

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092017\
 Data File : BF098659.D
 Acq On : 20 Sep 2017 14:49
 Operator : SJ/JU
 Sample : I5276-07
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SASP2P-ENV-118-3(5-10)

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF091117.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	4.792	456	460	478	rBV	279109	414642	16.27%	1.991%
2	5.228	530	534	554	rBV	2301521	2330273	91.44%	11.189%
3	6.275	708	712	728	rBV	2189931	2408507	94.51%	11.565%
4	6.392	728	732	735	rBV	2833104	2400298	94.19%	11.525%
5	6.616	767	770	781	rVB	597475	534018	20.96%	2.564%
6	6.775	793	797	810	rBV	1961052	1698848	66.67%	8.157%
7	7.192	863	868	871	rBV	1651977	1474985	57.88%	7.082%
8	7.904	985	989	1000	rBV	794237	725346	28.46%	3.483%
9	8.980	1167	1172	1175	rBV	3090577	2548319	100.00%	12.236%
10	9.380	1236	1240	1249	rBB	205383	184654	7.25%	0.887%
11	9.651	1282	1286	1290	rBV	892244	765588	30.04%	3.676%
12	10.445	1417	1421	1424	rBV	1629613	1355965	53.21%	6.511%
13	10.557	1438	1440	1448	rBV	27262	31671	1.24%	0.152%
14	11.133	1535	1538	1542	rBV	790298	706360	27.72%	3.392%
15	12.721	1803	1808	1811	rBV	2220497	1898680	74.51%	9.117%
16	12.951	1841	1847	1853	rBV	23596	32218	1.26%	0.155%
17	13.621	1956	1961	1968	rBV	80000	103439	4.06%	0.497%
18	13.774	1982	1987	1991	rBV	533451	536690	21.06%	2.577%
19	14.515	2111	2113	2121	rBV2	81853	105346	4.13%	0.506%
20	15.168	2219	2224	2230	rBV	474185	536596	21.06%	2.577%
21	16.009	2362	2367	2374	rBV4	18667	34002	1.33%	0.163%

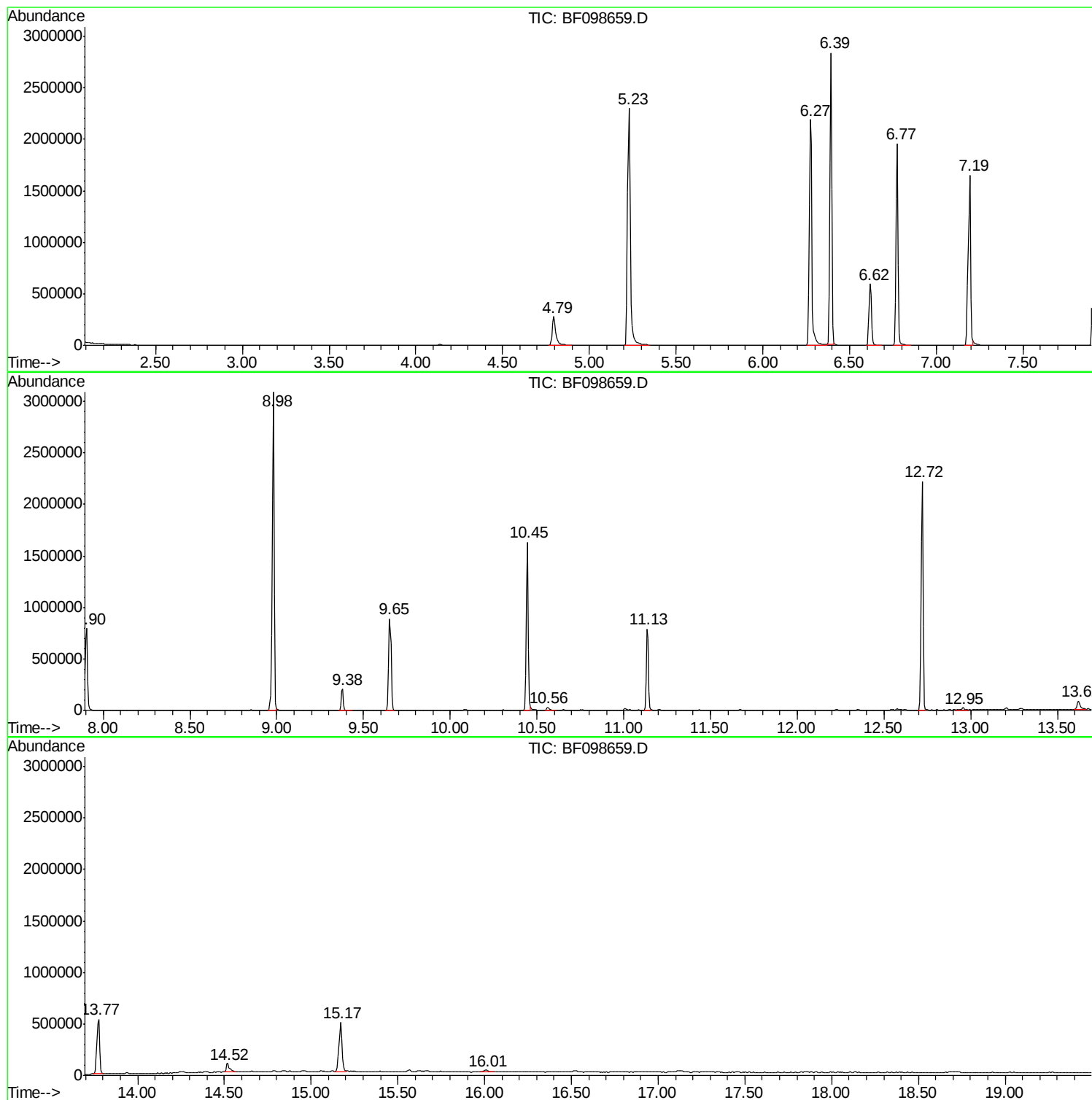
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF091117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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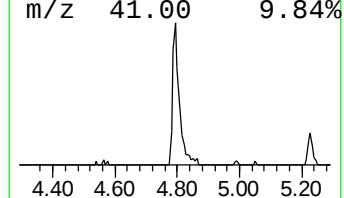
