

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF091918\
 Data File : BF109149.D
 Acq On : 19 Sep 2018 17:49
 Operator : JU/SJ
 Sample : J4983-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SAMPLE-01

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-BF091418.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	4.913	435	438	445	rVB	45351	56329	3.03%	0.308%
2	5.331	505	509	513	rBV	1690642	1542600	82.96%	8.444%
3	6.354	679	683	699	rBV2	1704276	1759935	94.64%	9.633%
4	6.489	702	706	709	rBV	1962679	1617890	87.01%	8.856%
5	6.719	742	745	749	rBV	920606	807616	43.43%	4.421%
6	6.878	768	772	775	rBV	1694767	1365292	73.42%	7.473%
7	7.278	836	840	843	rBV	1114645	937189	50.40%	5.130%
8	8.001	959	963	967	rBV	1168280	980971	52.75%	5.370%
9	9.077	1142	1146	1149	rBV	2259257	1859529	100.00%	10.179%
10	9.472	1209	1213	1220	rBV	247383	203253	10.93%	1.113%
11	9.754	1257	1261	1265	rBV	1153053	993851	53.45%	5.440%
12	10.536	1390	1394	1405	rBV	970483	881745	47.42%	4.826%
13	11.230	1508	1512	1515	rBV	1075375	887591	47.73%	4.858%
14	11.765	1598	1603	1606	rBV	30804	30413	1.64%	0.166%
15	12.812	1776	1781	1784	rBV	1622224	1519131	81.69%	8.315%
16	13.401	1878	1881	1887	rVB2	19479	22915	1.23%	0.125%
17	13.718	1932	1935	1943	rBV	102534	136191	7.32%	0.745%
18	13.854	1954	1958	1962	rBV	1071996	978987	52.65%	5.359%
19	13.971	1975	1978	1980	rVB3	24939	21629	1.16%	0.118%
20	14.042	1987	1990	1996	rVB	42101	43948	2.36%	0.241%
21	14.342	2038	2041	2046	rBV2	55845	59085	3.18%	0.323%
22	14.636	2088	2091	2096	rVB	73865	74039	3.98%	0.405%
23	14.706	2100	2103	2108	rVB	75852	75149	4.04%	0.411%
24	14.954	2142	2145	2150	rBV	91787	101441	5.46%	0.555%
25	15.259	2192	2197	2202	rBV	715242	868136	46.69%	4.752%
26	15.306	2202	2205	2211	rVB	95033	129285	6.95%	0.708%
27	15.712	2270	2274	2280	rBV	105143	162367	8.73%	0.889%
