

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF090917\
 Data File : BF098394.D
 Acq On : 9 Sep 2017 19:07
 Operator : SJ/JU
 Sample : I5140-03
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RL-5-13.75-27.5

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-BF081517.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	3.034	159	161	172	rBV	14905	37598	1.38%	0.140%
2	3.610	252	259	269	rBV	361004	554819	20.42%	2.062%
3	4.863	468	472	484	rVB	253804	392282	14.43%	1.458%
4	5.292	540	545	558	rBV	2453836	2531411	93.15%	9.406%
5	6.333	716	722	732	rBV	2322199	2717586	100.00%	10.098%
6	6.451	737	742	745	rVV	2670610	2514124	92.51%	9.342%
7	6.510	747	752	757	rVB5	21978	39174	1.44%	0.146%
8	6.675	776	780	787	rBV	821733	694734	25.56%	2.581%
9	6.833	802	807	810	rBV	2086384	1977362	72.76%	7.347%
10	7.104	848	853	862	rVB	222939	245711	9.04%	0.913%
11	7.245	869	877	880	rBV	1656648	1636377	60.21%	6.080%
12	7.275	880	882	887	rVB2	23683	28059	1.03%	0.104%
13	7.357	893	896	899	rBV2	34283	34456	1.27%	0.128%
14	7.716	952	957	960	rBV3	39186	48806	1.80%	0.181%
15	7.751	960	963	972	rVB6	17301	33467	1.23%	0.124%
16	7.863	979	982	986	rBV	30529	30128	1.11%	0.112%
17	7.957	991	998	1000	rVV	988343	894292	32.91%	3.323%
18	7.974	1000	1001	1007	rVB	215637	183228	6.74%	0.681%
19	8.316	1055	1059	1066	rVB6	14679	27451	1.01%	0.102%
20	8.604	1102	1108	1112	rBV2	21508	31222	1.15%	0.116%
21	8.669	1115	1119	1122	rVB	76295	67423	2.48%	0.251%
22	8.763	1131	1135	1139	rVB	49906	45841	1.69%	0.170%
23	8.880	1149	1155	1157	rBV2	30013	30314	1.12%	0.113%
24	9.033	1175	1181	1184	rBV	2750063	2639547	97.13%	9.808%
25	9.363	1234	1237	1240	rBV	28341	30545	1.12%	0.113%
26	9.427	1244	1248	1252	rVB	452100	365280	13.44%	1.357%
27	9.563	1267	1271	1275	rBV	47015	47099	1.73%	0.175%
28	9.704	1289	1295	1299	rBV	971089	896774	33.00%	3.332%
29	9.921	1326	1332	1335	rBV2	37081	51227	1.89%	0.190%
30	10.251	1385	1388	1395	rVB2	29818	34117	1.26%	0.127%
31	10.363	1401	1407	1410	rBV2	36019	42805	1.58%	0.159%
32	10.498	1423	1430	1433	rBV	1307764	1309849	48.20%	4.867%
33	10.615	1447	1450	1459	rVB3	67863	114487	4.21%	0.425%
34	11.057	1522	1525	1528	rBV2	38123	40194	1.48%	0.149%

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35	11.086	1528	1530	1537	rVB	51020	56177	2.07%	0.209%
36	11.186	1543	1547	1550	rBV	909343	817471	30.08%	3.038%
37	11.210	1550	1551	1555	rVV	145246	84806	3.12%	0.315%
38	11.262	1555	1560	1564	rVB	36371	48960	1.80%	0.182%
39	11.415	1583	1586	1593	rVB3	85060	94375	3.47%	0.351%
40	12.162	1711	1713	1720	rVB3	25488	36630	1.35%	0.136%
41	12.233	1723	1725	1727	rBV2	35842	31379	1.15%	0.117%
42	12.398	1749	1753	1756	rBV	151237	151064	5.56%	0.561%
43	12.621	1788	1791	1794	rVB	135570	117101	4.31%	0.435%
44	12.774	1812	1817	1820	rBV	1707406	1626235	59.84%	6.043%
45	13.668	1966	1969	1976	rVB	105804	126298	4.65%	0.469%
46	13.821	1990	1995	1998	rBV	814392	798446	29.38%	2.967%
47	13.845	1998	1999	2004	rVB	86842	56188	2.07%	0.209%
48	14.298	2074	2076	2079	rVB2	47278	45018	1.66%	0.167%
49	14.586	2123	2125	2128	rBV2	41248	43343	1.59%	0.161%
50	14.827	2163	2166	2173	rVB	78196	124265	4.57%	0.462%
51	14.898	2175	2178	2181	rBV	101335	96334	3.54%	0.358%
52	15.162	2221	2223	2228	rVB	70275	86061	3.17%	0.320%
53	15.227	2229	2234	2236	rBV	483648	624173	22.97%	2.319%
54	15.645	2300	2305	2309	rBV	201777	285841	10.52%	1.062%
55	16.097	2377	2382	2390	rVB	223881	398309	14.66%	1.480%
56	16.633	2467	2473	2480	rVB	163112	310812	11.44%	1.155%
57	17.262	2573	2580	2588	rVB3	123741	304774	11.21%	1.132%
58	17.997	2699	2705	2715	rVB2	65215	180744	6.65%	0.672%

