

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG013120\
 Data File : BG044266.D
 Acq On : 31 Jan 2020 19:05
 Operator : CG/JU
 Sample : L1331-02
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C-2_013020

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-BG013020.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	5.015	323	328	340	rVB	27126	49239	1.18%	0.220%
2	5.491	402	409	422	rBV	873450	1402153	33.73%	6.266%
3	7.083	672	680	698	rBV	944376	1712541	41.19%	7.652%
4	7.918	813	822	830	rBV	221986	389447	9.37%	1.740%
5	8.241	867	877	885	rBV	696146	1225304	29.47%	5.475%
6	9.069	1010	1018	1036	rBV	566904	1044565	25.12%	4.668%
7	10.714	1288	1298	1309	rBV	317852	612214	14.73%	2.736%
8	13.176	1710	1717	1732	rBV	1787128	2843202	68.39%	12.705%
9	14.005	1852	1858	1868	rBV	213252	323602	7.78%	1.446%
10	14.551	1944	1951	1963	rBV	674643	1071309	25.77%	4.787%
11	16.043	2198	2205	2225	rBV	1483539	2293099	55.15%	10.247%
12	17.295	2410	2418	2430	rBV	811013	1281899	30.83%	5.728%
13	19.915	2857	2864	2876	rBV	2978822	4157562	100.00%	18.578%
14	21.214	3080	3085	3096	rBV	141340	277041	6.66%	1.238%
15	21.543	3134	3141	3150	rBV	792234	1354668	32.58%	6.053%
16	21.754	3172	3177	3182	rBV	52054	73682	1.77%	0.329%
17	22.342	3272	3277	3282	rBV2	63797	96649	2.32%	0.432%
18	23.012	3385	3391	3397	rVB	78526	136993	3.30%	0.612%
19	23.194	3415	3422	3431	rBV	107672	199709	4.80%	0.892%
20	23.787	3516	3523	3530	rBV	75277	163894	3.94%	0.732%
21	24.639	3658	3668	3689	rBV2	408685	1526233	36.71%	6.820%
22	25.779	3853	3862	3874	rBV3	46575	143934	3.46%	0.643%

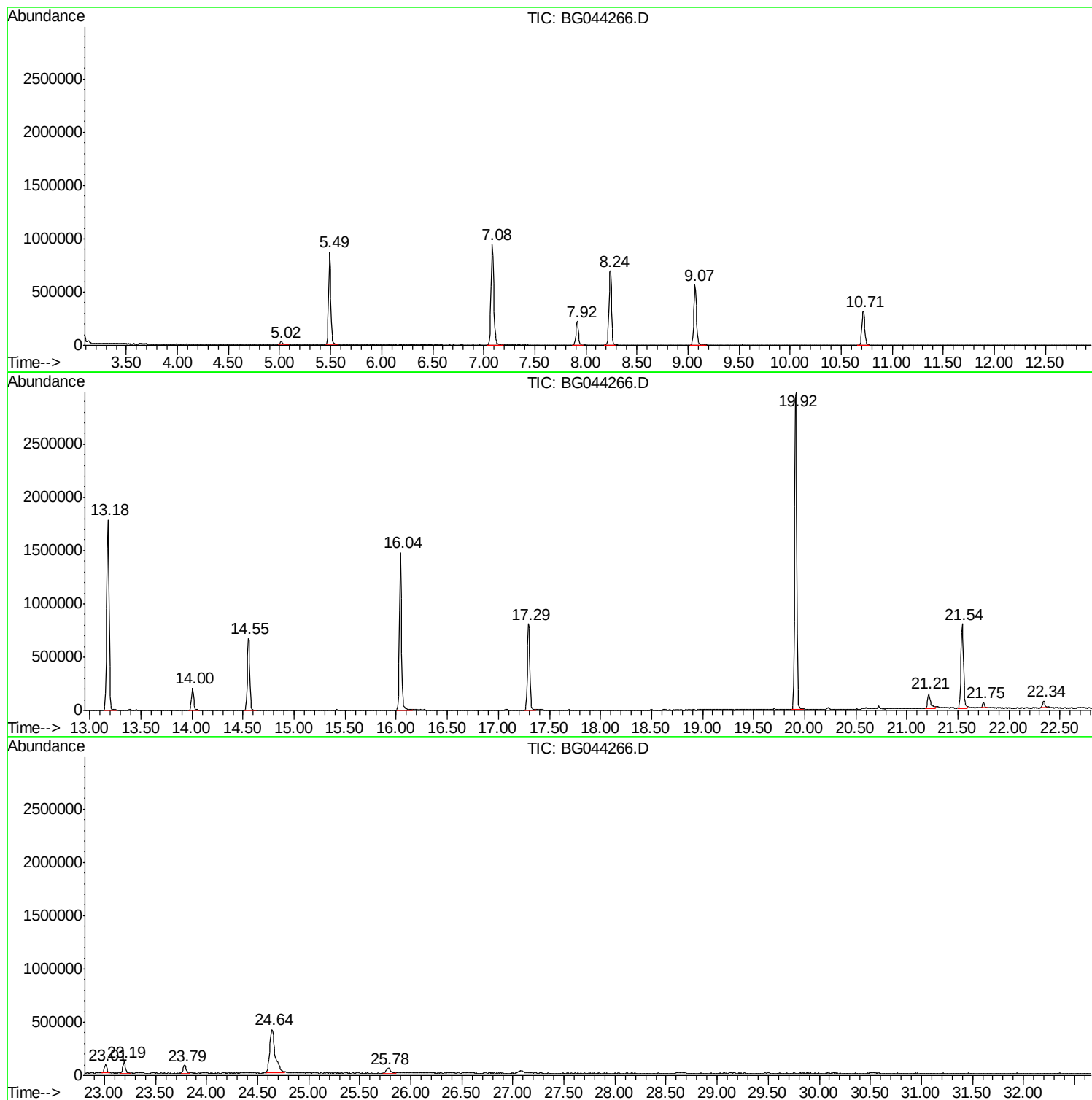
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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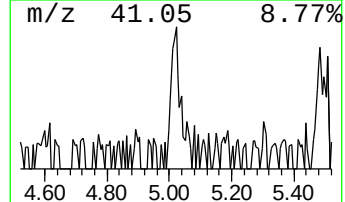
