

Data Path : Z:\HPCHEM1\BNA_F\Data\BF071217\
 Data File : BF096693.D
 Acq On : 12 Jul 2017 15:55
 Operator : SJ/JU
 Sample : I4126-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PP-NW-1

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-BF070117.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	4.863	467	472	485	rVB	347238	547731	15.66%	1.950%
2	5.287	538	544	547	rBV	3223538	3497861	100.00%	12.451%
3	6.316	713	719	722	rBV	2945327	3384499	96.76%	12.048%
4	6.440	735	740	743	rBV	3130257	3283004	93.86%	11.687%
5	6.669	775	779	782	rBV	782807	729578	20.86%	2.597%
6	6.828	801	806	809	rBV	2369505	2398657	68.57%	8.539%
7	7.234	869	875	877	rBV	2075511	2442363	69.82%	8.694%
8	7.945	992	996	1000	rBV	1117048	1114323	31.86%	3.967%
9	9.028	1174	1180	1183	rBV	3159756	3170903	90.65%	11.288%
10	9.422	1243	1247	1250	rBV	350710	313122	8.95%	1.115%
11	9.698	1289	1294	1298	rBV	1153974	1009876	28.87%	3.595%
12	10.145	1367	1370	1373	rVB	56479	43726	1.25%	0.156%
13	10.486	1423	1428	1431	rBV	1891271	1832970	52.40%	6.525%
14	10.616	1447	1450	1457	rBV	79355	88930	2.54%	0.317%
15	11.175	1541	1545	1549	rBV	1033469	937762	26.81%	3.338%
16	12.763	1810	1815	1818	rBV	2097209	2061042	58.92%	7.337%
17	13.669	1965	1969	1979	rVB2	153701	184421	5.27%	0.656%
18	13.804	1988	1992	1996	rBV	615901	559090	15.98%	1.990%
19	15.198	2225	2229	2235	rVB	419192	492159	14.07%	1.752%

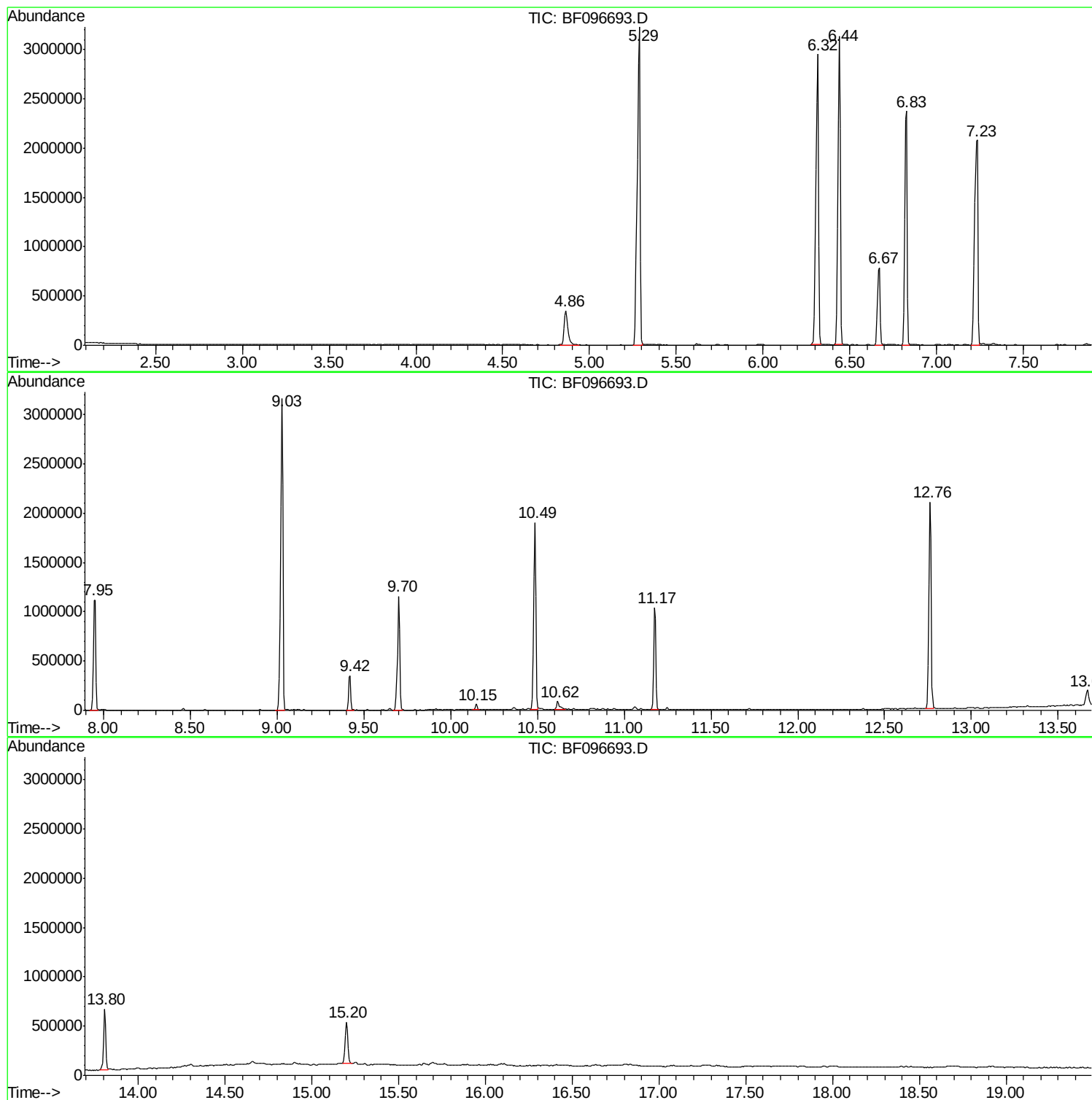
Sum of corrected areas: 28092017

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.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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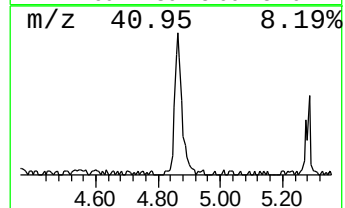
