.874	2001	2004	2007	rVB2	283127	290313	7.08%	0.714%
55	14.057	2030	2035	2038	rBV2	968379	1344755	32.79%	3.307%
56	14.080	2038	2039	2047	rVB	443835	388228	9.47%	0.955%
57	15.121	2213	2216	2219	rBV	324095	443038	10.80%	1.090%
58	15.433	2267	2269	2274	rVB	198762	239170	5.83%	0.588%
59	15.504	2278	2281	2285	rVB	303383	405593	9.89%	0.997%
60	15.568	2289	2292	2295	rVB	368388	447876	10.92%	1.101%

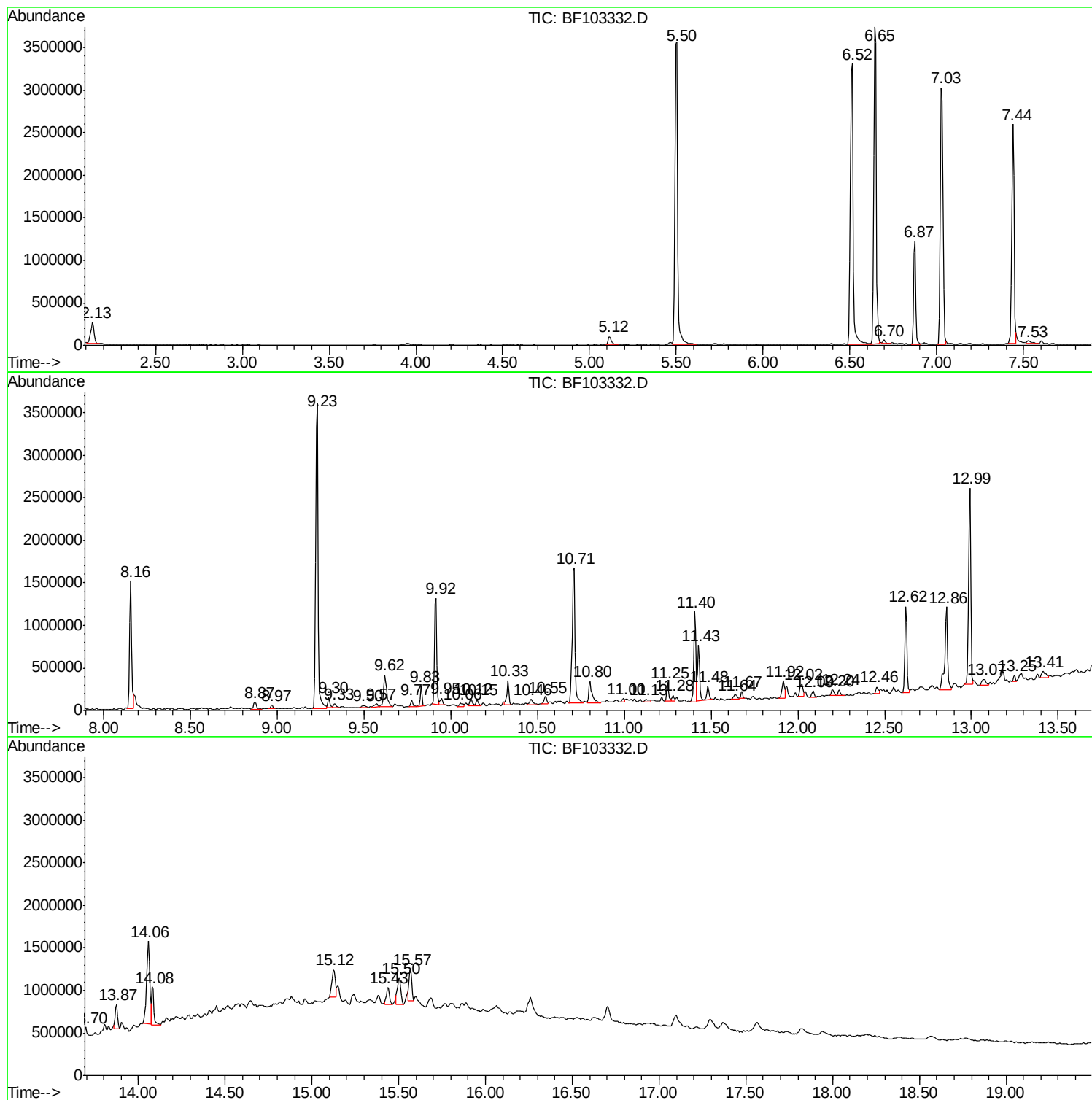
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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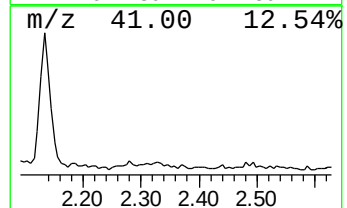
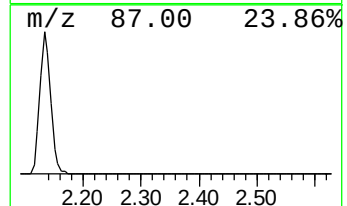
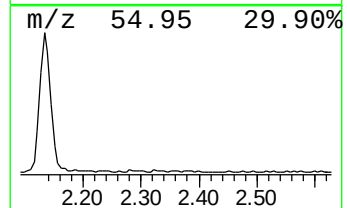
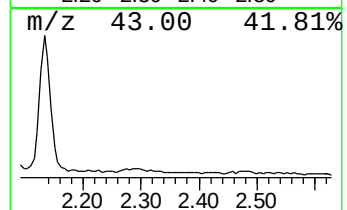
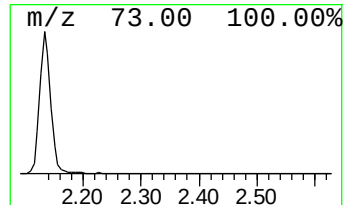
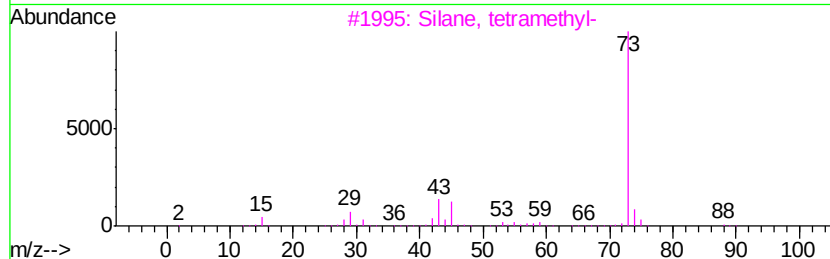
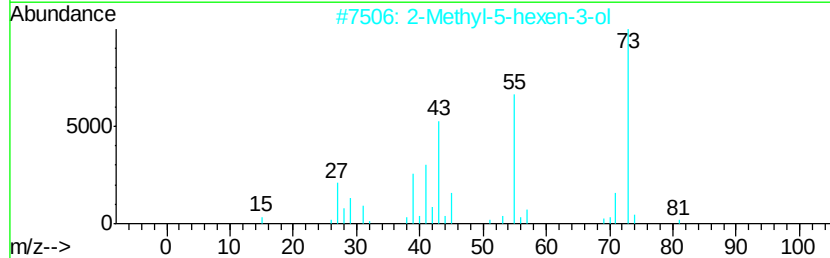
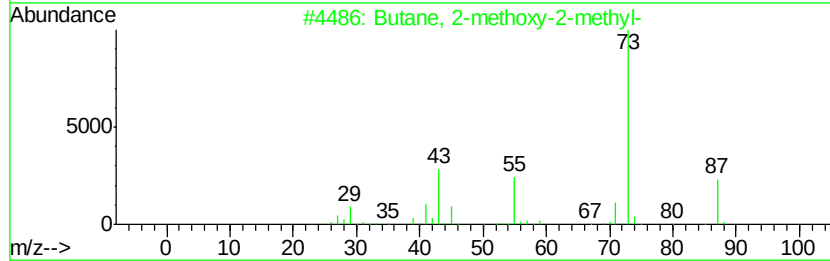
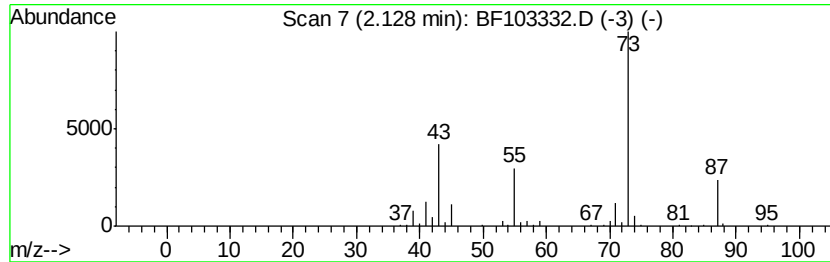
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.13	6.13 ng	335389	1,4-Dichlorobenzene-d4	6.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	38
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	35