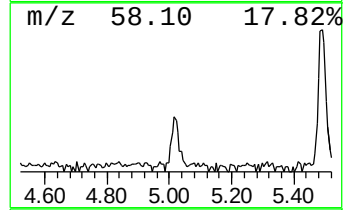
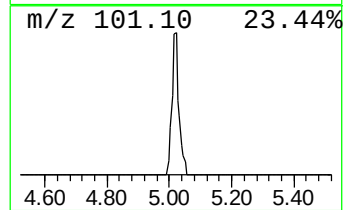
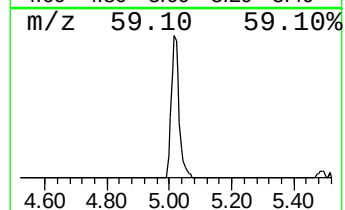
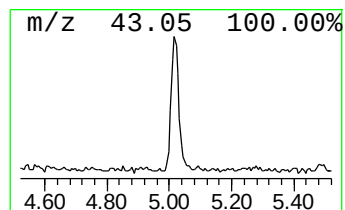
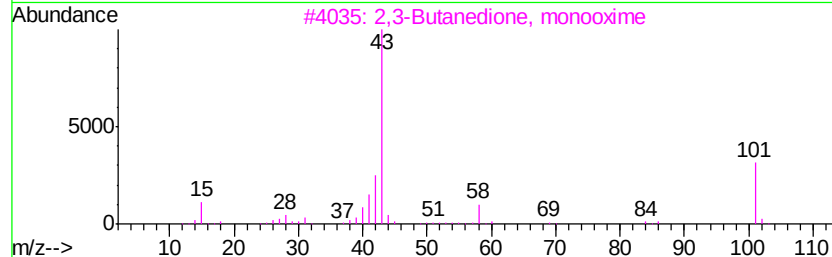
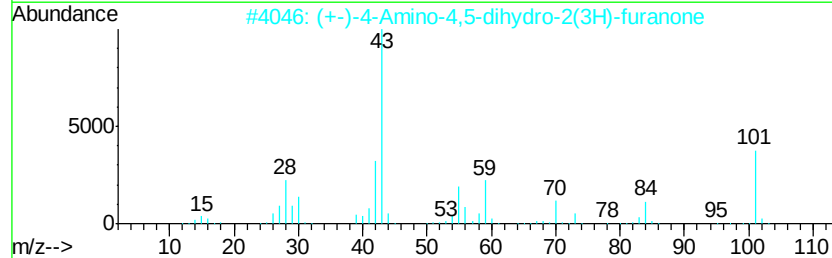
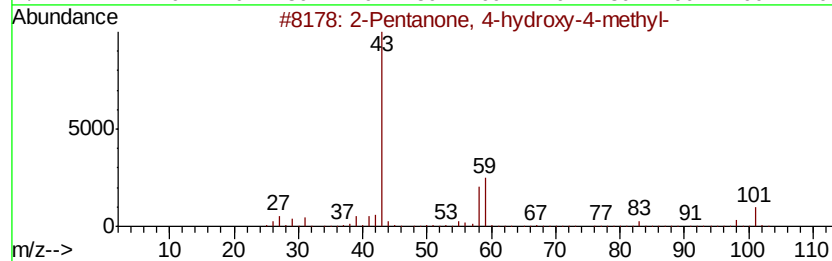
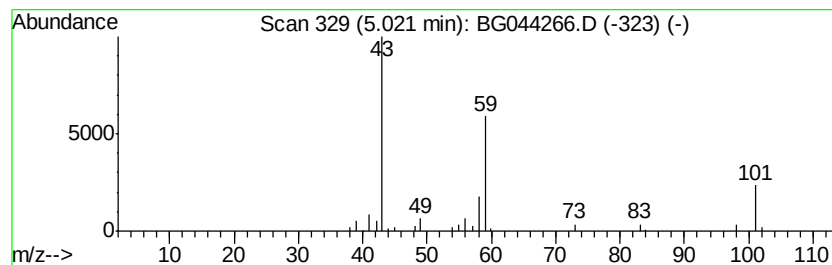
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.02	2.53 ng	49239	1,4-Dichlorobenzene-d4	7.91

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	17
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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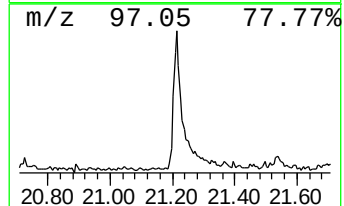
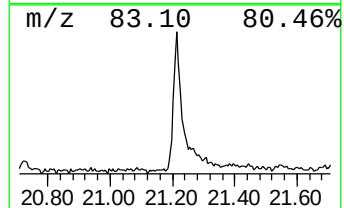
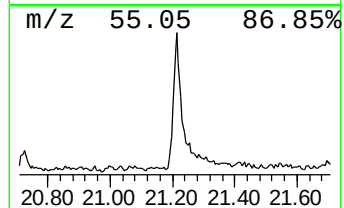
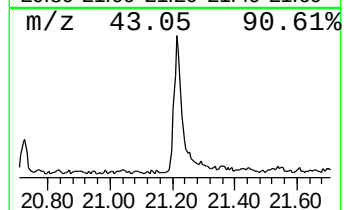
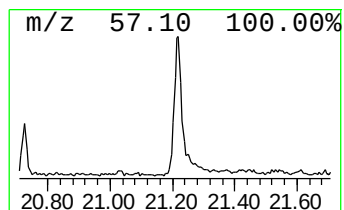
