

Data Path : Z:\HPCHEM1\BNA_F\Data\BF071417\
 Data File : BF096802.D
 Acq On : 15 Jul 2017 5:30
 Operator : SJ/JU
 Sample : I4159-01
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NORTH-DIESEL-SIDEWALL

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-BF071317.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	4.810	475	480	494	rVB	326346	530472	18.00%	2.396%
2	5.246	548	554	557	rBV	2664062	2947604	100.00%	13.316%
3	6.281	724	730	733	rBV	2478618	2872042	97.44%	12.975%
4	6.398	745	750	754	rBV	2760326	2917060	98.96%	13.178%
5	6.628	785	789	792	rBV	752283	629406	21.35%	2.843%
6	6.787	811	816	819	rBV	2223732	2067776	70.15%	9.341%
7	7.192	879	885	888	rBV	1899334	1790597	60.75%	8.089%
8	7.904	1003	1006	1010	rBV	805373	744176	25.25%	3.362%
9	8.986	1185	1190	1193	rBV	2456473	2154172	73.08%	9.732%
10	9.381	1253	1257	1260	rBV	295101	233956	7.94%	1.057%
11	9.657	1300	1304	1308	rBV	699662	583718	19.80%	2.637%
12	10.104	1377	1380	1383	rVB	52653	41545	1.41%	0.188%
13	10.445	1433	1438	1441	rBV	1190254	1140494	38.69%	5.152%
14	10.575	1458	1460	1468	rBV	52009	57832	1.96%	0.261%
15	11.133	1551	1555	1558	rVB	595420	482171	16.36%	2.178%
16	12.722	1820	1825	1828	rBV	1818029	1697062	57.57%	7.667%
17	13.627	1975	1979	1985	rVB	231285	265203	9.00%	1.198%
18	13.763	1998	2002	2005	rBV	616199	552888	18.76%	2.498%
19	14.521	2128	2131	2134	rBV	82319	74146	2.52%	0.335%
20	15.145	2233	2237	2240	rVB	303820	353148	11.98%	1.595%

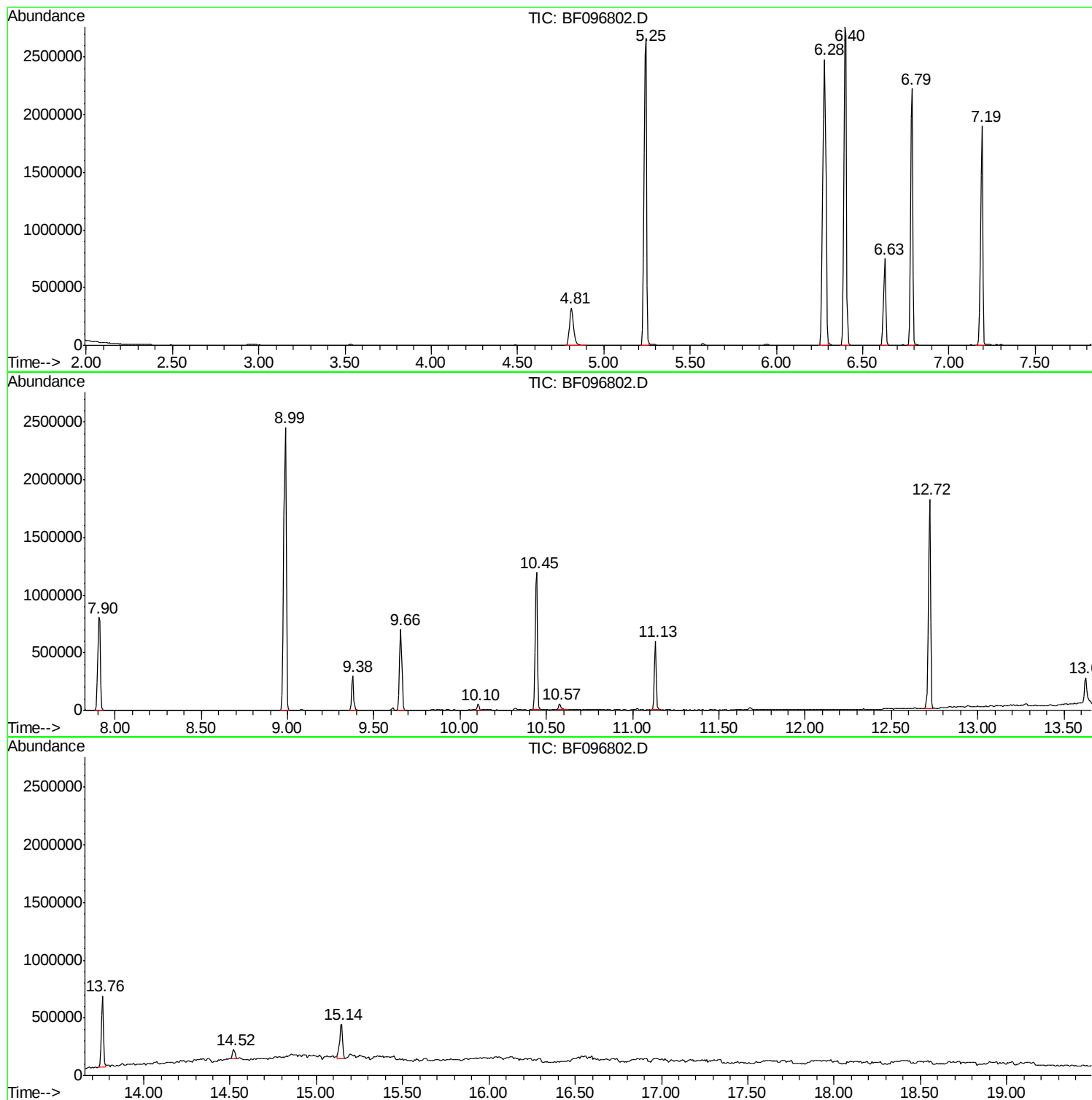
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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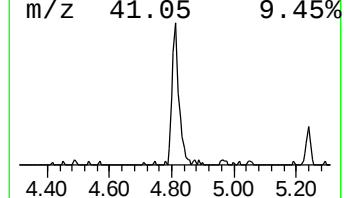