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5		Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9



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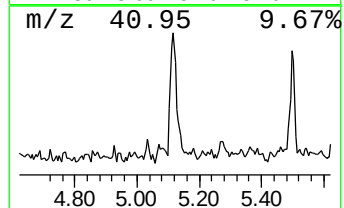
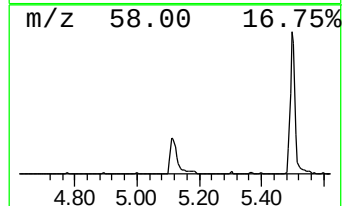
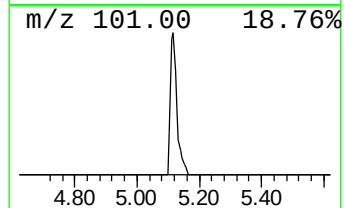
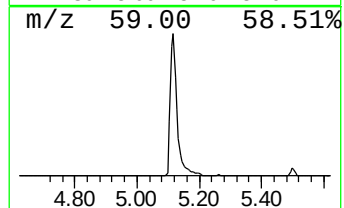
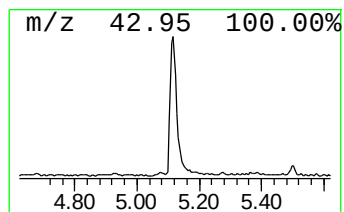
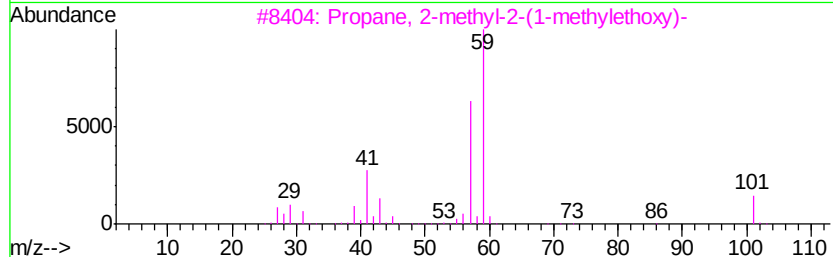
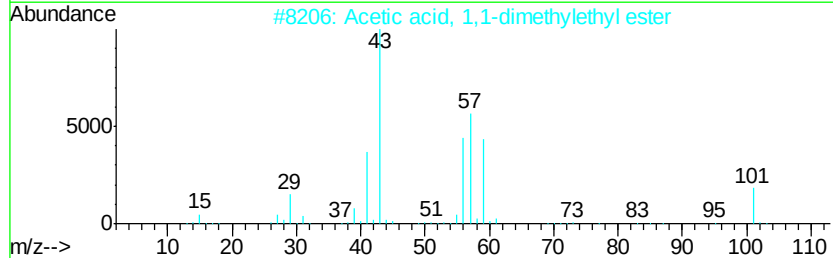
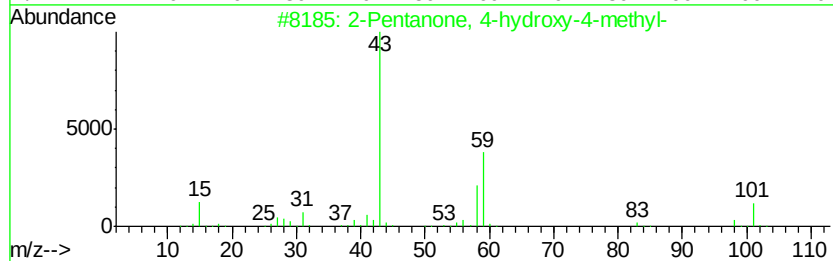
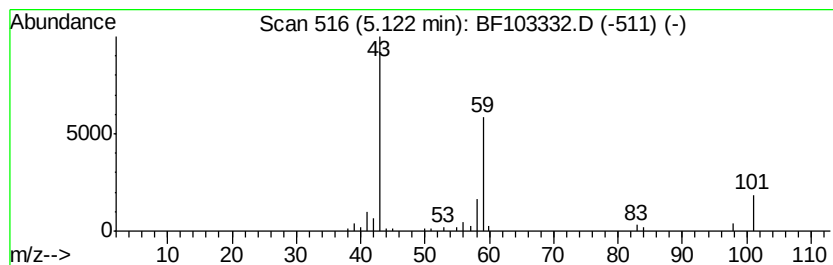
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	2.45 ng	134184	1,4-Dichlorobenzene-d4	6.87

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
4			Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	23
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16



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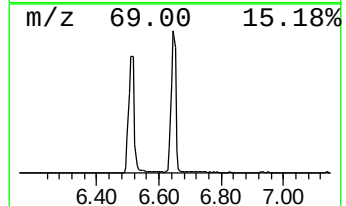
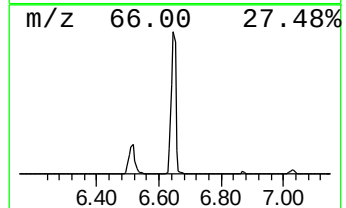
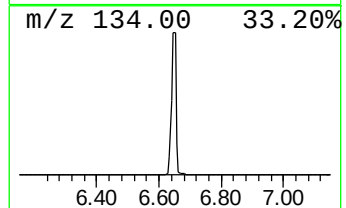
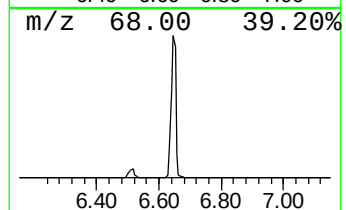
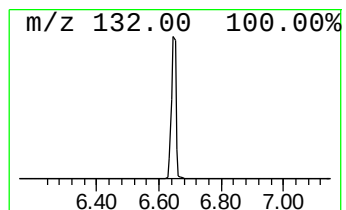
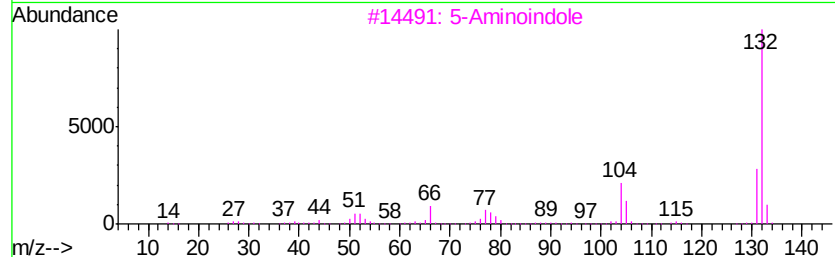
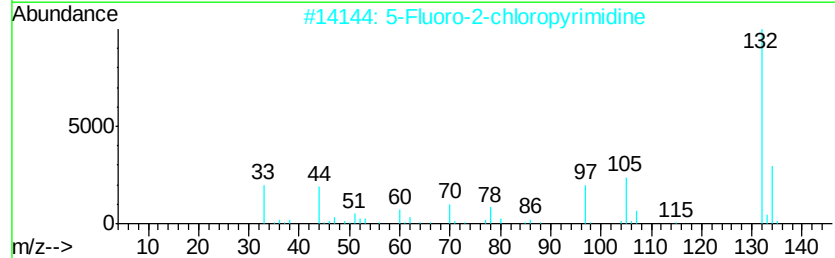
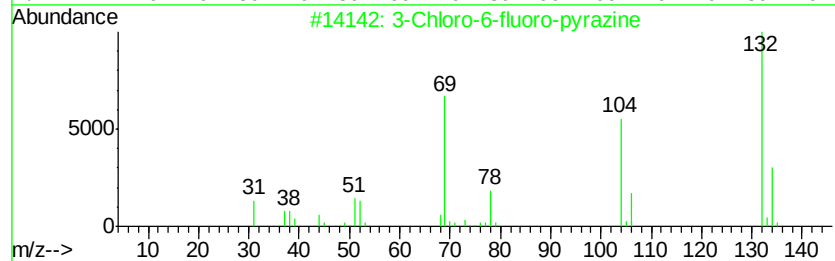
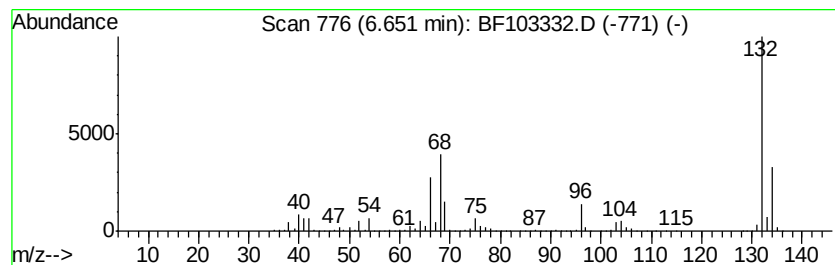
