

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG030220\
 Data File : BG044638.D
 Acq On : 2 Mar 2020 18:29
 Operator : CG/JU
 Sample : L1740-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GP-08-022720

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-BG022420.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	5.372	383	389	404	rBV	652716	1046610	30.93%	6.518%
2	6.982	655	663	681	rBV	665145	1153504	34.08%	7.184%
3	7.805	795	803	811	rBV	164564	277724	8.21%	1.730%
4	8.128	850	858	871	rBV	527529	909540	26.88%	5.665%
5	8.962	991	1000	1013	rBV	386901	722045	21.34%	4.497%
6	10.602	1272	1279	1295	rBV	216694	412682	12.19%	2.570%
7	13.075	1693	1700	1716	rBV	1215567	1897064	56.06%	11.815%
8	13.915	1836	1843	1855	rBV	45383	80133	2.37%	0.499%
9	14.456	1928	1935	1945	rBV	431628	668536	19.75%	4.164%
10	15.942	2182	2188	2203	rBV	988229	1534217	45.33%	9.555%
11	17.200	2395	2402	2416	rBV	603376	893107	26.39%	5.562%
12	19.820	2842	2848	2861	rBV	2513808	3384247	100.00%	21.078%
13	21.113	3063	3068	3079	rBV2	115108	224300	6.63%	1.397%
14	21.436	3116	3123	3135	rBV	652003	1035632	30.60%	6.450%
15	21.642	3154	3158	3163	rBV2	29057	41586	1.23%	0.259%
16	22.212	3250	3255	3264	rBV	49932	90411	2.67%	0.563%
17	22.864	3356	3366	3373	rBV	73450	136666	4.04%	0.851%
18	23.616	3487	3494	3501	rVB2	69078	137309	4.06%	0.855%
19	24.450	3626	3636	3658	rBV2	400791	1230437	36.36%	7.663%
20	25.543	3813	3822	3833	rVB2	42702	117781	3.48%	0.734%
21	26.800	4027	4036	4045	rBV3	20132	62475	1.85%	0.389%

