

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051316\
 Data File : BF086985.D
 Acq On : 13 May 2016 22:21
 Operator : UM/SJ
 Sample : H2992-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 90THAVE-SB-26(0.5-10.0)

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:\HPCHEM1\BNA_F\METHODS\8270-BF050916.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	1.736	11	13	48	rVB	4319147	8666127	100.00%	17.554%
2	4.742	274	276	289	rBV	701565	1196019	13.80%	2.423%
3	5.199	314	316	319	rBV	2960710	4082639	47.11%	8.270%
4	5.610	350	352	362	rVB	64767	100737	1.16%	0.204%
5	6.273	407	410	413	rBV	3502602	4157455	47.97%	8.421%
6	6.387	417	420	422	rBV	3889285	4516606	52.12%	9.149%
7	6.605	437	439	442	rBV	643699	894106	10.32%	1.811%
8	6.765	451	453	456	rBV	3281458	3109927	35.89%	6.299%
9	7.187	487	490	492	rBV	2990457	3162644	36.49%	6.406%
10	7.896	550	552	555	rBV	1186188	1190728	13.74%	2.412%
11	8.982	644	647	649	rBV	4625443	4923084	56.81%	9.972%
12	9.382	679	682	684	rBV	417192	414813	4.79%	0.840%
13	9.645	702	705	708	rBV	1464642	1374559	15.86%	2.784%
14	10.091	742	744	746	rVB	131246	104845	1.21%	0.212%
15	10.434	771	774	777	rBV	2791939	2996037	34.57%	6.069%
16	10.559	783	785	792	rVB	143925	184870	2.13%	0.374%
17	11.119	832	834	837	rBV	1262845	1317660	15.20%	2.669%
18	12.708	970	973	976	rBV	3884701	4483840	51.74%	9.083%
19	13.622	1050	1053	1061	rBV	66781	163199	1.88%	0.331%
20	13.748	1061	1064	1067	rBV	1423861	1340432	15.47%	2.715%
21	15.108	1180	1183	1186	rBV	878401	987577	11.40%	2.000%

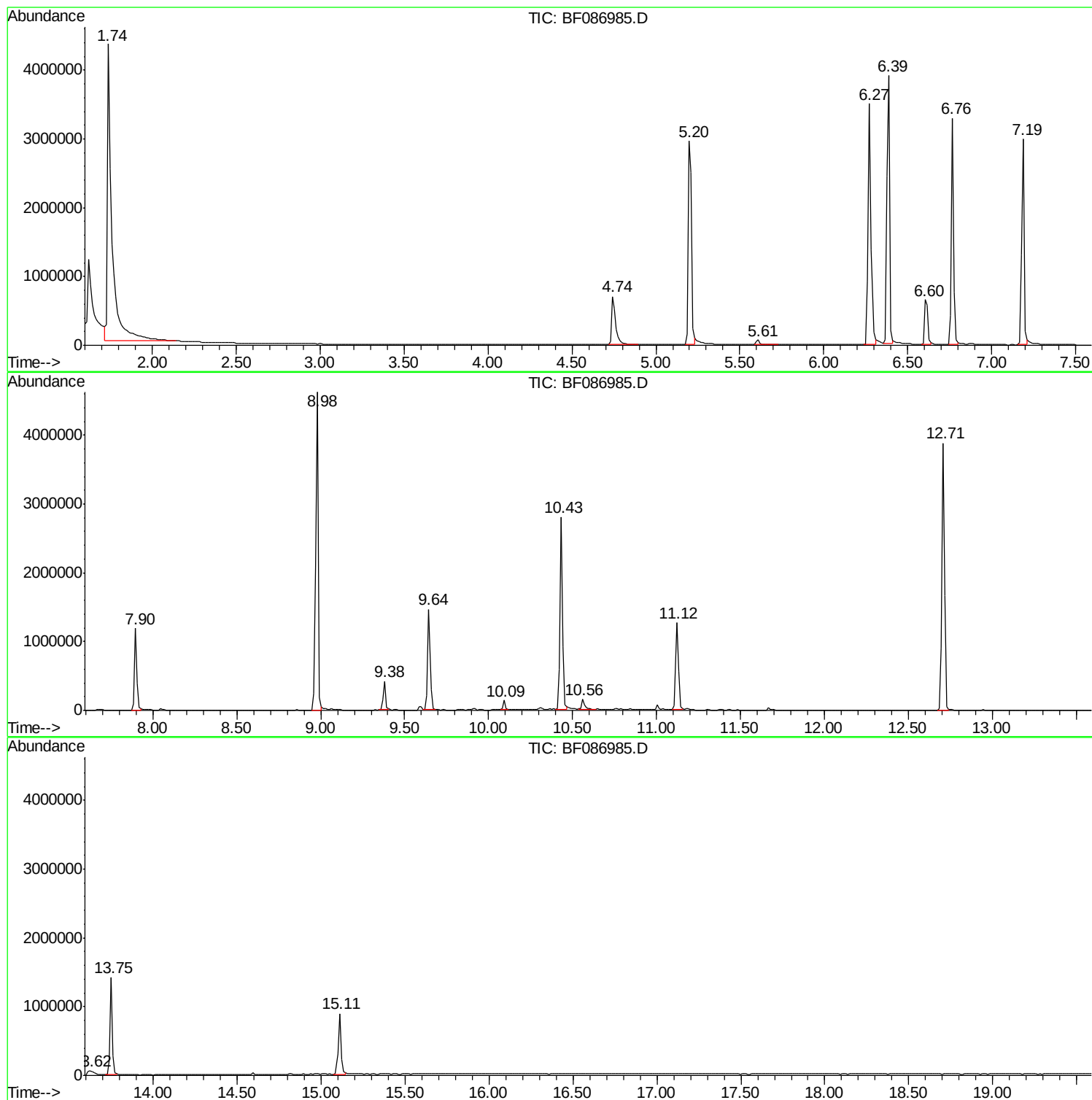
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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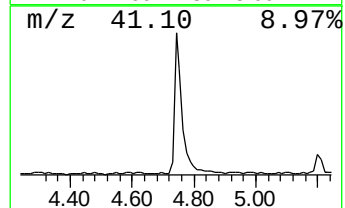
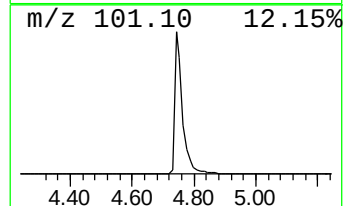