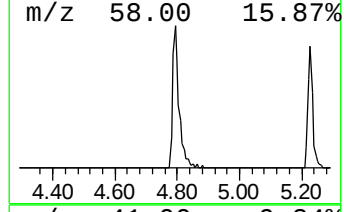
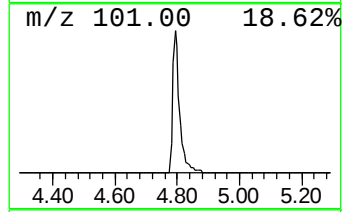
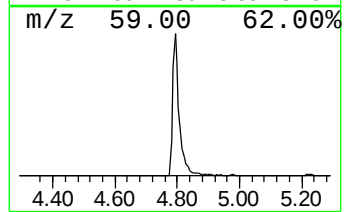
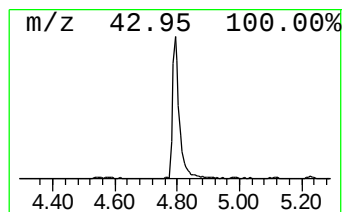
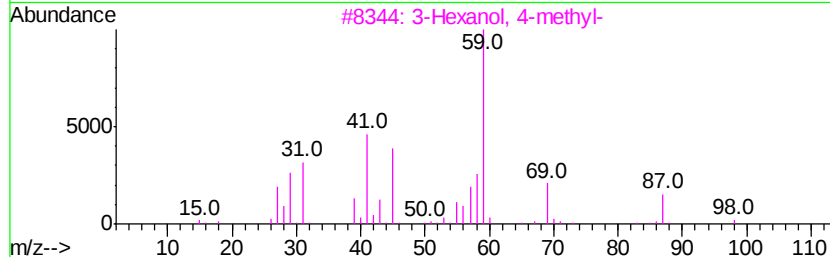
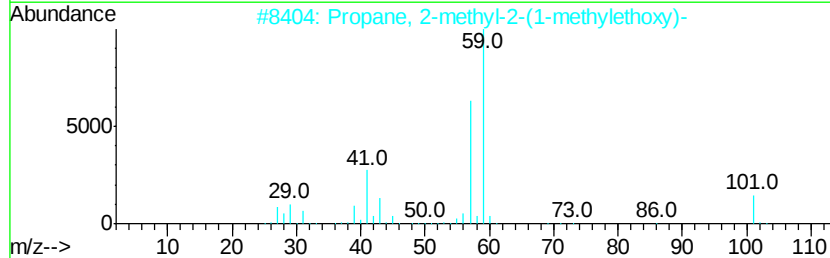
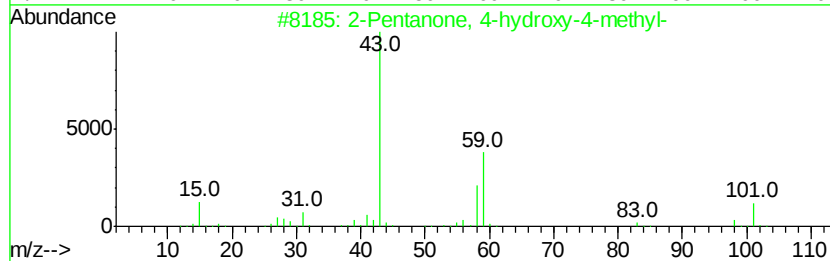
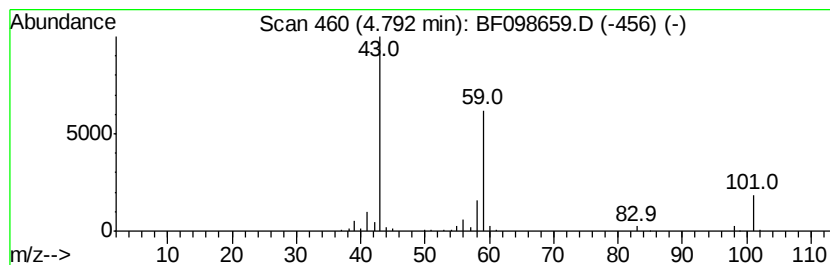
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.79	15.53 ng	414642	1,4-Dichlorobenzene-d4	6.62

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	42
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25



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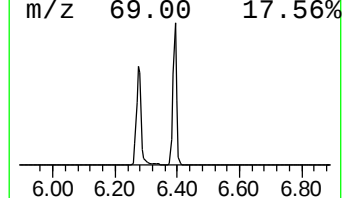
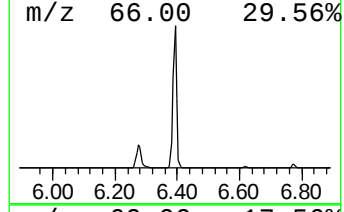
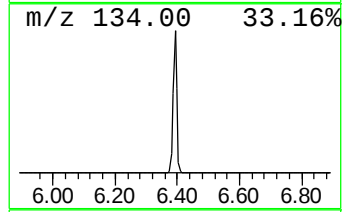
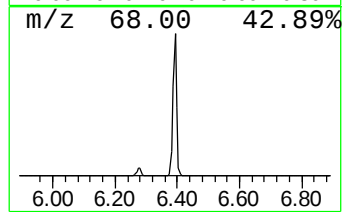
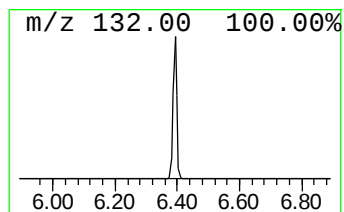