28	16.183	2349	2354	2365	rVB2	79739	152599	8.21%	0.835%

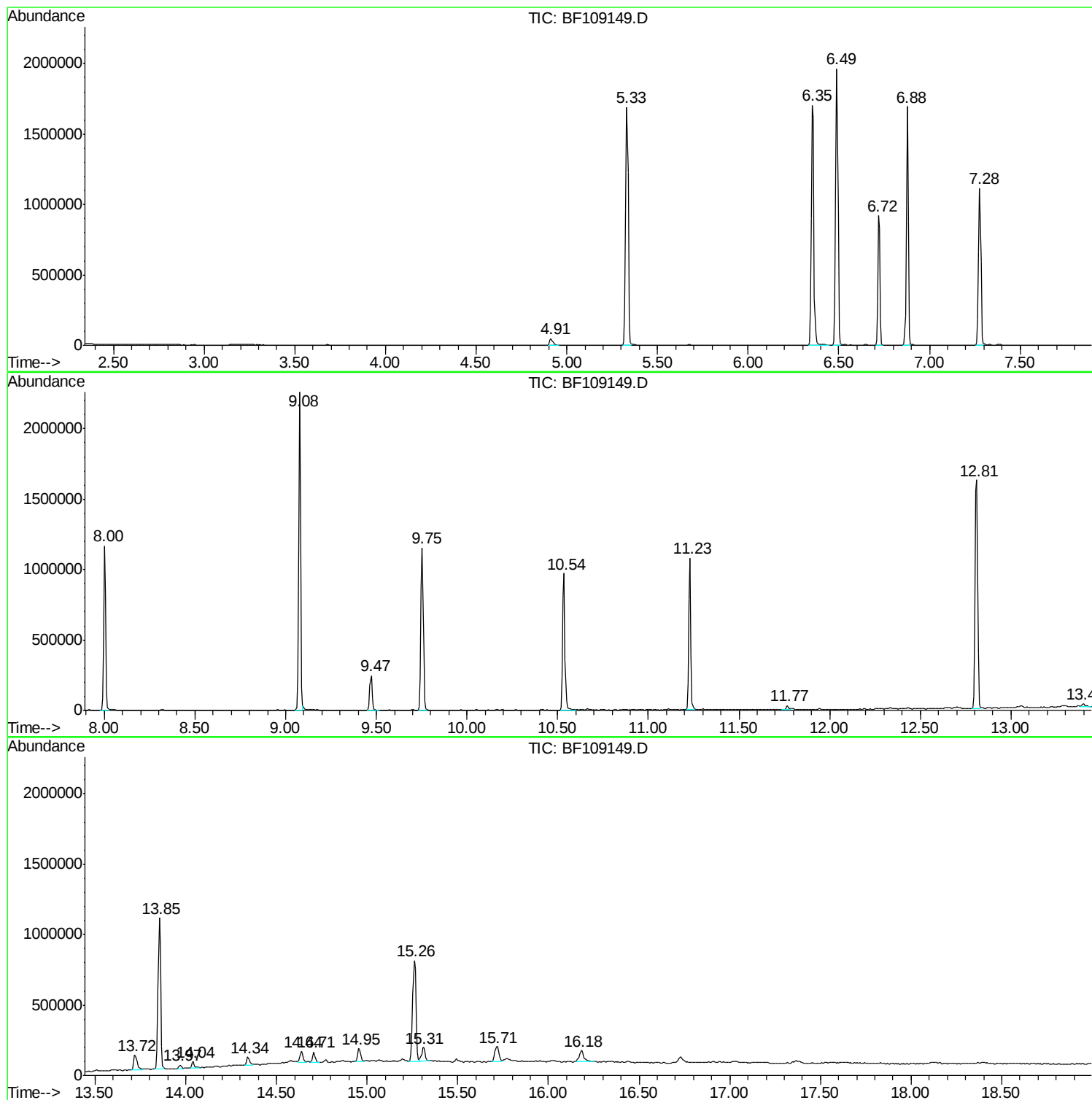
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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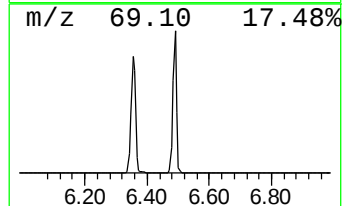
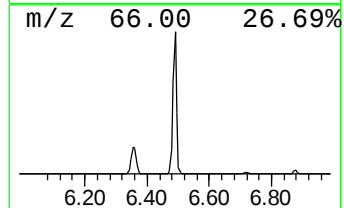
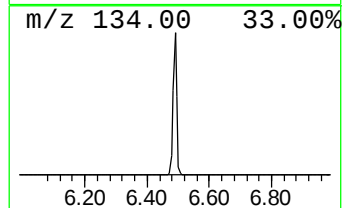
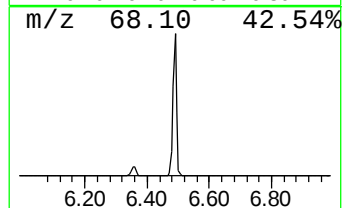
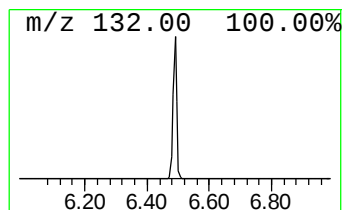
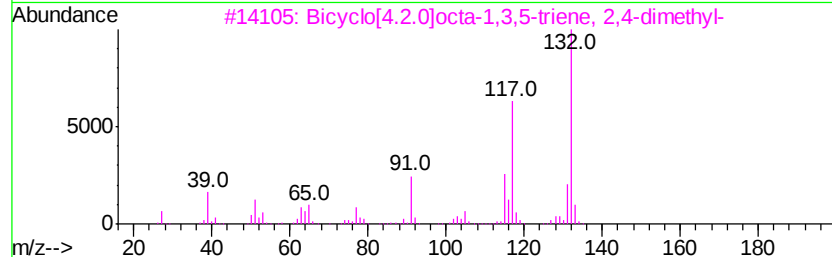
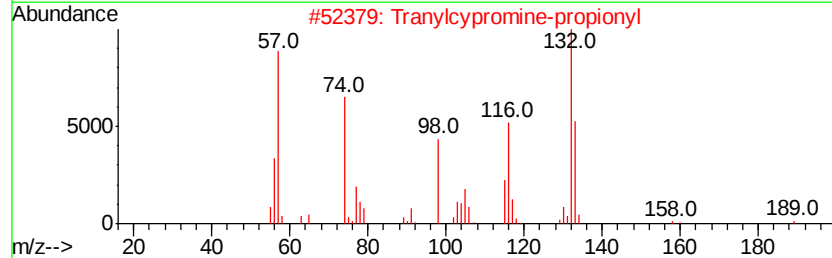
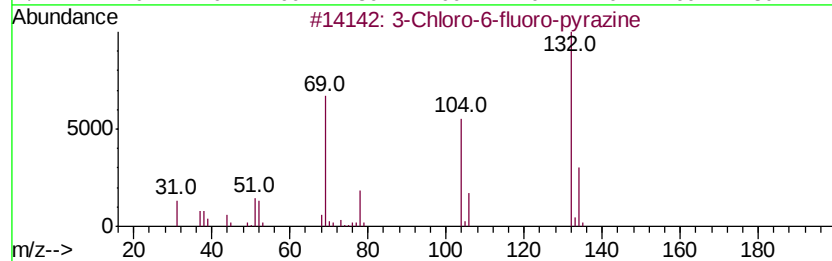
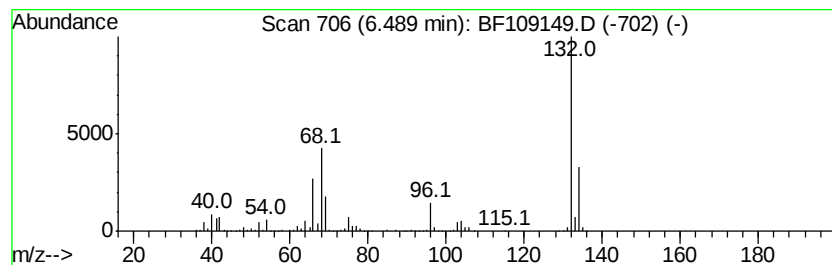
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.49 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	40.07 ng	1617890	1,4-Dichlorobenzene-d4	6.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
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		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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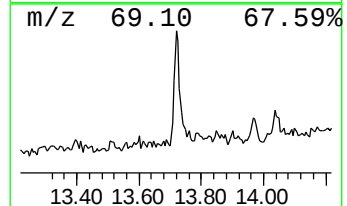