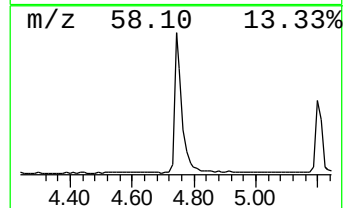
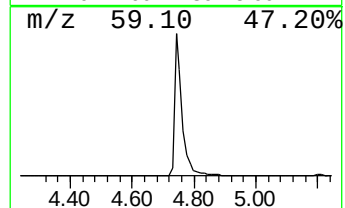
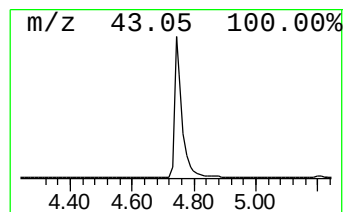
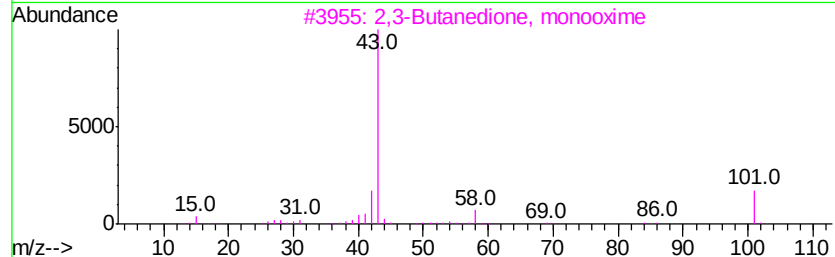
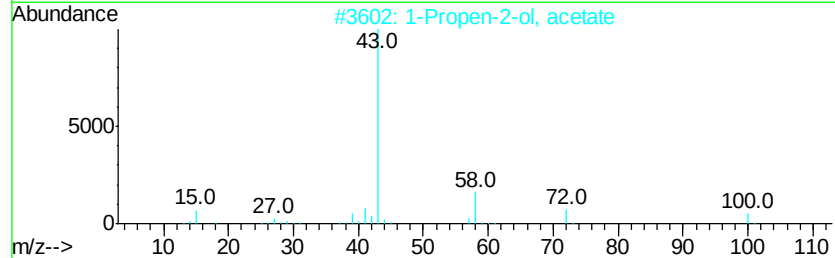
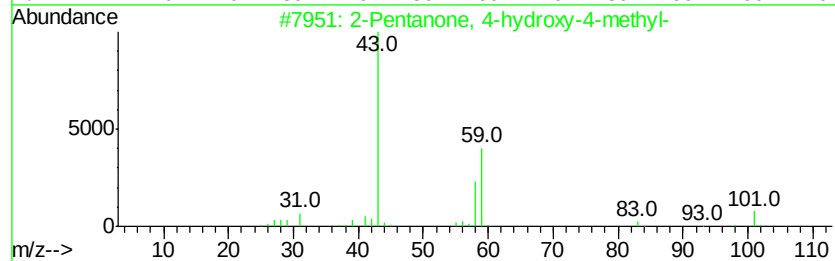
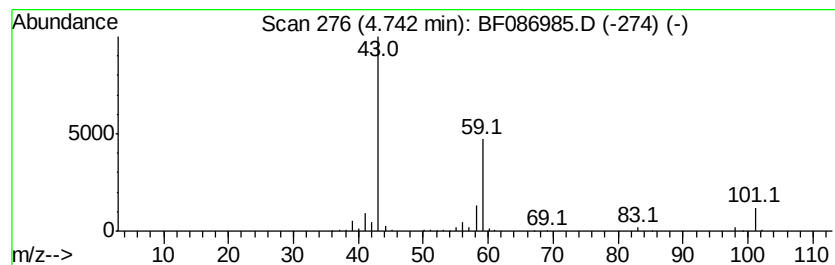
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.74	26.75 ng	1196020	1,4-Dichlorobenzene-d4	6.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		Diacetamide	101	C4H7NO2	000625-77-4	9
5		2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9



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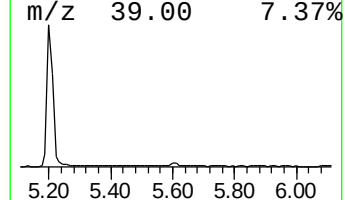
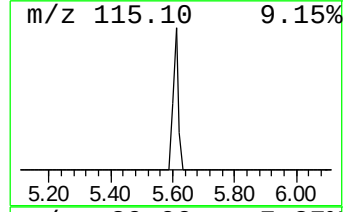
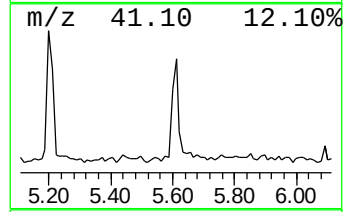
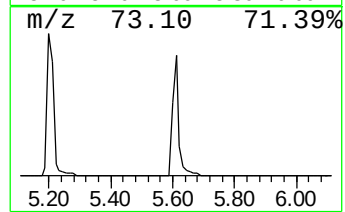
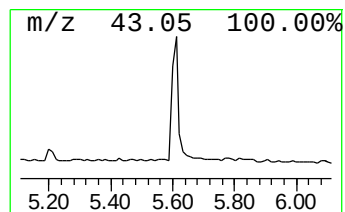
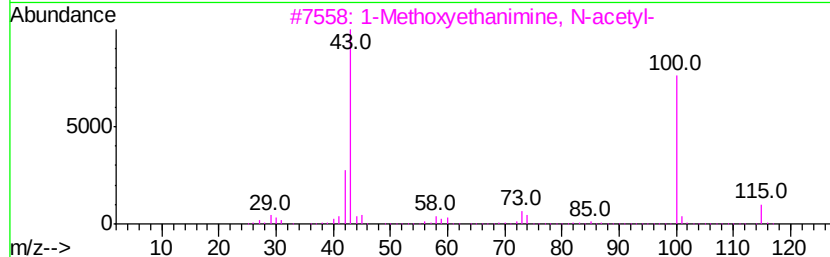
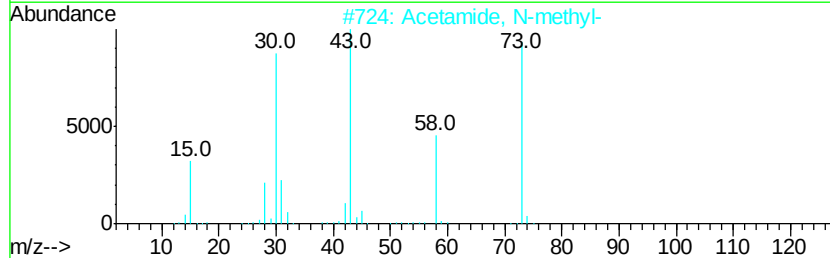
