

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090420\
 Data File : BG046539.D
 Acq On : 4 Sep 2020 13:55
 Operator : CG/JU
 Sample : L3894-09
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 AL-FALAH-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:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG082520.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	3.135	3	8	26	rVB	240375	472082	11.45%	2.362%
2	5.050	328	334	344	rVB	36761	58555	1.42%	0.293%
3	5.520	406	414	426	rBV	881536	1441858	34.98%	7.213%
4	7.101	675	683	701	rBV	837058	1506313	36.54%	7.536%
5	7.929	817	824	832	rVB	251149	410744	9.96%	2.055%
6	9.081	1012	1020	1036	rBV	546642	1014683	24.61%	5.076%
7	10.726	1292	1300	1314	rBV	317505	579170	14.05%	2.897%
8	13.182	1711	1718	1727	rBV	1586592	2464177	59.78%	12.327%
9	14.010	1853	1859	1871	rBV	114922	187322	4.54%	0.937%
10	14.557	1943	1952	1965	rBV2	532590	839740	20.37%	4.201%
11	16.043	2199	2205	2227	rBV	1092354	1817077	44.08%	9.090%
12	17.301	2412	2419	2434	rBV	694436	1070370	25.97%	5.355%
13	18.200	2567	2572	2580	rBV2	37729	66305	1.61%	0.332%
14	19.915	2858	2864	2877	rBV	2997831	4122290	100.00%	20.622%
15	21.126	3064	3070	3073	rBV2	31269	52032	1.26%	0.260%
16	21.214	3080	3085	3094	rBV3	264796	457336	11.09%	2.288%
17	21.284	3094	3097	3103	rVB	43157	66695	1.62%	0.334%
18	21.549	3135	3142	3154	rBV2	776877	1318944	32.00%	6.598%
19	21.966	3206	3213	3219	rBV3	42063	91770	2.23%	0.459%
20	22.342	3272	3277	3287	rBV3	49915	110016	2.67%	0.550%
21	22.935	3373	3378	3385	rVV3	38235	82534	2.00%	0.413%
22	23.012	3385	3391	3398	rVB2	63212	123986	3.01%	0.620%
23	23.781	3517	3522	3530	rVB2	61796	120583	2.93%	0.603%
24	24.651	3659	3670	3689	rBV2	461854	1423591	34.53%	7.122%