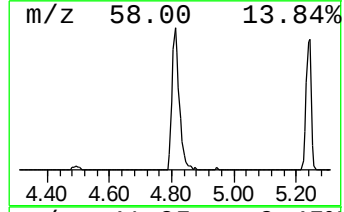
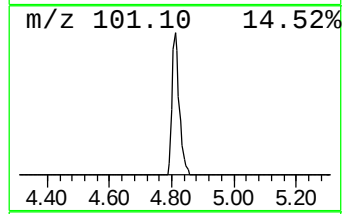
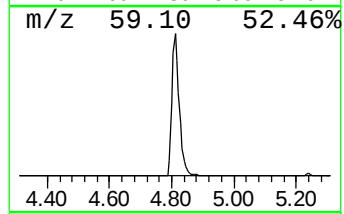
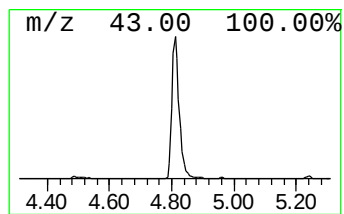
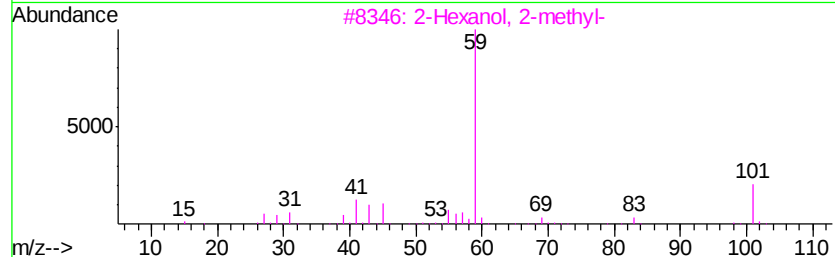
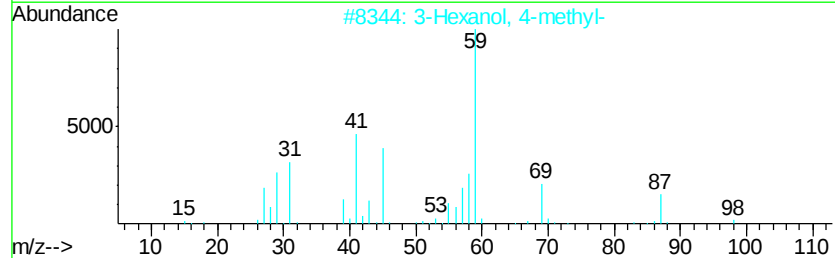
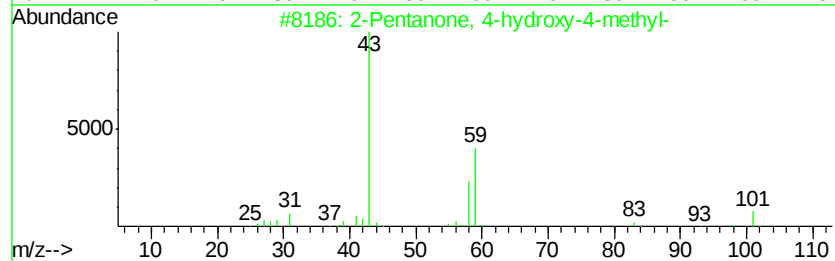
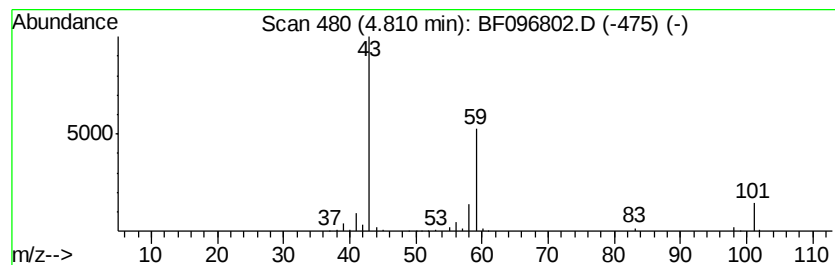
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.81	16.86 ng	530472	1,4-Dichlorobenzene-d4	6.63

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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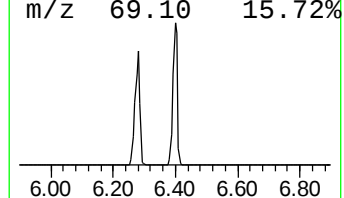
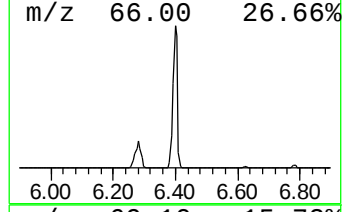
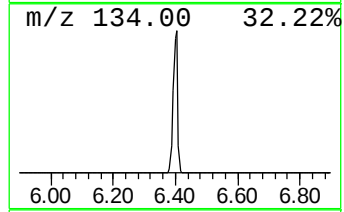
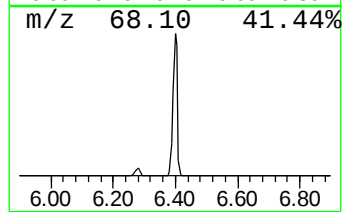
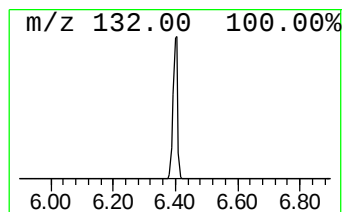
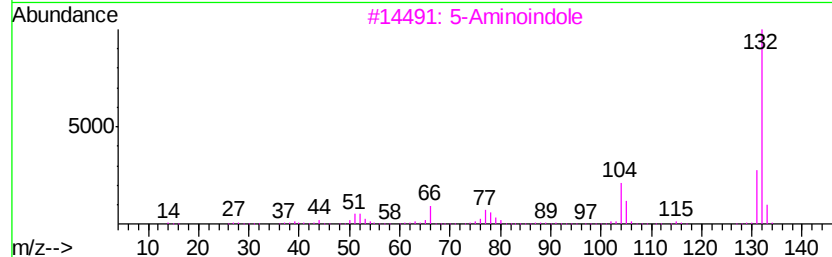
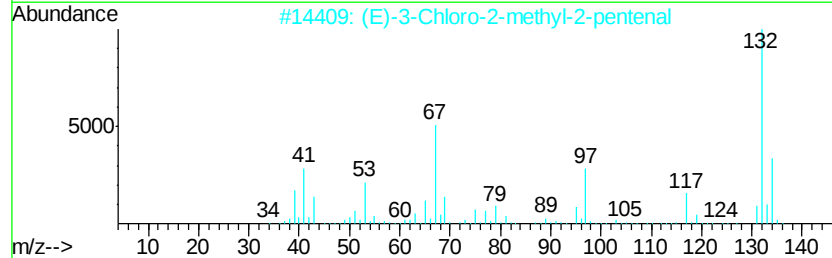