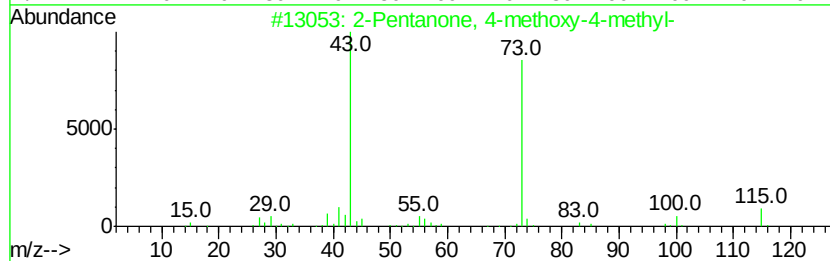
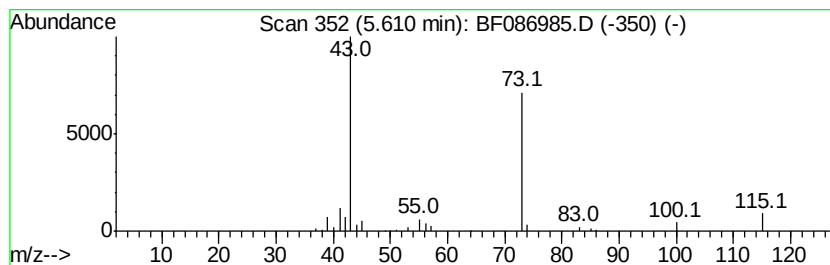
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-methoxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.61	2.25 ng	100737	1,4-Dichlorobenzene-d4	6.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	83
2		Acetamide, N-methyl-	73	C3H7NO	000079-16-3	49
3		1-Methoxyethanimine, N-acetyl-	115	C5H9NO2	099028-43-0	33
4		Pent-3-enylamine	85	C5H11N	087156-70-5	9
5		Guanidine, methyl-	73	C2H7N3	000471-29-4	9



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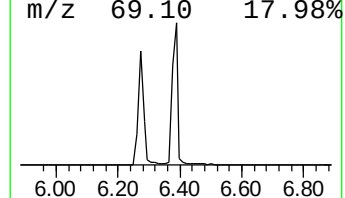
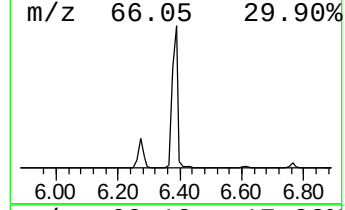
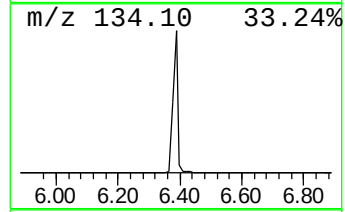
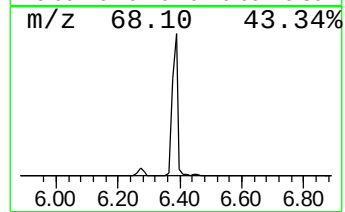
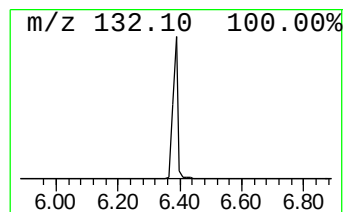
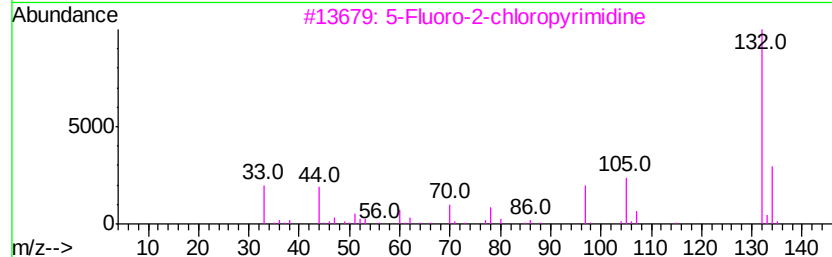
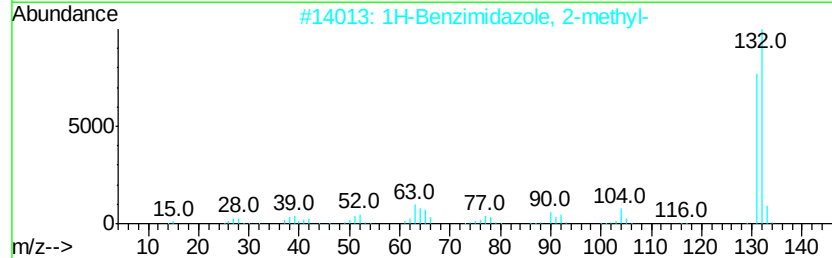
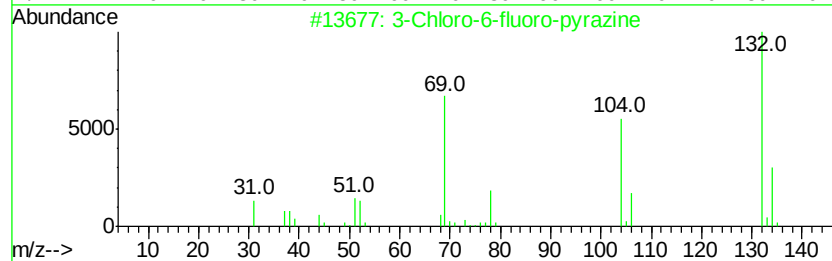