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	70.78 ng	3870620	1,4-Dichlorobenzene-d4	6.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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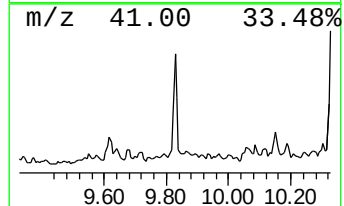
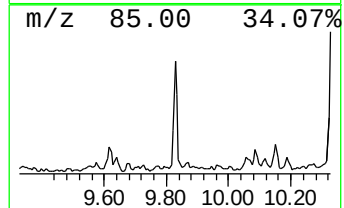
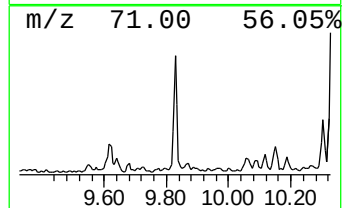
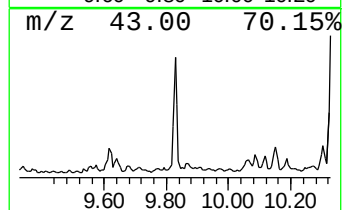
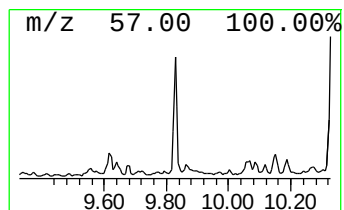
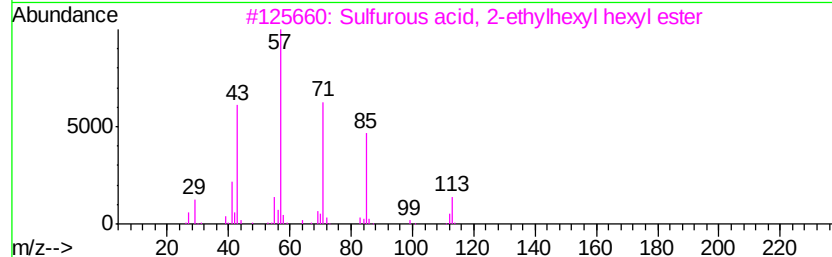
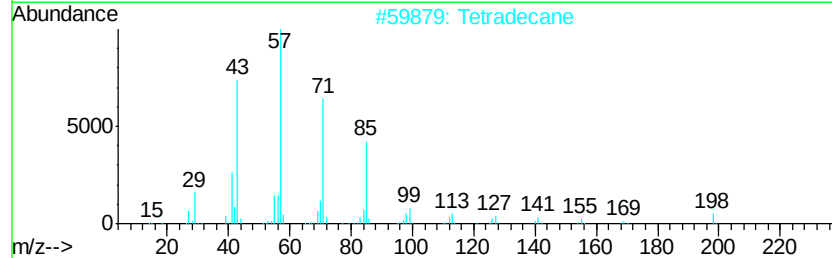
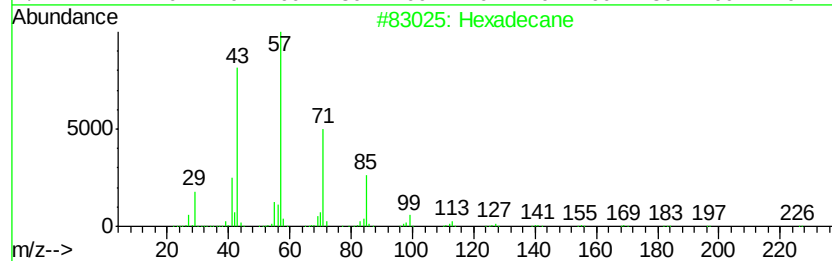
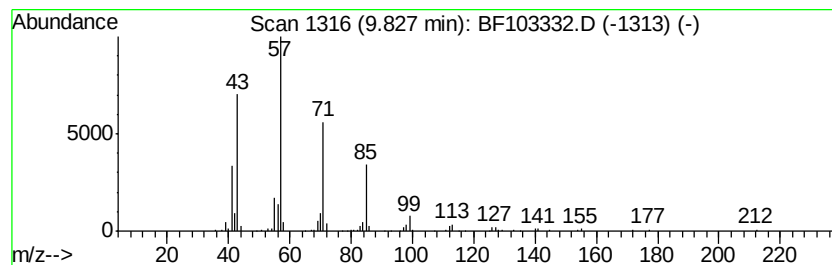
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.83	2.92 ng	178440	Acenaphthene-d10	9.92

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	91
2	Tetradecane		198	C14H30	000629-59-4	80
3	Sulfurous acid, 2-ethylhexyl hex...		278	C14H30O3S	1000309-20-2	80
4	Decane, 2,3,5-trimethyl-		184	C13H28	062238-11-3	72
5	Tridecane		184	C13H28	000629-50-5	59



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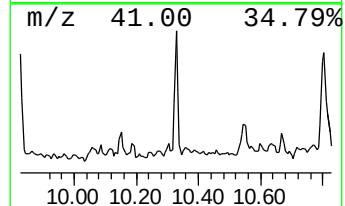
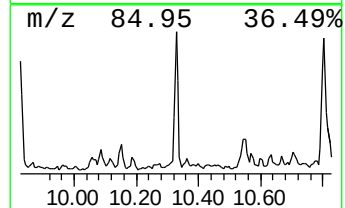
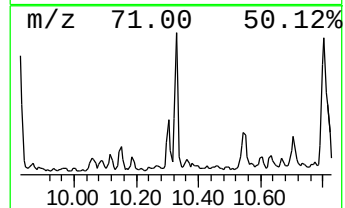
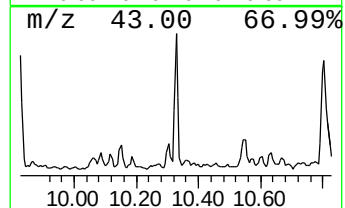