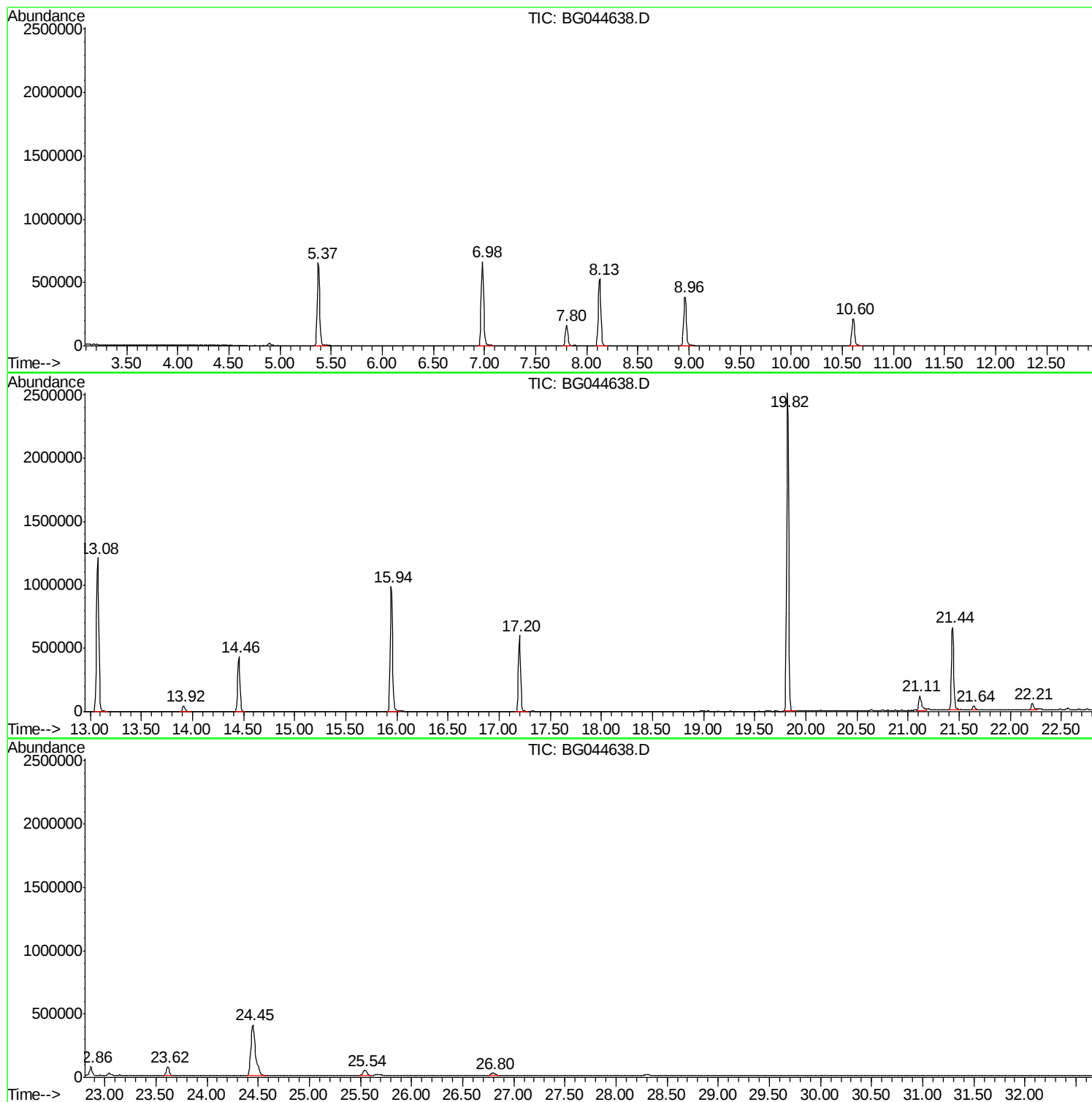
Sum of corrected areas: 16056006

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG022420.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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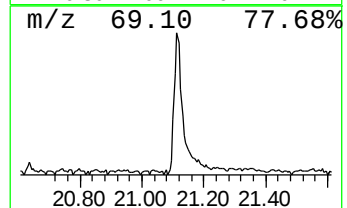
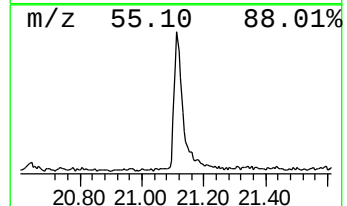
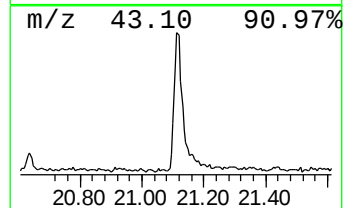
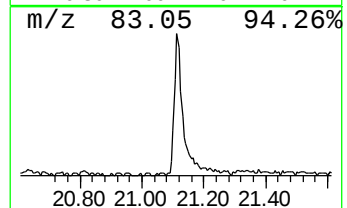
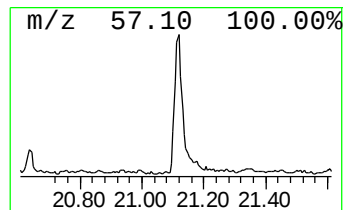
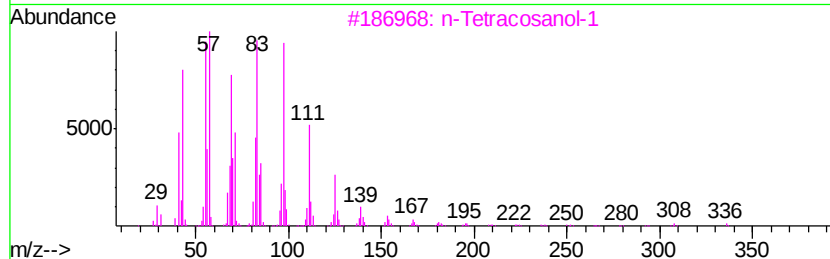
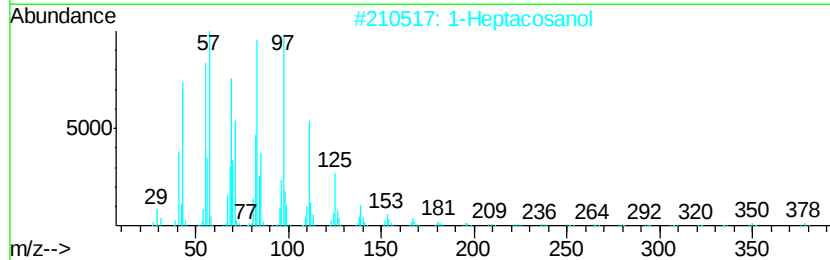
