

Data Path : Z:\HPCHEM1\BNA F\DATA\BF010416\
 Data File : BF084249.D
 Acq On : 4 Jan 2016 19:57
 Operator : UM/SJ
 Sample : G4986-03
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PJ-SB-4-1.5-2.0

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-BF123115.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	2.155	4	6	16	rVB4	68485	178366	3.26%	0.332%
2	5.081	260	262	267	rVB	1415436	1651750	30.20%	3.076%
3	5.504	296	299	301	rBV	5797161	5391055	98.56%	10.040%
4	6.533	386	389	391	rBV	5255913	5194927	94.97%	9.675%
5	6.659	398	400	403	rVB	5337700	5337614	97.58%	9.941%
6	6.887	418	420	423	rVB	1968681	1692187	30.94%	3.151%
7	7.047	431	434	436	rBV	5088660	4327320	79.11%	8.059%
8	7.459	467	470	472	rBV	3714215	4047312	73.99%	7.538%
9	7.550	475	478	480	rBV	49625	66571	1.22%	0.124%
10	8.179	530	533	535	rBV	2606483	2341981	42.82%	4.362%
11	9.253	624	627	630	rBV	6147274	5469979	100.00%	10.187%
12	9.653	660	662	664	rBV	1372481	1203838	22.01%	2.242%
13	9.870	676	681	682	rBV	43299	65029	1.19%	0.121%
14	9.928	684	686	689	rVB	1938026	2339586	42.77%	4.357%
15	10.362	720	724	726	rVB	125242	143515	2.62%	0.267%
16	10.590	740	744	747	rBV	48258	96856	1.77%	0.180%
17	10.716	753	755	758	rBV2	3248706	3709170	67.81%	6.908%
18	10.842	764	766	769	rBV	173904	261501	4.78%	0.487%
19	11.288	803	805	806	rBV	144239	127078	2.32%	0.237%
20	11.413	814	816	819	rBV	2406546	2171578	39.70%	4.044%
21	13.002	953	955	957	rBV	4943444	4293521	78.49%	7.996%
22	13.482	995	997	1003	rBV3	33934	83934	1.53%	0.156%
23	13.905	1031	1034	1040	rBV	255725	370795	6.78%	0.691%
24	14.054	1044	1047	1049	rBV	1469518	1564294	28.60%	2.913%
25	14.534	1087	1089	1091	rBV2	44425	79319	1.45%	0.148%
26	15.494	1170	1173	1175	rBV	1094584	1424270	26.04%	2.653%
