

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060319\
 Data File : BG041042.D
 Acq On : 3 Jun 2019 20:03
 Operator : HP/JU
 Sample : K3097-13
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 0433-HR-13(HACKENSACK)

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:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG052919.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	3.124	3	6	15	rVB	121368	224472	6.32%	0.754%
2	3.206	15	20	35	rVB	433998	655739	18.47%	2.202%
3	5.656	430	437	450	rBV	1283705	2063206	58.13%	6.928%
4	7.266	702	711	724	rBV	1404042	2495060	70.29%	8.378%
5	7.648	767	776	785	rBV	1280547	2269606	63.94%	7.621%
6	8.118	848	856	866	rVB	260651	439387	12.38%	1.475%
7	8.441	903	911	920	rBV	852411	1460006	41.13%	4.902%
8	9.299	1048	1057	1065	rBV	812490	1479878	41.69%	4.969%
9	10.938	1328	1336	1349	rVB	350705	635767	17.91%	2.135%
10	13.370	1741	1750	1759	rBV	1774056	2733798	77.02%	9.180%
11	14.187	1884	1889	1897	rVB	214996	317170	8.94%	1.065%
12	14.751	1978	1985	1992	rBV2	573693	908344	25.59%	3.050%
13	16.232	2230	2237	2245	rBV	1870920	2901210	81.73%	9.742%
14	17.495	2443	2452	2457	rBV	741239	1147247	32.32%	3.852%
15	17.530	2457	2458	2464	rVB	37410	43153	1.22%	0.145%
16	17.624	2467	2474	2478	rBV	25797	37703	1.06%	0.127%
17	18.341	2592	2596	2604	rBV3	125389	206445	5.82%	0.693%
18	19.263	2749	2753	2759	rBV2	33091	49091	1.38%	0.165%
19	19.534	2794	2799	2805	rBV	250580	356949	10.06%	1.199%
20	19.610	2808	2812	2818	rVB6	42264	76555	2.16%	0.257%
21	19.863	2850	2855	2857	rVV2	57541	94072	2.65%	0.316%
22	19.898	2857	2861	2868	rVV	227714	349000	9.83%	1.172%
23	19.963	2868	2872	2878	rVB2	26914	41241	1.16%	0.138%
24	20.086	2887	2893	2905	rVB	2755901	3549579	100.00%	11.919%
25	20.215	2909	2915	2921	rBV6	41779	100226	2.82%	0.337%
26	20.386	2940	2944	2949	rVB	109095	156192	4.40%	0.524%
27	20.486	2956	2961	2965	rBV	93457	134293	3.78%	0.451%