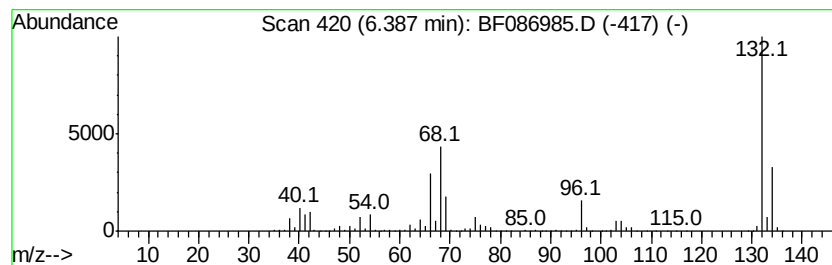
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Peak Number 4 unknown6.39 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.39	101.03 ng	4516610	1,4-Dichlorobenzene-d4	6.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3	Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	1H	Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3	5	Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4	(5	METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5	Benzene,	1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9



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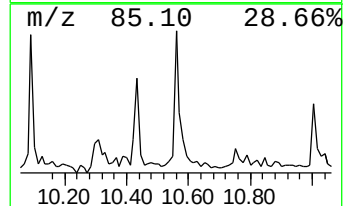
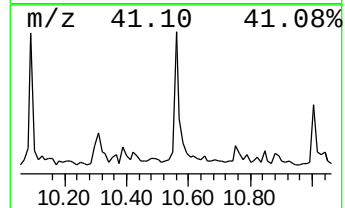
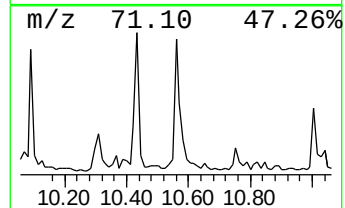
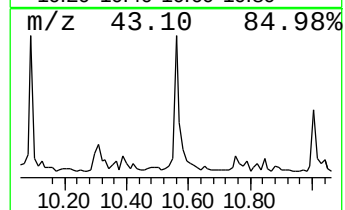
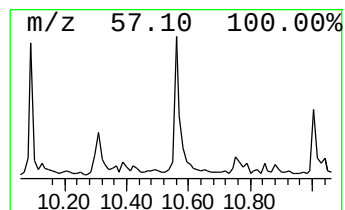
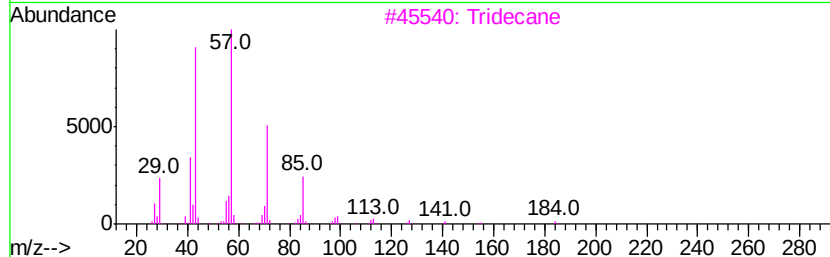
