

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111117\
 Data File : BF100438.D
 Acq On : 11 Nov 2017 14:09
 Operator : SJ/JU
 Sample : I6407-01 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 3091-4235

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-BF110817.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	5.122	512	516	527	rBV	404885	443090	18.16%	1.924%
2	5.516	579	583	587	rBV	2221055	2222475	91.09%	9.652%
3	6.528	749	755	762	rBV	2372804	2394798	98.16%	10.400%
4	6.669	775	779	783	rBV	2288031	2286121	93.70%	9.928%
5	6.904	815	819	822	rBV	1220280	954512	39.12%	4.145%
6	7.063	841	846	849	rVB	1859643	1753680	71.88%	7.616%
7	7.463	909	914	917	rBV	1616874	1373857	56.31%	5.967%
8	8.186	1033	1037	1040	rBV	1525734	1203228	49.32%	5.226%
9	9.257	1215	1219	1223	rBV	2606320	2439739	100.00%	10.596%
10	9.651	1282	1286	1294	rBV	126488	119023	4.88%	0.517%
11	9.804	1308	1312	1317	rBV	26865	25902	1.06%	0.112%
12	9.939	1331	1335	1346	rBV2	1243838	1171067	48.00%	5.086%
13	10.727	1465	1469	1479	rBV	1340695	1175986	48.20%	5.107%
14	11.427	1584	1588	1591	rBV	1215078	1028734	42.17%	4.468%
15	11.451	1591	1592	1596	rVV	70689	38574	1.58%	0.168%
16	11.504	1598	1601	1604	rVB	51177	44687	1.83%	0.194%
17	11.939	1671	1675	1683	rBV3	40090	56292	2.31%	0.244%
18	12.639	1791	1794	1798	rVB	138859	115938	4.75%	0.504%
19	12.868	1831	1833	1838	rVB	136704	113578	4.66%	0.493%
20	13.010	1853	1857	1861	rBV	2079102	1836163	75.26%	7.974%
21	13.898	2005	2008	2013	rBV2	79591	96106	3.94%	0.417%
22	14.068	2032	2037	2040	rBV	1009257	957015	39.23%	4.156%
23	15.109	2212	2214	2224	rVB2	77470	149706	6.14%	0.650%
24	15.486	2275	2278	2282	rBV	61393	78311	3.21%	0.340%
25	15.551	2285	2289	2300	rVB	504684	750456	30.76%	3.259%
26	17.503	2616	2621	2629	rVB	99156	196960	8.07%	0.855%

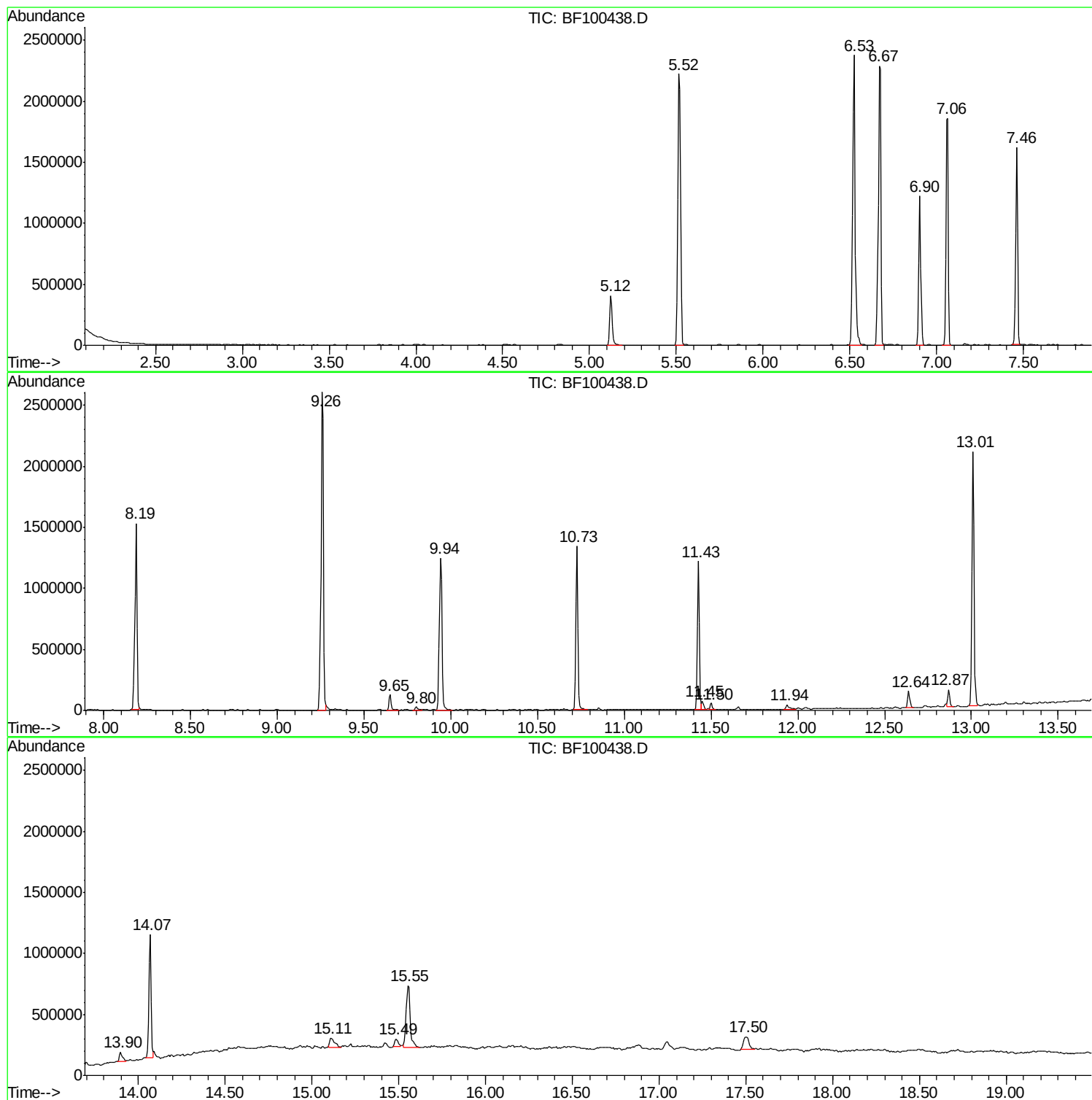