28	20.544	2965	2971	2974	rVB2	41294	74409	2.10%	0.250%
29	20.679	2991	2994	2998	rBV	34631	51521	1.45%	0.173%
30	20.726	2999	3002	3005	rBV3	28623	39815	1.12%	0.134%
31	21.367	3102	3111	3119	rBV5	77645	225865	6.36%	0.758%
32	21.437	3119	3123	3127	rVB3	26677	37509	1.06%	0.126%
33	21.772	3171	3180	3185	rBV2	773968	1580312	44.52%	5.306%
34	21.819	3185	3188	3194	rVB	147501	233145	6.57%	0.783%

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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

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35	21.925	3201	3206	3210	rVB5	37361	64427	1.82%	0.216%
36	21.972	3210	3214	3219	rBV7	23445	54631	1.54%	0.183%
37	23.253	3427	3432	3438	rBV5	41177	80339	2.26%	0.270%
38	23.447	3462	3465	3477	rVB4	43868	91659	2.58%	0.308%
39	24.040	3558	3566	3570	rBV2	98652	293773	8.28%	0.986%
40	24.299	3607	3610	3618	rVB3	24166	57094	1.61%	0.192%
41	24.786	3688	3693	3702	rVB2	60041	162454	4.58%	0.545%
42	24.939	3711	3719	3732	rBV	93189	284818	8.02%	0.956%
43	25.110	3732	3748	3758	rBV2	404834	1325782	37.35%	4.452%
44	26.226	3934	3938	3947	rVB9	29734	75300	2.12%	0.253%
45	28.923	4387	4397	4409	rBV5	26887	122953	3.46%	0.413%

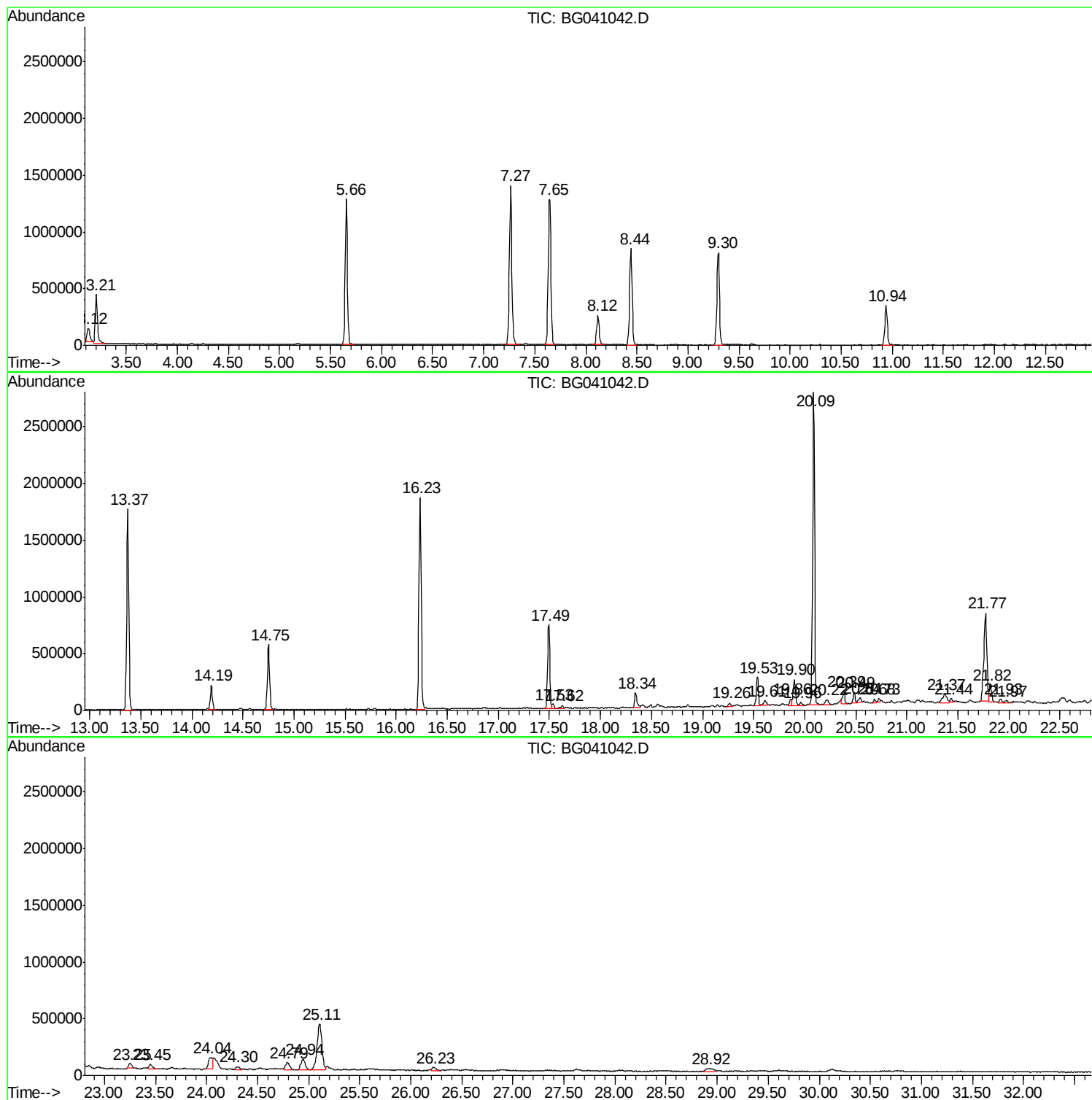
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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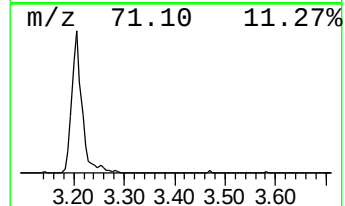
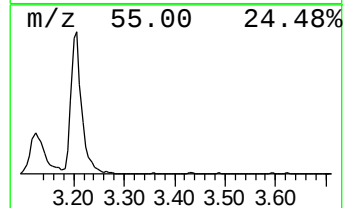
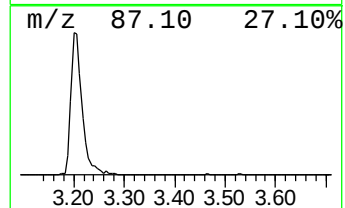
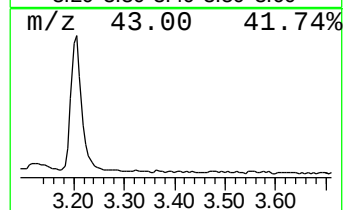
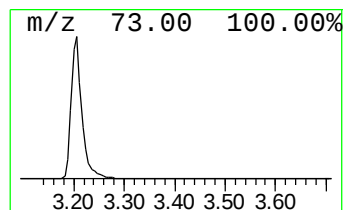
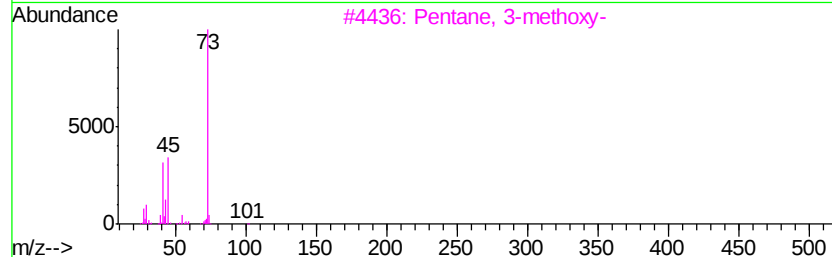
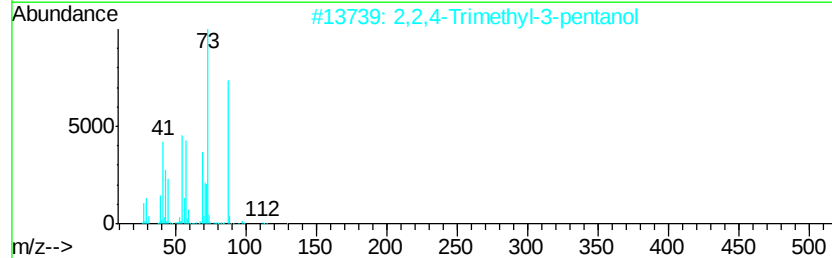
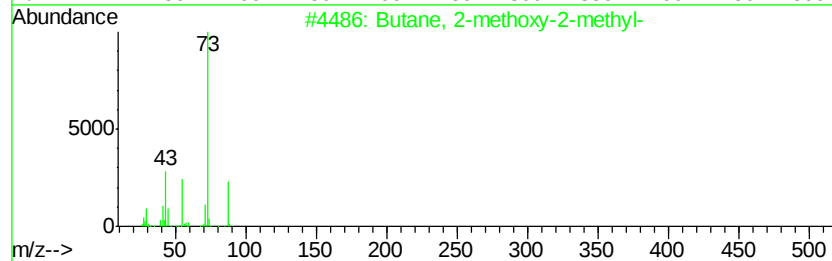
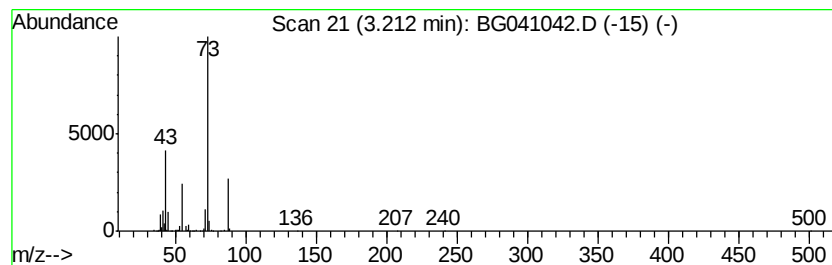
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.21	29.85 ng	655739	1,4-Dichlorobenzene-d4	8.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	38
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	9