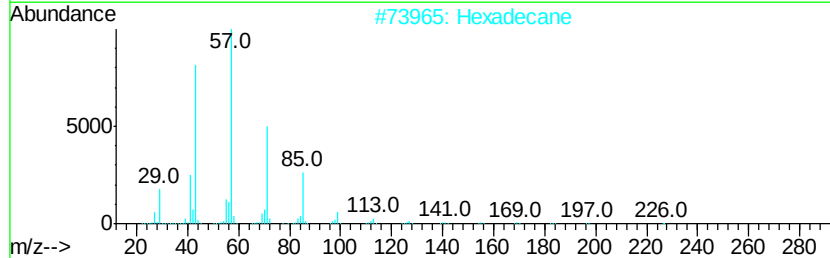
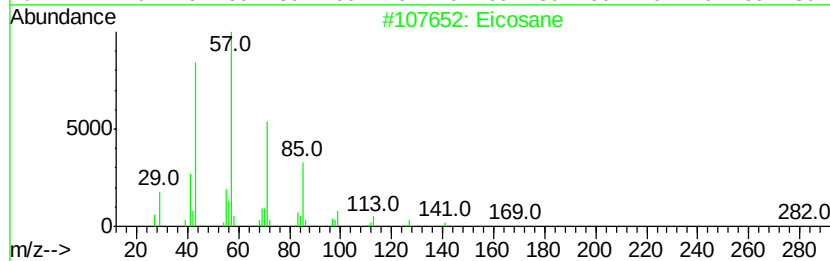
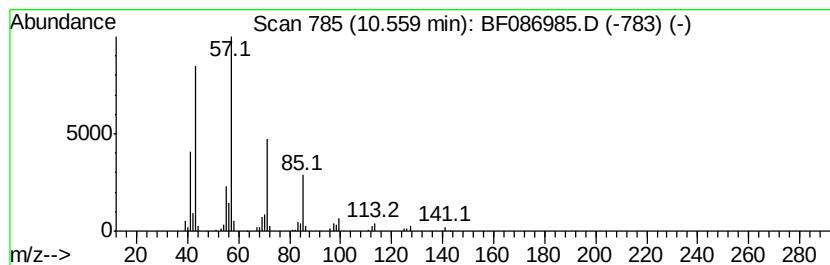
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 6 Eicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.56	2.81 ng	184870	Phenanthrene-d10	11.12

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	91
2	Hexadecane		226	C16H34	000544-76-3	91
3	Tridecane		184	C13H28	000629-50-5	90
4	Octadecane		254	C18H38	000593-45-3	86
5	Pentadecane		212	C15H32	000629-62-9	86



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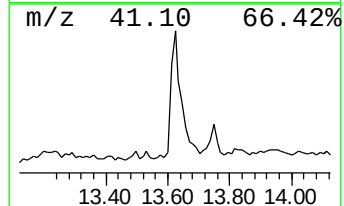
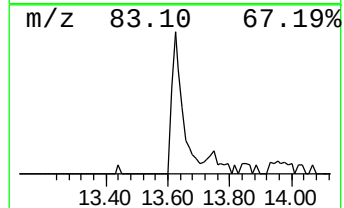
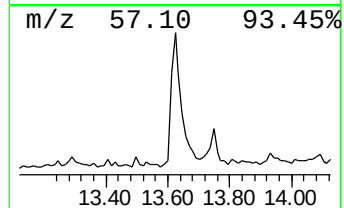
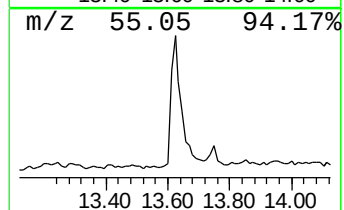
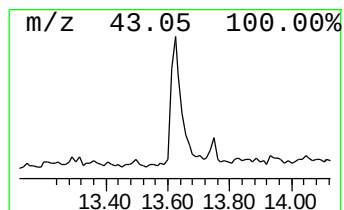
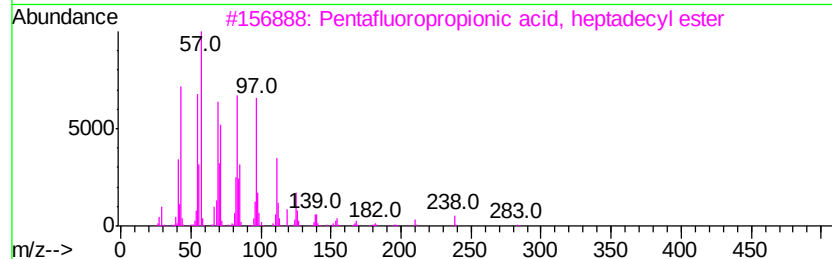
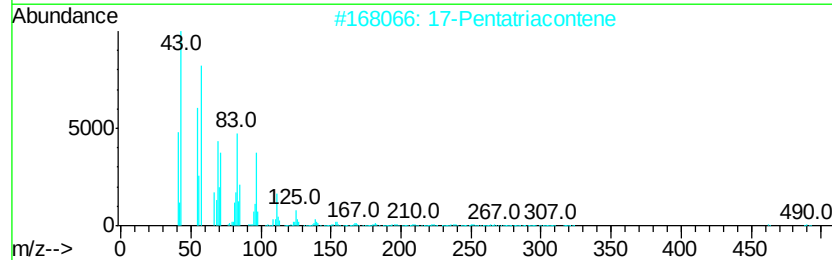
