

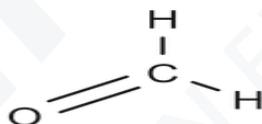
Certificate of Analysis

Formaldehyde GC Standard

Product No.: NKC0277

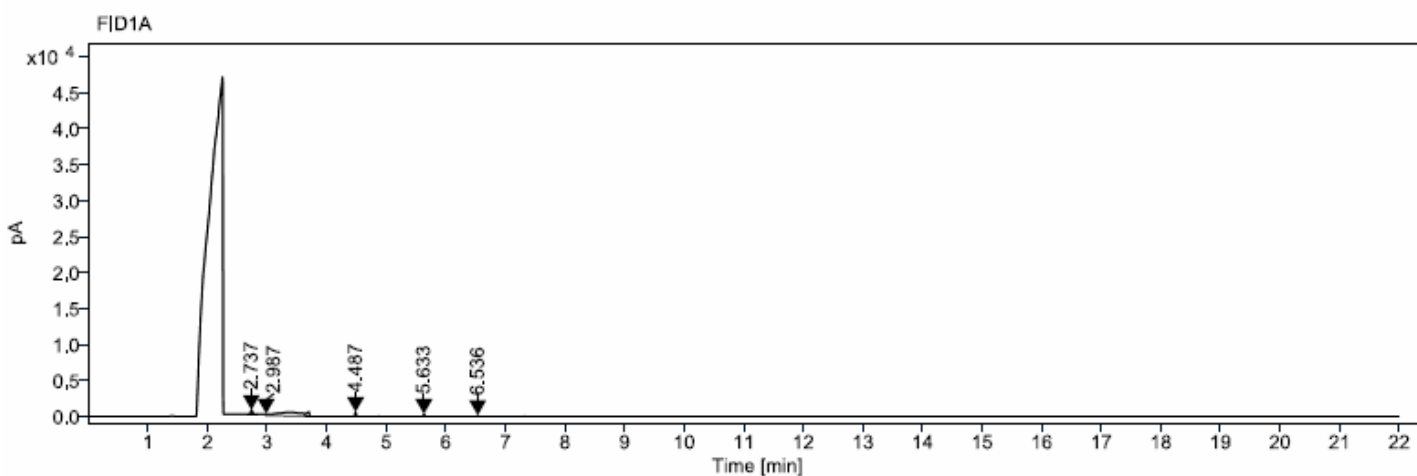
Product Information	
Cas No.:	50-00-0
Molecular Formula:	HCHO
Molecular Weight:	30.03 g/mol.
Grade:	GC Standard.
Storage:	Room Temperature
MFG Date:	Feb-2026.
EXP Date:	Feb-2031.
Batch No.:	NKCF23M067.

Test	Specification	Result
Description	A Clear, Colorless Liquid	A Clear, Colorless Liquid
Purity By GC	37-41%	38.72%
WEIGHT PER ML AT 20° C	1.090 - 1.120 g/ml	1.1004 g/ml
Identification by MASS	Conforms to Molecular mass.	Complies
Identification by IR	Conforms to structure.	Complies
Identification by 1H NMR	Conforms to structure.	Complies.



- The product complies with the prescribed standards of quality
- The product has been tested by the Quality Control Laboratory of N K Chem to the above specifications
- This is Electronic Generated Specification do not require signature

Data file:	NKCF23M067.dx	Project Name:	April-2026
Sample name:	NKCF23M067	Operator:	ADMIN (ADMIN)
Instrument ID:	GCHS-01	Injection date:	2026-04-02 14:30:37+05:30
Inj. volume (uL):	1.000		
Vial No.:	102		
Acq. method:	ALS_250.amx		
Processing method:	*GC_LC Area Percent_DefaultMethod.pmx		