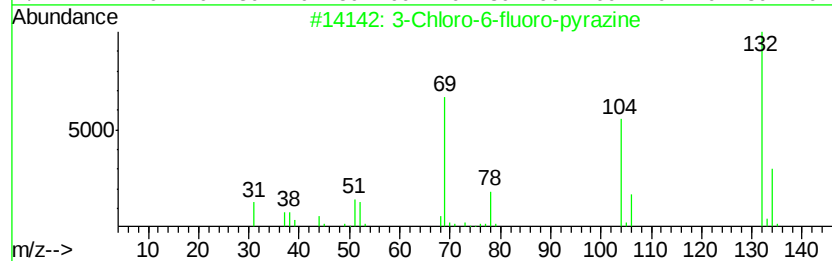
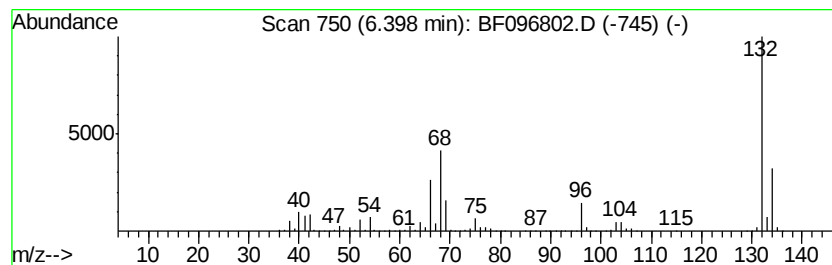
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Peak Number 2 unknown6.40 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.40	92.69 ng	2917060	1,4-Dichlorobenzene-d4	6.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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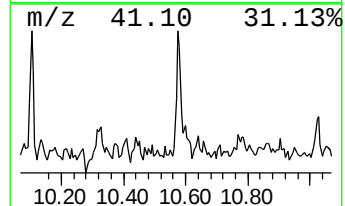
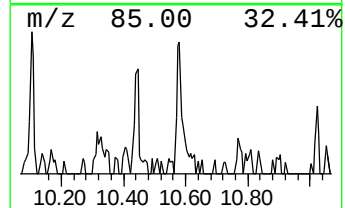
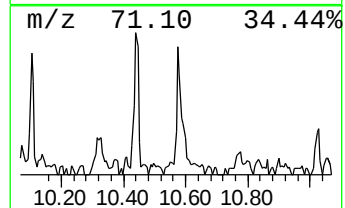
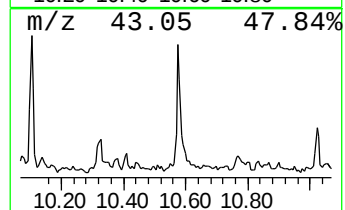
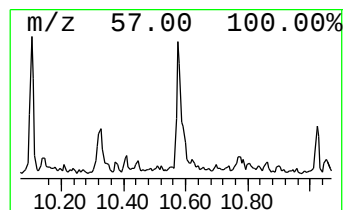
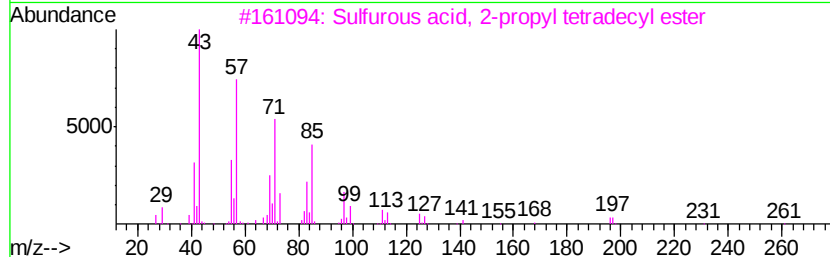
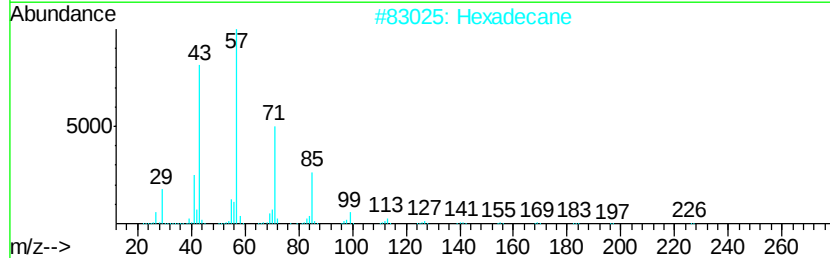
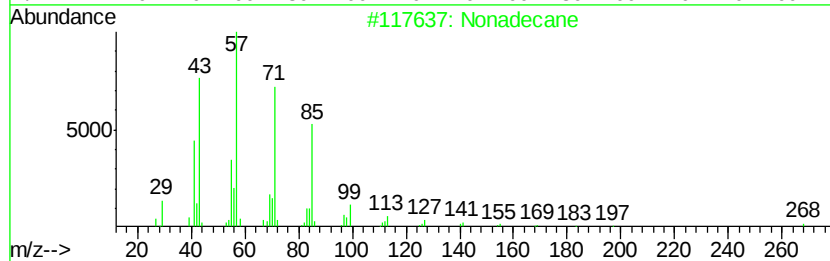
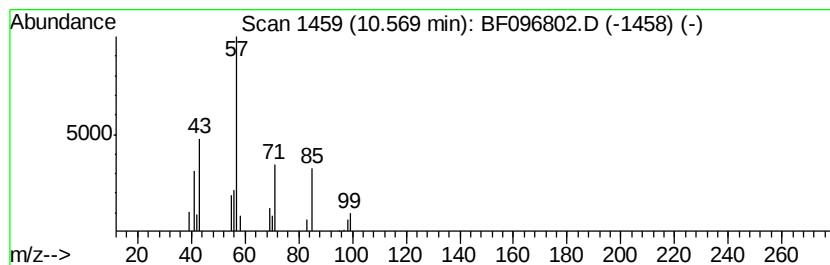
