

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081518\
 Data File : BF108189.D
 Acq On : 16 Aug 2018 2:08
 Operator : JU/SJ
 Sample : J4440-04
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 080918-TIPC-F-U-S

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_F\METHODS\8270-BF080818.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	2.372	3	6	24	rVB	1649949	2196335	30.68%	5.206%
2	5.248	492	495	504	rVB	79267	84156	1.18%	0.199%
3	5.636	557	561	564	rBV	2211943	1884338	26.32%	4.467%
4	6.625	725	729	732	rBV	1576874	1256781	17.55%	2.979%
5	6.783	751	756	759	rBV	3908346	3561513	49.74%	8.442%
6	7.013	791	795	799	rBV	1546165	1308343	18.27%	3.101%
7	7.172	817	822	826	rBV	4978476	4829843	67.46%	11.449%
8	7.577	885	891	894	rBV	3743983	4308187	60.17%	10.212%
9	8.295	1009	1013	1017	rBV	1710035	1628412	22.74%	3.860%
10	9.377	1190	1197	1200	rBV	6851184	7159895	100.00%	16.972%
11	9.760	1258	1262	1267	rBV	480836	388443	5.43%	0.921%
12	10.060	1309	1313	1317	rVB2	1929992	1724905	24.09%	4.089%
13	10.465	1380	1382	1386	rVB2	82684	73183	1.02%	0.173%
14	10.718	1422	1425	1428	rBV	220422	166100	2.32%	0.394%
15	10.854	1443	1448	1457	rBV	2422536	2363582	33.01%	5.603%
16	10.942	1460	1463	1470	rVB2	82177	96512	1.35%	0.229%
17	11.554	1563	1567	1571	rBV	1690150	1500979	20.96%	3.558%
18	13.142	1832	1837	1841	rBV	4676299	4518794	63.11%	10.712%
19	14.042	1985	1990	2005	rBV2	107030	280879	3.92%	0.666%
20	14.206	2014	2018	2022	rBV	1363372	1191100	16.64%	2.823%
21	15.342	2206	2211	2216	rBV	81830	101972	1.42%	0.242%
22	15.771	2277	2284	2290	rBV	876937	1266794	17.69%	3.003%
23	16.236	2358	2363	2368	rVB2	77573	120890	1.69%	0.287%
24	16.794	2453	2458	2463	rVB	59712	101318	1.42%	0.240%
25	17.447	2564	2569	2577	rVB3	35275	72746	1.02%	0.172%