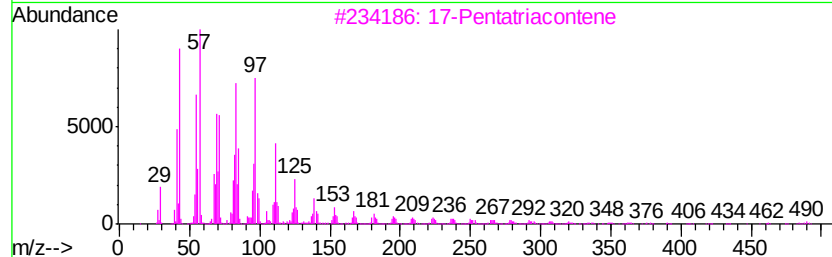
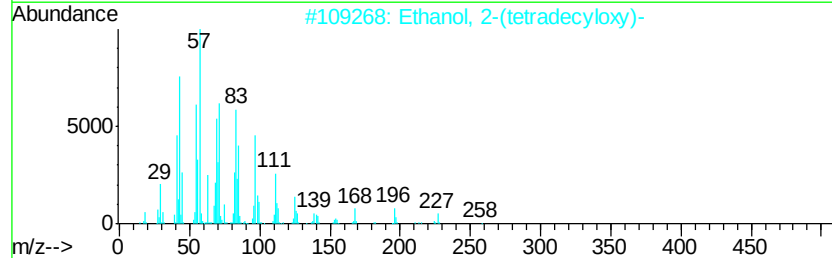
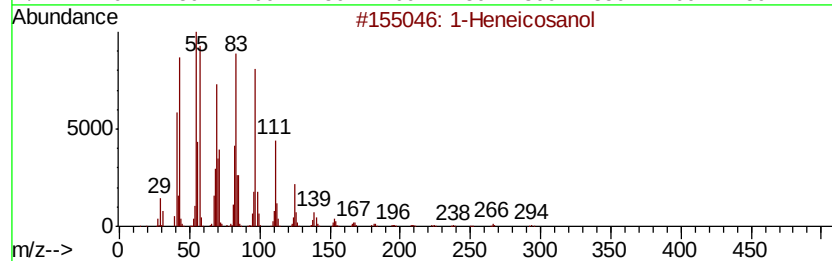
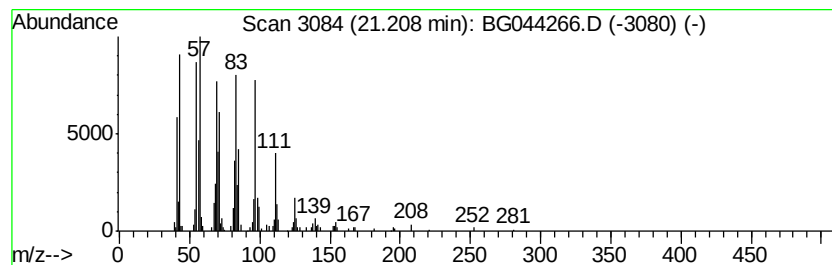
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1-Heneicosanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.21	4.09 ng	277041	Chrysene-d12	21.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Heneicosanol	312	C21H44O	015594-90-8	93
2		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
3		17-Pentatriacontene	491	C35H70	006971-40-0	91
4		Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	90
5		Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	81



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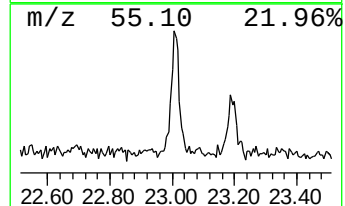
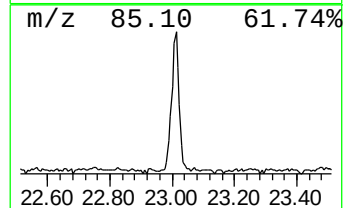
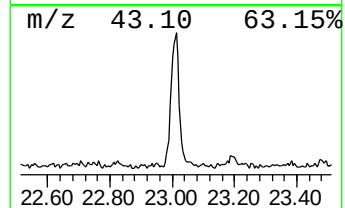
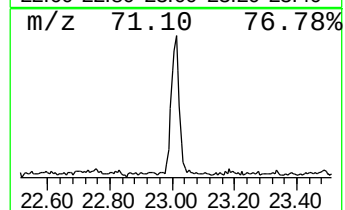