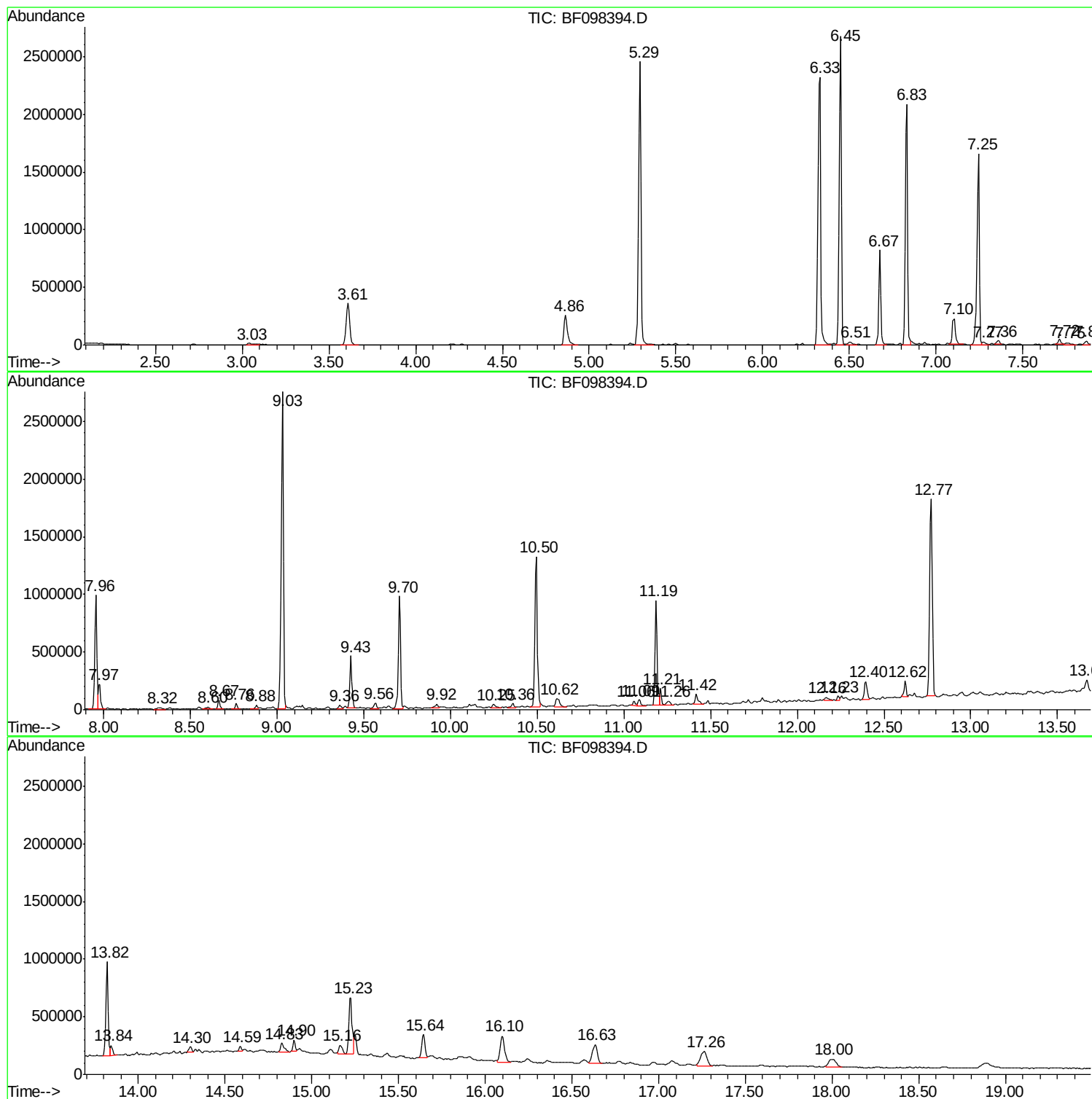
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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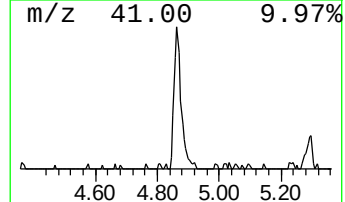
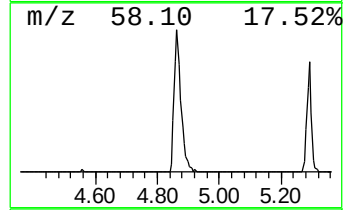
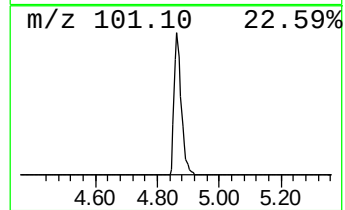
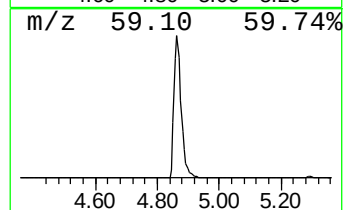
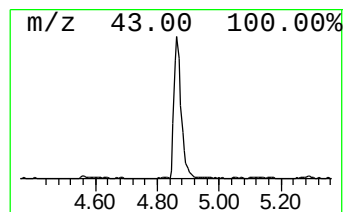
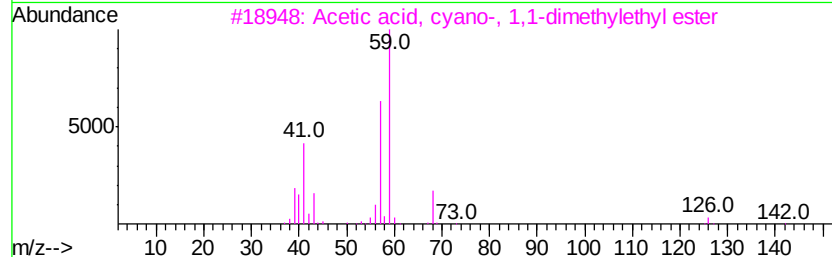
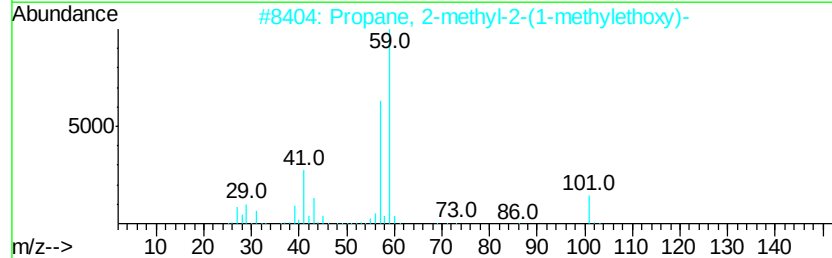
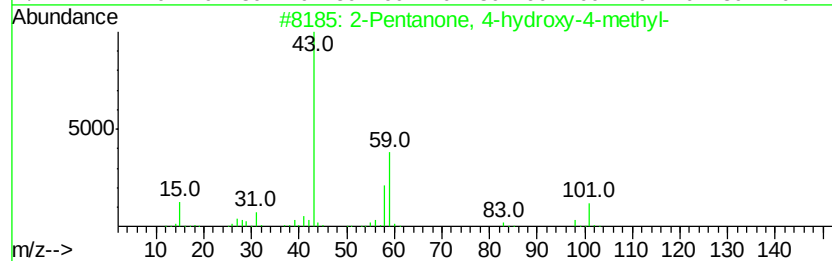
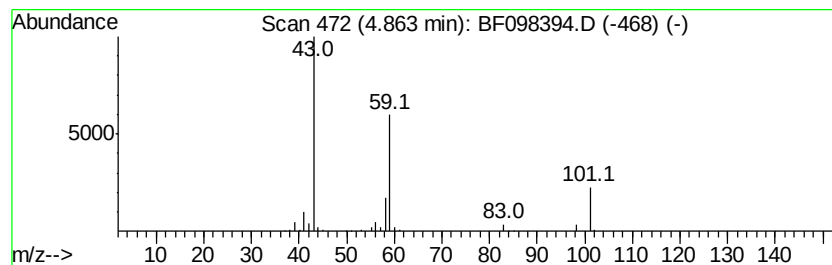
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.86	11.29 ng	392282	1,4-Dichlorobenzene-d4	6.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	50
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		Acetone	58	C3H6O	000067-64-1	9