27	15.962	1212	1214	1217	rBV3	31589	61570	1.13%	0.115%

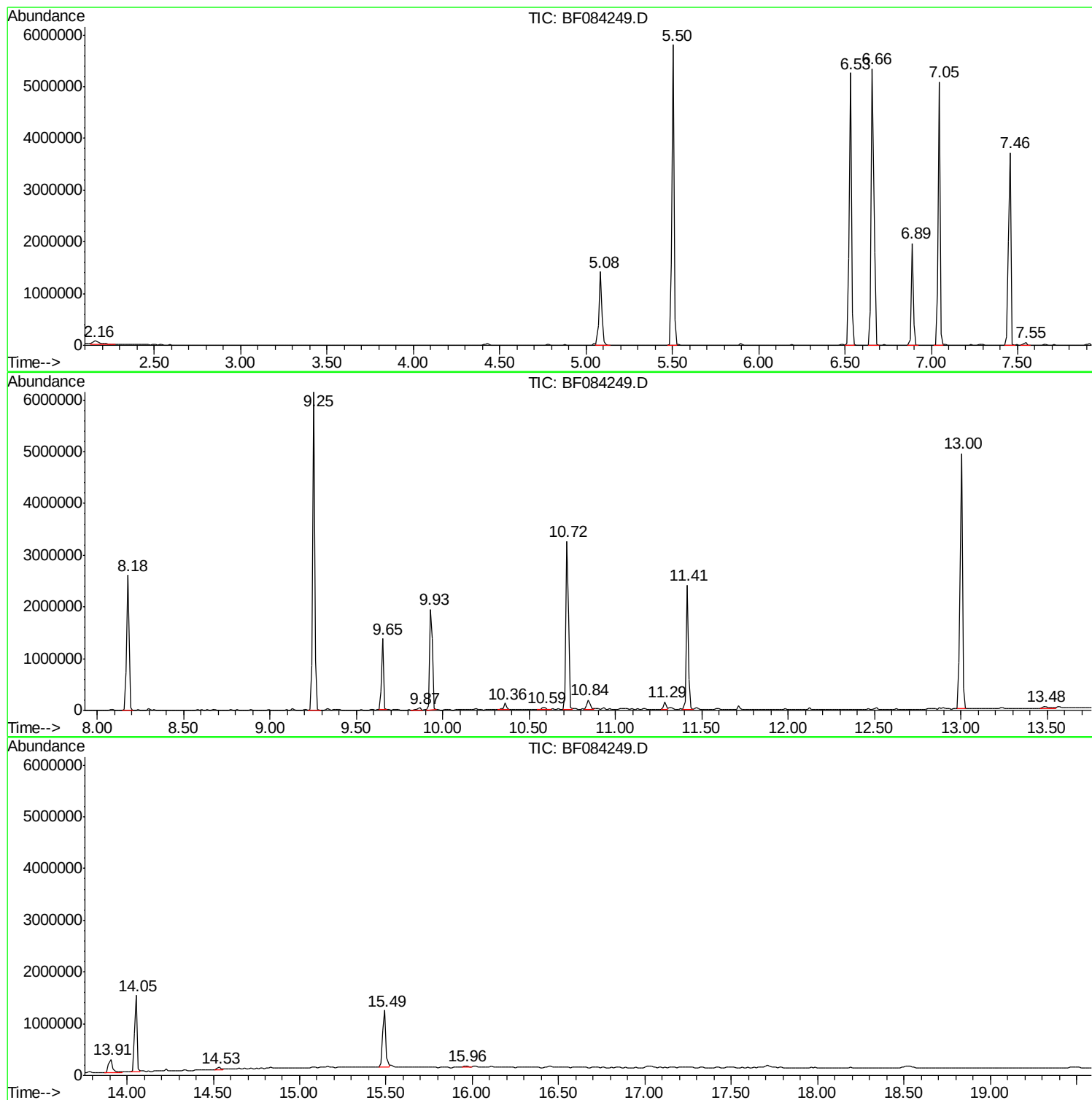
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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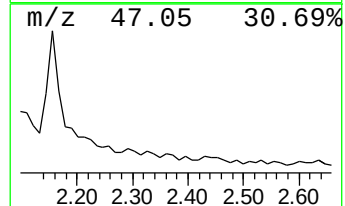
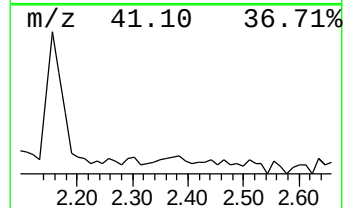
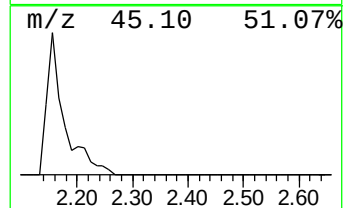
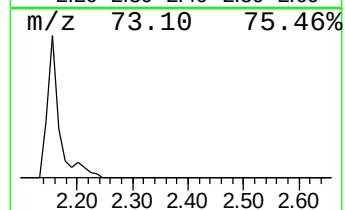
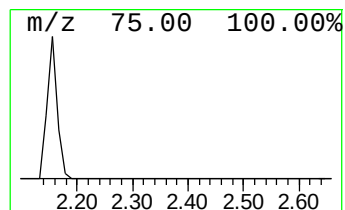
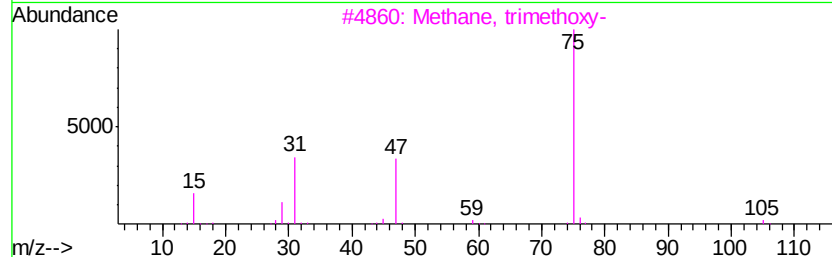
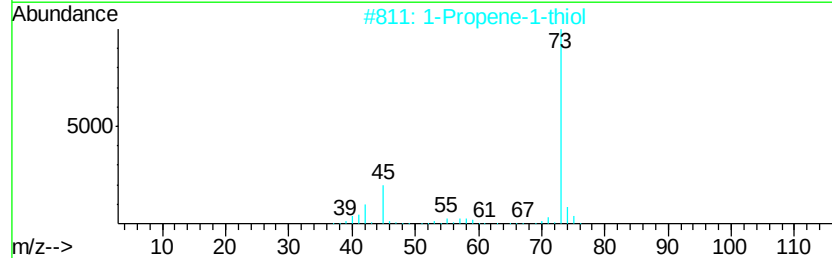
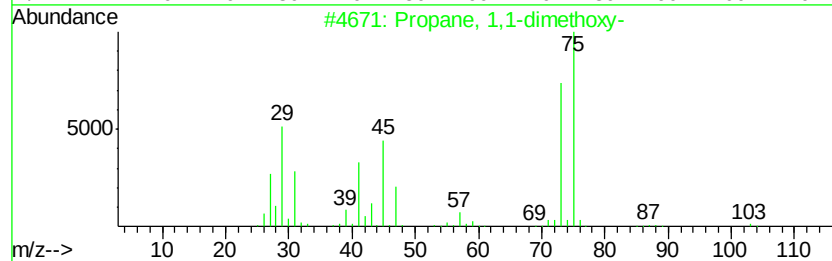
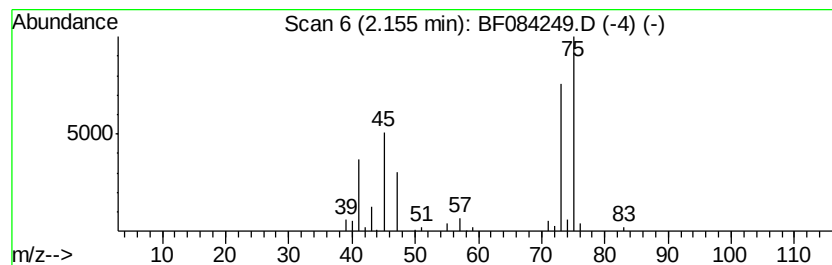
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Propane, 1,1-dimethoxy- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.16	2.11 ng	178366	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	64
2		1-Propene-1-thiol	74	C3H6S	000925-89-3	9
3		Methane, trimethoxy-	106	C4H10O3	000149-73-5	9