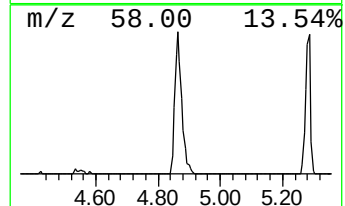
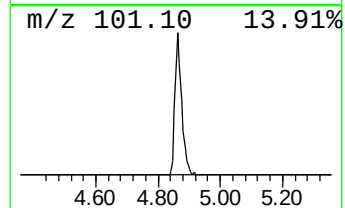
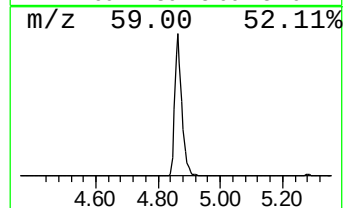
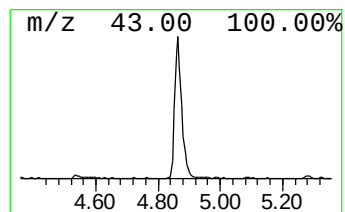
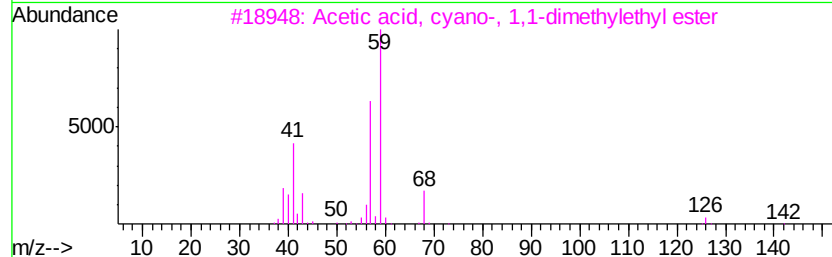
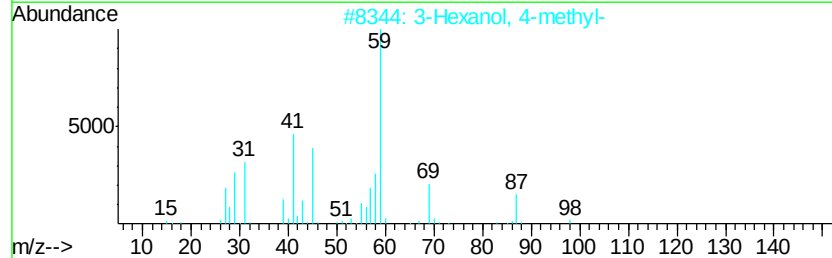
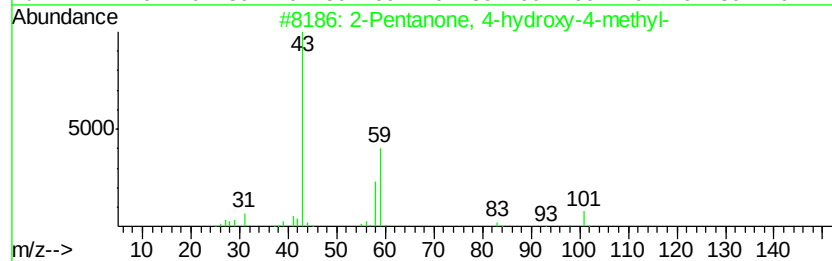
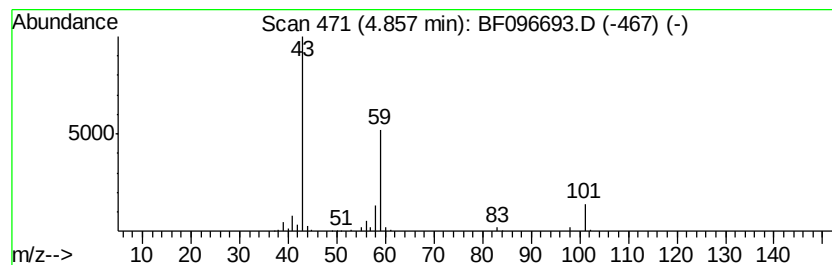
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.86	15.02 ng	547731	1,4-Dichlorobenzene-d4	6.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
5			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	9



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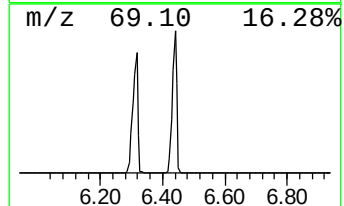
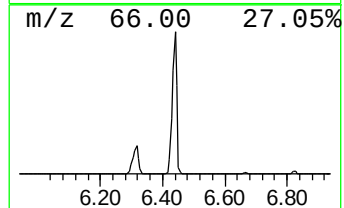
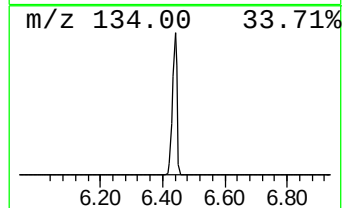