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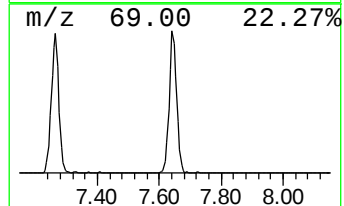
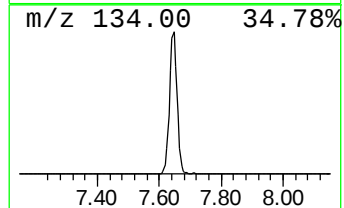
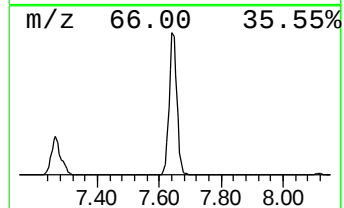
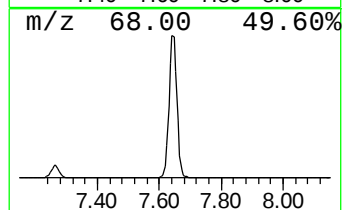
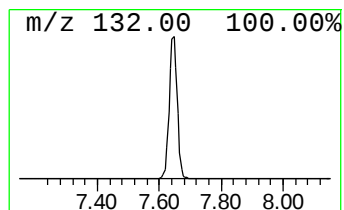
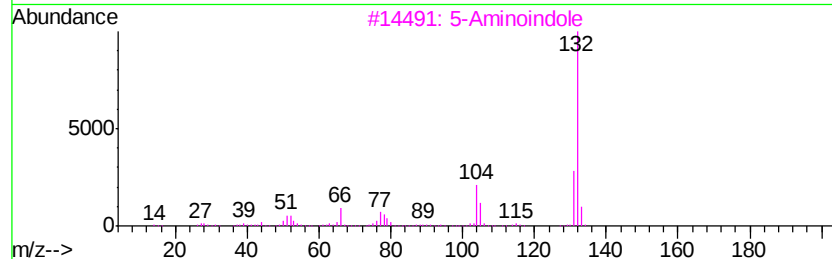
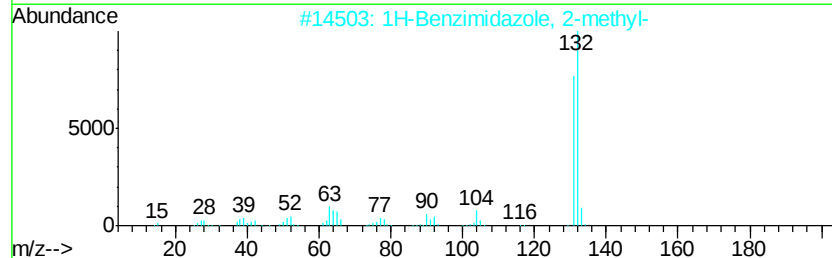
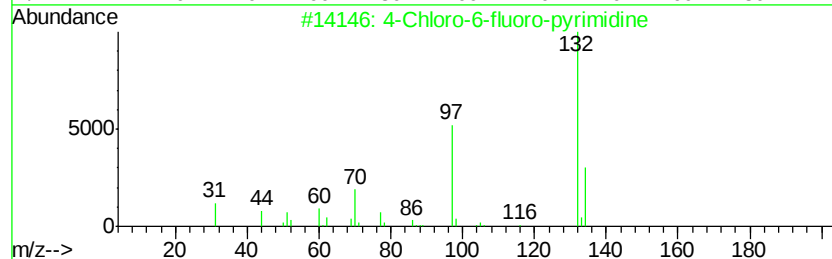
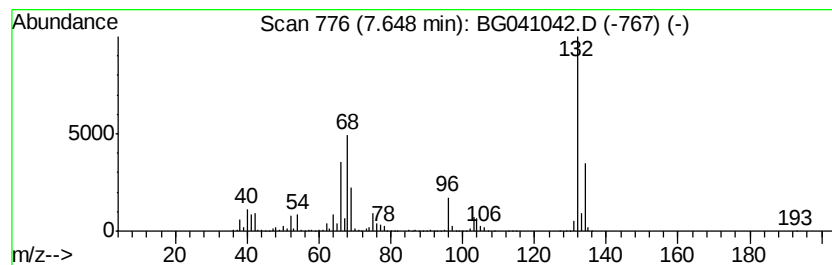
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TIC Library : C:\Database\NIST11.L
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Peak Number 3 unknown7.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.65	103.31 ng	2269610	1,4-Dichlorobenzene-d4	8.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		5-Aminoindole	132	C8H8N2	005192-03-0	22
4		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	14



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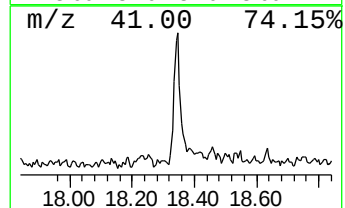
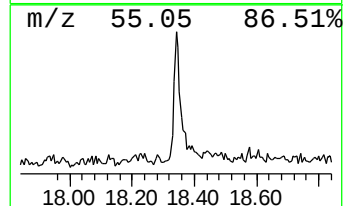