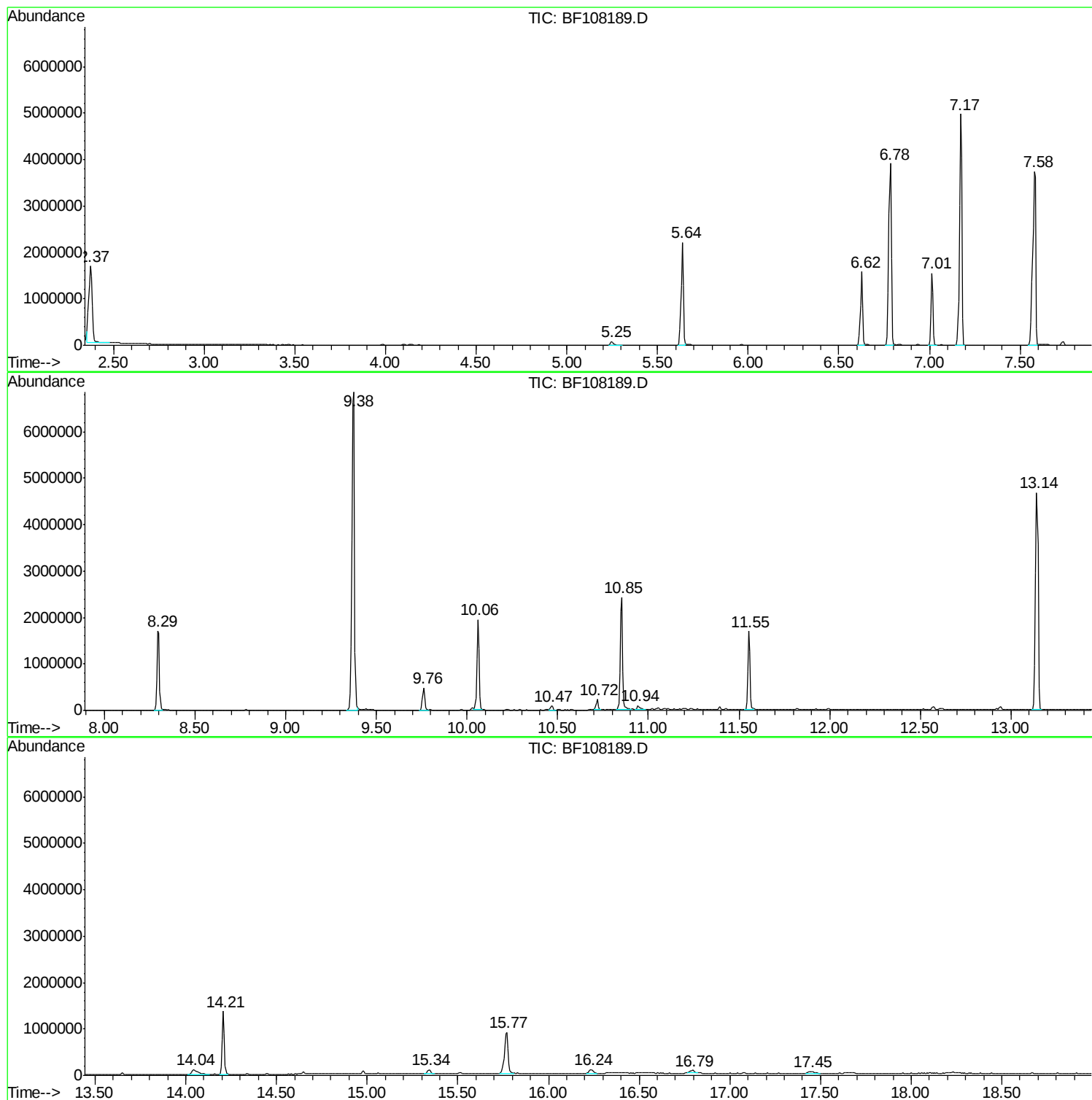
Sum of corrected areas: 42186000

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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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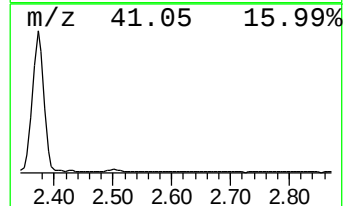
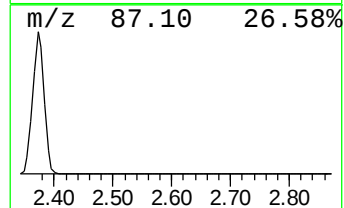
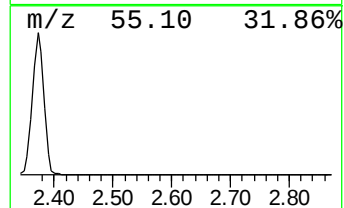
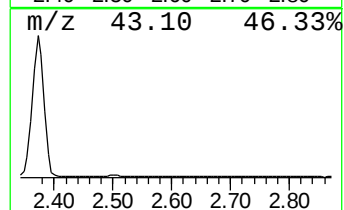
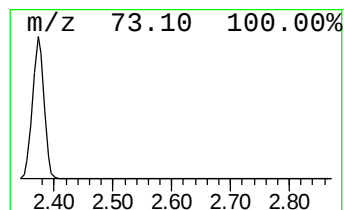
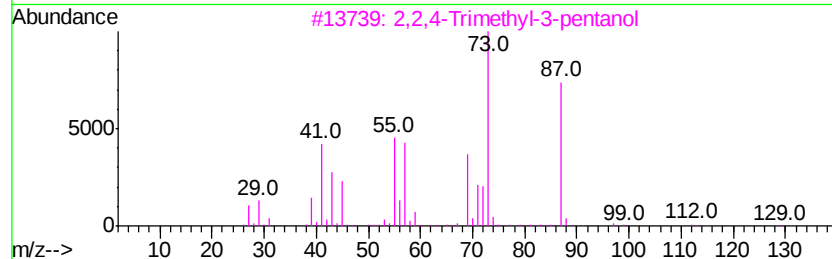
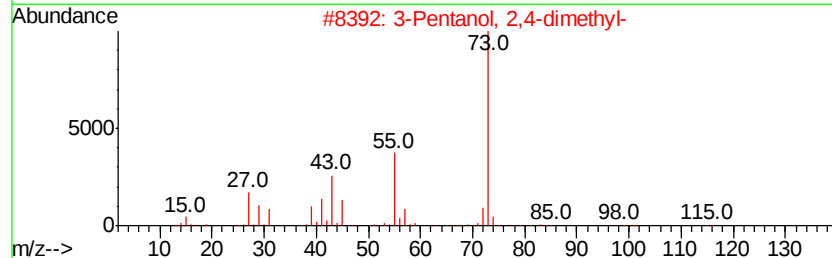
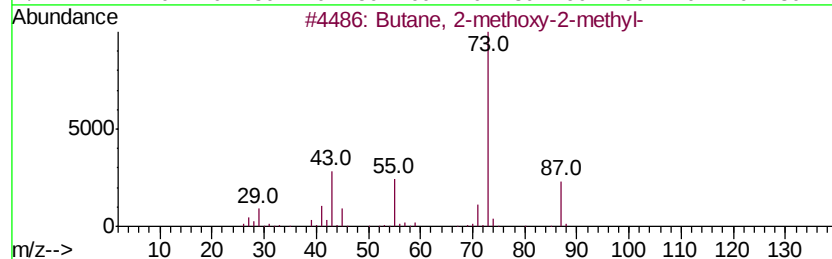
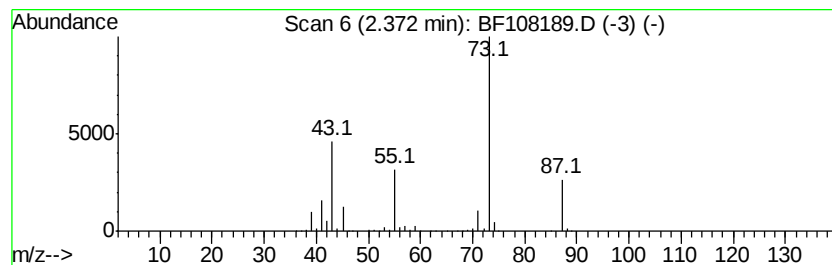
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.37	33.57 ng	2196340	1,4-Dichlorobenzene-d4	7.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	43
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23