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Instrument :
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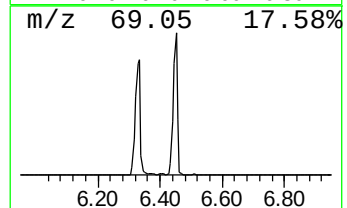
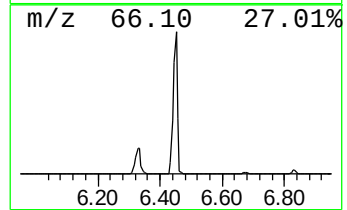
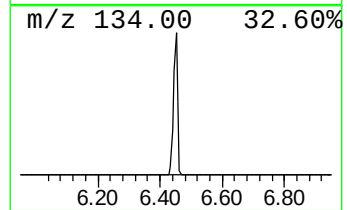
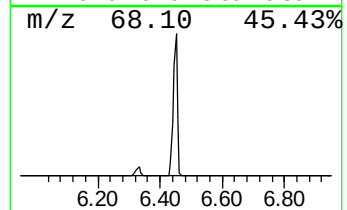
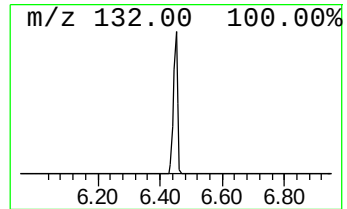
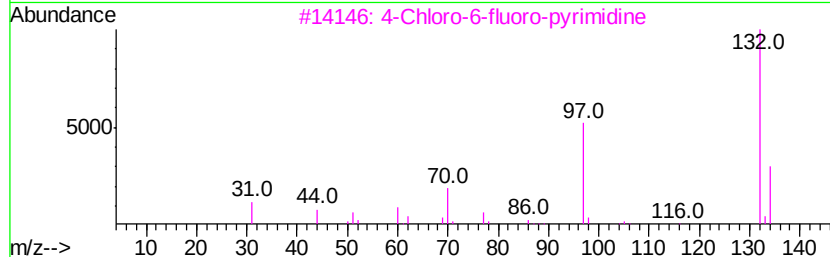
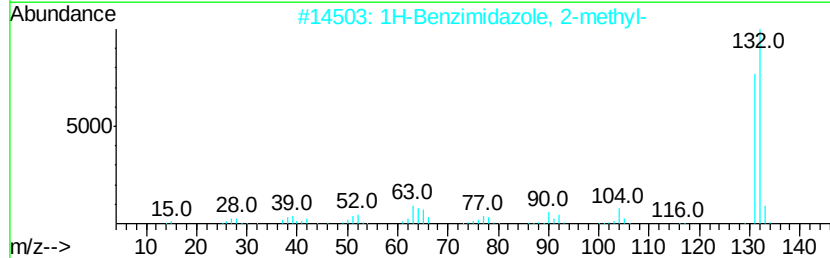
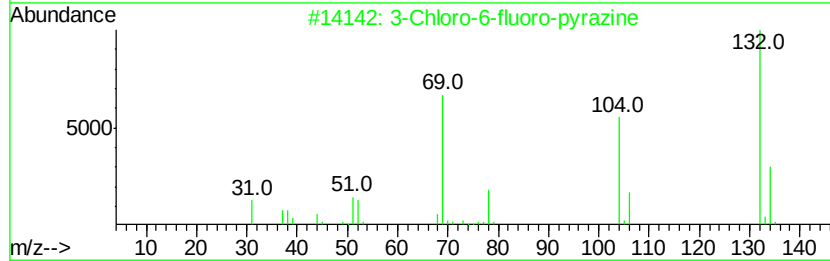
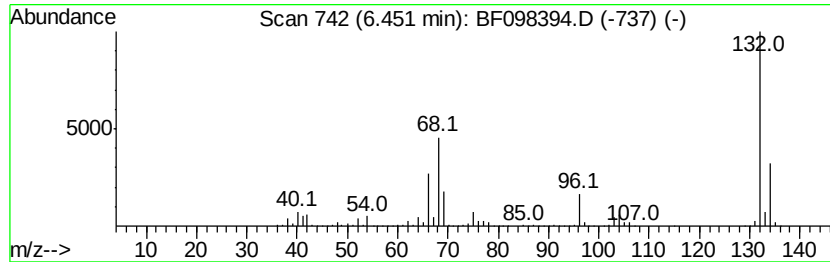
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 unknown6.45 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.45	72.38 ng	2514120	1,4-Dichlorobenzene-d4	6.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
3		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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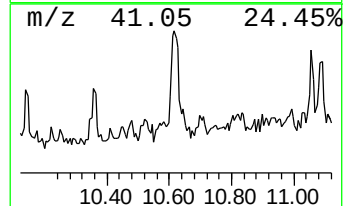
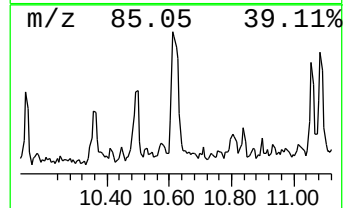
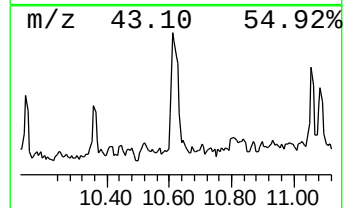
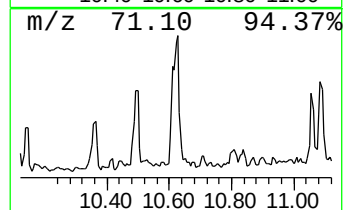
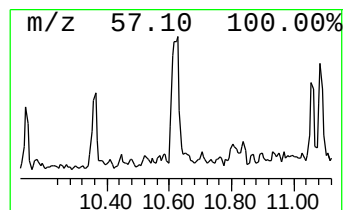
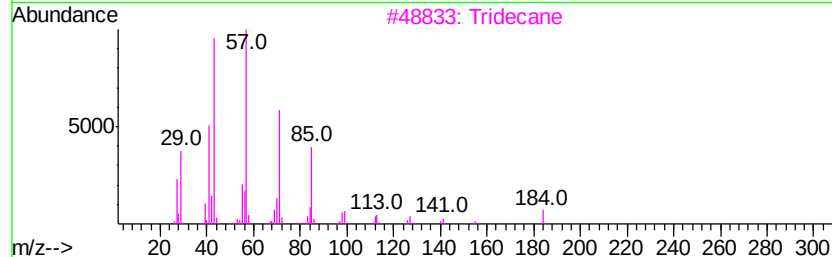
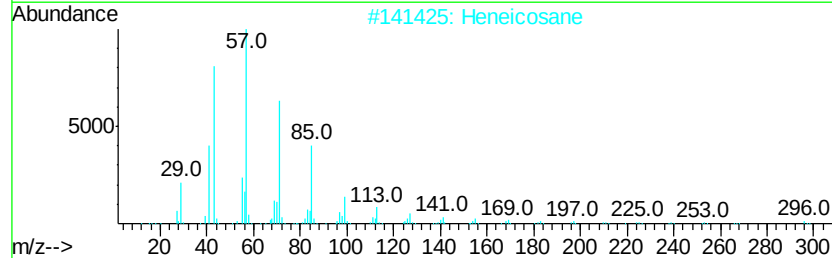
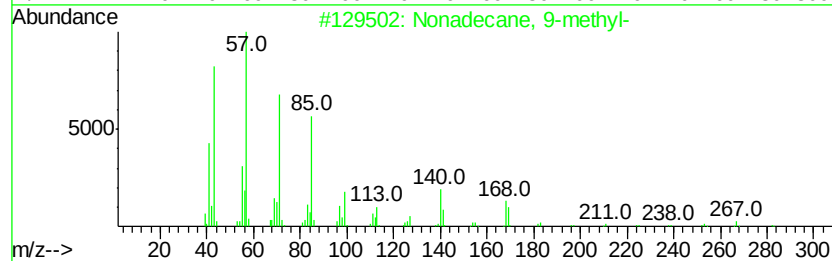
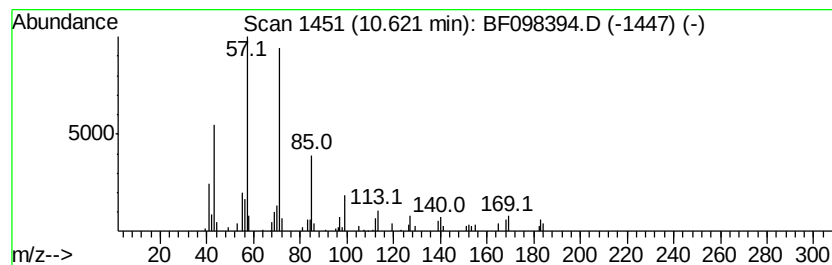
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Nonadecane, 9-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.62	2.80 ng	114487	Phenanthrene-d10	11.19	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	86
2	Heneicosane	296	C21H44	000629-94-7	83
3	Tridecane	184	C13H28	000629-50-5	81
4	Tetradecane, 4-methyl-	212	C15H32	025117-24-2	80
5	Tetracosane	338	C24H50	000646-31-1	80