Sum of corrected areas: 23025998

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.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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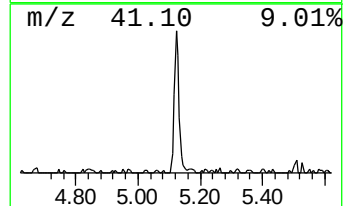
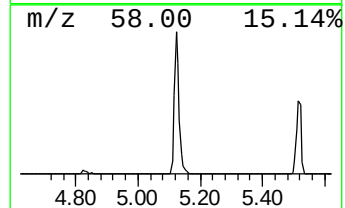
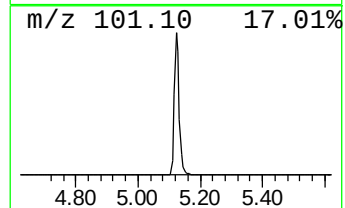
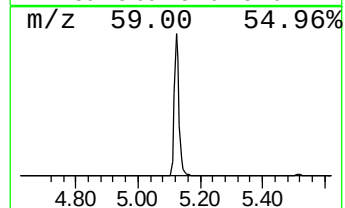
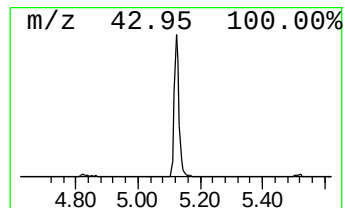
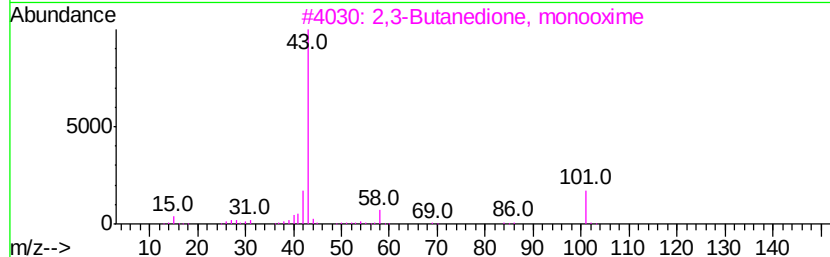
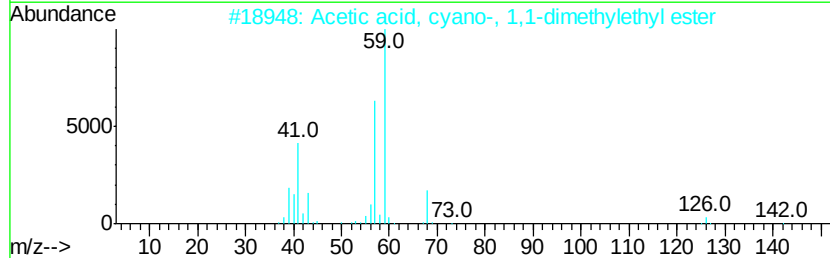
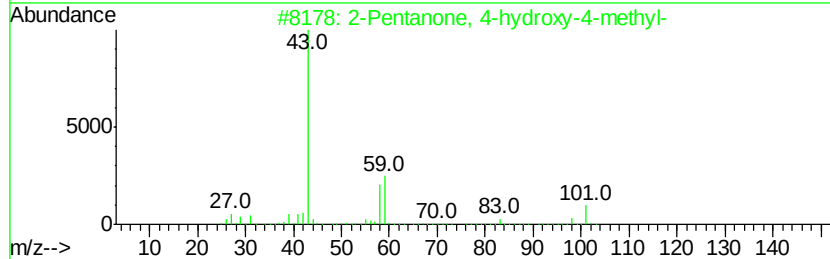
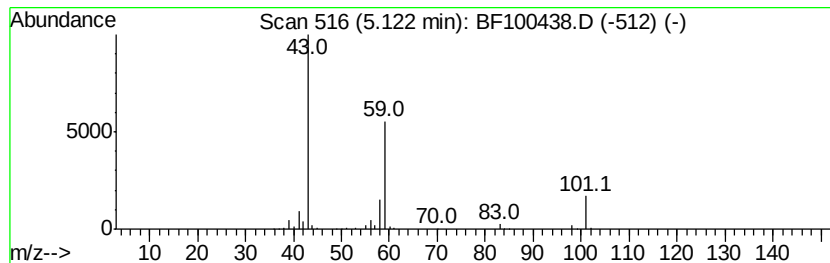
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	9.28 ng	443090	1,4-Dichlorobenzene-d4	6.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			Propanoic acid, 2-nitro-, methyl...	133	C4H7NO4	006118-50-9	9
5			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9



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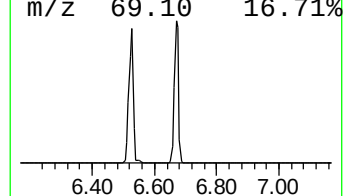