4		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	9
5		1,3-Dioxolane, 2-(1-methylethyl)-	116	C6H12O2	000822-83-3	9



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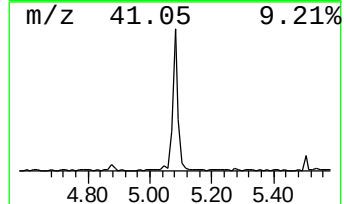
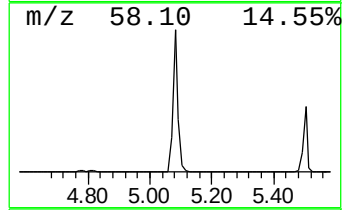
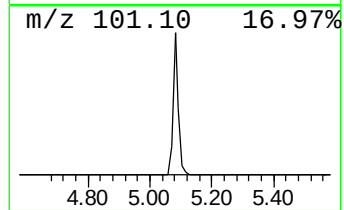
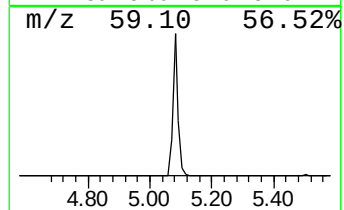
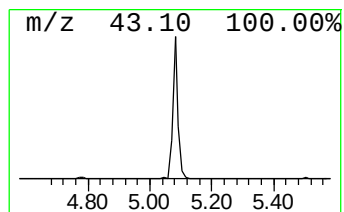
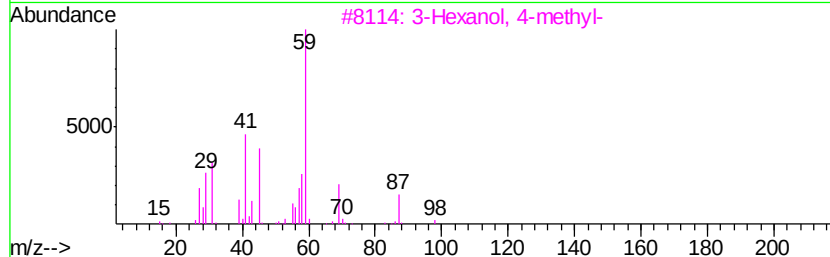
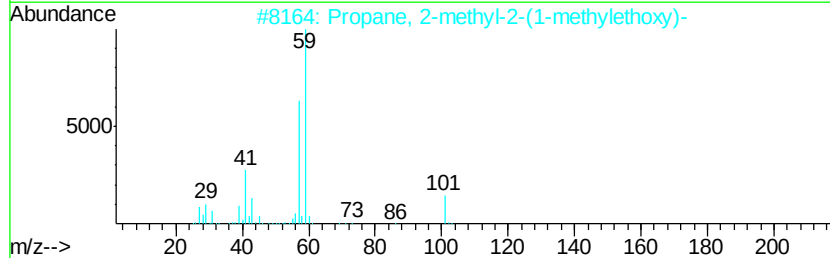
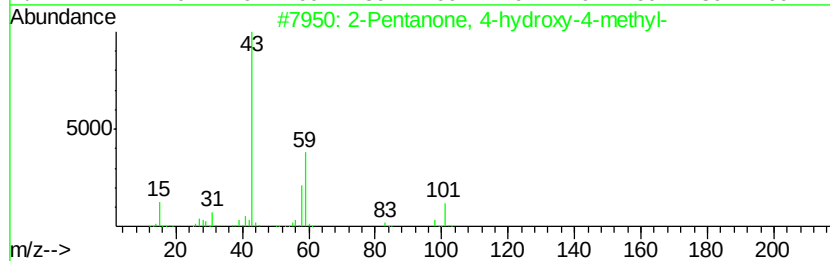
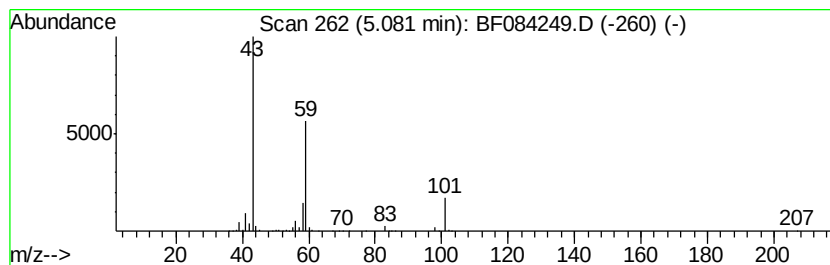
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.08	19.52 ng	1651750	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	36
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	32
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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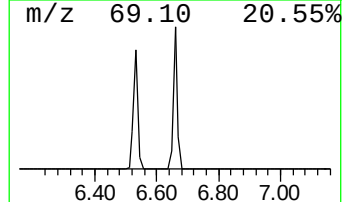