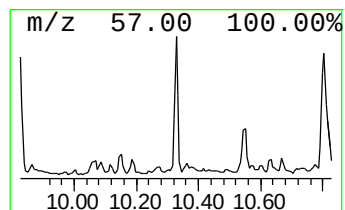
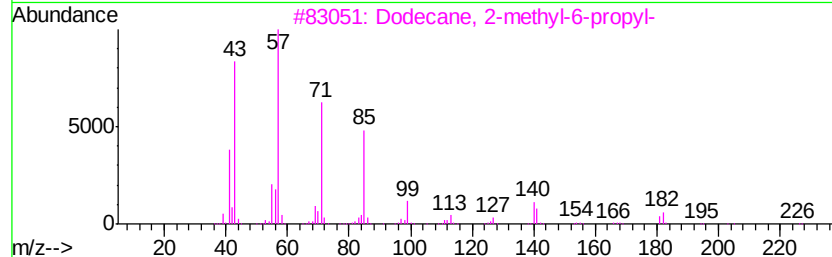
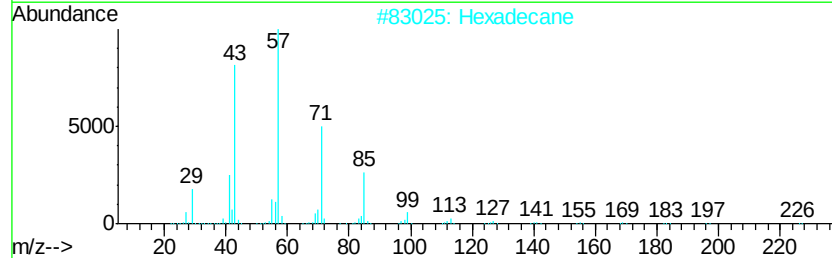
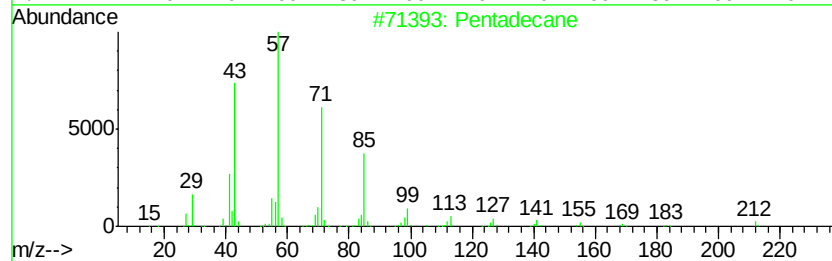
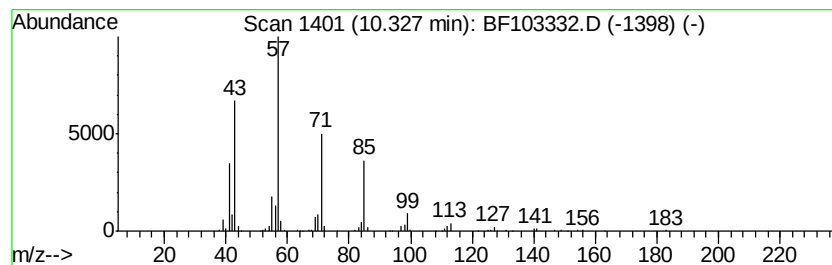
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Pentadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.33	3.79 ng	231439	Acenaphthene-d10	9.92

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane		212	C15H32	000629-62-9	90
2	Hexadecane		226	C16H34	000544-76-3	86
3	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	86
4	Decane, 2,3,5-trimethyl-		184	C13H28	062238-11-3	72
5	Sulfurous acid, 2-ethylhexyl hex...		278	C14H30O3S	1000309-20-2	72



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Acq On : 1 Mar 2018 13:15
Operator : SJ/JU
Sample : J1720-15
Misc :
ALS Vial : 10 Sample Multiplier: 1

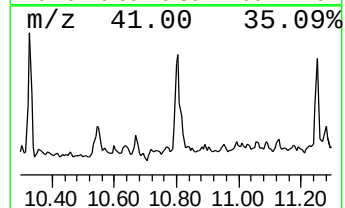
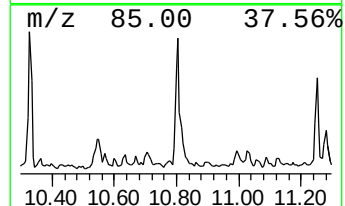
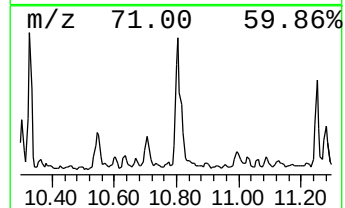
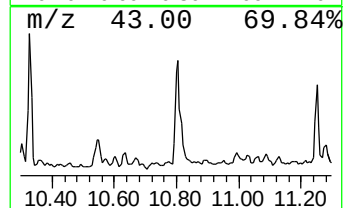
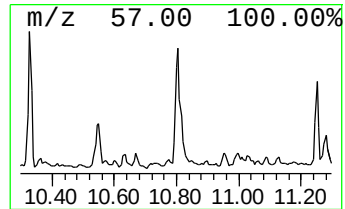
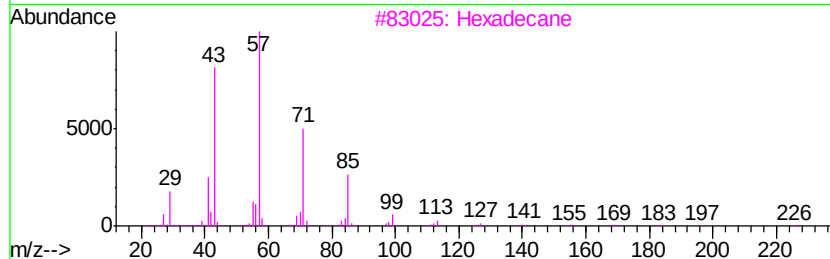
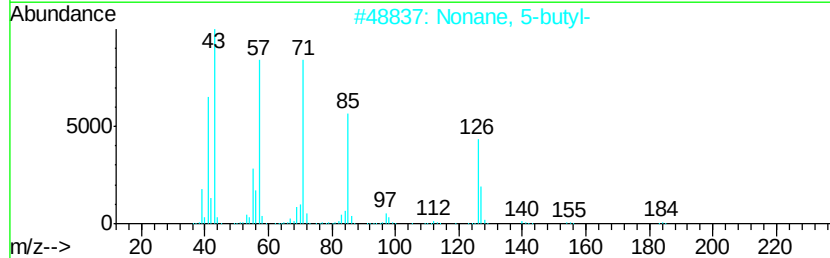
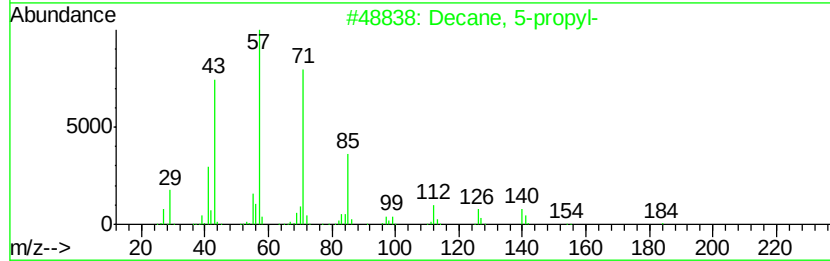
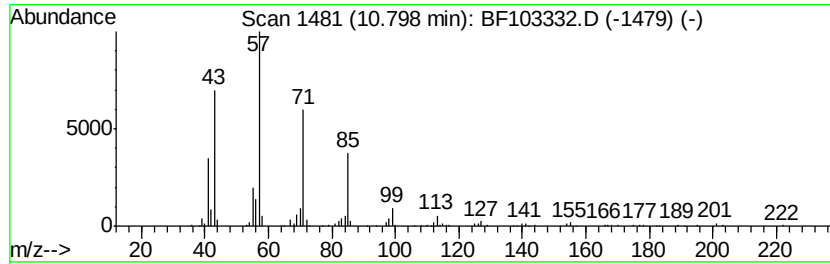
