

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052225\
 Data File : BF142506.D
 Acq On : 22 May 2025 13:11
 Operator : RC/JU
 Sample : Q2074-08
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-18

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:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052025.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF142506.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.116	486	497	505	rBV	137232	207101	9.41%	1.270%
2	5.510	558	564	569	rBV	1389343	1835699	83.45%	11.254%
3	6.516	728	735	749	rBV	1416255	1951034	88.69%	11.962%
4	6.898	794	800	805	rBV	508449	645890	29.36%	3.960%
5	7.457	888	895	900	rBV	860646	1133311	51.52%	6.948%
6	7.710	933	938	943	rBB	27060	33193	1.51%	0.204%
7	8.181	1012	1018	1034	rBV	667002	862162	39.19%	5.286%
8	8.392	1049	1054	1061	rVB2	17080	23042	1.05%	0.141%
9	9.251	1194	1200	1205	rBV	1729083	2199760	100.00%	13.486%
10	9.934	1311	1316	1321	rBV	780223	964739	43.86%	5.915%
11	10.645	1432	1437	1444	rBV	183266	227090	10.32%	1.392%
12	10.722	1444	1450	1455	rBV	1243747	1578941	71.78%	9.680%
13	10.957	1479	1490	1495	rBV2	18470	26902	1.22%	0.165%
14	11.422	1564	1569	1576	rBV	793753	996003	45.28%	6.106%
15	11.945	1651	1658	1670	rBV2	129838	213476	9.70%	1.309%
16	12.198	1697	1701	1710	rVB4	16989	23521	1.07%	0.144%
17	12.633	1771	1775	1784	rBV	17344	30592	1.39%	0.188%
18	12.727	1787	1791	1800	rVV3	15545	26618	1.21%	0.163%
19	13.010	1833	1839	1844	rBV	1564594	2032369	92.39%	12.460%
20	13.904	1986	1991	2002	rBV	51974	104377	4.74%	0.640%
21	14.063	2009	2018	2028	rVB	435606	593491	26.98%	3.639%
22	15.557	2265	2272	2285	rBV	367513	601587	27.35%	3.688%

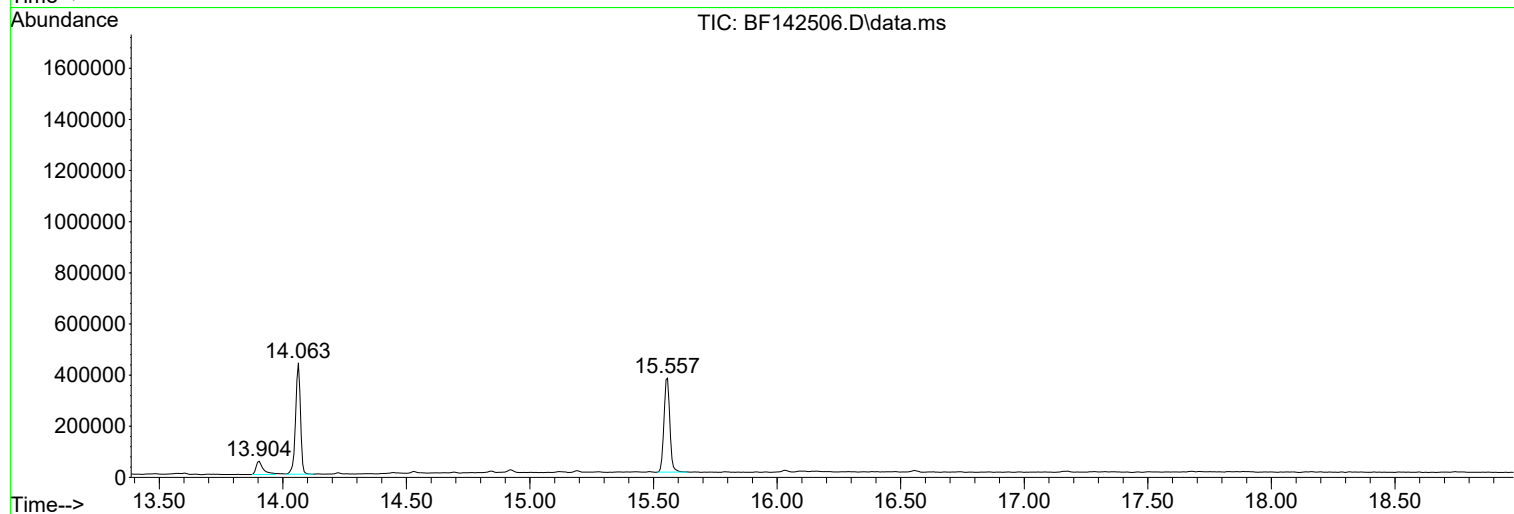
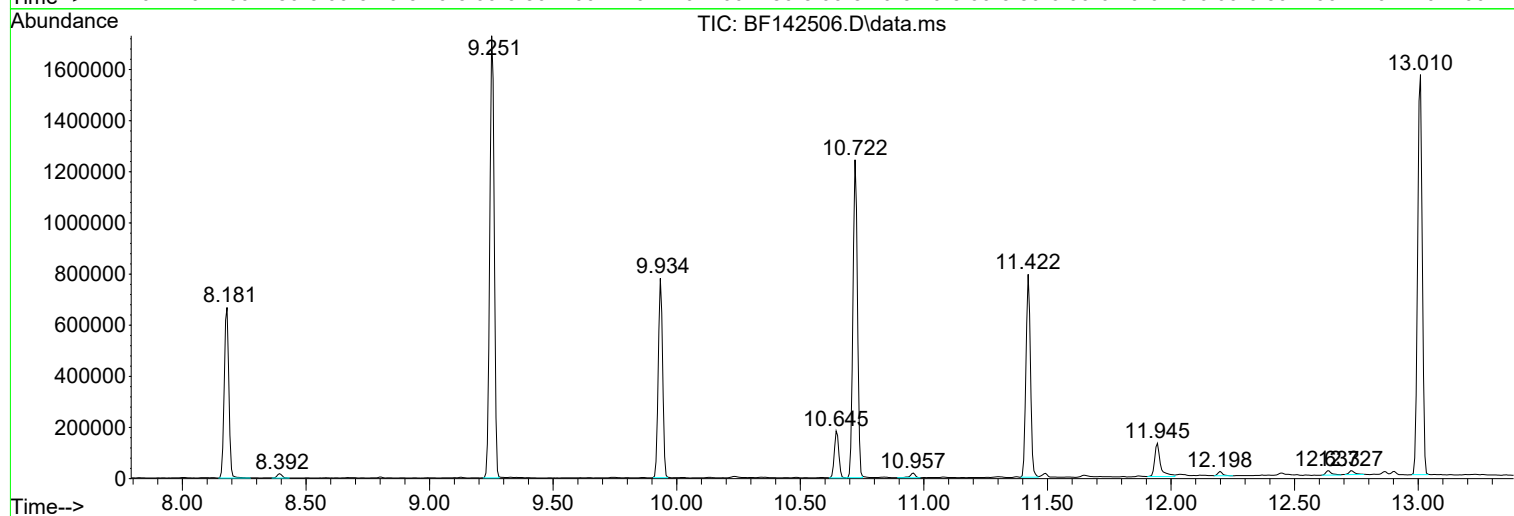
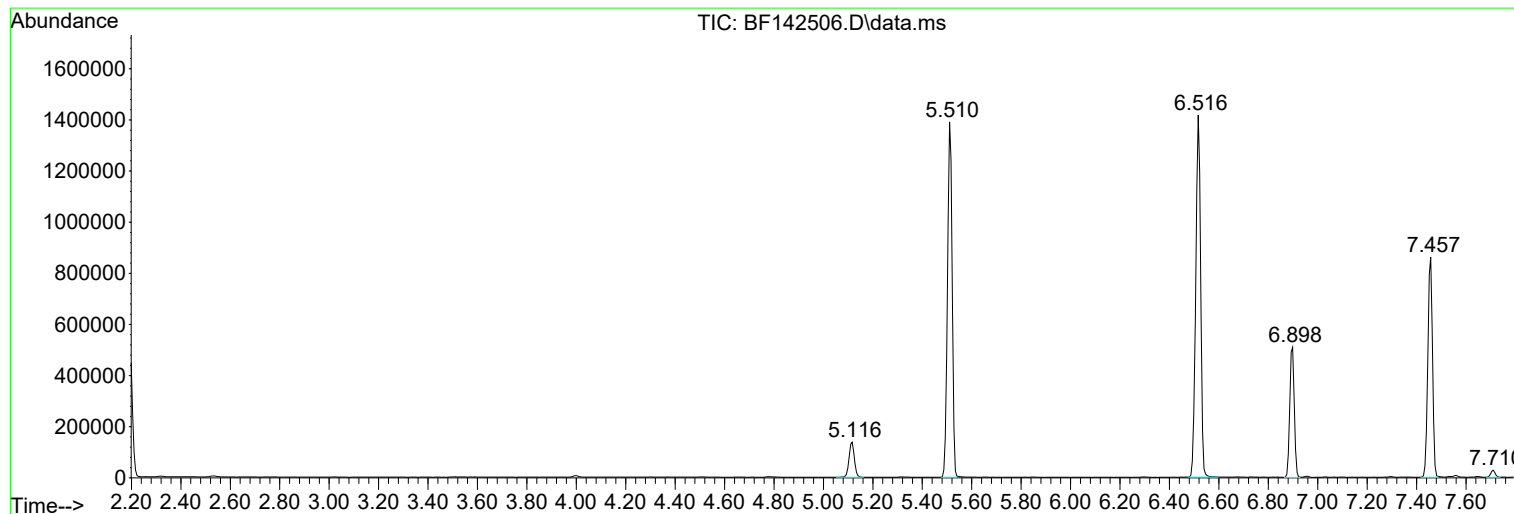