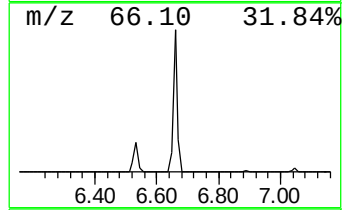
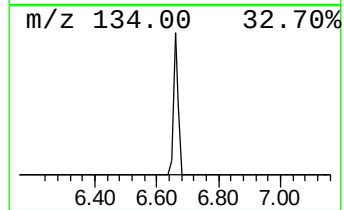
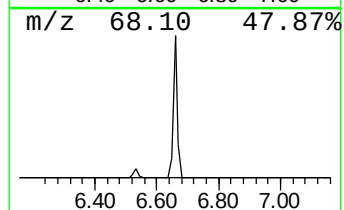
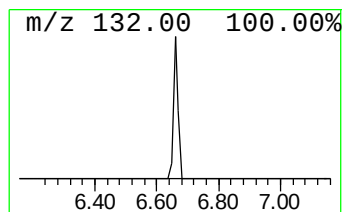
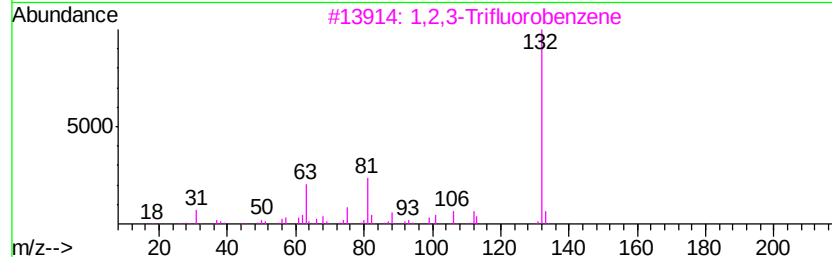
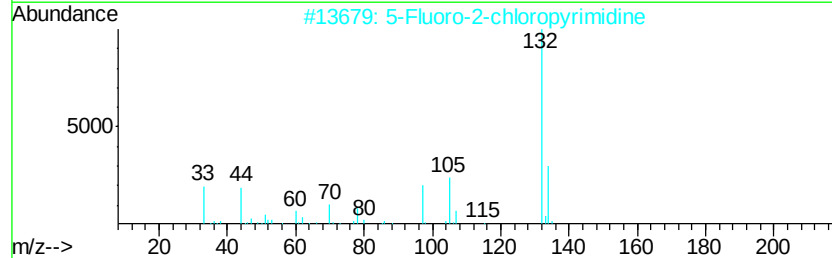
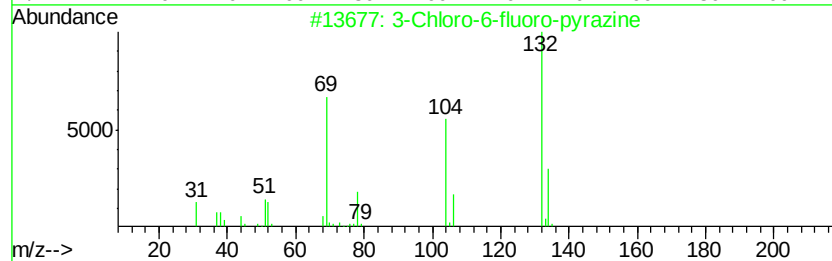
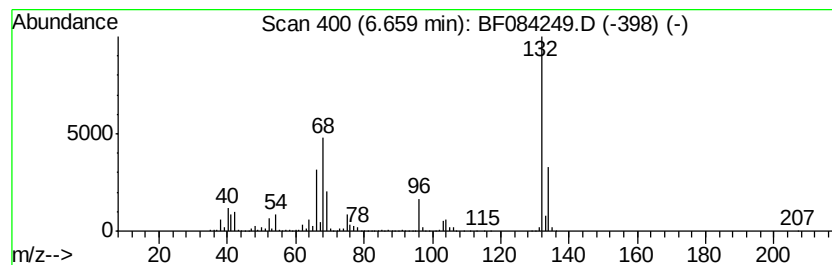
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 unknown6.66 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.66	63.09 ng	5337610	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		1,2,3-Trifluorobenzene	132	C6H3F3	001489-53-8	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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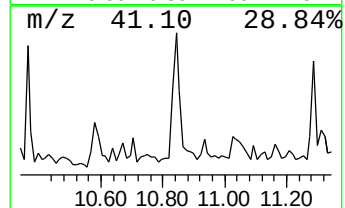
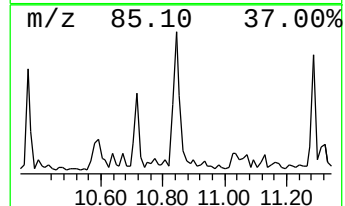
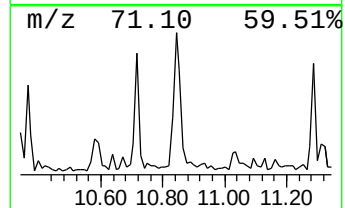
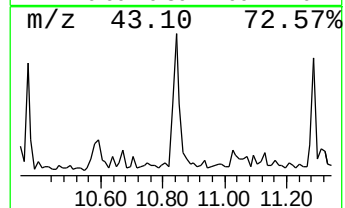
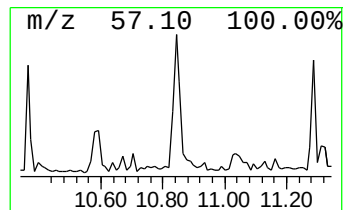