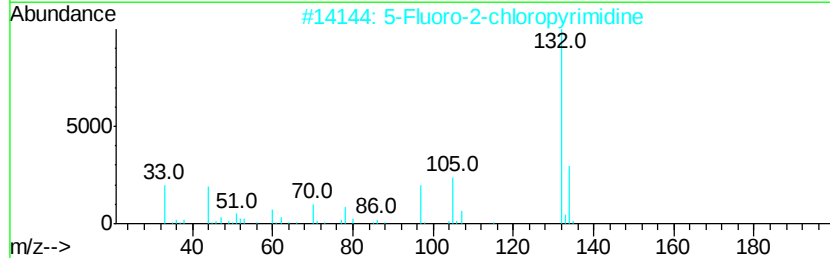
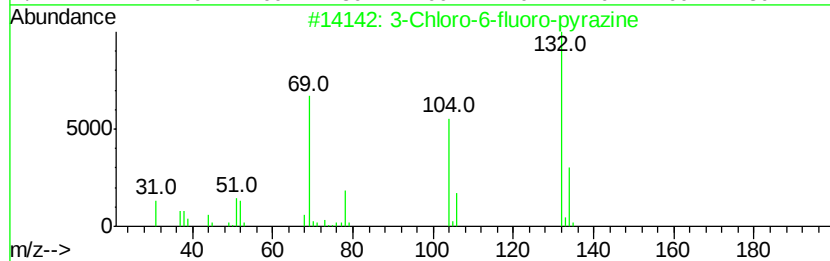
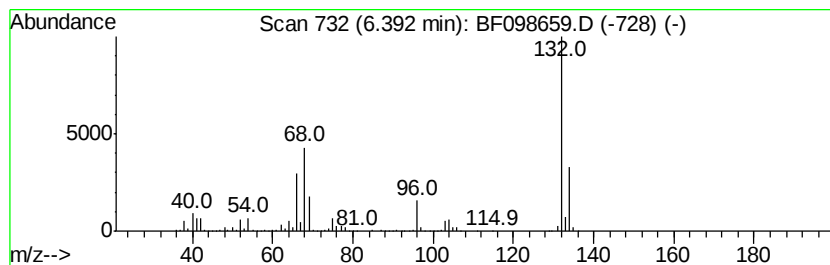
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Peak Number 2 unknown6.39 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.39	89.90 ng	2400300	1,4-Dichlorobenzene-d4	6.62

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3	Tranvlcv		promine-propionyl	189	C12H15NO	1000123-86-3	10
4	4		Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5	4		Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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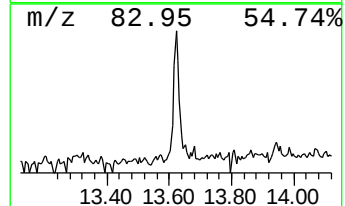
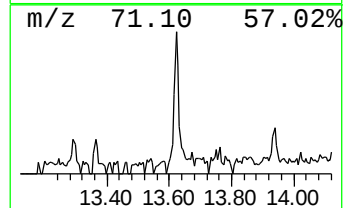
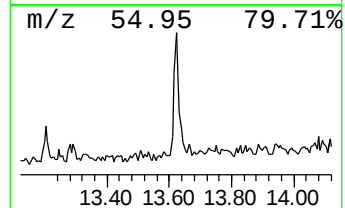
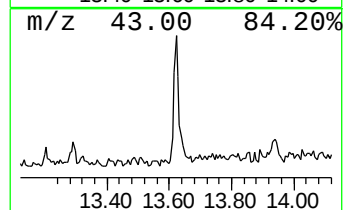
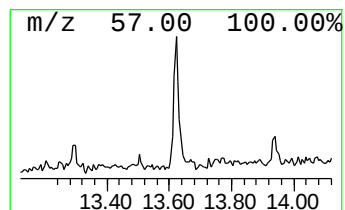
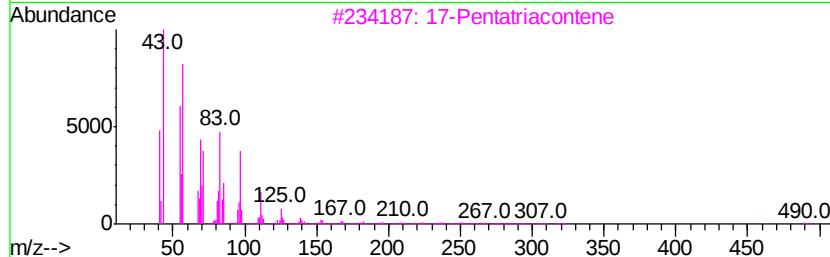