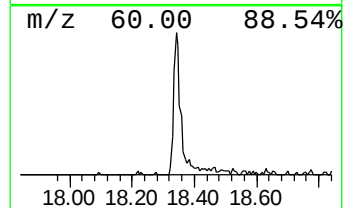
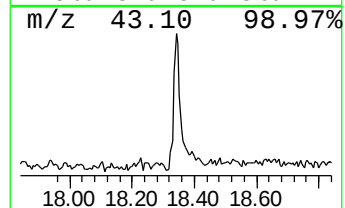
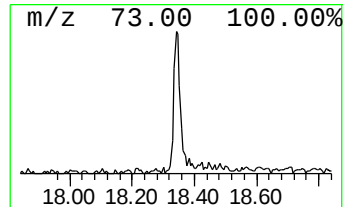
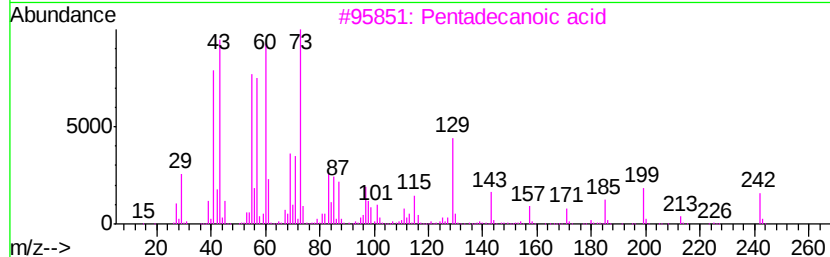
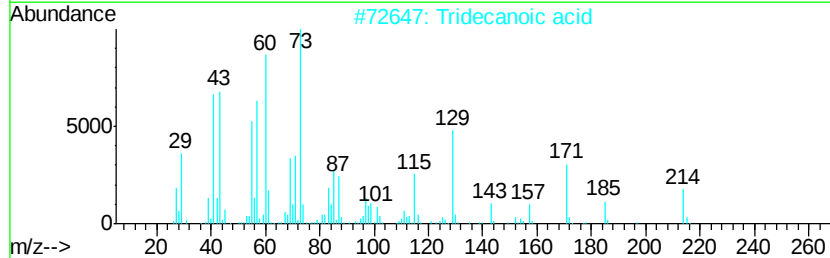
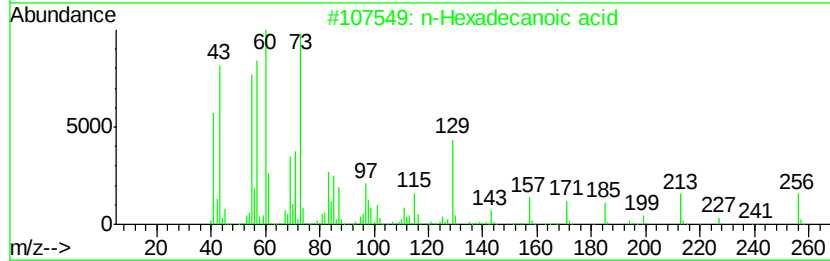
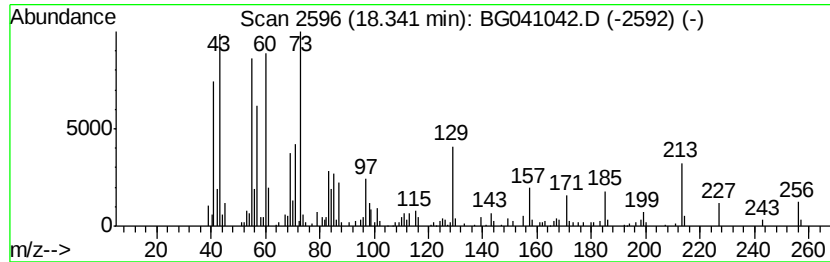
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.34	3.60 ng	206445	Phenanthrene-d10	17.49	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2	Tridecanoic acid	214	C13H26O2	000638-53-9	91
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	83
4	Undecanoic acid	186	C11H22O2	000112-37-8	80
5	Dodecanoic acid	200	C12H24O2	000143-07-7	72



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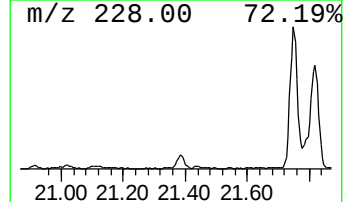
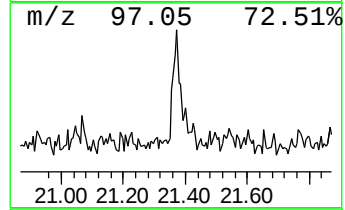
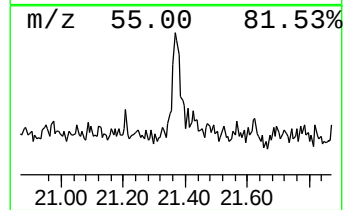
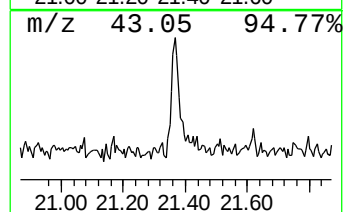
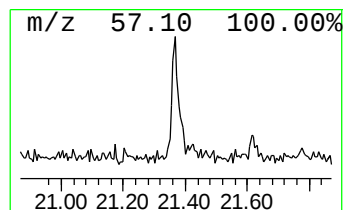
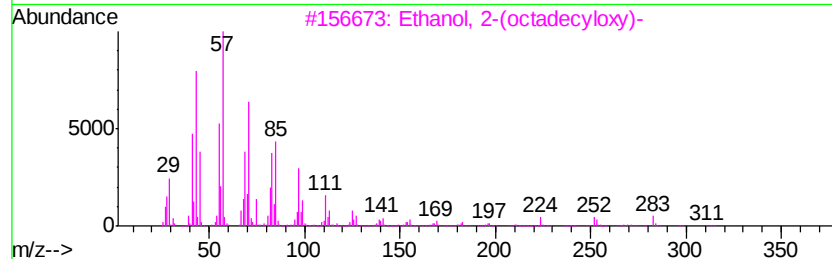