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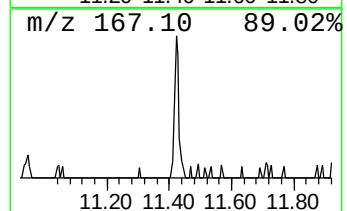
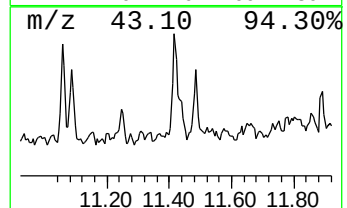
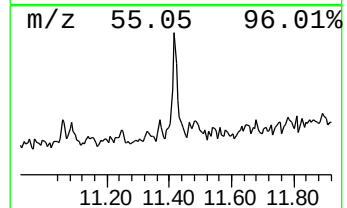
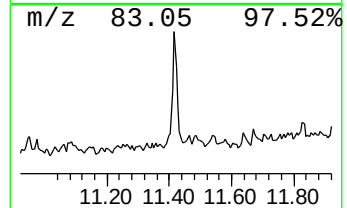
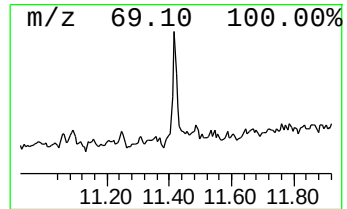
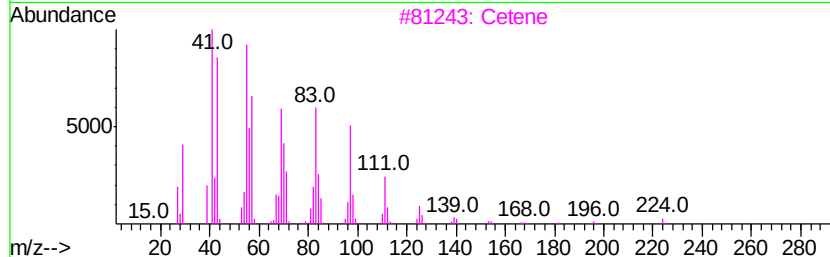
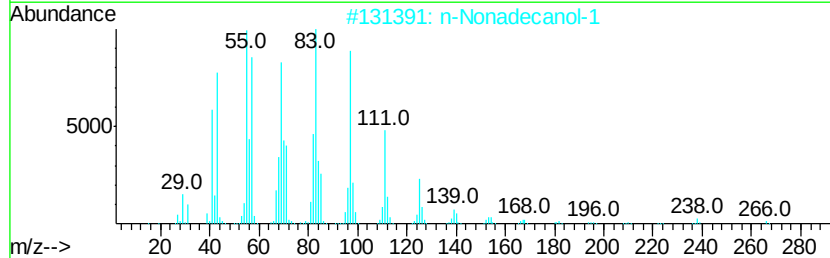
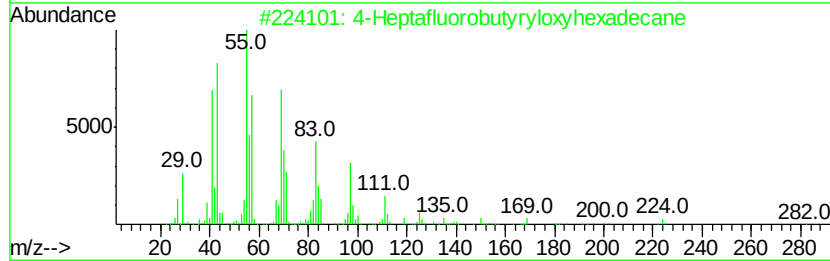
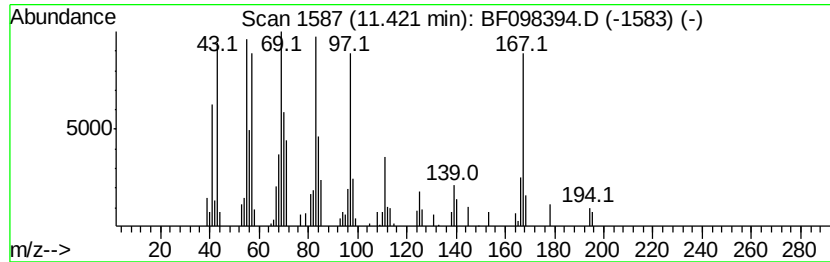
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 4-Heptafluorobutyryloxyhexa... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.42	2.31 ng	94375	Phenanthrene-d10	11.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Heptafluorobutyryloxyhexadecane	438	C20H33F7O2	1000282-97-2	91
2		n-Nonadecanol-1	284	C19H40O	001454-84-8	90
3		Cetene	224	C16H32	000629-73-2	64
4		3-Eicosene, (E)-	280	C20H40	074685-33-9	62
5		9-Eicosene, (E)-	280	C20H40	074685-29-3	49