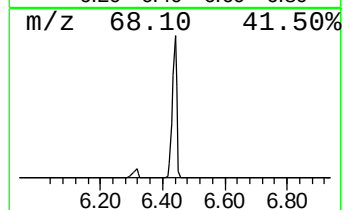
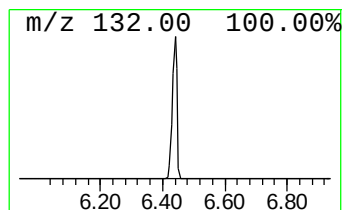
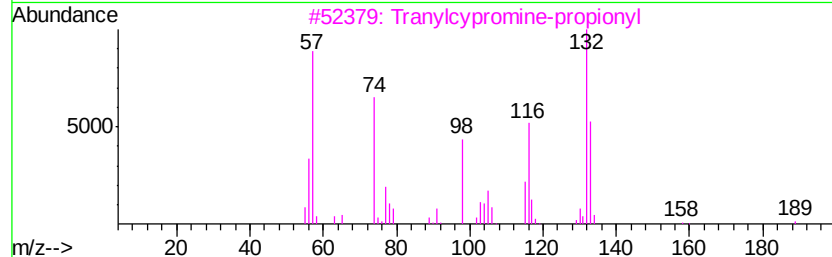
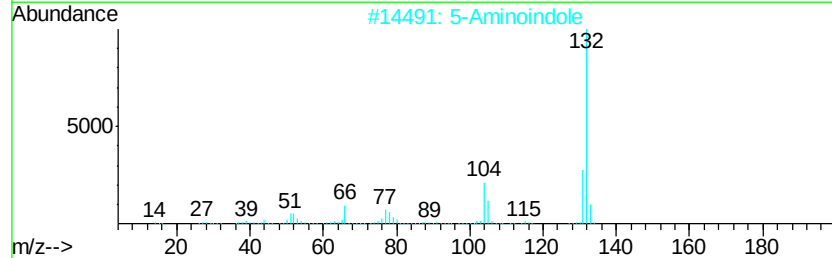
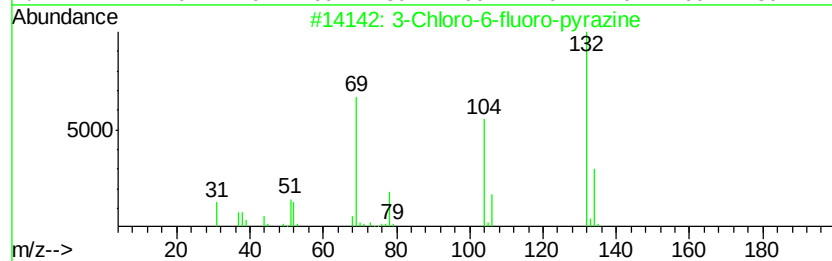
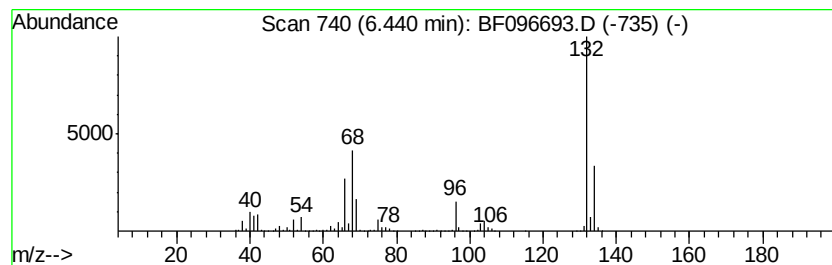
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Peak Number 2 unknown6.44 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.44	90.00 ng	3283000	1,4-Dichlorobenzene-d4	6.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4			1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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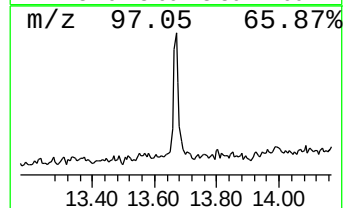
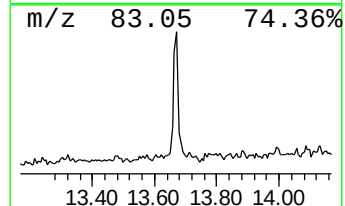
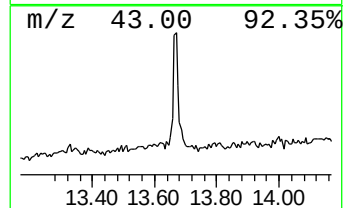
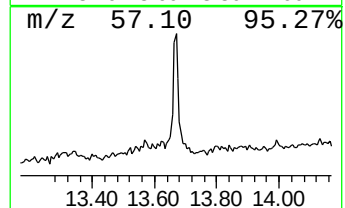
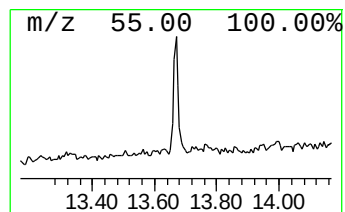