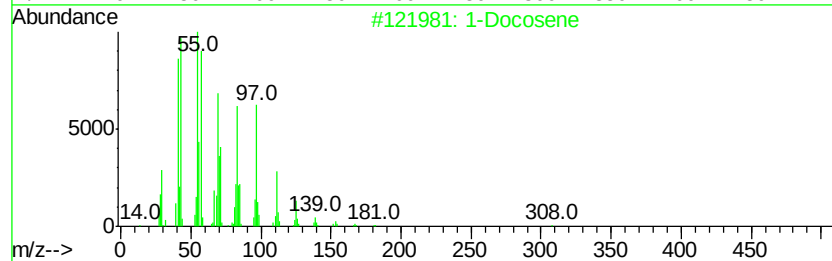
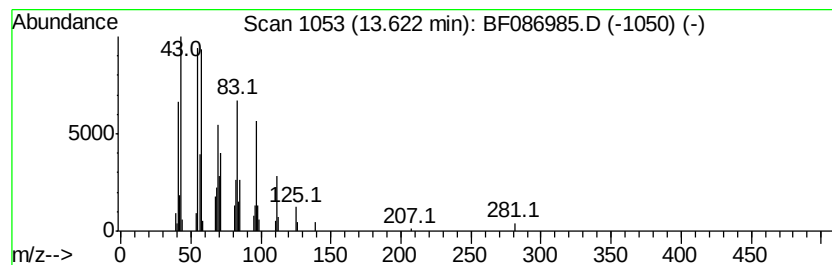
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 1-Docosene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.62	2.44 ng	163199	Chrysene-d12	13.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Docosene	308	C22H44	001599-67-3	94
2		17-Pentatriacontene	491	C35H70	006971-40-0	90
3		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	87
4		5-Eicosene, (E)-	280	C20H40	074685-30-6	86
5		1-Nonadecene	266	C19H38	018435-45-5	83



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

TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.74	26.8	ng	1196020	1	6.62	894106	20.0
2-Pentanone, 4-me...	5.61	2.3	ng	100737	1	6.62	894106	20.0
unknown6.39	6.39	101.0	ng	4516610	1	6.62	894106	20.0
Eicosane	10.56	2.8	ng	184870	4	11.12	1317660	20.0
1-Docosene	13.62	2.4	ng	163199	5	13.75	1340430	20.0