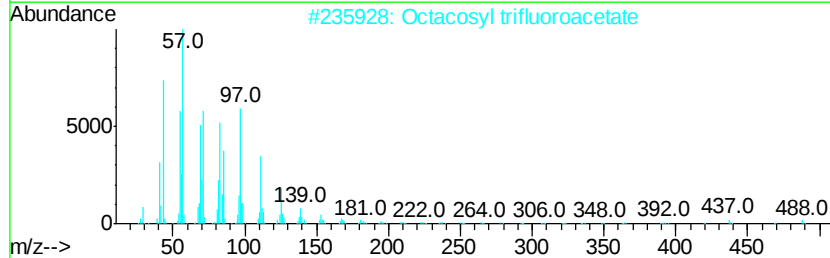
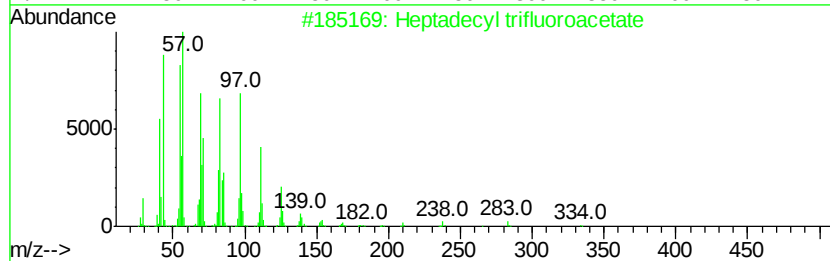
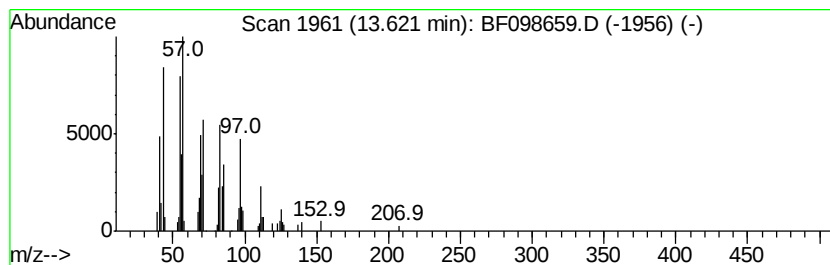
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Peak Number 4 Heptadecyl trifluoroacetate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.62	3.85 ng	103439	Chrysene-d12	13.77

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	87
2		Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	87
3		17-Pentatriacontene	491	C35H70	006971-40-0	87
4		Heneicosyl heptafluorobutyrte	508	C25H43F7O2	1000351-83-8	87
5		5-Eicosene, (E)-	280	C20H40	074685-30-6	87



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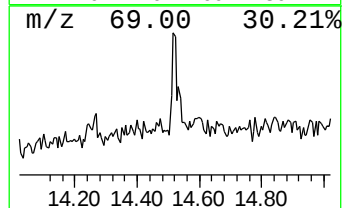
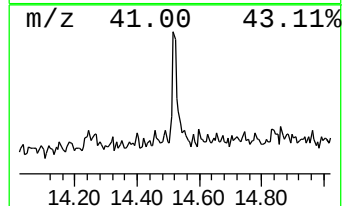
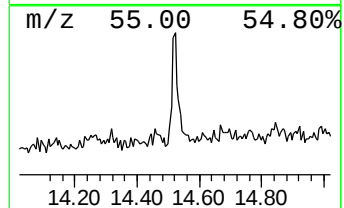
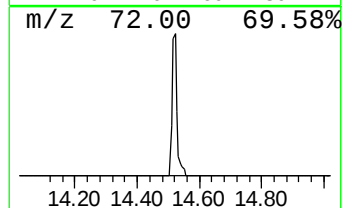
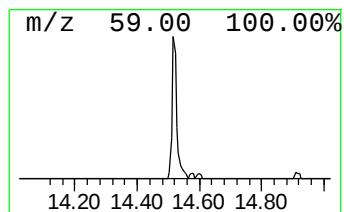