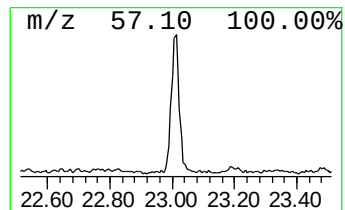
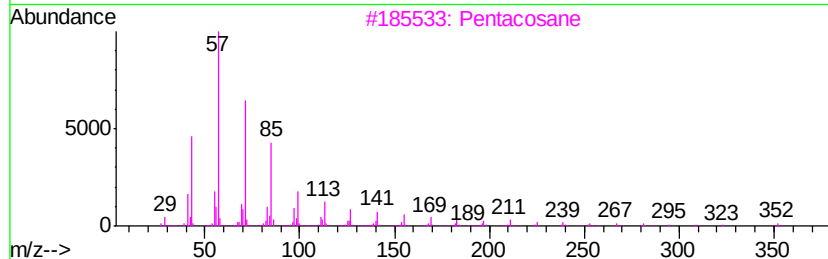
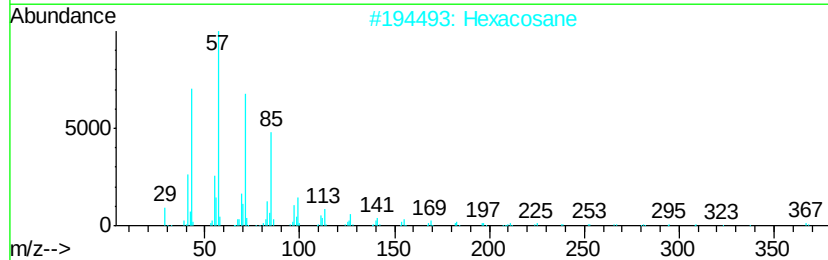
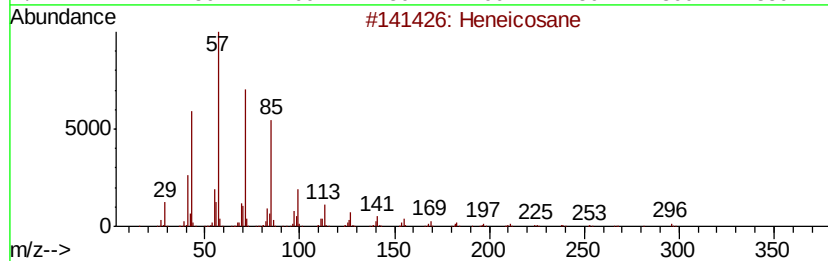
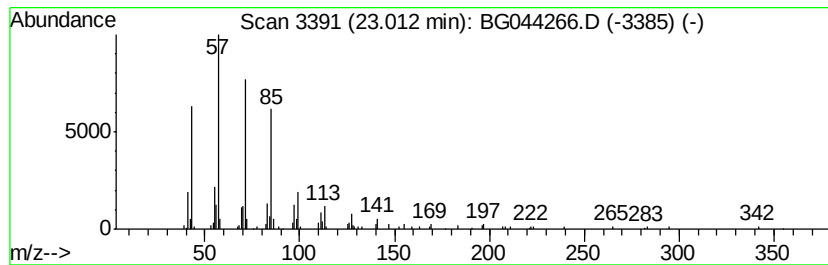
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Peak Number 4 Heneicosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.01	2.02 ng	136993	Chrysene-d12	21.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Hexacosane	366	C26H54	000630-01-3	91
3		Pentacosane	352	C25H52	000629-99-2	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Heptadecane	240	C17H36	000629-78-7	91



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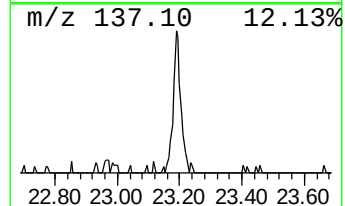
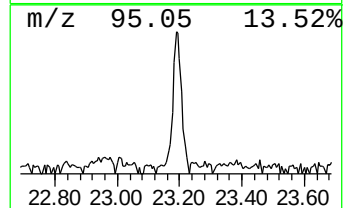
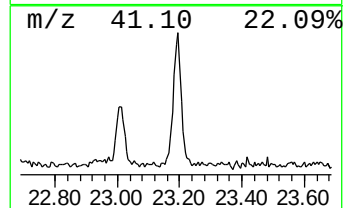
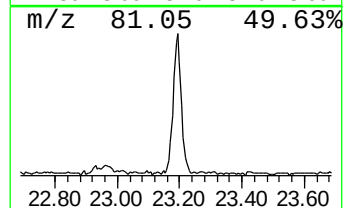
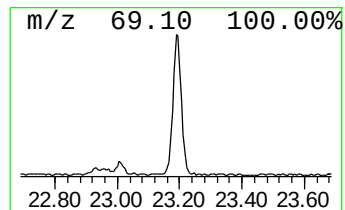
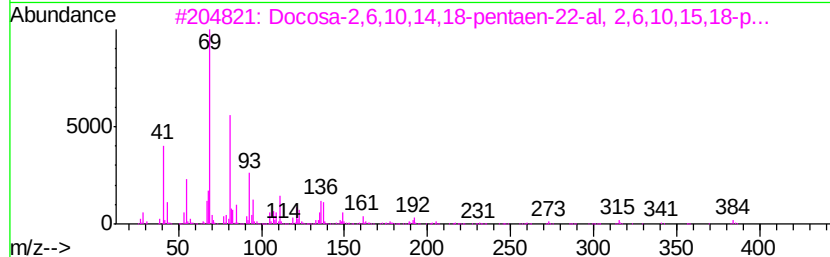