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Peak Number 4 Nonadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.57	2.40 ng	57832	Phenanthrene-d10	11.13

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	80
2	Hexadecane		226	C16H34	000544-76-3	72
3	Sulfurous acid, 2-propyl tetradec...		320	C17H36O3S	1000309-12-5	64
4	Sulfurous acid, 2-ethylhexyl iso...		278	C14H30O3S	1000309-19-0	64
5	Dodecane, 2,6,11-trimethyl-		212	C15H32	031295-56-4	59



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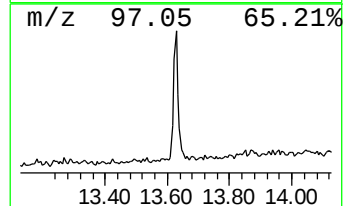
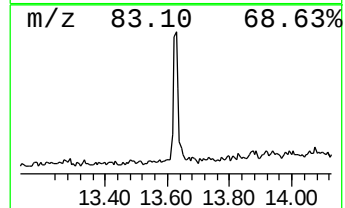
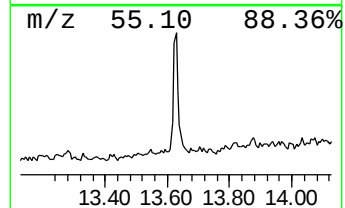
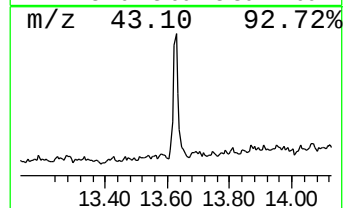
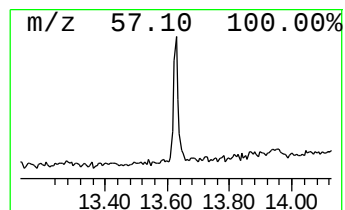
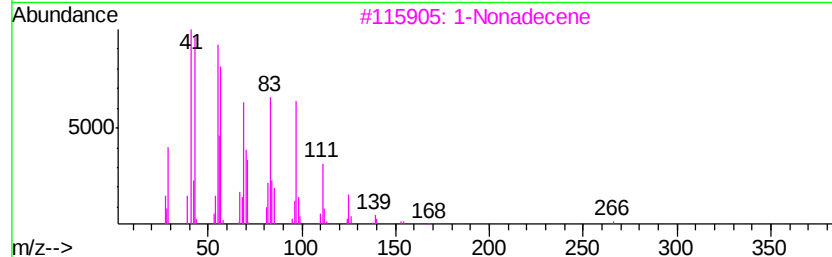
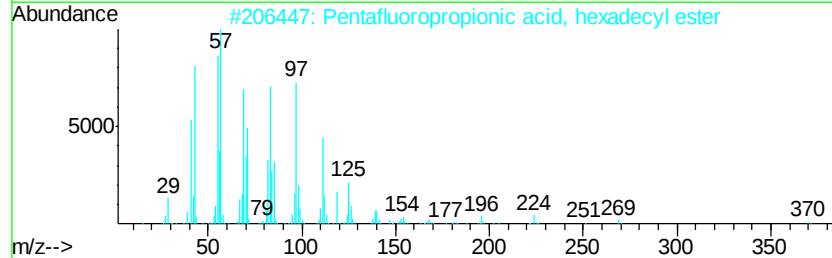
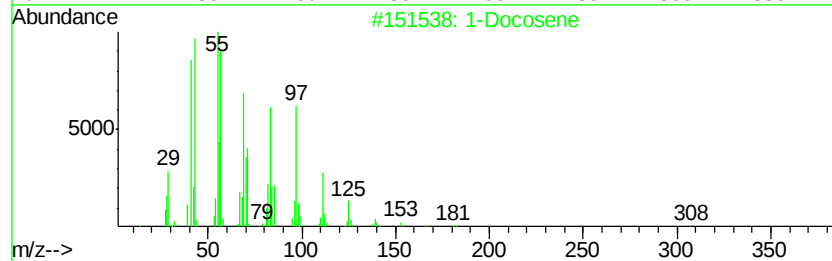
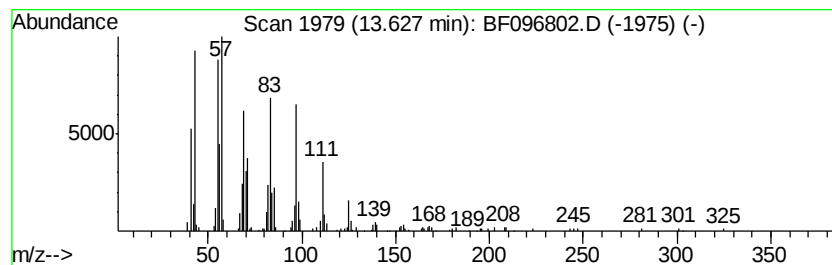