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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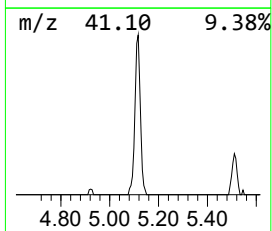
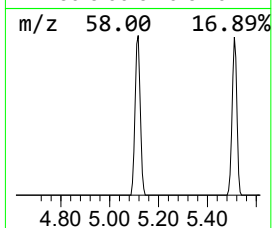
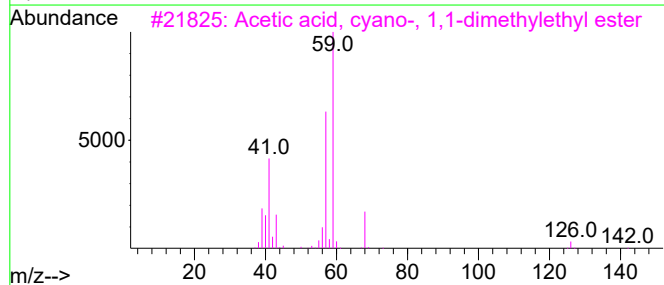
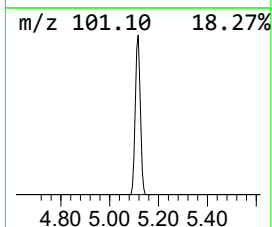
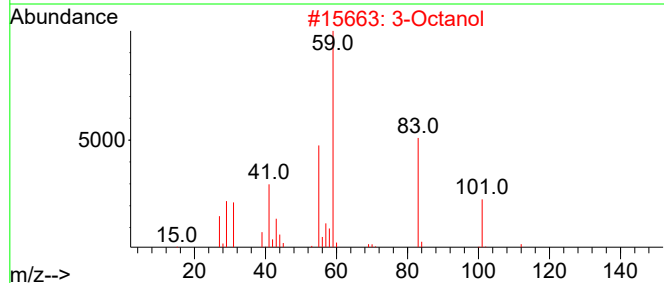
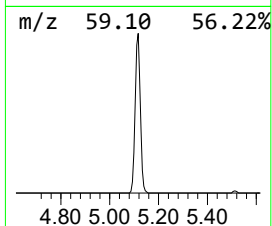
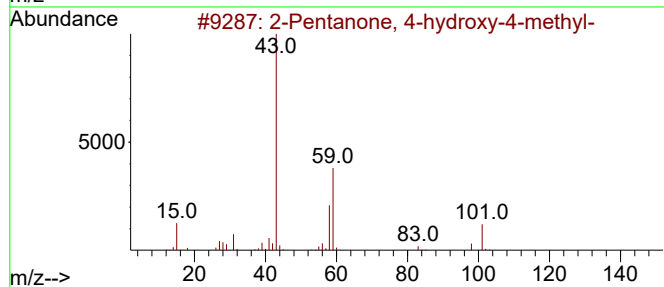
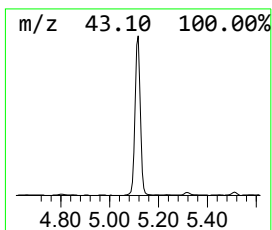
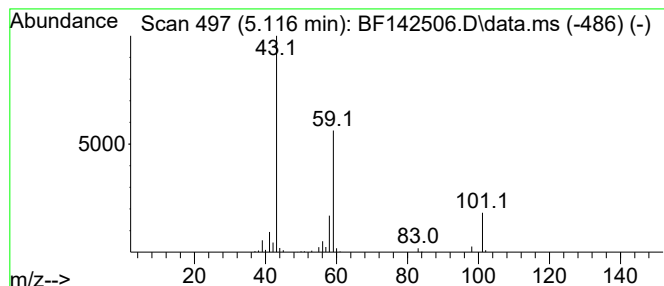
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.116	6.41 ng	207101	1,4-Dichlorobenzene-d4	6.898

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59		
2	3-Octanol	130	C8H18O	000589-98-0	36		
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25		
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12		
5	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	9		



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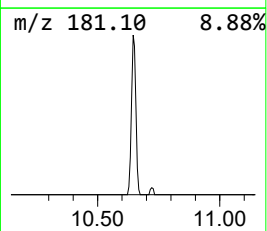