Signal: FID1A

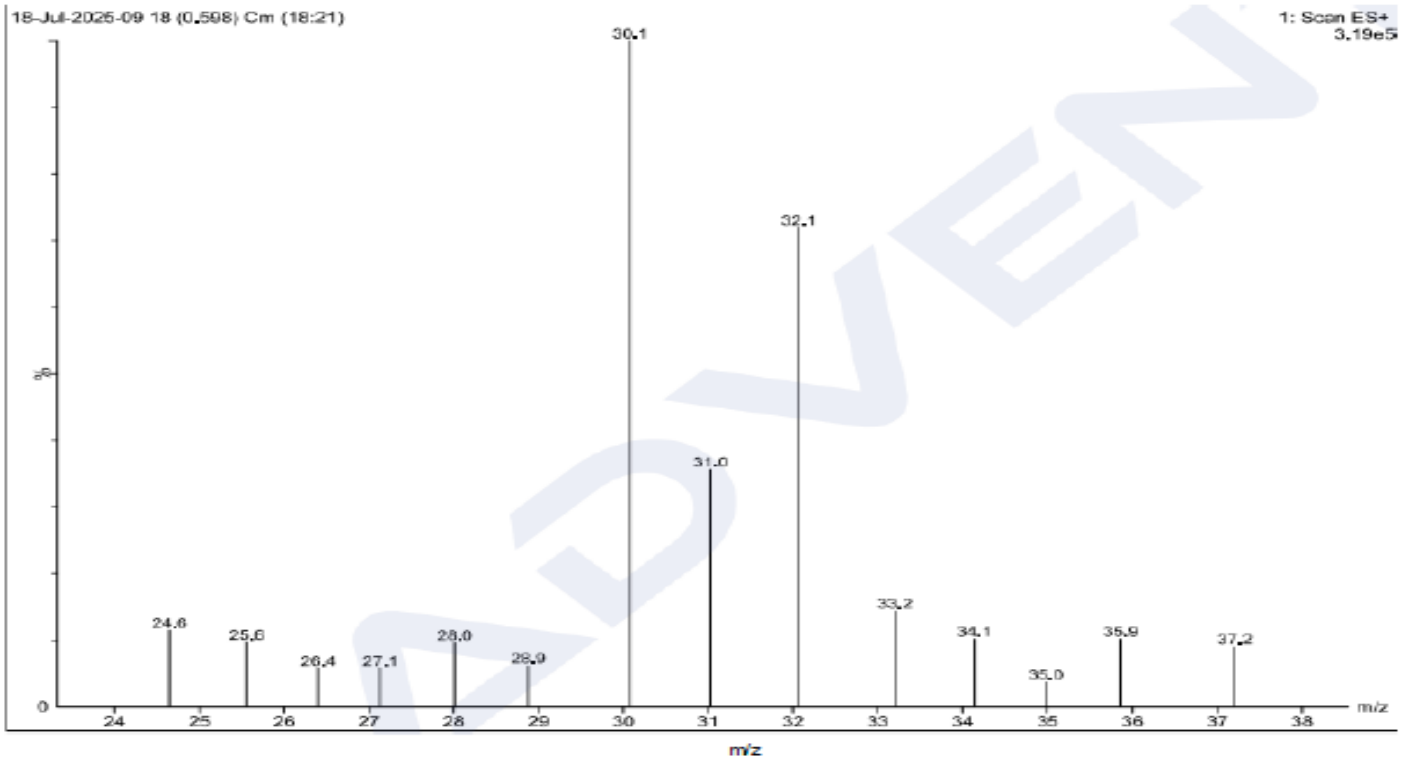
Sr.No.	RT [min]	Area	Height	Peak Area Percent
1	2,737	1191,71	604,814	6,35
2	2,987	16392,33	240,401	87,37
3	4,487	725,09	606,183	3,86
4	5,633	326,22	330,467	1,74
5	6,536	127,58	127,389	0,68

K M Pharma Solution

Empower 3
SOFTWARE

SAMPLE INFORMATION

Sample Name:	NKCF23M067	Acquired By:	System
Sample Type:	Unknown	Date Acquired:	03-04-2026 13:59:54 IST
Injection Volume:	6.00 ul	Acq. Method Set:	Mass Analysis_30_200
Run Time:	1.0 Minutes	Date Processed:	06-04-2026 12:21:39 IST,



Name: SampleName: NKCF23M067 Date Acquired: 03-04-2026 13:59:54 IST Channel Description 5: QDa Negative(-)
Scan (30.00-1200.00)Da, Centroid, CV=10

Reported by User: System
Report Method: Mass Analysis
Report Method ID: 2817
Page: 1 of 1

Project Name: 2026\April

Sr No	M/Z	Fragment
1	30.1	M+1

Conclusion: The signal of the mass and their interpretation are consistent with the structural formula.