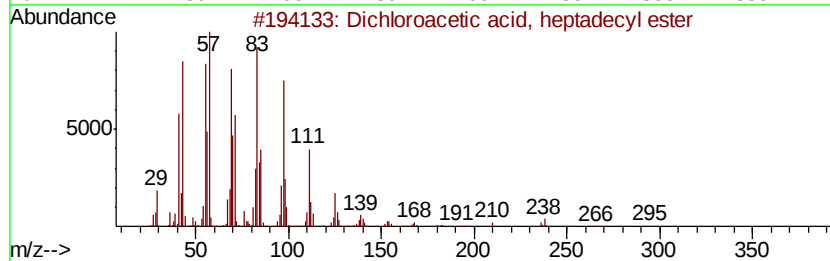
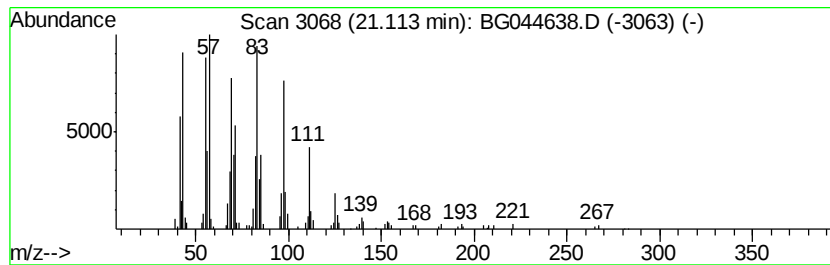
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG022420.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Dichloroacetic acid, heptad... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.11	4.33 ng	224300	Chrysene-d12	21.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	95
2		1-Heptacosanol	396	C27H56O	002004-39-9	94
3		n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4		1-Heneicosanol	312	C21H44O	015594-90-8	93
5		Behenic alcohol	326	C22H46O	000661-19-8	91