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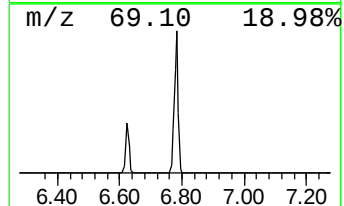
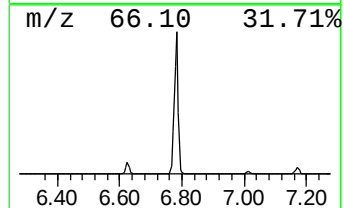
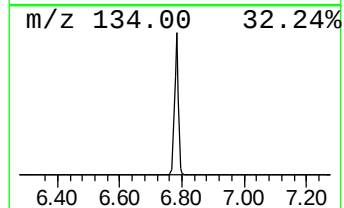
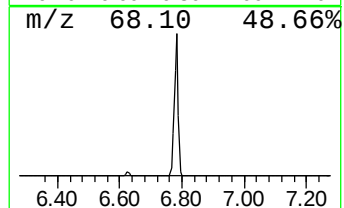
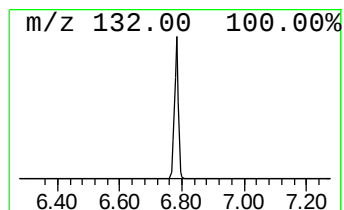
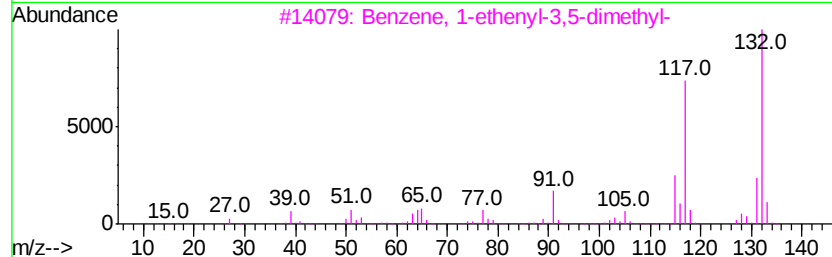
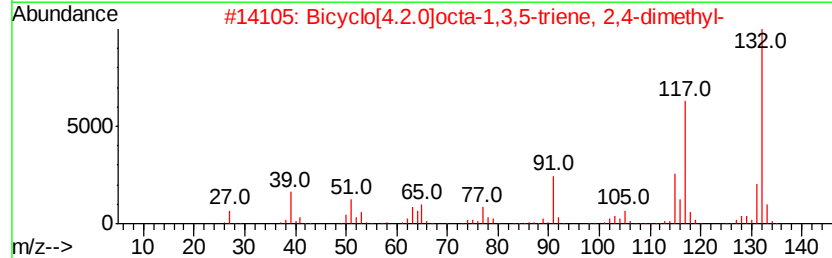
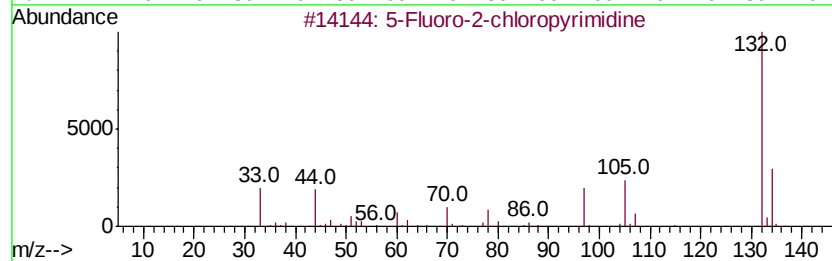
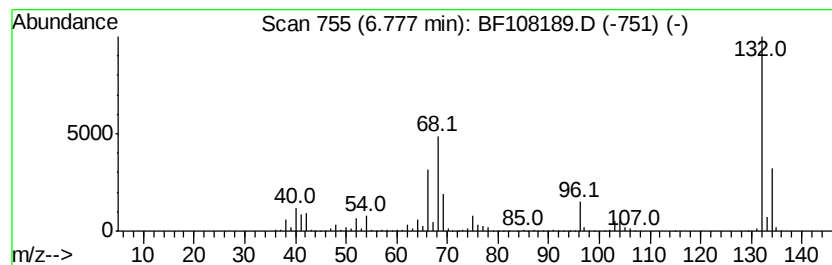
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Peak Number 2 unknown6.78 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.78	54.44 ng	3561510	1,4-Dichlorobenzene-d4	7.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
3		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	9



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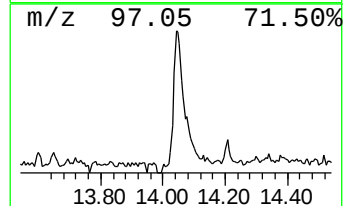
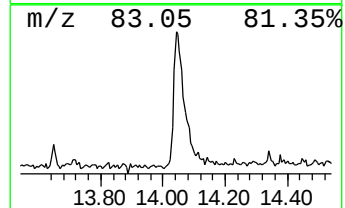
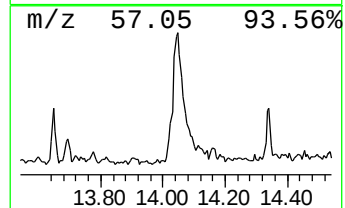