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Decane, 5-propyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.80	7.35 ng	356843	Phenanthrene-d10	11.40	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 5-propyl-	184	C13H28	017312-62-8	91
2	Nonane, 5-butyl-	184	C13H28	017312-63-9	90
3	Hexadecane	226	C16H34	000544-76-3	90
4	Bacchotricuneatin c	342	C20H22O5	066563-30-2	87
5	Dodecane, 4-methyl-	184	C13H28	006117-97-1	81



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030118\
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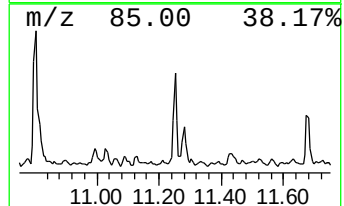
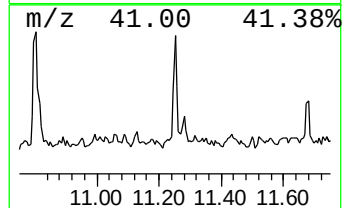
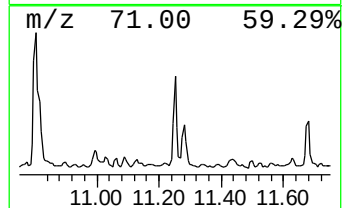
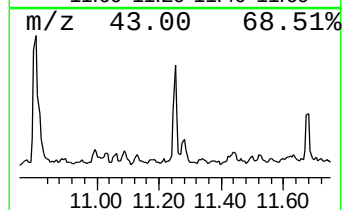
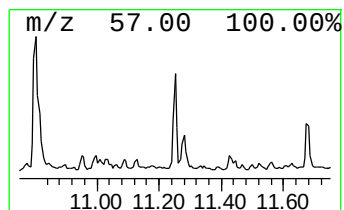
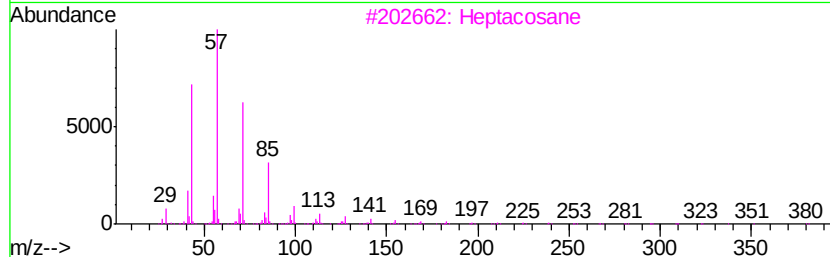
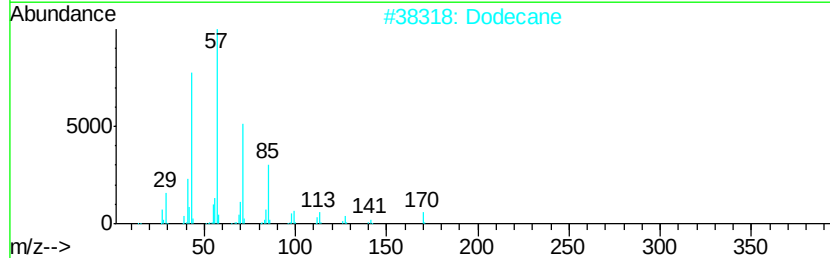
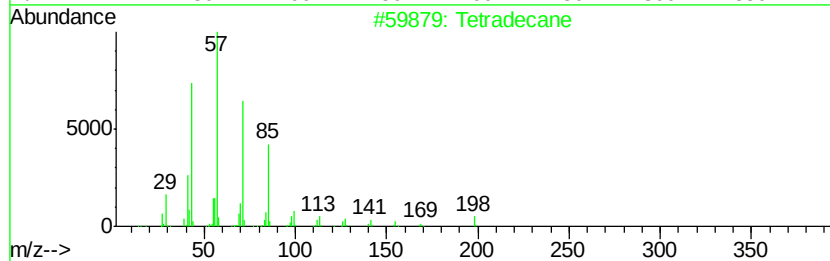
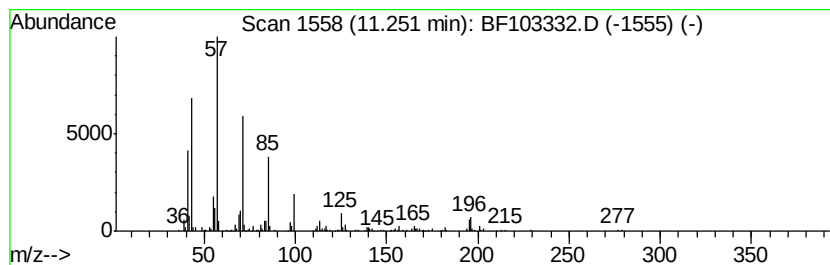
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tetradecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.25	3.79 ng	184165	Phenanthrene-d10	11.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	74