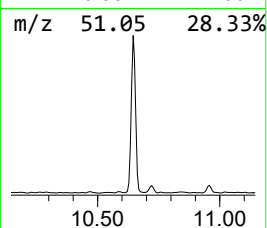
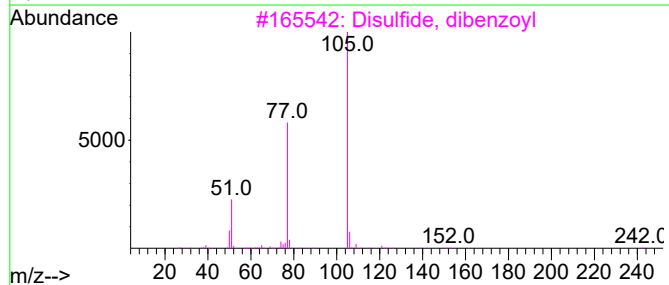
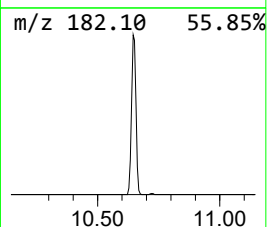
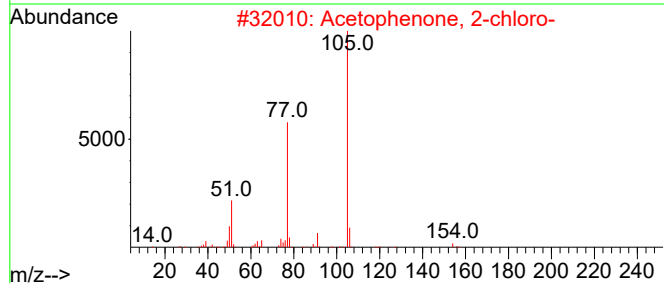
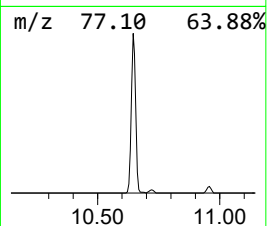
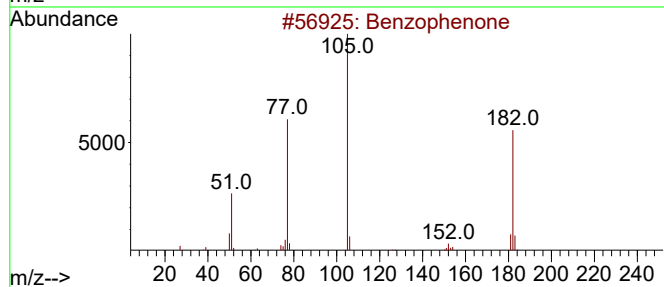
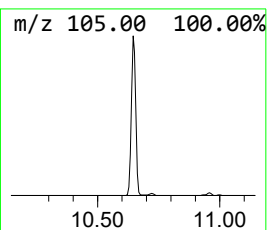
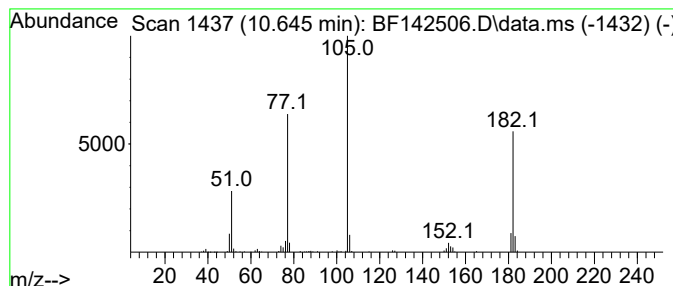
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Peak Number 2 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.645	4.71 ng	227090	Acenaphthene-d10	9.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	49
3			Disulfide, dibenzoyl	274	C14H10O2S2	000644-32-6	49
4			Pentafluoroethyl phenyl ketone	224	C9H5F5O	000394-52-5	47
5			N-(1-Cyanovinyl)benzamide	172	C10H8N2O	1000186-19-1	47



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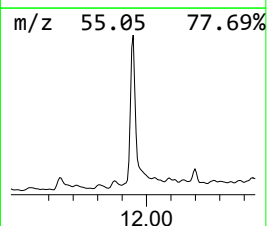
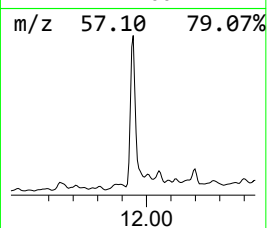