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Peak Number 5 1-Docosene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.63	9.59 ng	265203	Chrysene-d12	13.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Docosene	308	C22H44	001599-67-3	94
2		Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	94
3		1-Nonadecene	266	C19H38	018435-45-5	91
4		1-Octadecanol	270	C18H38O	000112-92-5	91
5		10-Heneicosene (c,t)	294	C21H42	095008-11-0	91



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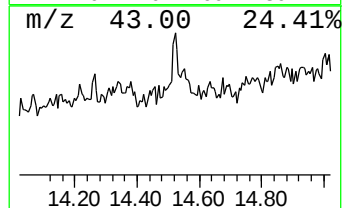
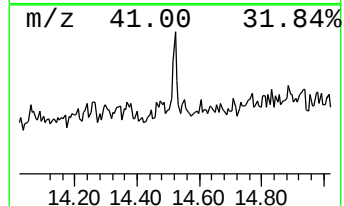
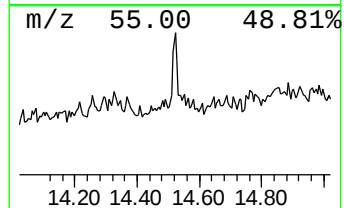
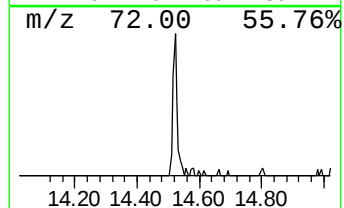
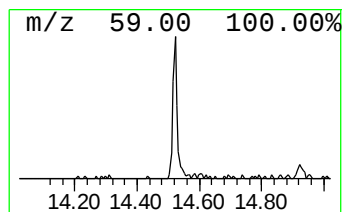
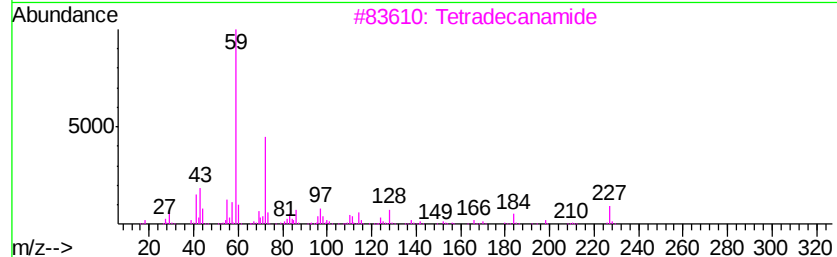
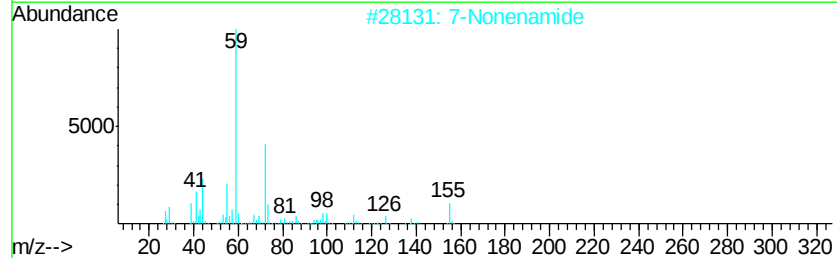