2		Dodecane	170	C12H26	000112-40-3	74
3		Heptacosane	380	C27H56	000593-49-7	72
4		Tetracosane	338	C24H50	000646-31-1	72
5		Undecane	156	C11H24	001120-21-4	72



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF030118\
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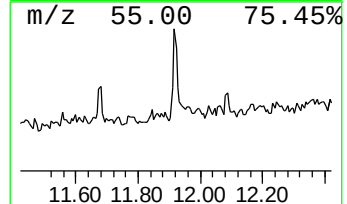
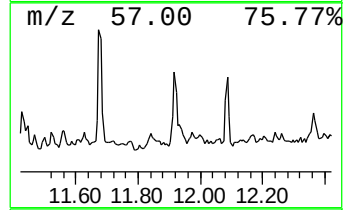
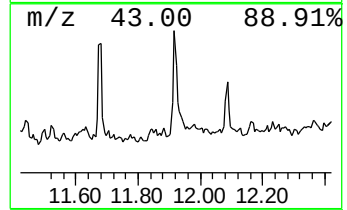
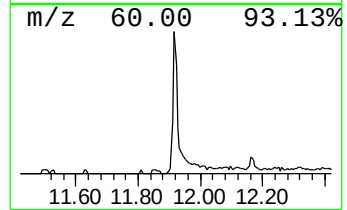
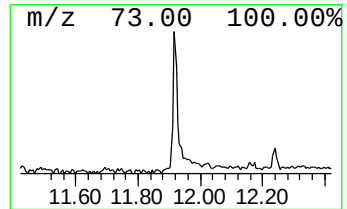
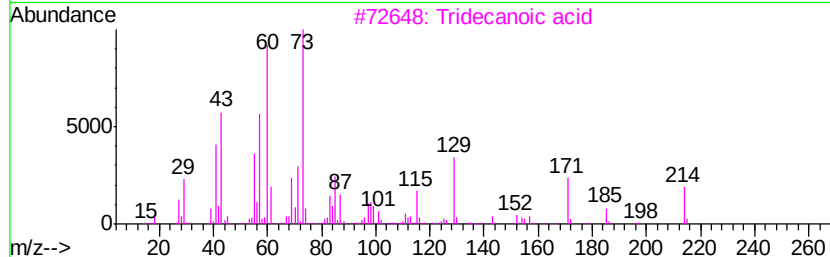
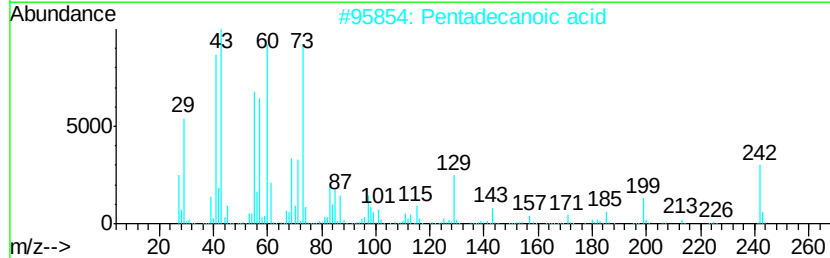
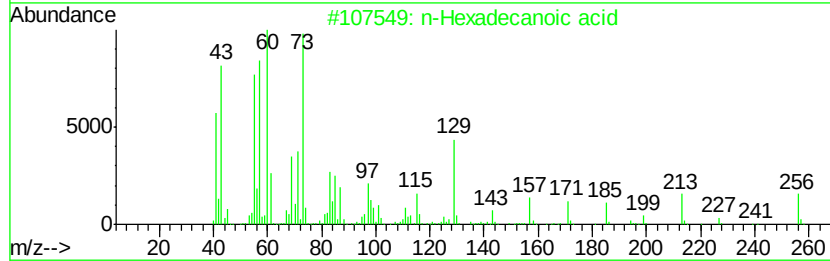
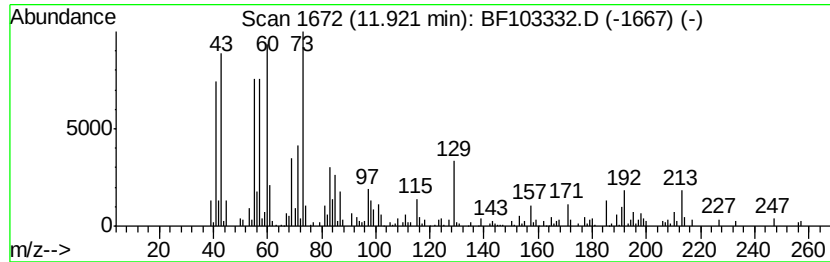
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.92	4.94 ng	239867	Phenanthrene-d10	11.40	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	68
3	Tridecanoic acid	214	C13H26O2	000638-53-9	52
4	n-Decanoic acid	172	C10H20O2	000334-48-5	49
5	Tetradecane	198	C14H30	000629-59-4	25



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030118\
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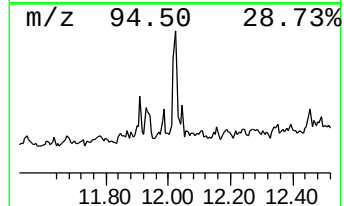
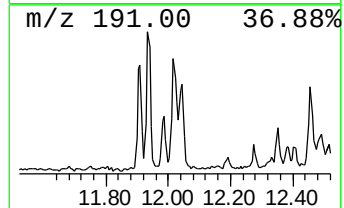
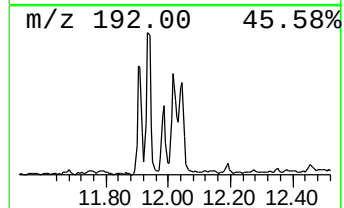