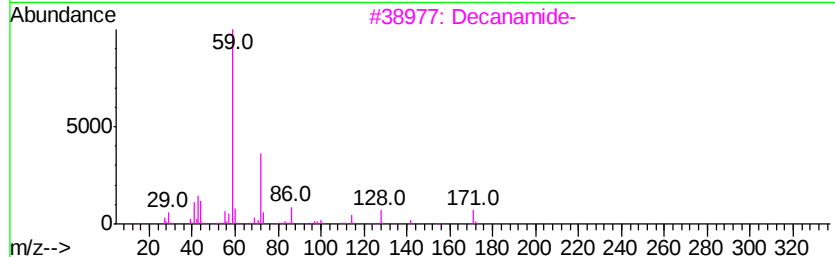
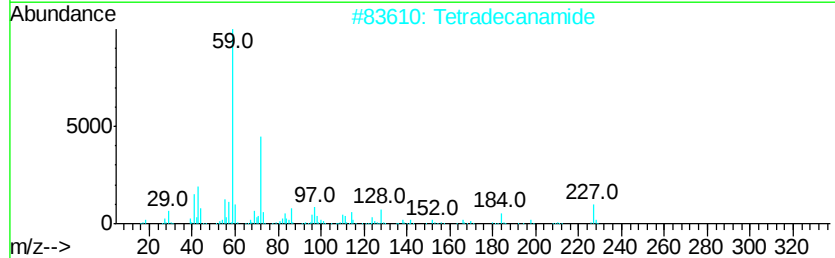
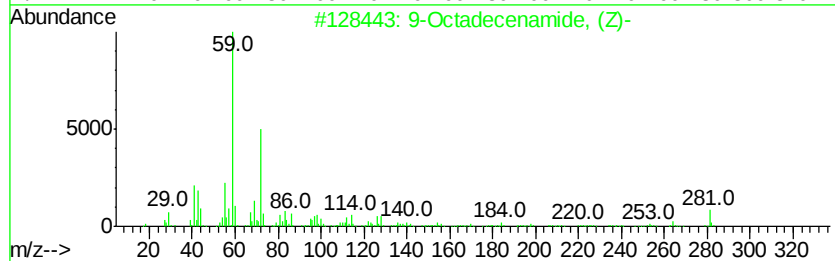
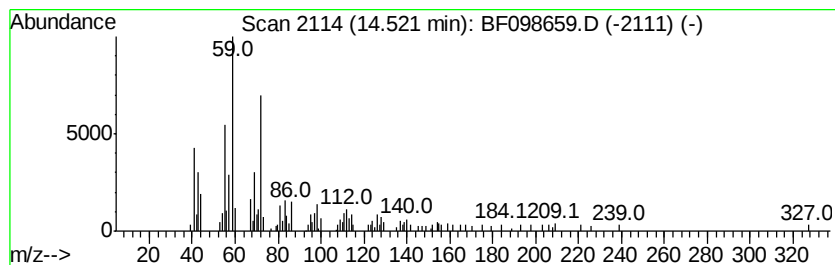
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Peak Number 5 9-Octadecenamide, (Z)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.52	3.93 ng	105346	Perylene-d12	15.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	86
2		Tetradecanamide	227	C14H29NO	000638-58-4	64
3		Decanamide-	171	C10H21NO	002319-29-1	64
4		7-Nonenamide	155	C9H17NO	090949-53-4	64
5		Octadecanamide	283	C18H37NO	000124-26-5	59



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.79	15.5	ng	414642	1	6.62	534018	20.0
unknown6.39	6.39	89.9	ng	2400300	1	6.62	534018	20.0
Heptadecyl triflu...	13.62	3.9	ng	103439	5	13.77	536690	20.0
9-Octadecenamide,...	14.52	3.9	ng	105346	6	15.17	536596	20.0