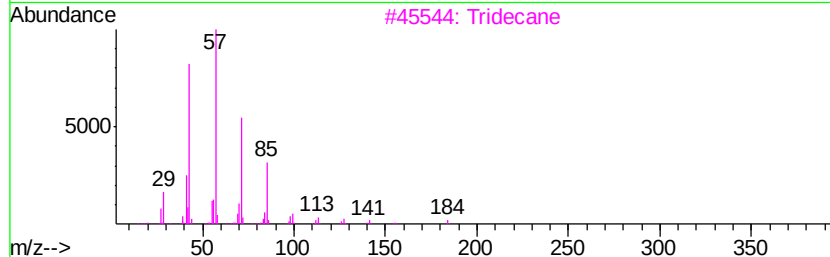
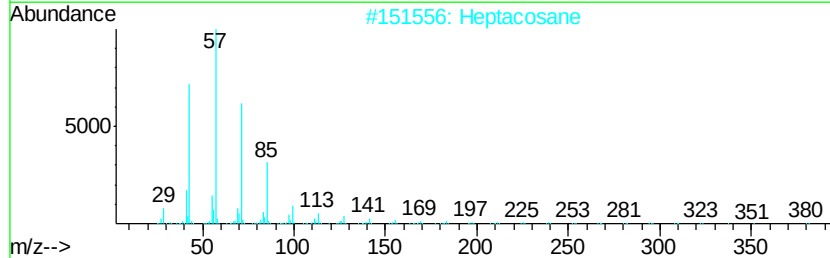
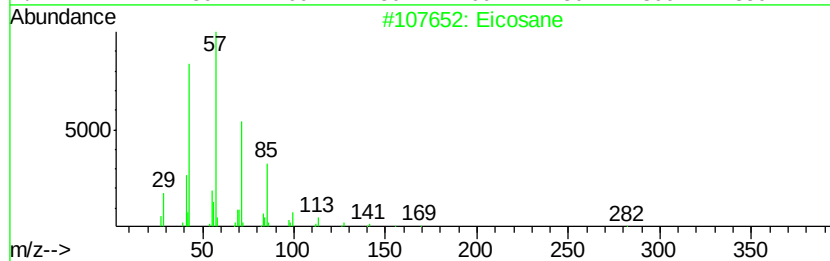
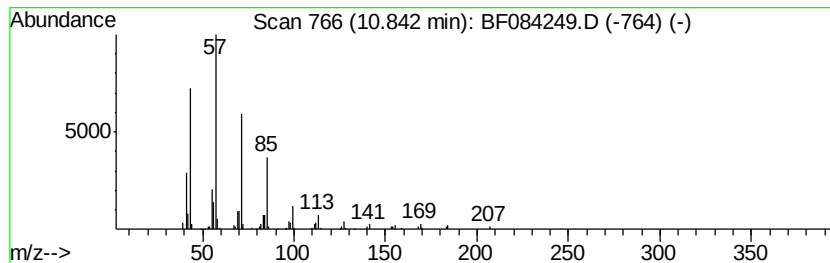
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Peak Number 5 Eicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.84	2.41 ng	261501	Phenanthrene-d10	11.41	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	91
2	Heptacosane	380	C27H56	000593-49-7	90
3	Tridecane	184	C13H28	000629-50-5	90
4	Heptadecane	240	C17H36	000629-78-7	90
5	Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	86



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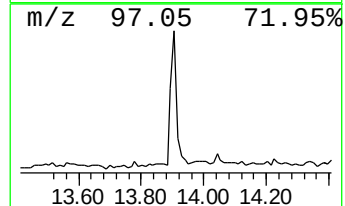
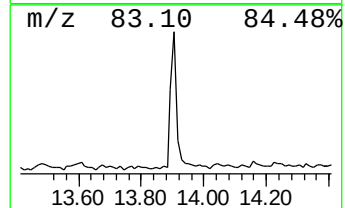
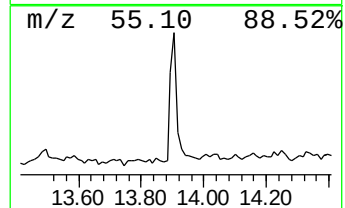
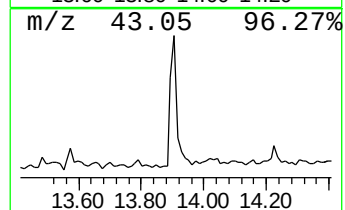
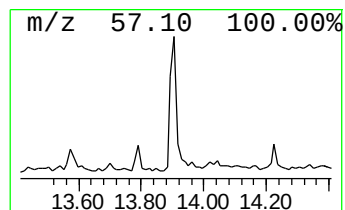
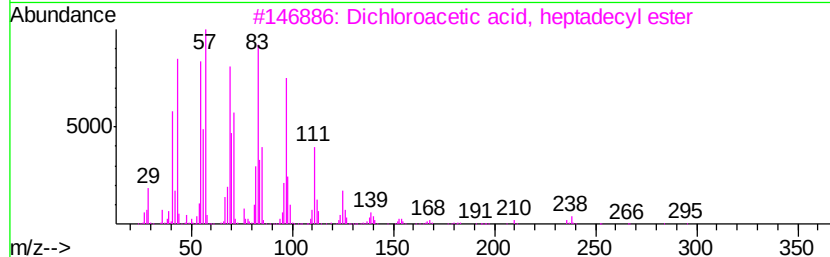
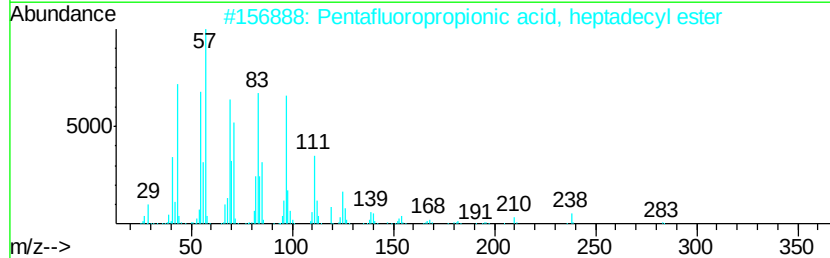