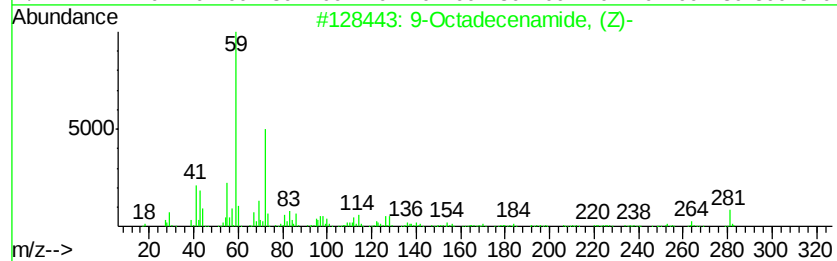
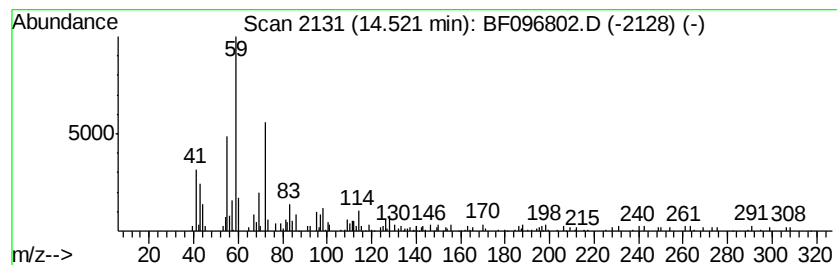
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 9-Octadecenamide, (Z)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.52	4.20 ng	74146	Perylene-d12	15.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	78
2		7-Nonenamide	155	C9H17NO	090949-53-4	72
3		Tetradecanamide	227	C14H29NO	000638-58-4	72
4		Benzeneethanamine, 2-fluoro-.bet...	229	C11H16FN03	061338-98-5	59
5		Hexadecanamide	255	C16H33NO	000629-54-9	53



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.81	16.9	ng	530472	1	6.63	629406	20.0
unknown6.40	6.40	92.7	ng	2917060	1	6.63	629406	20.0
Nonadecane	10.57	2.4	ng	57832	4	11.13	482171	20.0
1-Docosene	13.63	9.6	ng	265203	5	13.76	552888	20.0
9-Octadecenamide,...	14.52	4.2	ng	74146	6	15.14	353148	20.0