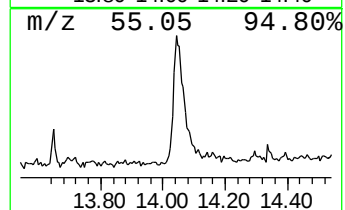
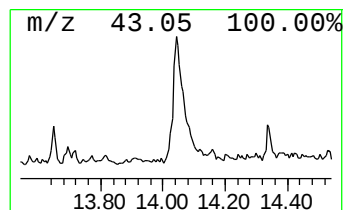
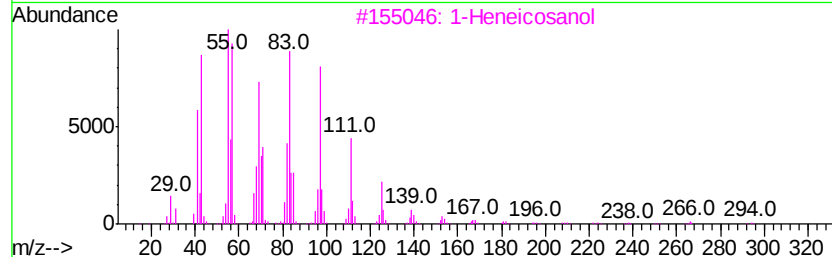
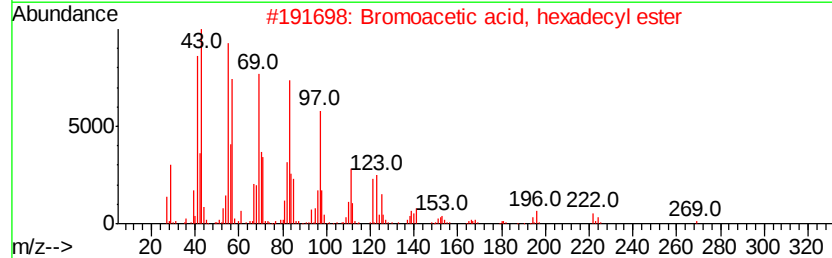
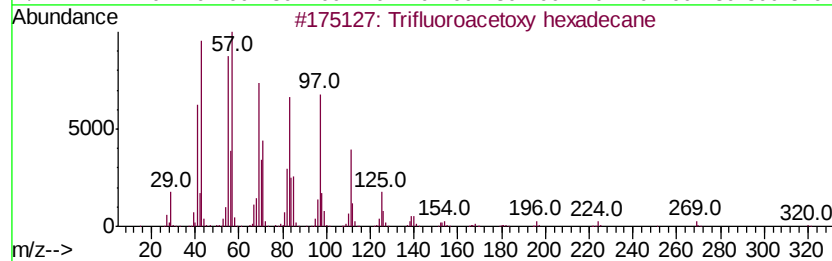
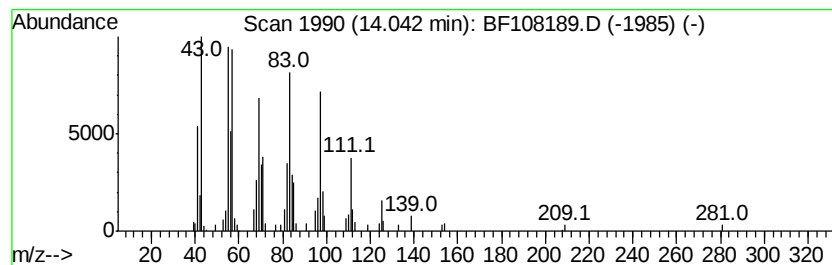
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Trifluoroacetoxy hexadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.04	4.72 ng	280879	Chrysene-d12	14.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
2		Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	94
3		1-Heneicosanol	312	C21H44O	015594-90-8	93
4		n-Nonadecanol-1	284	C19H40O	001454-84-8	93
5		n-Tetracosanol-1	354	C24H50O	000506-51-4	91



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.37	33.6	ng	2196340	1	7.01	1308340	20.0
unknown6.78	6.78	54.4	ng	3561510	1	7.01	1308340	20.0
Trifluoroacetoxy ...	14.04	4.7	ng	280879	5	14.21	1191100	20.0