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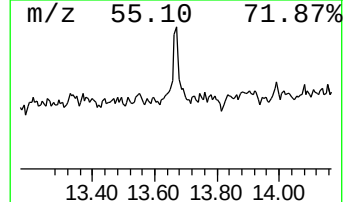
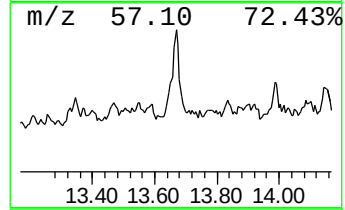
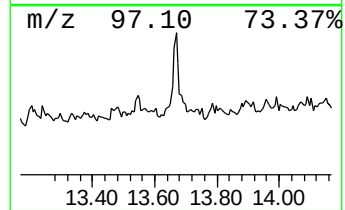
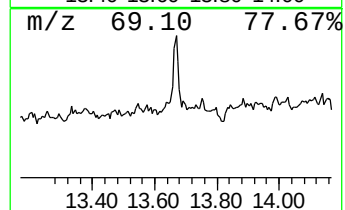
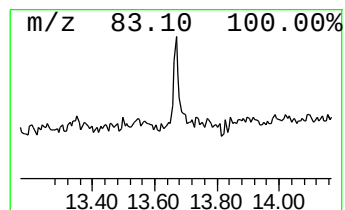
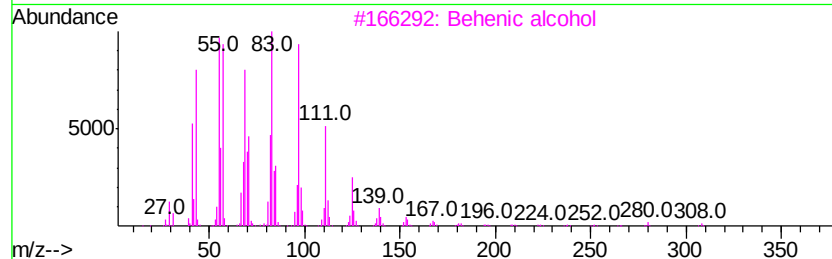
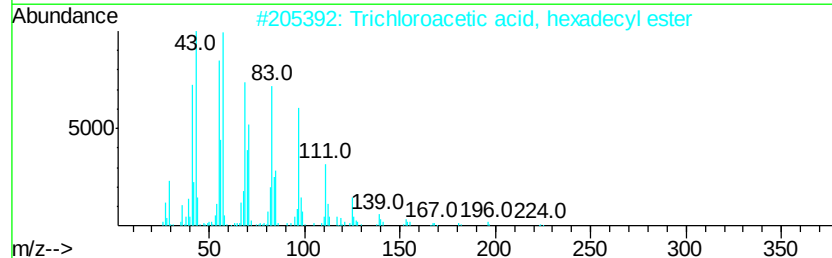
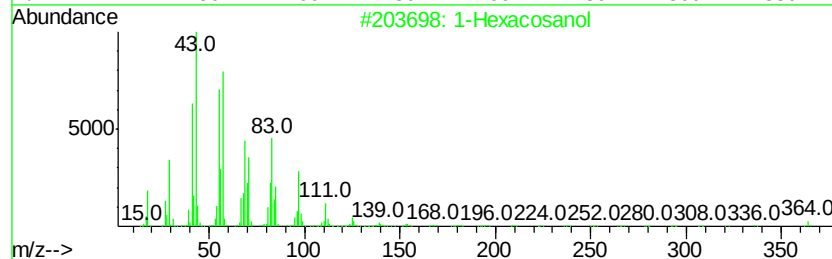
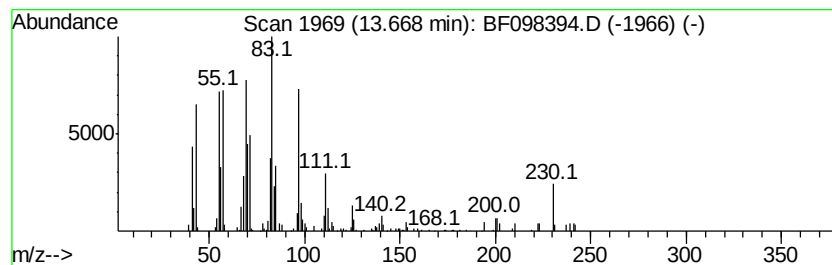
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 RL-5-13.75-27.5

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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 1-Hexacosanol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.67	3.16 ng	126298	Chrysene-d12	13.82	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosanol	382	C26H54O	000506-52-5	83
2	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	81
3	Behenic alcohol	326	C22H46O	000661-19-8	72
4	3-Eicosene, (E)-	280	C20H40	074685-33-9	72
5	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	70



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090917\
Data File : BF098394.D
Acq On : 9 Sep 2017 19:07
Operator : SJ/JU
Sample : I5140-03
Misc :
ALS Vial : 19 Sample Multiplier: 1

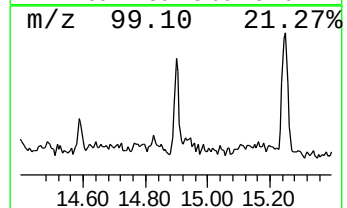
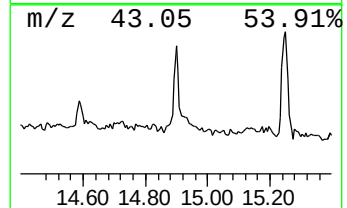
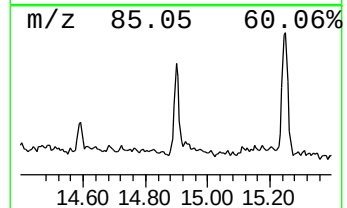
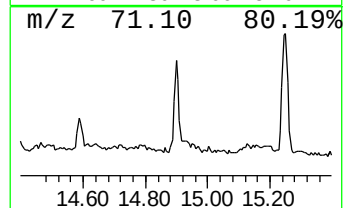
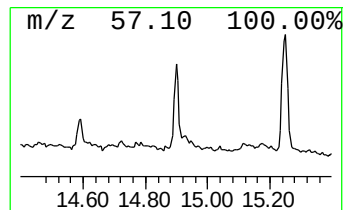
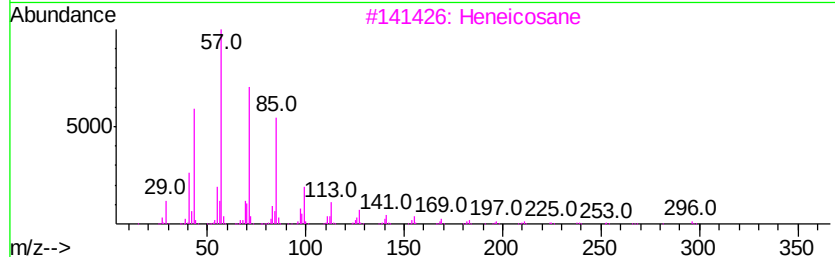
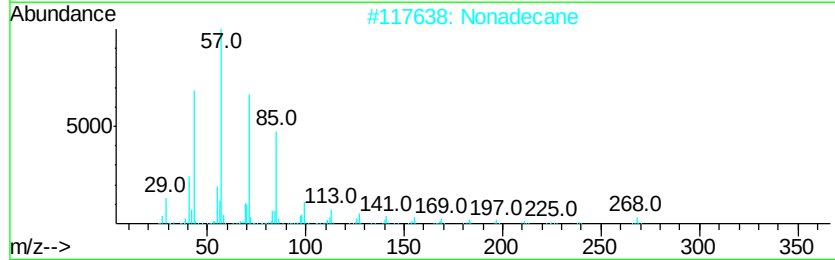
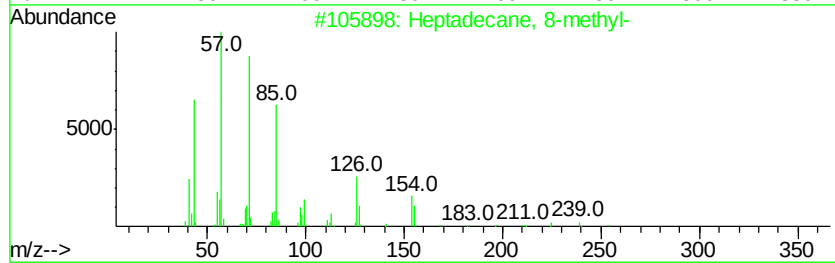
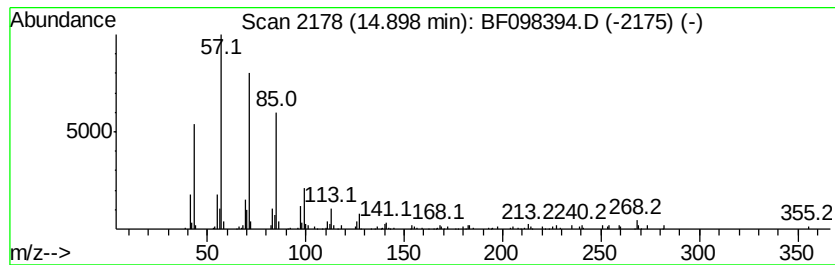
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RL-5-13.75-27.5