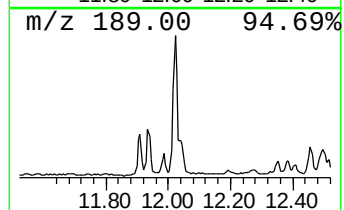
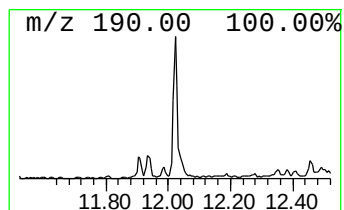
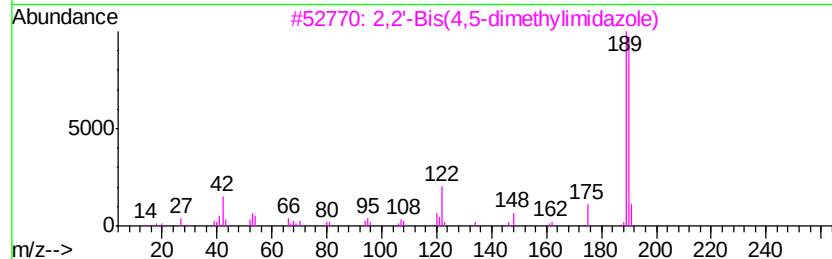
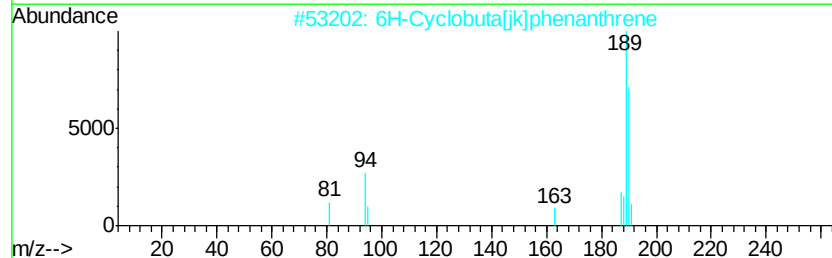
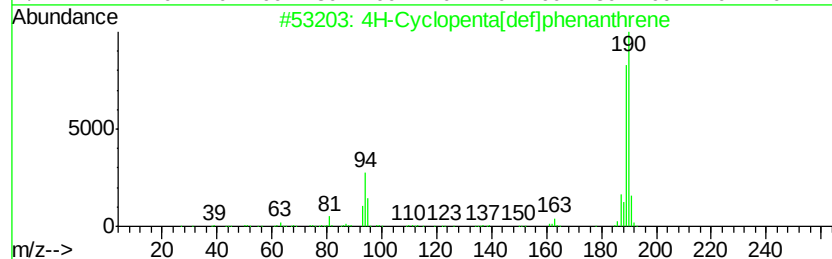
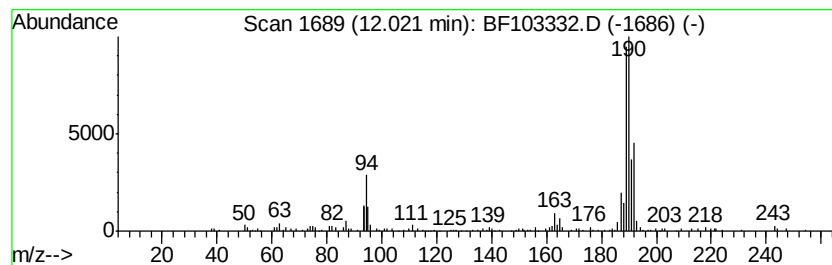
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.02	3.28 ng	159360	Phenanthrene-d10	11.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
2		6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	72
3		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	37
4		Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	28
5		Carbonic acid, ethyl 2-formyl-4,...	262	C10H8Cl2O4	1000331-34-4	17



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF030118\
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Acq On : 1 Mar 2018 13:15
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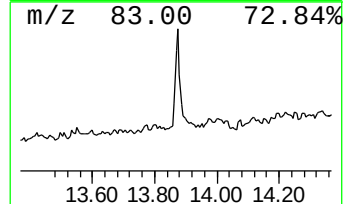
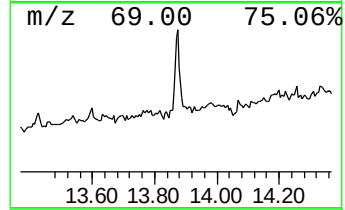
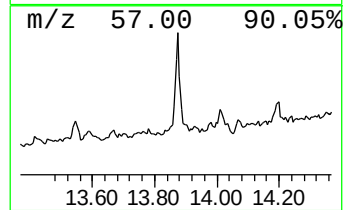
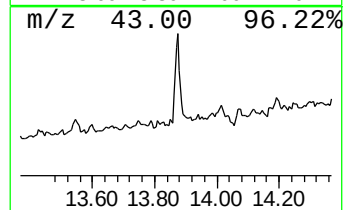
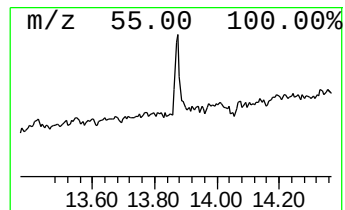
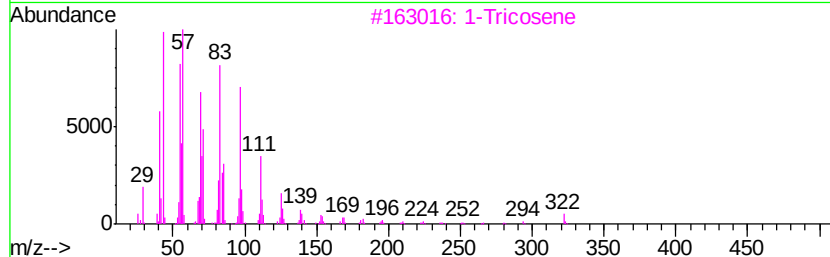
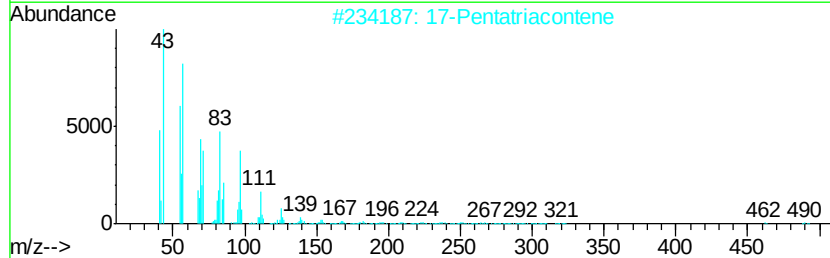
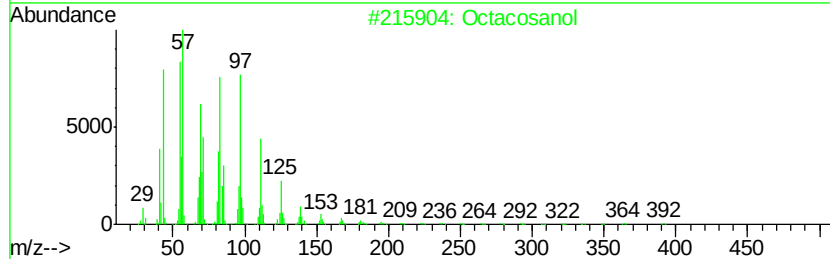