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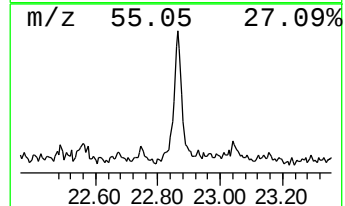
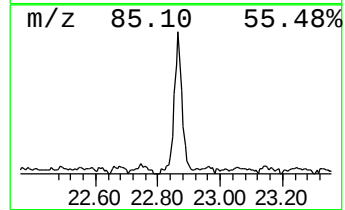
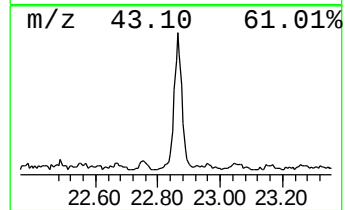
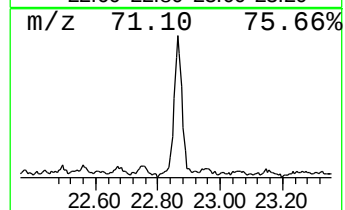
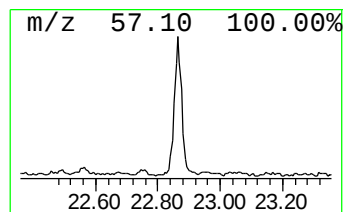
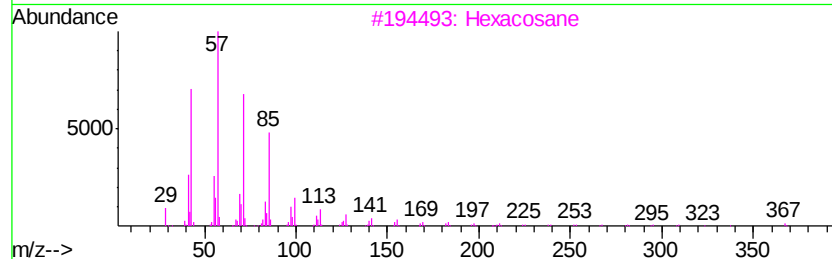
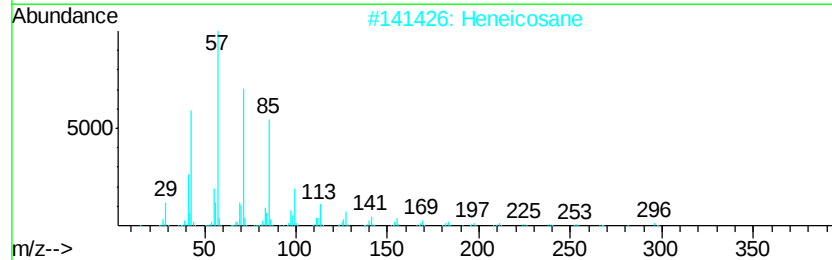
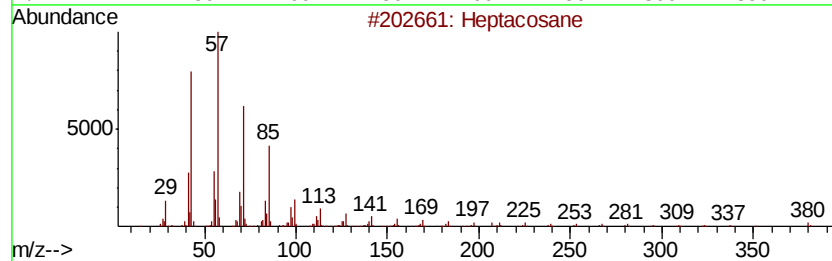
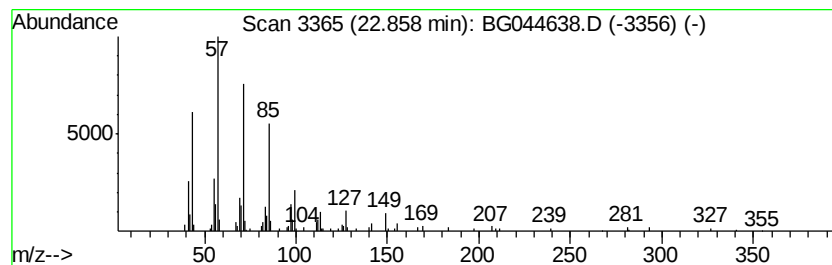
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Peak Number 3 Heptacosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.86	2.64 ng	136666	Chrysene-d12	21.44

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane		380	C27H56	000593-49-7	87
2	Heneicosane		296	C21H44	000629-94-7	87
3	Hexacosane		366	C26H54	000630-01-3	87
4	Octacosane		394	C28H58	000630-02-4	83
5	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	83