25	25.773	3854	3861	3869	rVB5	34325	91350	2.22%	0.457%

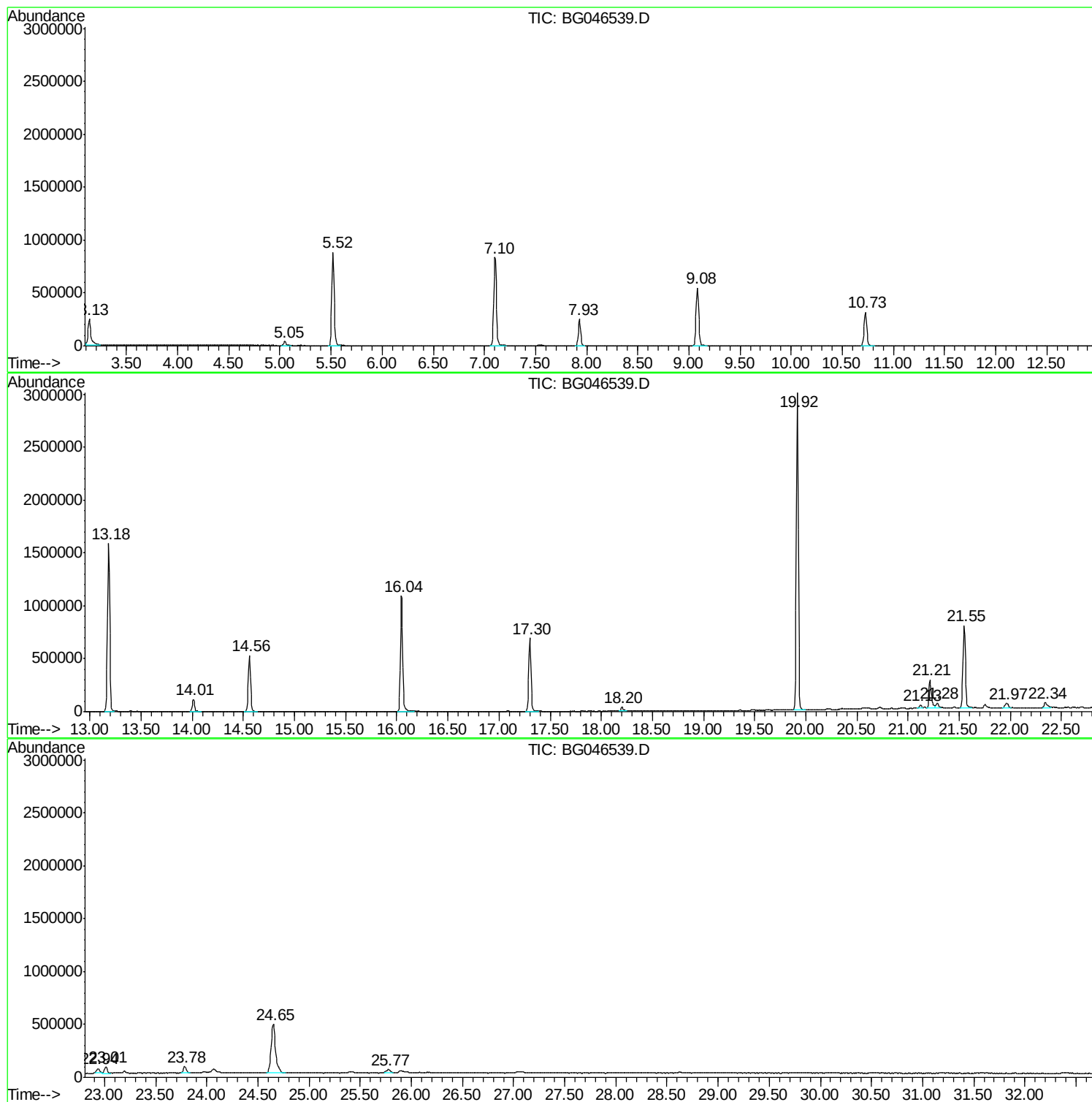
Sum of corrected areas: 19989523

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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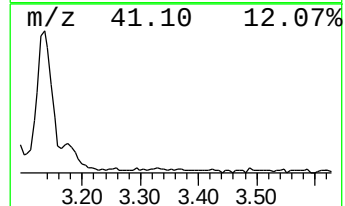
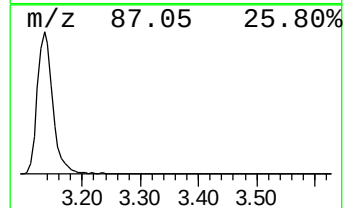
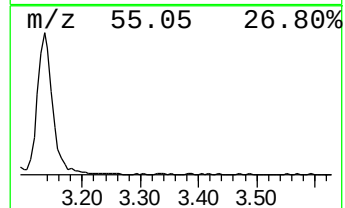
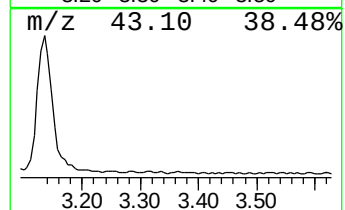
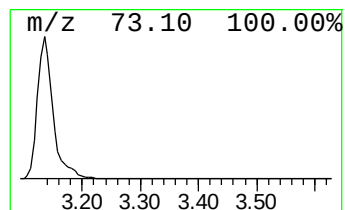
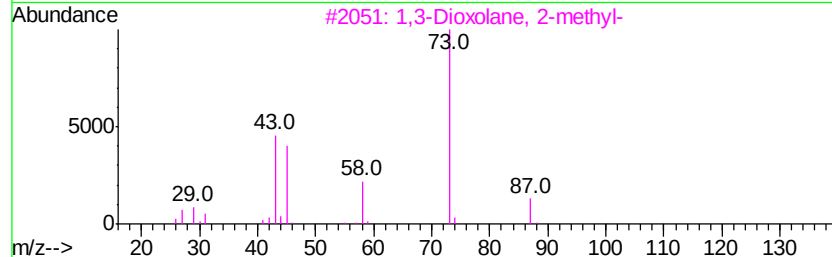
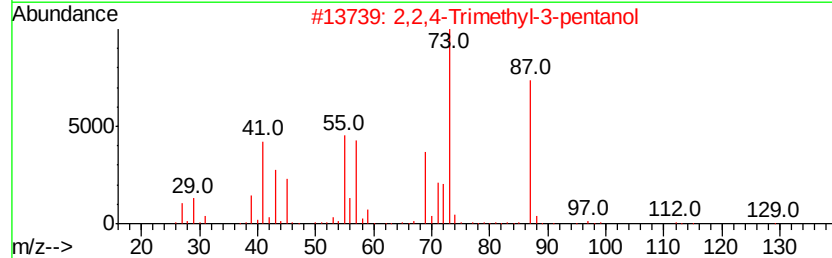
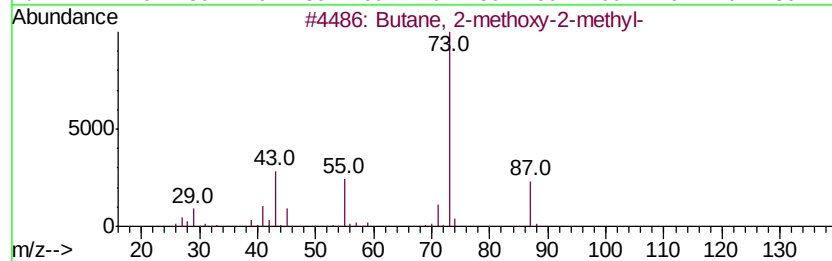
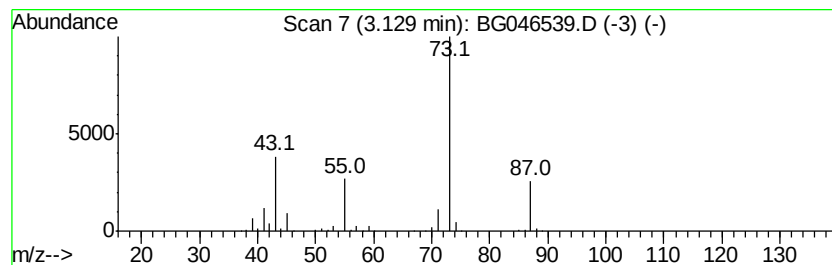
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.13	22.99 ng	472082	1,4-Dichlorobenzene-d4	7.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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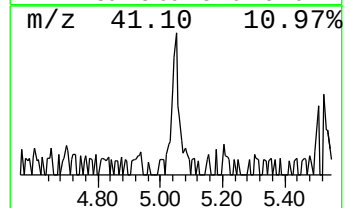