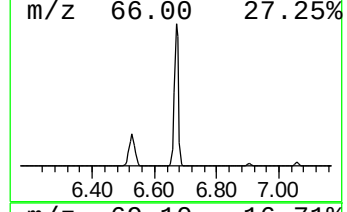
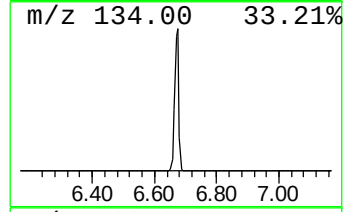
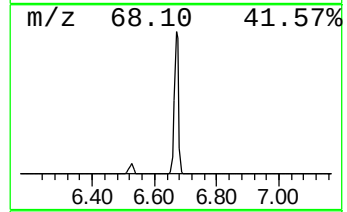
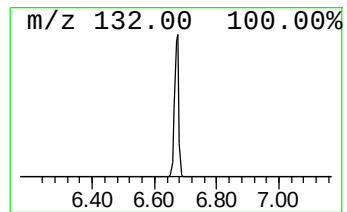
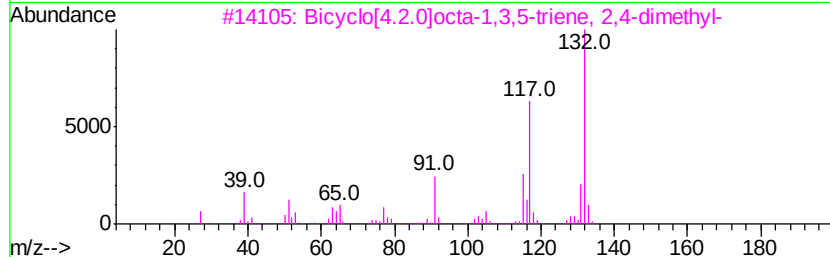
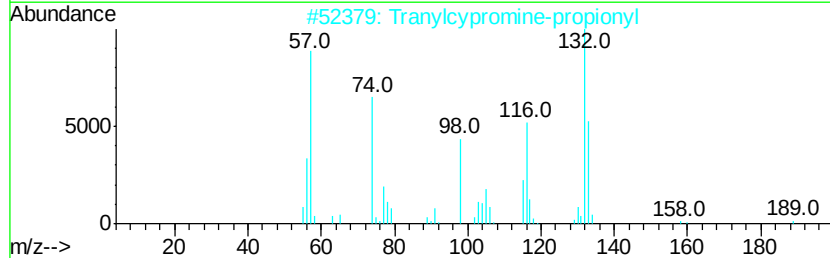
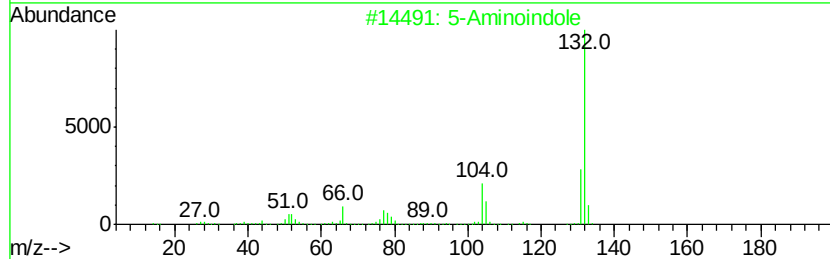
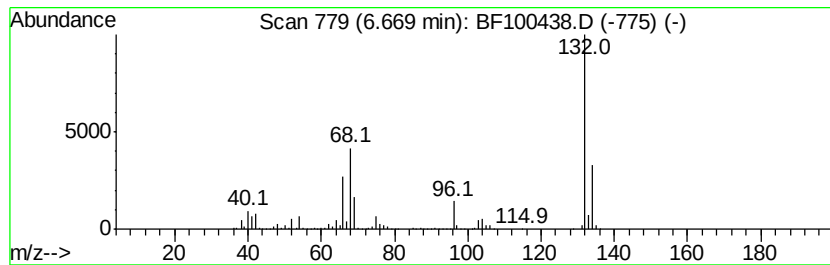
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Peak Number 2 unknown6.67 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.67	47.90 ng	2286120	1,4-Dichlorobenzene-d4	6.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Aminoindole	132	C8H8N2	005192-03-0	12
2			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4			2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5			1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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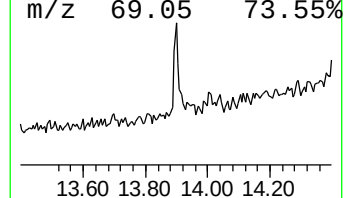
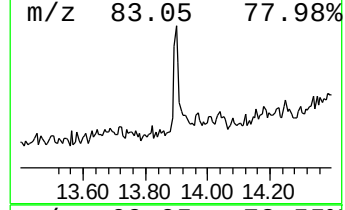