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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 Heptadecane, 8-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
14.90	3.09 ng	96334	Perylene-d12		15.23
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 8-methyl-	254	C18H38	013287-23-5	91
2	Nonadecane	268	C19H40	000629-92-5	87
3	Heneicosane	296	C21H44	000629-94-7	86
4	2-Bromo dodecane	248	C12H25Br	013187-99-0	86
5	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	86



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090917\
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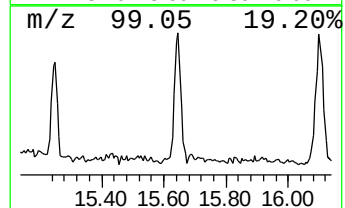
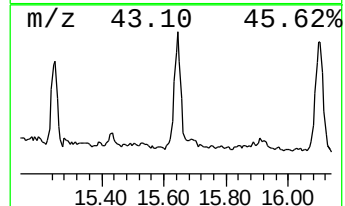
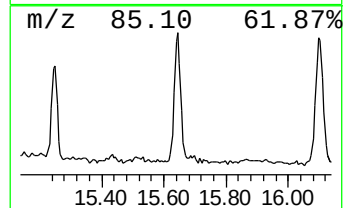
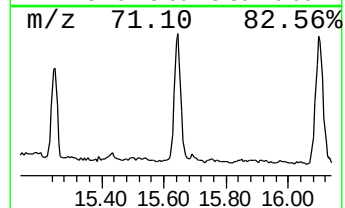
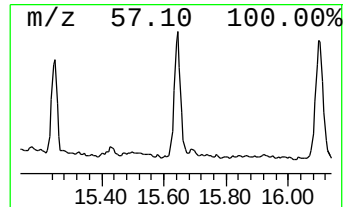
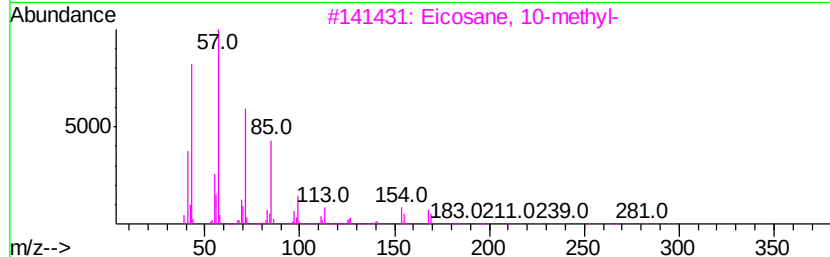
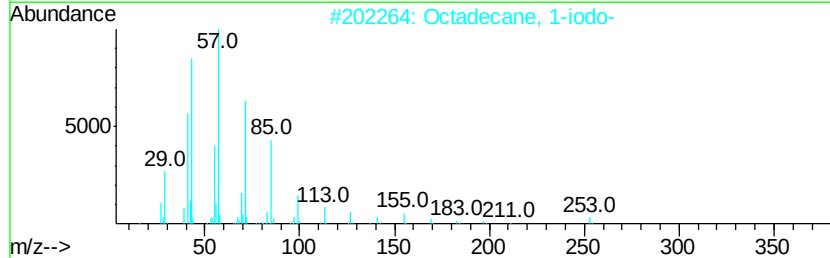
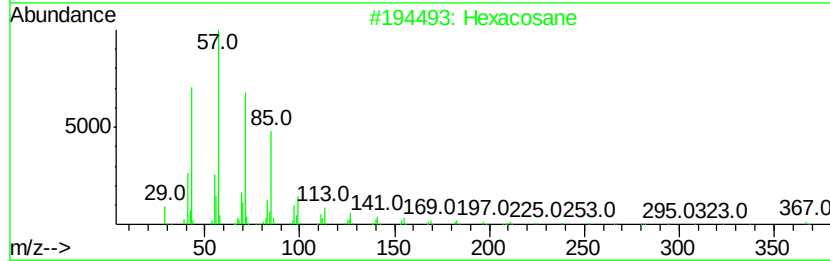
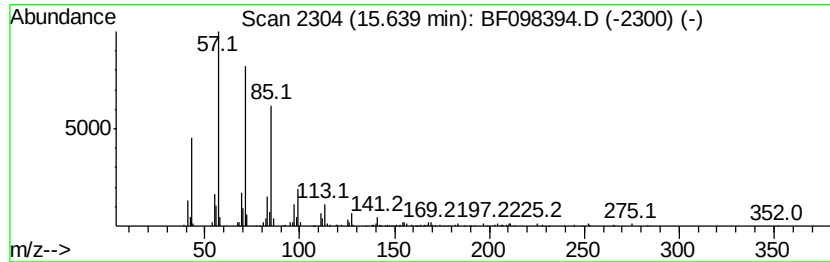
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.64	9.16 ng	285841	Perylene-d12	15.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexacosane	366 C26H54	000630-01-3	91
2	Octadecane, 1-iodo-	380 C18H37I	000629-93-6	86
3	Eicosane, 10-methyl-	296 C21H44	054833-23-7	86
4	Heptacosane	380 C27H56	000593-49-7	83
5	Heneicosane	296 C21H44	000629-94-7	76



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090917\
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Operator : SJ/JU
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ALS Vial : 19 Sample Multiplier: 1