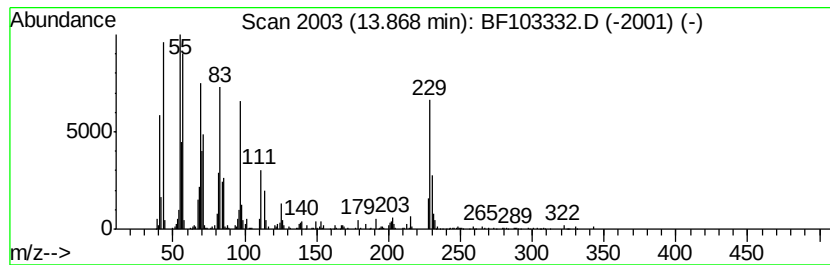
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octacosanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	4.32 ng	290313	Chrysene-d12	14.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosanol	410	C28H58O	000557-61-9	70
2		17-Pentatriacontene	491	C35H70	006971-40-0	62
3		1-Tricosene	322	C23H46	018835-32-0	62
4		1-Hexacosanol	382	C26H54O	000506-52-5	62
5		1-Tricosanol	340	C23H48O	003133-01-5	58



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Misc :
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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
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2-Pentanone, 4-hy...	5.12	2.5 ng		134184	1	6.87	1093730	20.0
unknown6.65	6.65	70.8 ng		3870620	1	6.87	1093730	20.0
Hexadecane	9.83	2.9 ng		178440	3	9.92	1220610	20.0
Pentadecane	10.33	3.8 ng		231439	3	9.92	1220610	20.0
Decane, 5-propyl-	10.80	7.3 ng		356843	4	11.40	971177	20.0
Tetradecane	11.25	3.8 ng		184165	4	11.40	971177	20.0
n-Hexadecanoic acid	11.92	4.9 ng		239867	4	11.40	971177	20.0
4H-Cyclopenta[def...	12.02	3.3 ng		159360	4	11.40	971177	20.0
Octacosanol	13.87	4.3 ng		290313	5	14.06	1344760	20.0