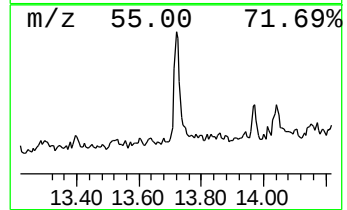
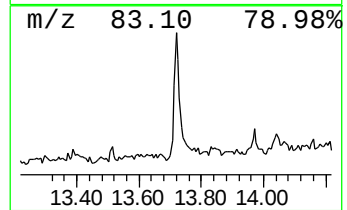
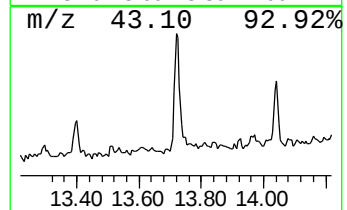
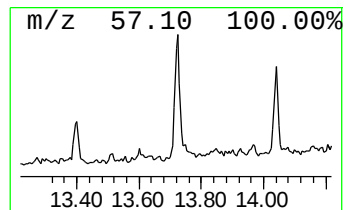
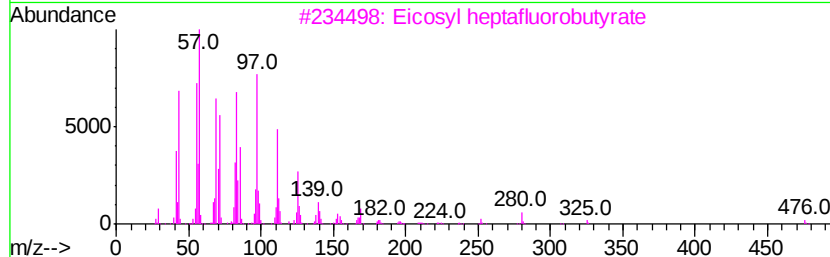
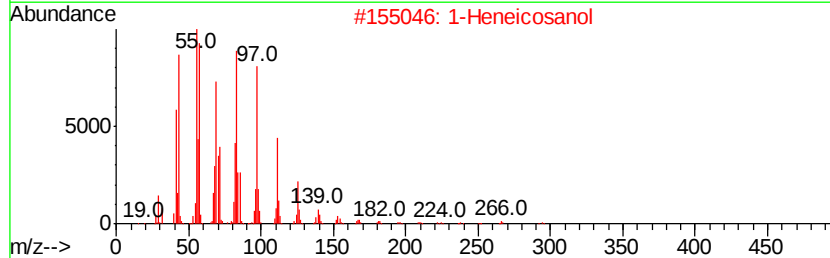
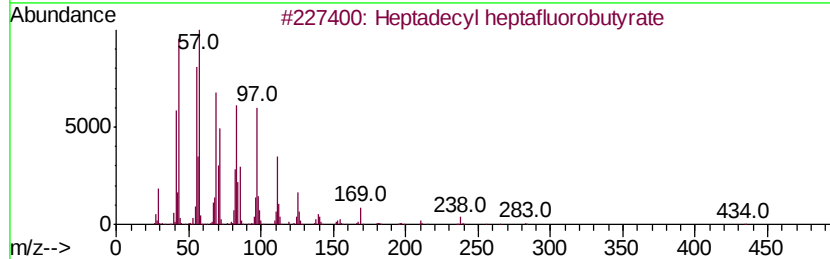
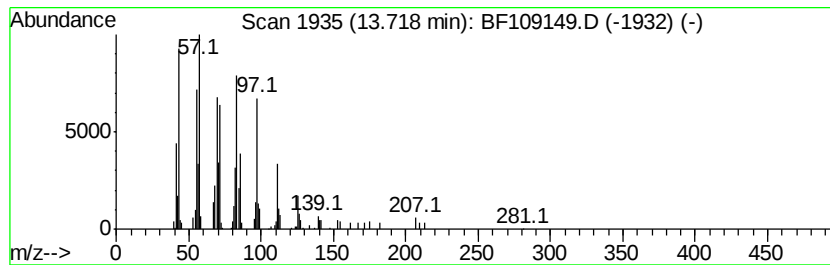
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Heptadecyl heptafluorobutyrate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	2.78 ng	136191	Chrysene-d12	13.85

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecyl heptafluorobutvrate	452	C21H35F7O2	959085-66-6	94
2		1-Heneicosanol	312	C21H44O	015594-90-8	93
3		Eicosyl heptafluorobutvrate	494	C24H41F7O2	1000351-83-5	91
4		Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	91
5		Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91