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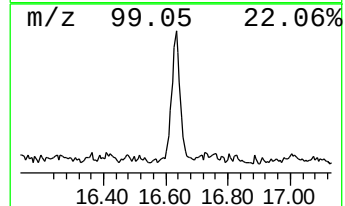
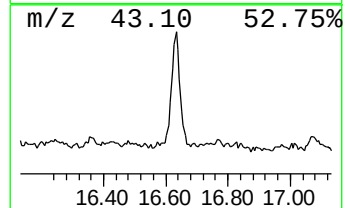
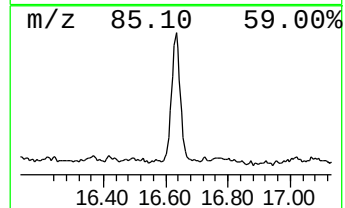
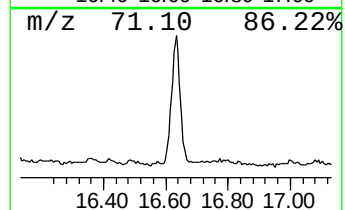
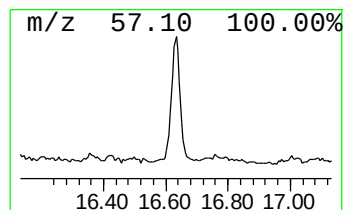
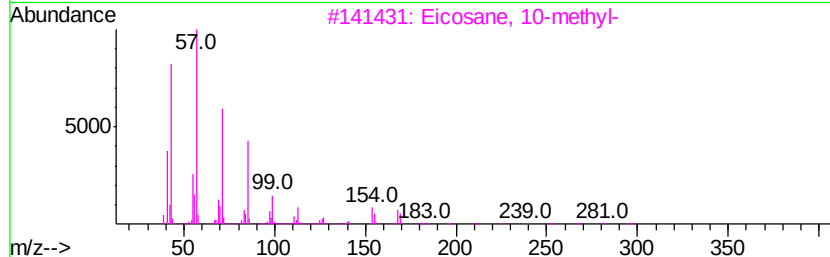
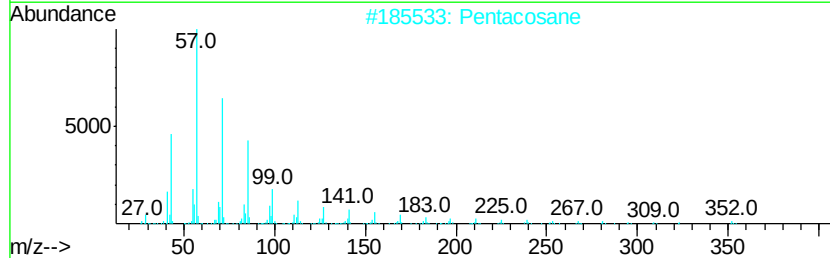
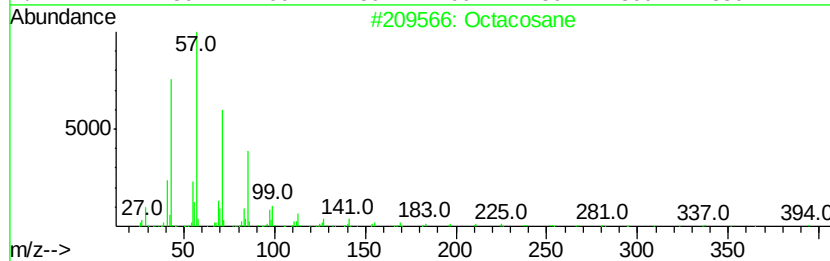
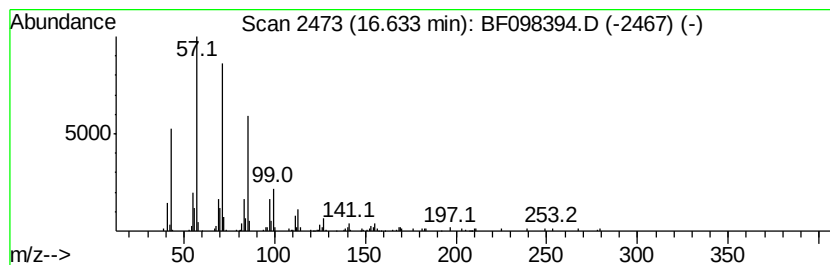
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.63	9.96 ng	310812	Perylene-d12	15.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Pentacosane	352	C25H52	000629-99-2	90
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	90
4		Hexacosane	366	C26H54	000630-01-3	87
5		Heneicosane	296	C21H44	000629-94-7	87



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090917\
Data File : BF098394.D
Acq On : 9 Sep 2017 19:07
Operator : SJ/JU
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Misc :
ALS Vial : 19 Sample Multiplier: 1

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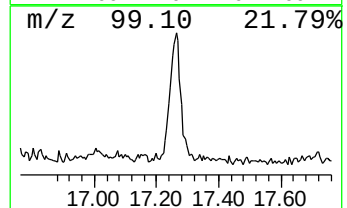
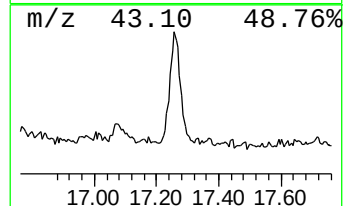
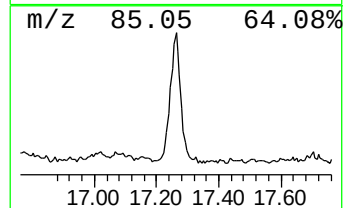
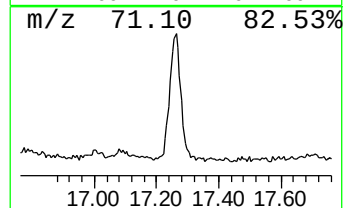
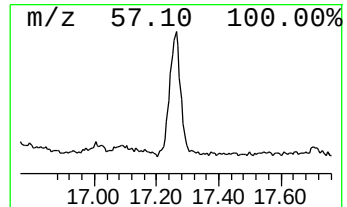
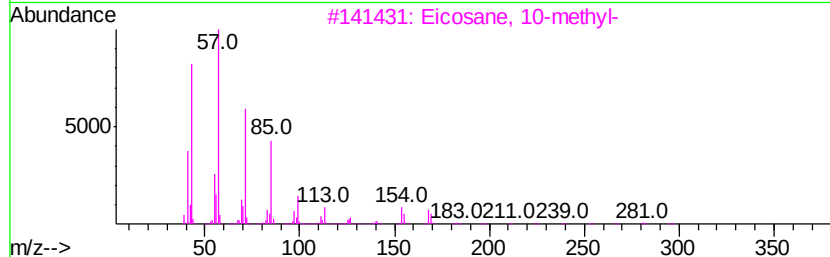
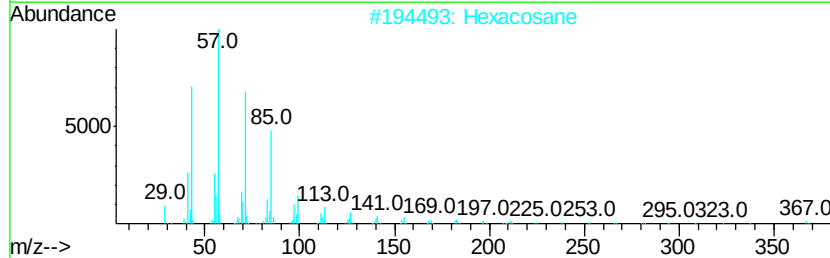
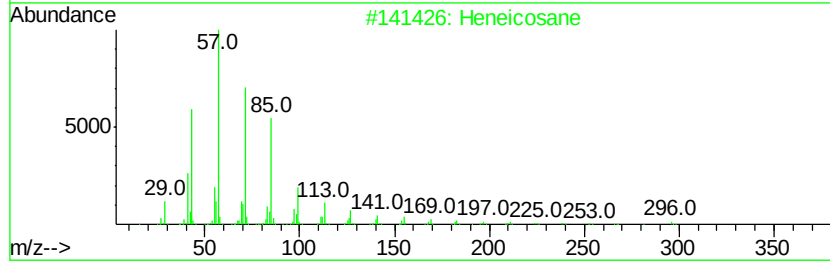
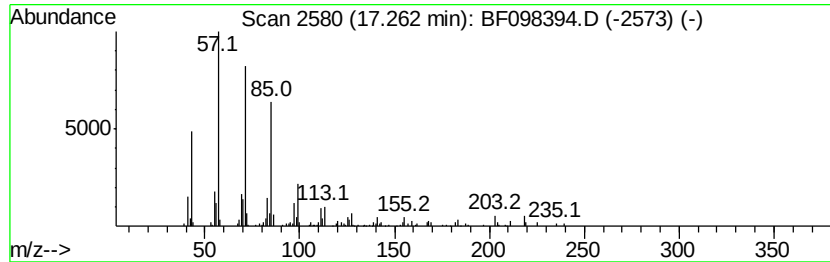
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.26	9.77 ng	304774	Perylene-d12	15.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	87
2		Hexacosane	366	C26H54	000630-01-3	87
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	86
4		2-methyloctacosane	408	C29H60	1000376-72-8	86
5		Dodecane, 3-methyl-	184	C13H28	017312-57-1	80