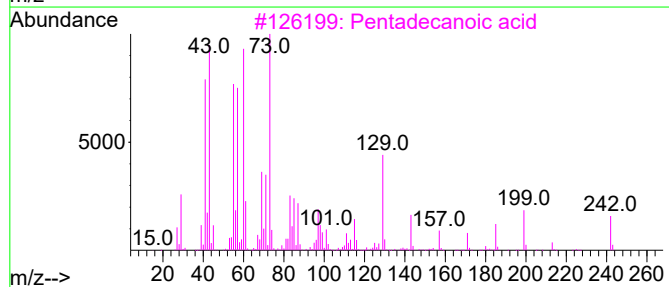
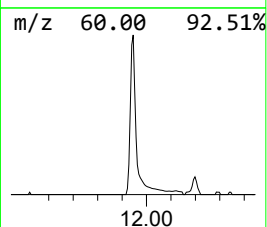
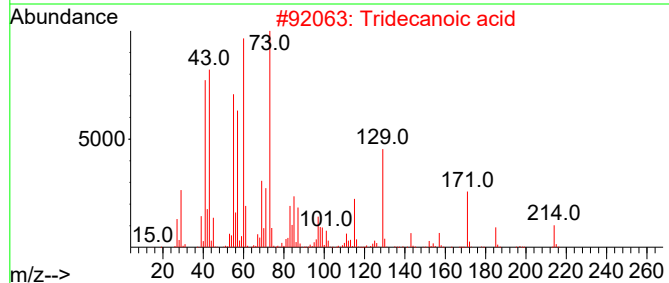
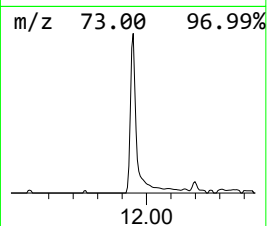
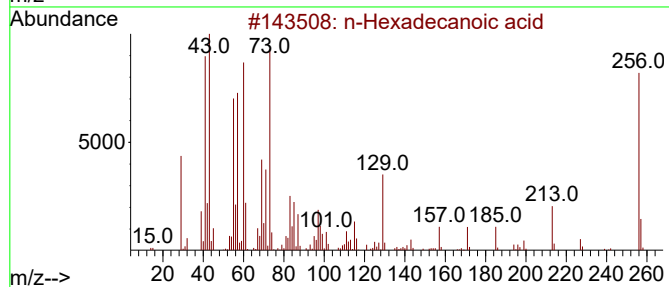
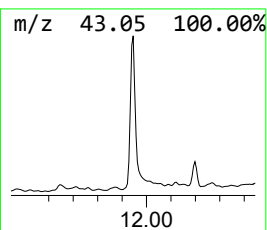
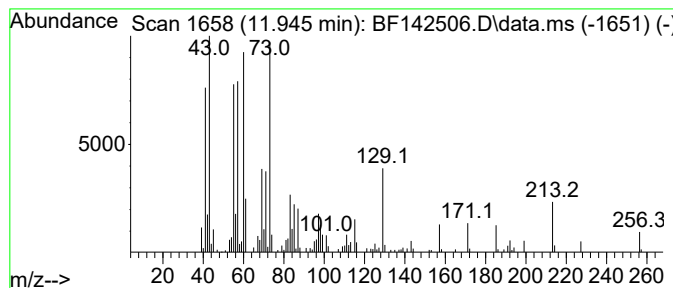
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Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.945	4.29 ng	213476	Phenanthrene-d10	11.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Tridecanoic acid	214	C13H26O2	000638-53-9	91
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5			n-Decanoic acid	172	C10H20O2	000334-48-5	58



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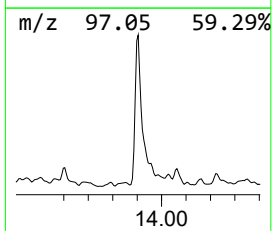
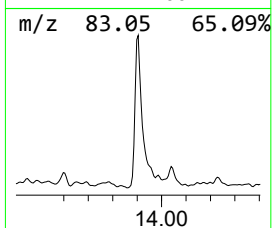
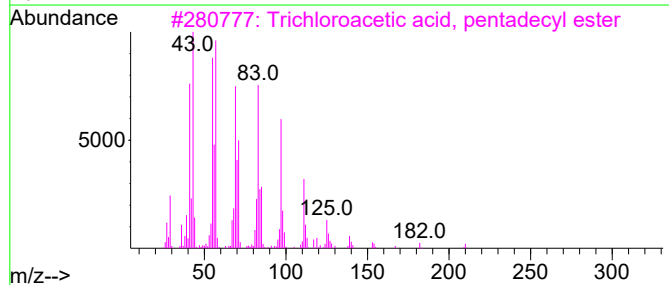
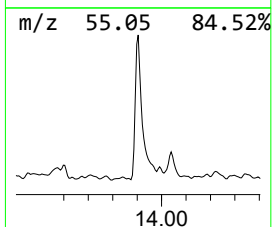
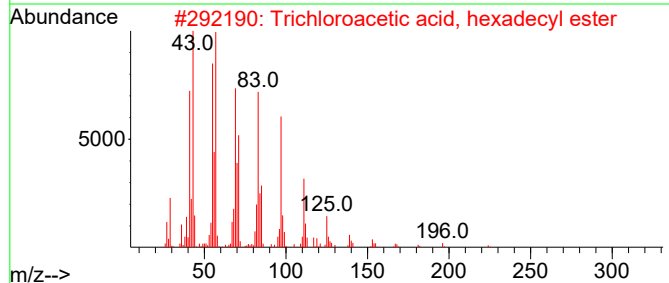