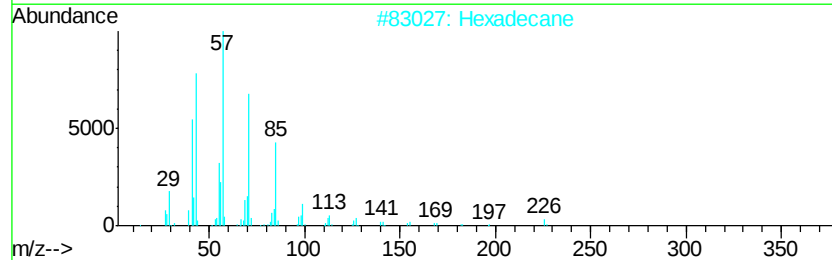
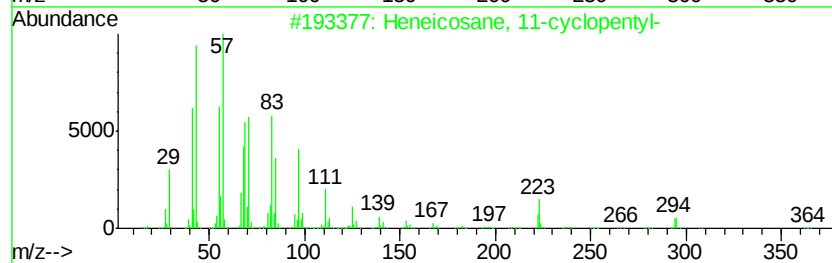
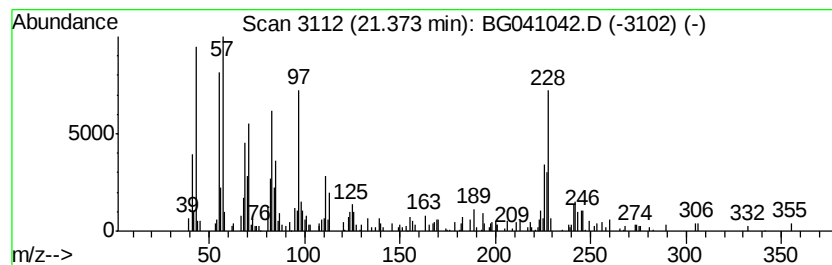
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Peak Number 6 Heneicosane, 11-cyclopentyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.37	2.86 ng	225865	Chrysene-d12	21.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane, 11-cvclopentvl-	364	C26H52	006703-81-7	64
2		Hexadecane	226	C16H34	000544-76-3	56
3		Ethanol, 2-(octadecvloxy)-	314	C20H42O2	002136-72-3	52
4		Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	52
5		Oxalic acid, propyl tetradecyl e...	328	C19H36O4	1000309-26-7	52



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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...	3.21	29.9	ng	655739	1	8.12	439387	20.0
unknown7.65	7.65	103.3	ng	2269610	1	8.12	439387	20.0
n-Hexadecanoic acid	18.34	3.6	ng	206445	4	17.49	1147250	20.0
Heneicosane, 11-c...	21.37	2.9	ng	225865	5	21.77	1580310	20.0