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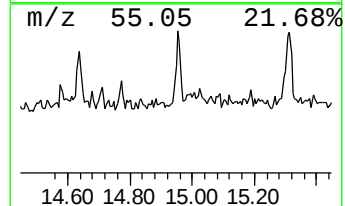
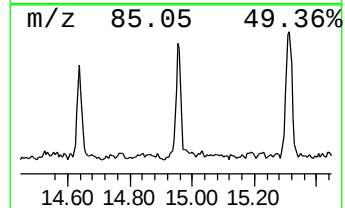
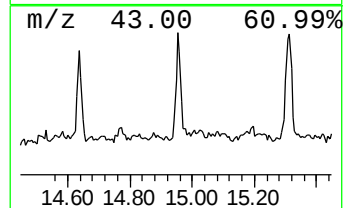
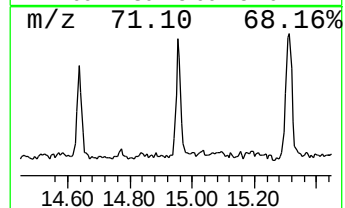
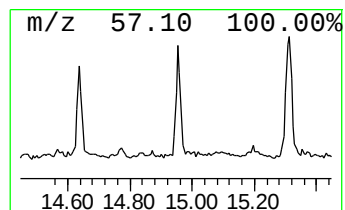
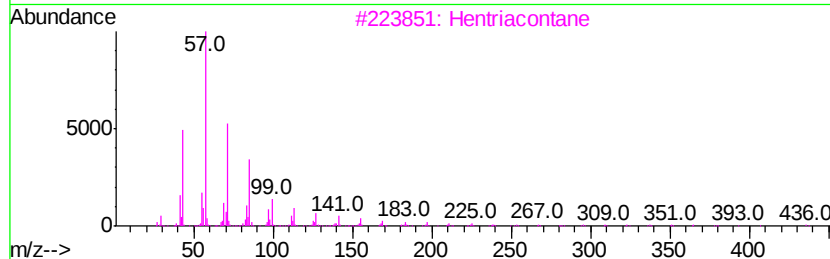
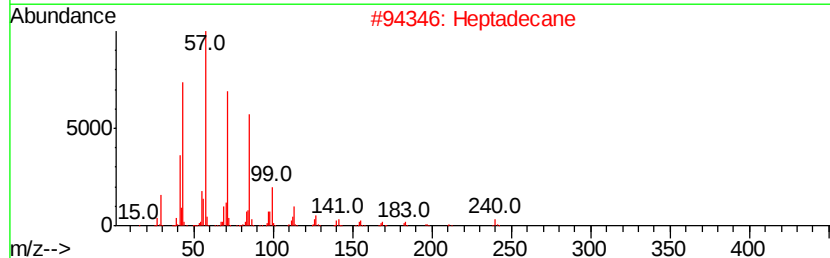
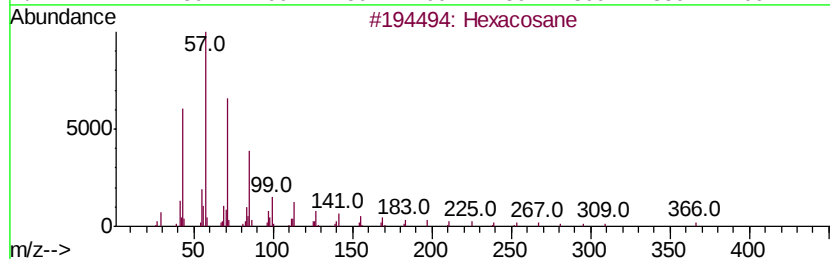
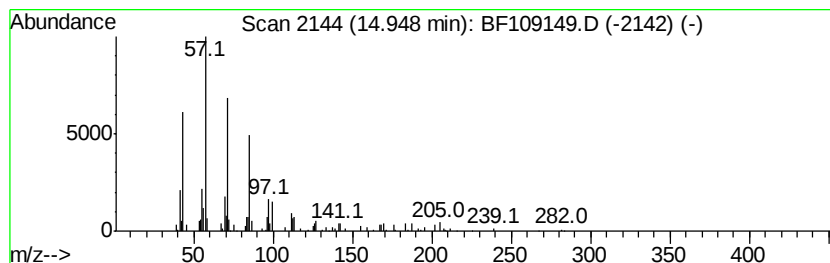
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Hexacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.95	2.34 ng	101441	Perylene-d12	15.26

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane		366	C26H54	000630-01-3	91
2	Heptadecane		240	C17H36	000629-78-7	90
3	Hentriacontane		437	C31H64	000630-04-6	90
4	Heneicosane		296	C21H44	000629-94-7	87
5	Methoxyacetic acid, 2-tetradecyl...		286	C17H34O3	1000282-04-8	86



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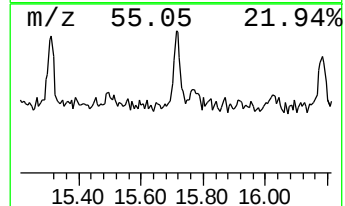