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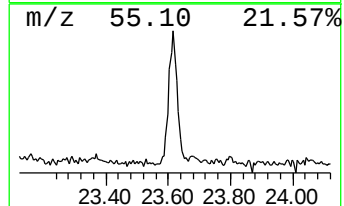
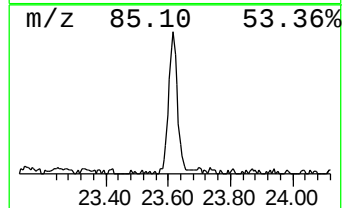
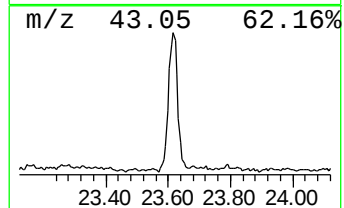
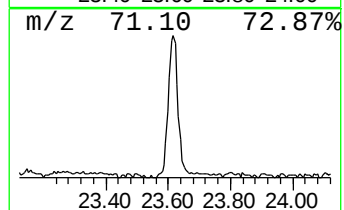
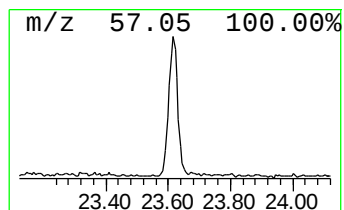
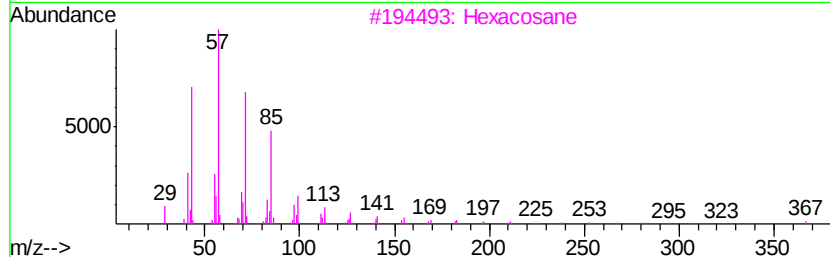
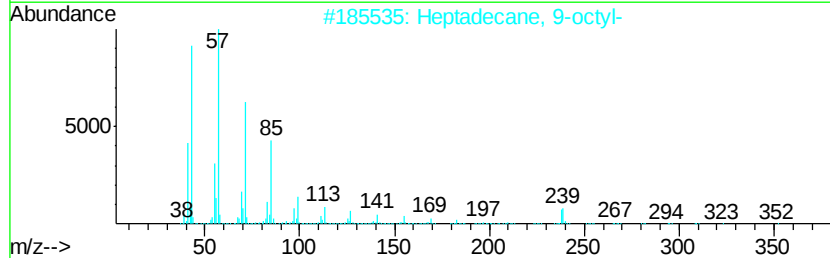
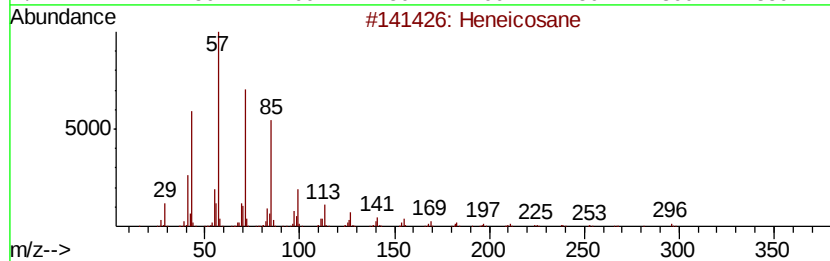
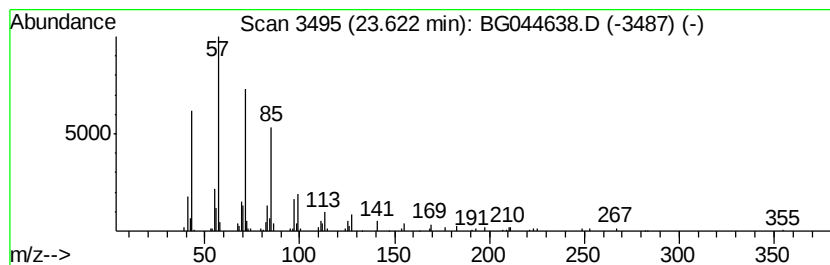
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Peak Number 4 Heneicosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.62	2.23 ng	137309	Perylene-d12	24.45

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	91
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3			Hexacosane	366	C26H54	000630-01-3	90
4			Eicosane	282	C20H42	000112-95-8	90
5			Heptadecane	240	C17H36	000629-78-7	87



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Dichloroacetic ac...	21.11	4.3	ng	224300	5	21.44	1035630	20.0
Heptacosane	22.86	2.6	ng	136666	5	21.44	1035630	20.0
Heneicosane	23.62	2.2	ng	137309	6	24.45	1230440	20.0