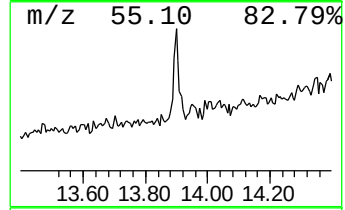
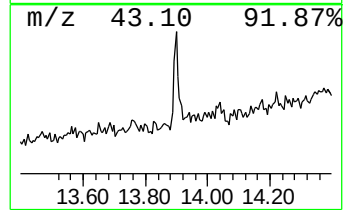
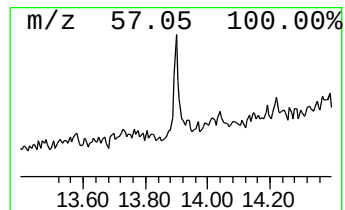
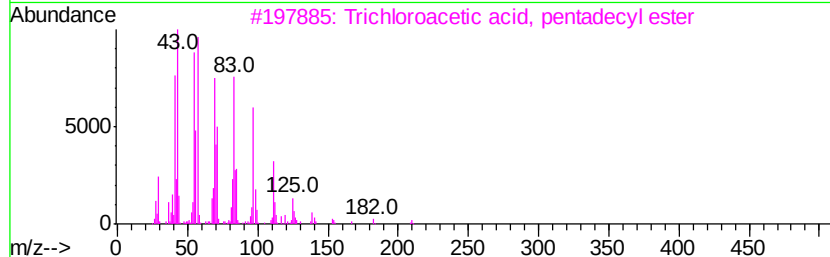
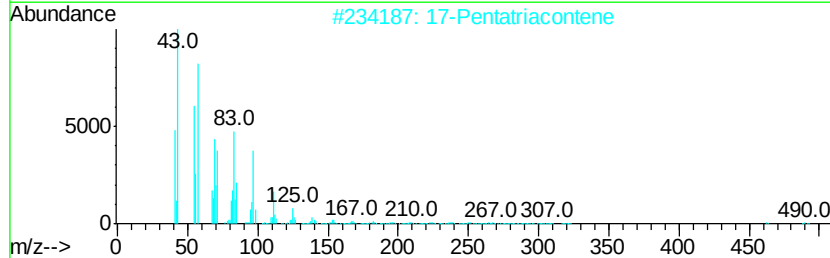
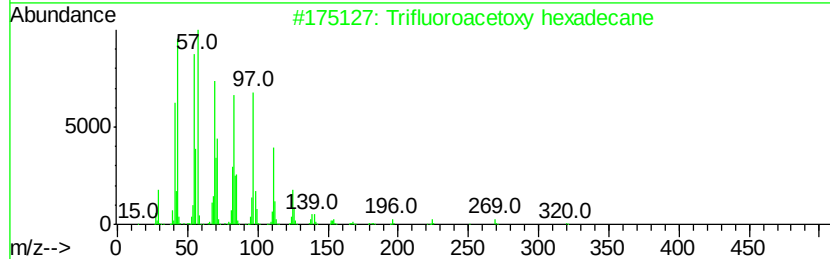
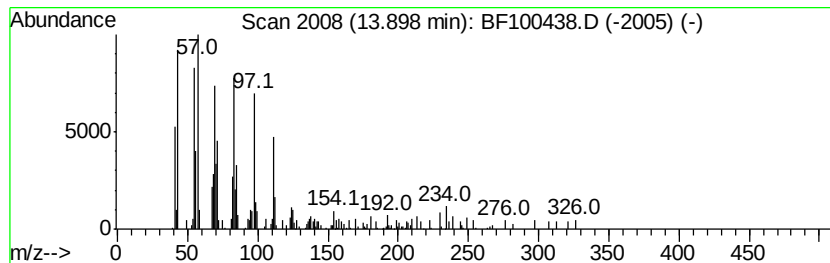
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Peak Number 4 Trifluoroacetoxy hexadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.90	2.01 ng	96106	Chrysene-d12	14.07	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	91
2	17-Pentatriacontene	491	C35H70	006971-40-0	87
3	Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	87
4	Cyclodocosane, ethyl-	336	C24H48	1000151-22-6	83
5	Octacosanol	410	C28H58O	000557-61-9	76



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.12	9.3	ng	443090	1	6.90	954512	20.0
unknown6.67	6.67	47.9	ng	2286120	1	6.90	954512	20.0
Trifluoroacetoxy ...	13.90	2.0	ng	96106	5	14.07	957015	20.0