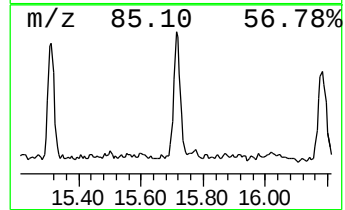
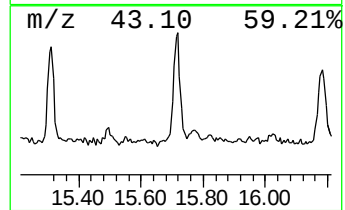
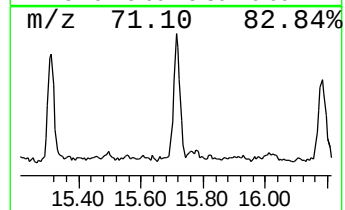
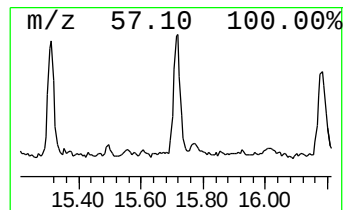
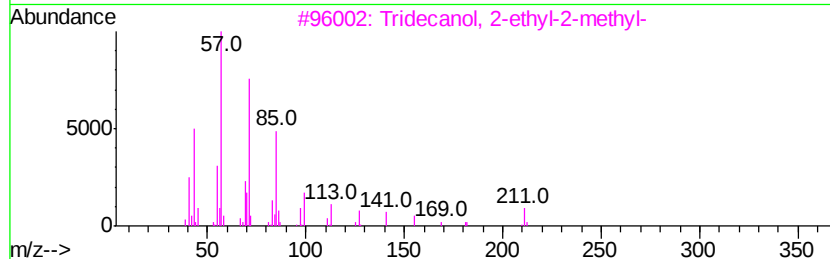
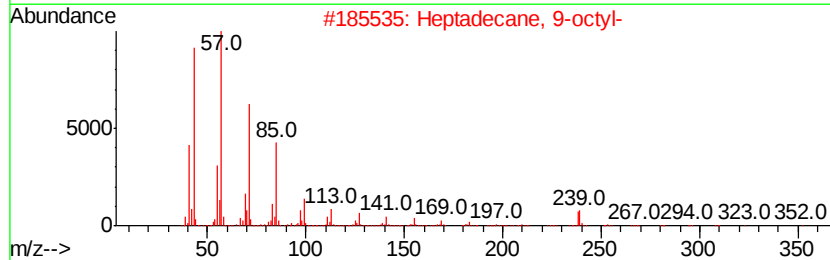
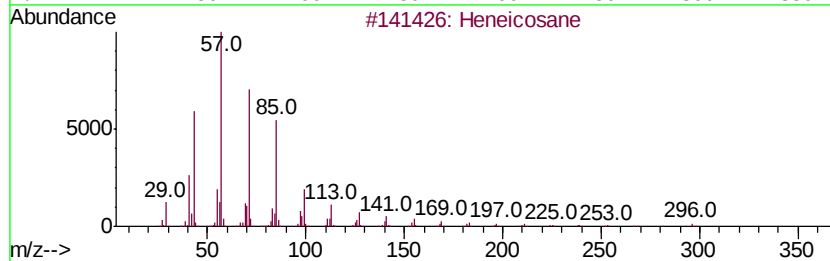
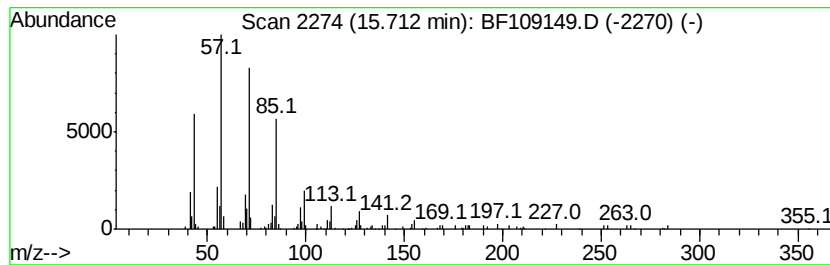
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Heneicosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.71	3.74 ng	162367	Perylene-d12	15.26

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
3	Tridecanol, 2-ethyl-2-methyl-		242	C16H34O	1000115-66-1	91
4	Triacontane		422	C30H62	000638-68-6	91
5	Hexacosane		366	C26H54	000630-01-3	90



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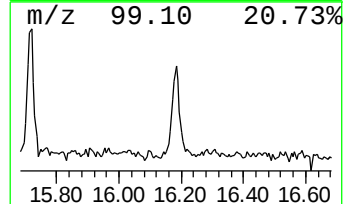
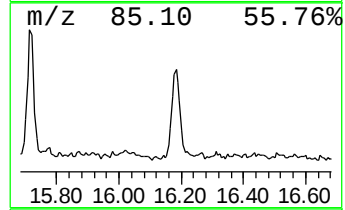
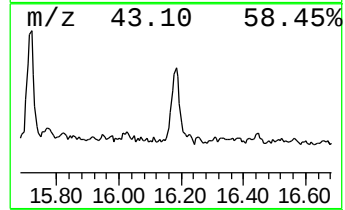
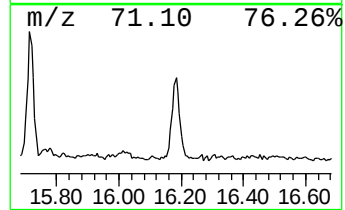