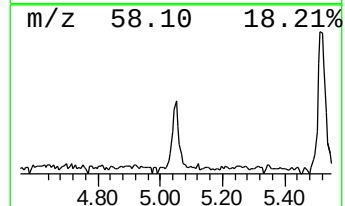
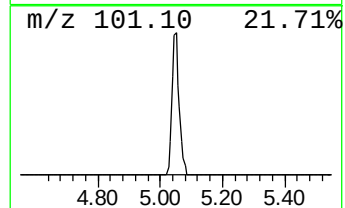
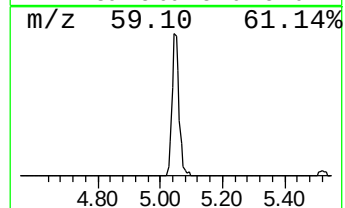
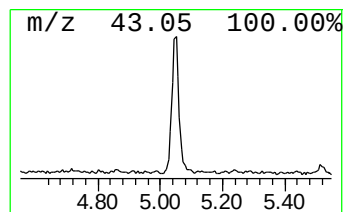
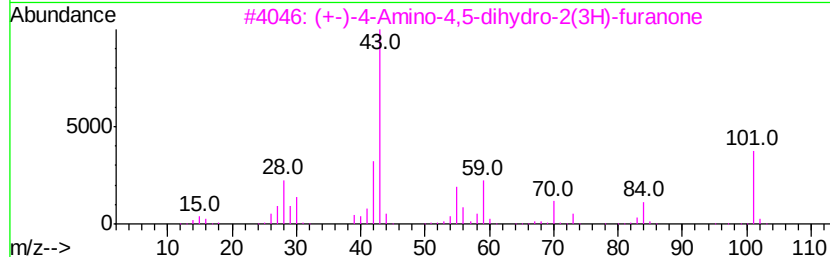
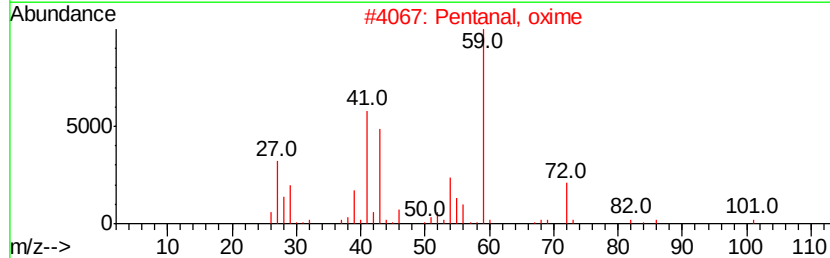
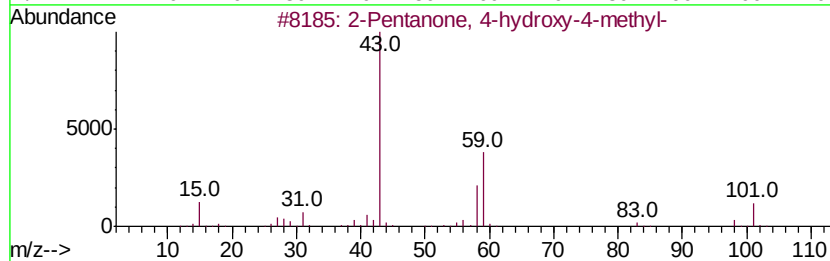
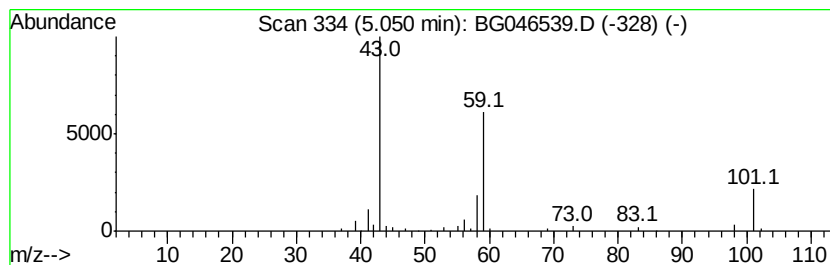
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	2.85 ng	58555	1,4-Dichlorobenzene-d4	7.93

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2		Pentanal, oxime	101	C5H11NO	000628-79-5	23
3		(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Silane, ethyldimethyl-	88	C4H12Si	000758-21-4	9



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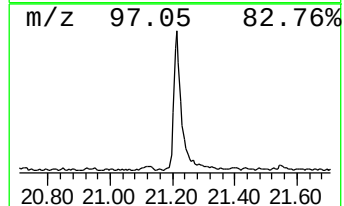
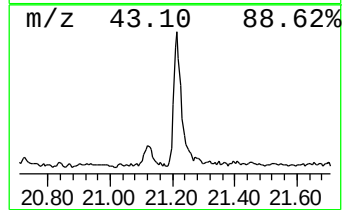
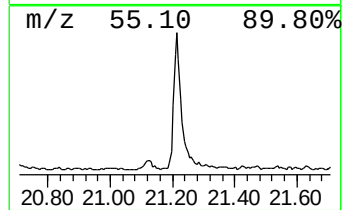
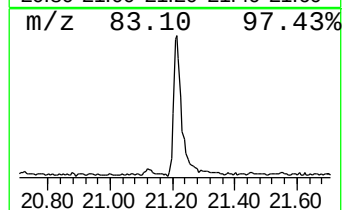