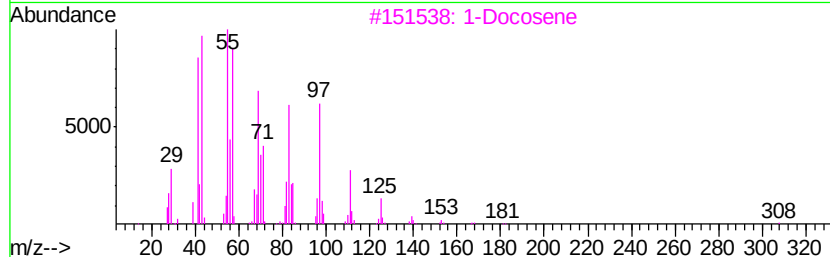
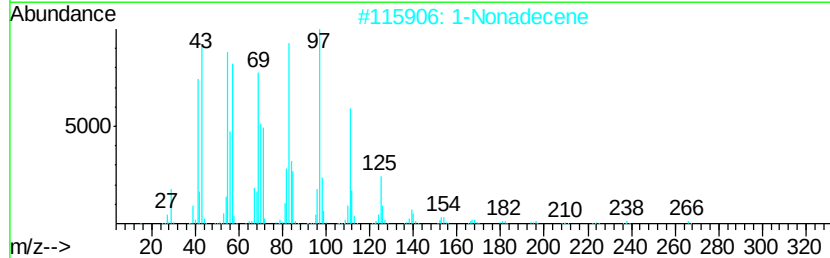
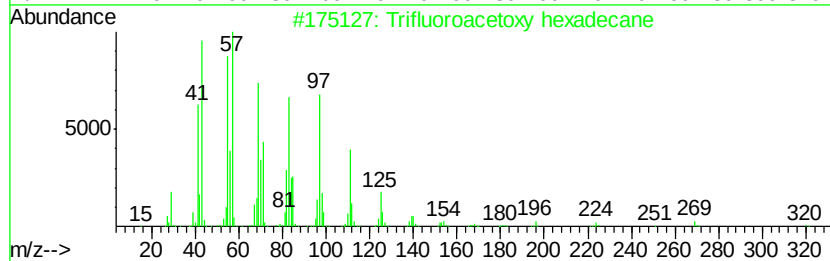
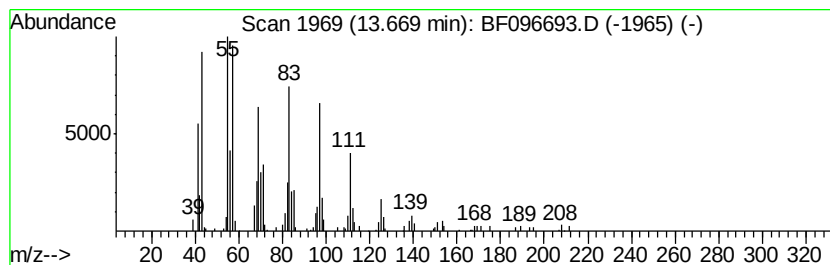
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Peak Number 4 Trifluoroacetoxy hexadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.67	6.60 ng	184421	Chrysene-d12	13.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	95
2		1-Nonadecene	266	C19H38	018435-45-5	94
3		1-Docosene	308	C22H44	001599-67-3	93
4		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	93
5		1-Heneicosanol	312	C21H44O	015594-90-8	91



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.86	15.0	ng	547731	1	6.67	729578	20.0
unknown6.44	6.44	90.0	ng	3283000	1	6.67	729578	20.0
Trifluoroacetoxy ...	13.67	6.6	ng	184421	5	13.80	559090	20.0