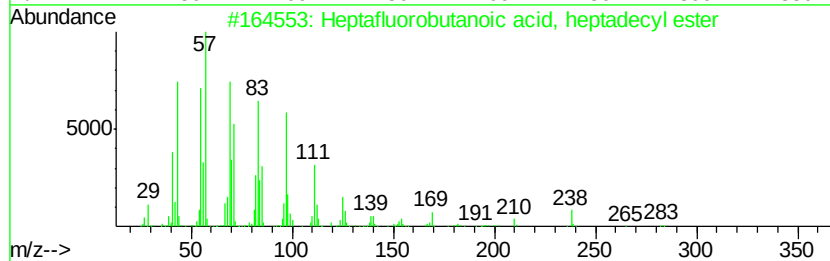
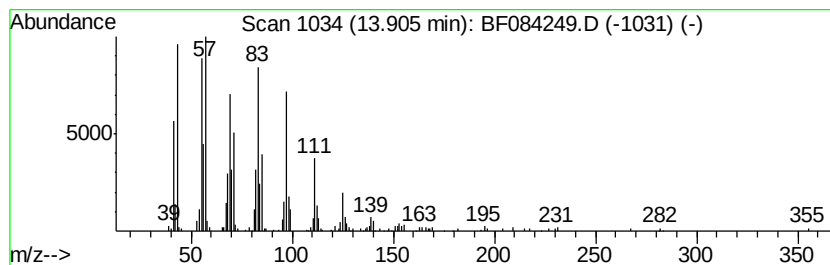
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Heptafluorobutanoic acid, h... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.91	4.74 ng	370795	Chrysene-d12	14.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	94
2		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	93
3		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	91
4		1-Tricosene	322	C23H46	018835-32-0	91
5		1-Heneicosyl formate	340	C22H44O2	077899-03-7	90



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 1,1-dime...	2.16	2.1	ng	178366	1	6.89	1692190	20.0
2-Pentanone, 4-hy...	5.08	19.5	ng	1651750	1	6.89	1692190	20.0
unknown6.66	6.66	63.1	ng	5337610	1	6.89	1692190	20.0
Eicosane	10.84	2.4	ng	261501	4	11.41	2171580	20.0
Heptafluorobutano...	13.91	4.7	ng	370795	5	14.05	1564290	20.0