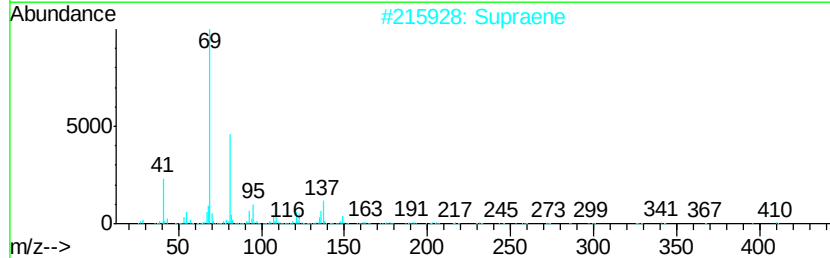
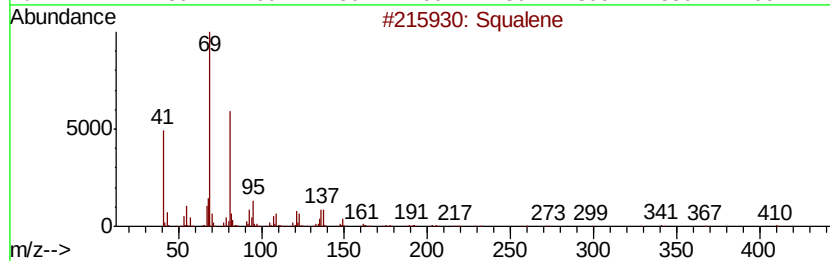
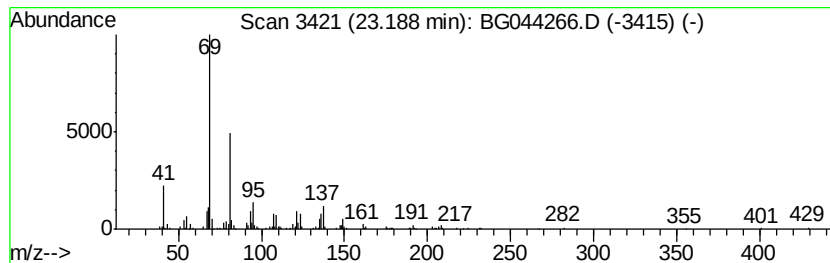
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Peak Number 5 Squalene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.19	2.62 ng	199709	Perylene-d12	24.64
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Squalene		410 C30H50	000111-02-4 91
2	Supraene		410 C30H50	007683-64-9 91
3	Docosa-2,6,10,14,18-pentaen-22-a...		384 C27H44O	1000163-04-7 83
4	7,11-Dimethyldodeca-2,6,10-trien...		208 C14H24O	125529-10-4 80
5	(.+/-)-Lavandulol, pentafluorop...		300 C13H17F5O2	1000352-63-1 72



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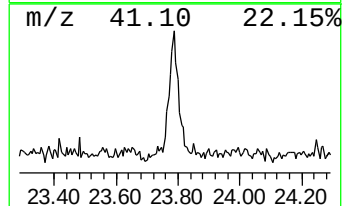
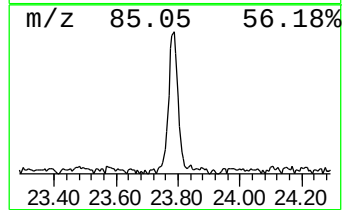
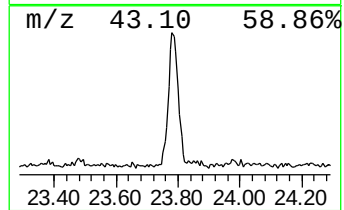
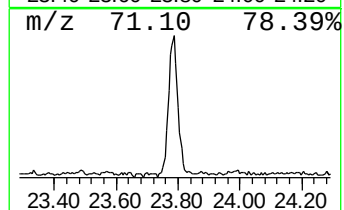