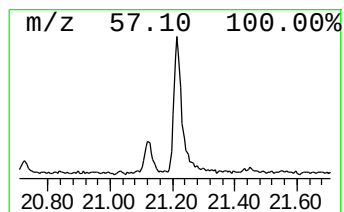
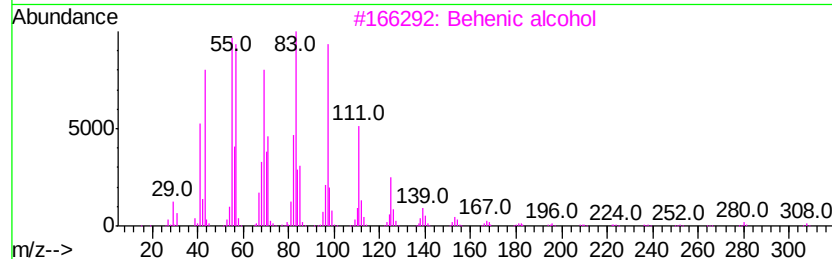
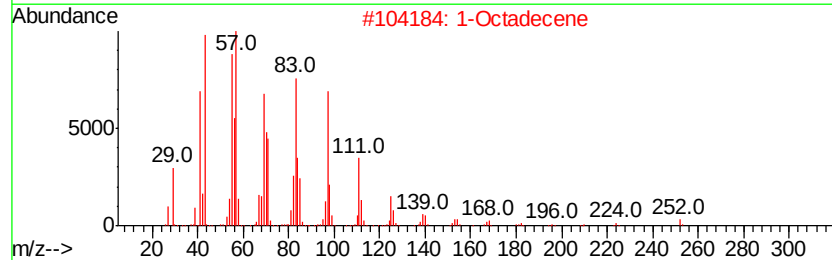
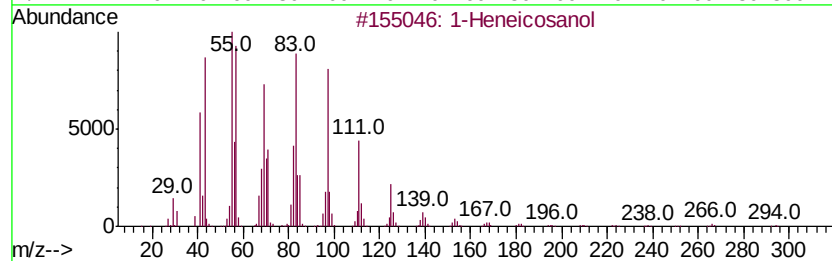
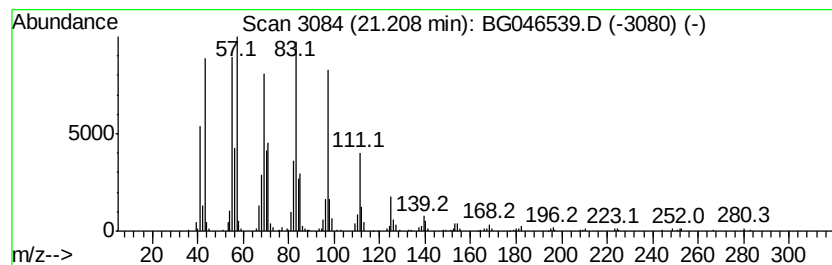
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Peak Number 3 1-Heneicosanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.
21.21	6.93 ng	457336	Chrysene-d12		21.55
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	95
2	1-Octadecene	252	C18H36	000112-88-9	95
3	Behenic alcohol	326	C22H46O	000661-19-8	95
4	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
5	n-Tetracosanol-1	354	C24H50O	000506-51-4	94



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.13	23.0	ng	472082	1	7.93	410744	20.0
2-Pentanone, 4-hy...	5.05	2.9	ng	58555	1	7.93	410744	20.0
1-Heneicosanol	21.21	6.9	ng	457336	5	21.55	1318940	20.0