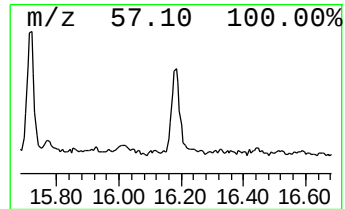
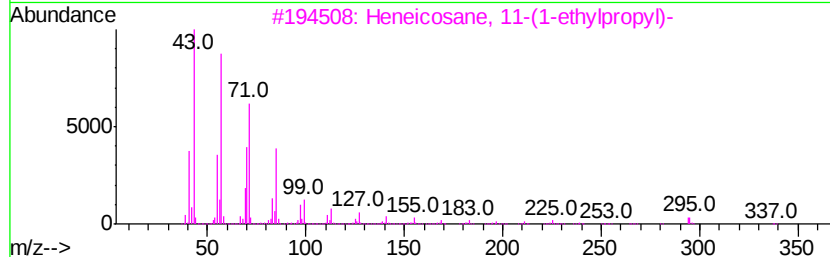
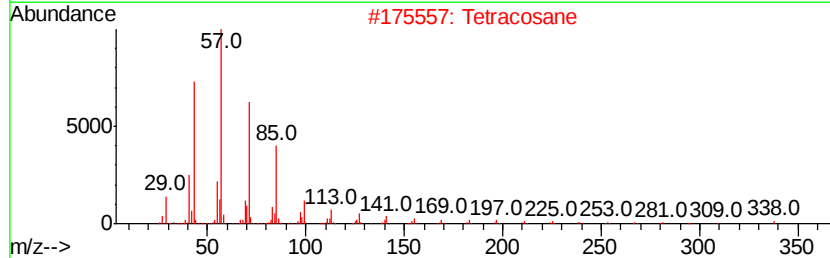
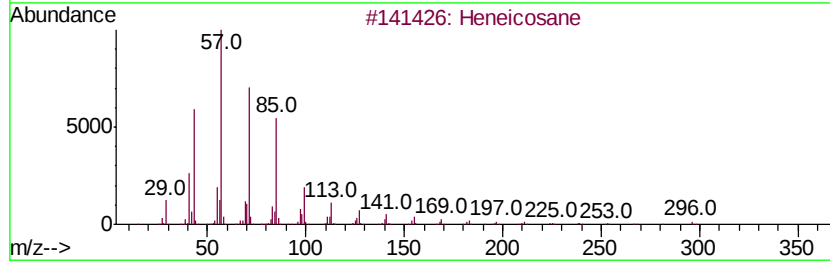
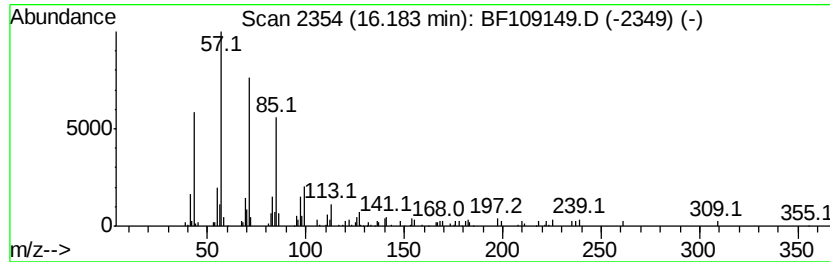
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Tetracosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.18	3.52 ng	152599	Perylene-d12	15.26

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	90
2	Tetracosane		338	C24H50	000646-31-1	90
3	Heneicosane, 11-(1-ethylpropyl)-		366	C26H54	055282-11-6	90
4	Octacosane		394	C28H58	000630-02-4	87
5	Hexadecane		226	C16H34	000544-76-3	83



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.49	6.49	40.1	ng	1617890	1	6.72	807616	20.0
Heptadecyl heptaf...	13.72	2.8	ng	136191	5	13.85	978987	20.0
Hexacosane	14.95	2.3	ng	101441	6	15.26	868136	20.0
Heneicosane	15.71	3.7	ng	162367	6	15.26	868136	20.0
Tetracosane	16.18	3.5	ng	152599	6	15.26	868136	20.0