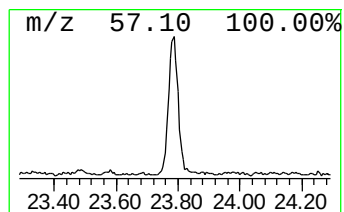
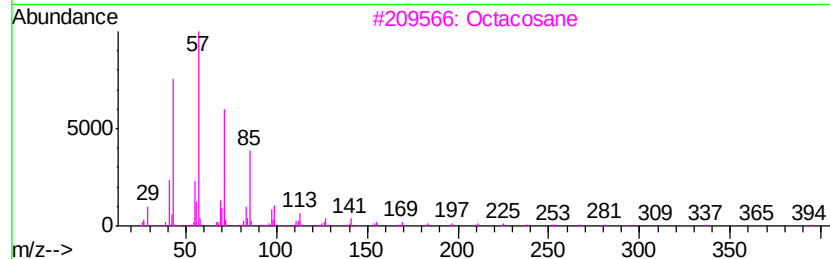
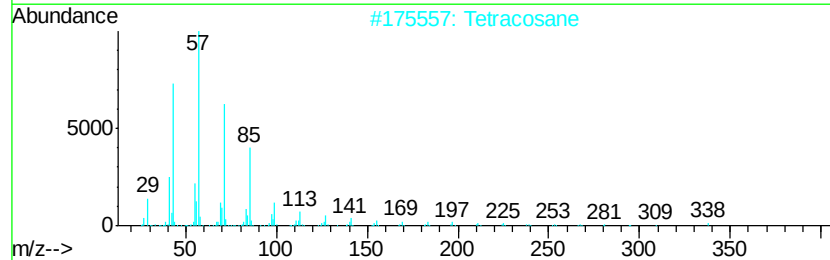
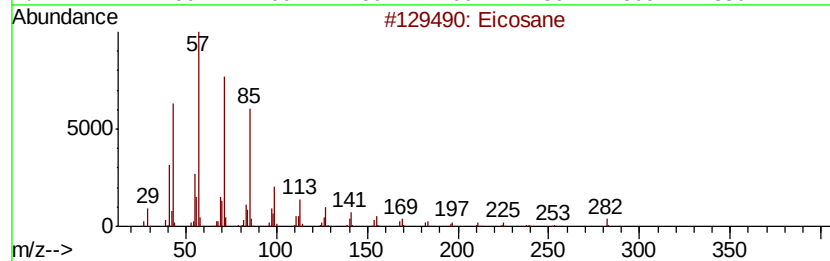
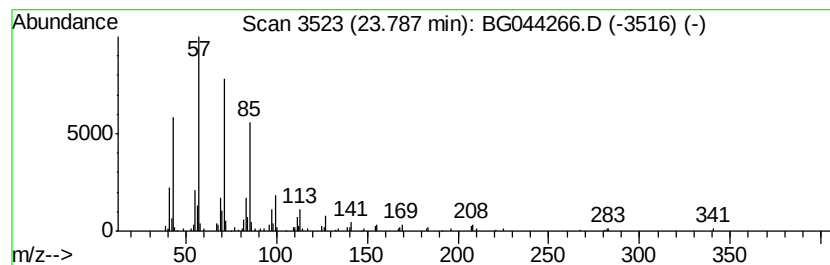
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Peak Number 6 Eicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.79	2.15 ng	163894	Perylene-d12	24.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	94
2		Tetracosane	338	C24H50	000646-31-1	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Docosane, 7-hexyl-	394	C28H58	055373-86-9	91



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.02	2.5	ng	49239	1	7.91	389447	20.0
1-Heneicosanol	21.21	4.1	ng	277041	5	21.54	1354670	20.0
Heneicosane	23.01	2.0	ng	136993	5	21.54	1354670	20.0
Squalene	23.19	2.6	ng	199709	6	24.64	1526230	20.0
Eicosane	23.79	2.1	ng	163894	6	24.64	1526230	20.0