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090917\
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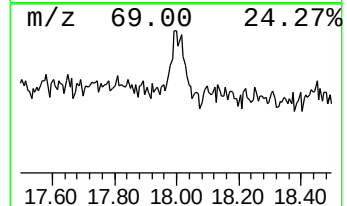
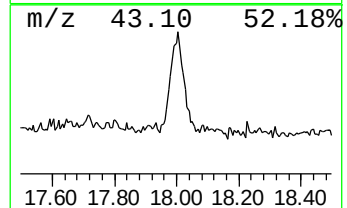
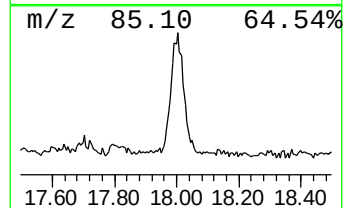
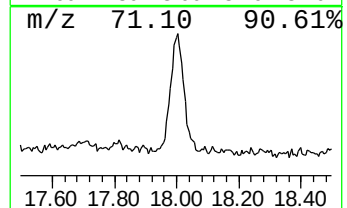
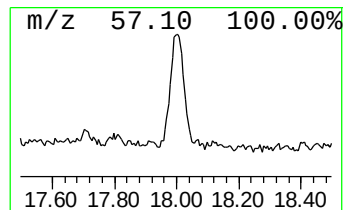
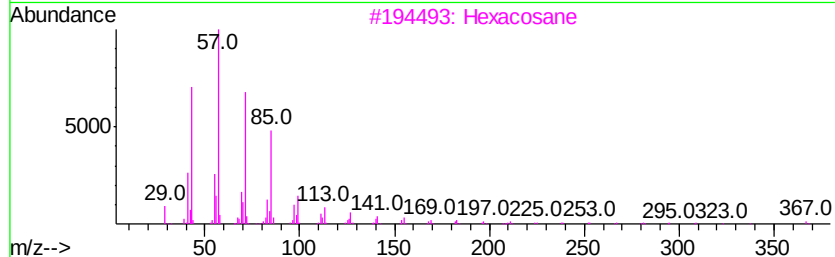
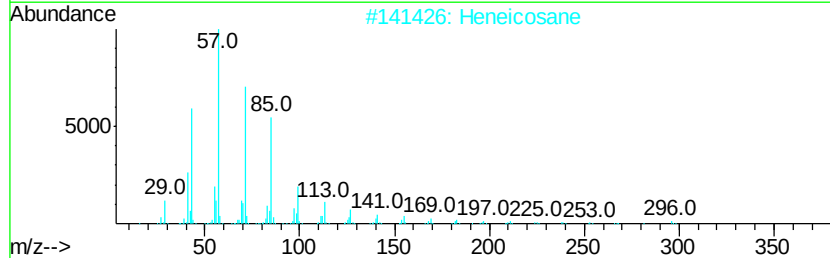
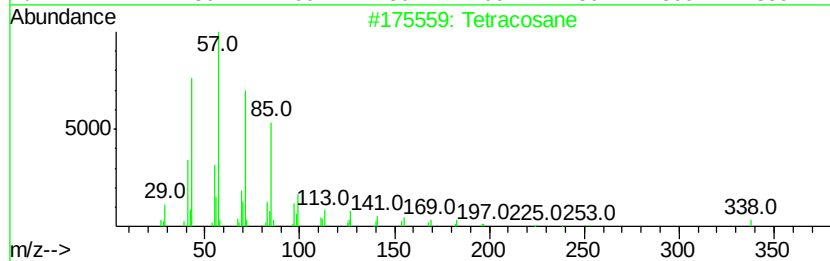
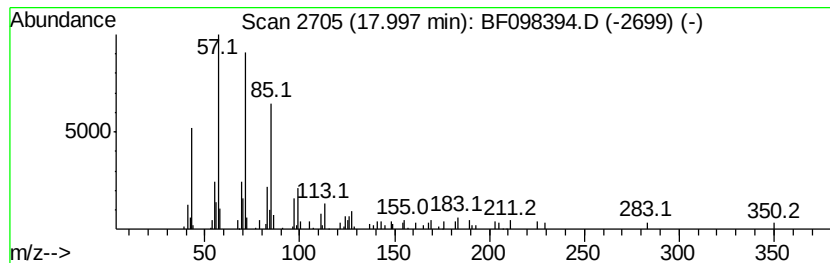
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tetracosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.00	5.79 ng	180744	Perylene-d12	15.23	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	91
2	Heneicosane	296	C21H44	000629-94-7	90
3	Hexacosane	366	C26H54	000630-01-3	90
4	Tricosane, 2-methyl-	338	C24H50	001928-30-9	90
5	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	90



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF090917\
Data File : BF098394.D
Acq On : 9 Sep 2017 19:07
Operator : SJ/JU
Sample : I5140-03
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.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
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unknown6.45	6.45	72.4	ng	2514120	1	6.67	694734	20.0
Nonadecane, 9-met...	10.62	2.8	ng	114487	4	11.19	817471	20.0
4-Heptafluorobuty...	11.42	2.3	ng	94375	4	11.19	817471	20.0
1-Hexacosanol	13.67	3.2	ng	126298	5	13.82	798446	20.0
Heptadecane, 8-me...	14.90	3.1	ng	96334	6	15.23	624173	20.0
Hexacosane	15.64	9.2	ng	285841	6	15.23	624173	20.0
Octacosane	16.63	10.0	ng	310812	6	15.23	624173	20.0
Heneicosane	17.26	9.8	ng	304774	6	15.23	624173	20.0
Tetracosane	18.00	5.8	ng	180744	6	15.23	624173	20.0