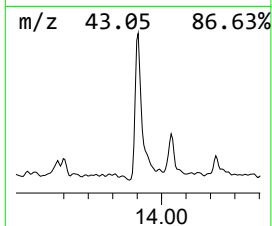
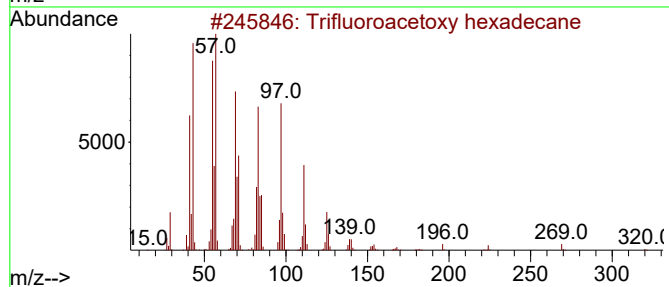
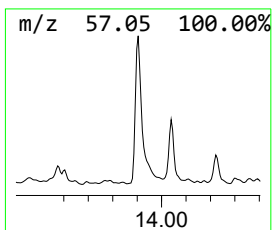
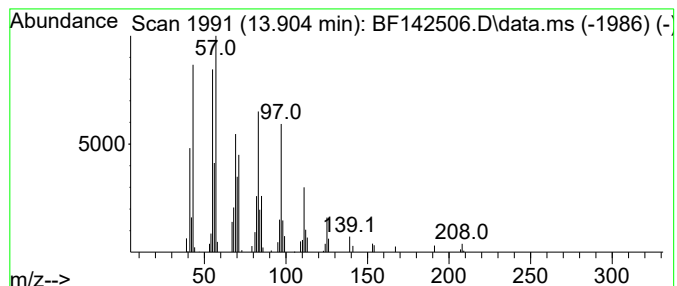
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Peak Number 4 Trifluoroacetoxy hexadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.904	3.52 ng	104377	Chrysene-d12	14.063

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	95
2			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	95
3			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	94
4			1-Docosene	308	C22H44	001599-67-3	94
5			17-Pentatriacontene	491	C35H70	006971-40-0	91



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.116	6.4	ng	207101	1	6.898	645890	20.0
Benzophenone	10.645	4.7	ng	227090	3	9.934	964739	20.0
n-Hexadecanoic ...	11.945	4.3	ng	213476	4	11.422	996003	20.0
Trifluoroacetox...	13.904	3.5	ng	104377	5	14.063	593491	20.0