NKCF23M067
CDC13 PROTON



7.285
5.037
5.010
4.990
4.983
4.979
4.970
4.967
4.955
4.942
4.930
4.925
4.918
4.912
4.905
4.886
4.813
4.777
4.766
4.759
4.741
4.593
3.496
3.452
3.433
3.423
3.419
3.418
3.412
3.378
3.080
1.897

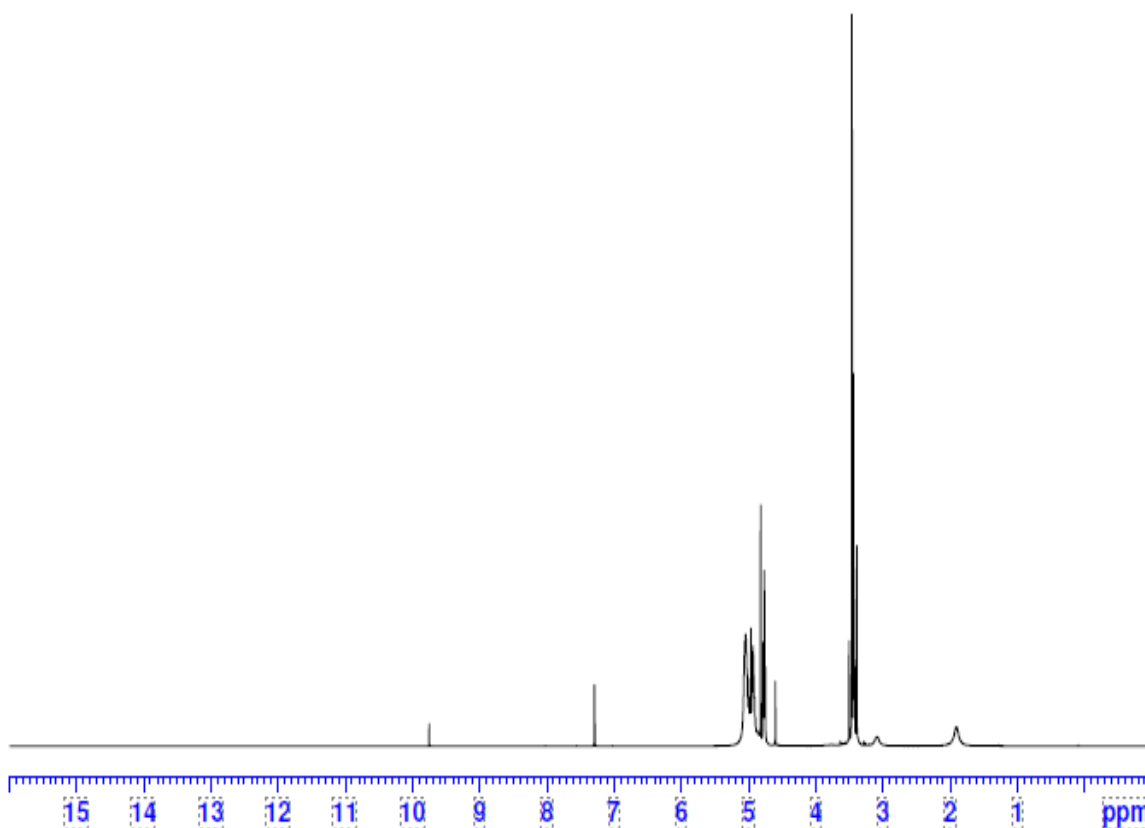
Current Data Parameters
NAME 117-K.M.PHARMA
EXPNO 1
PROCNO 1

F2 - Acquisition Parameters

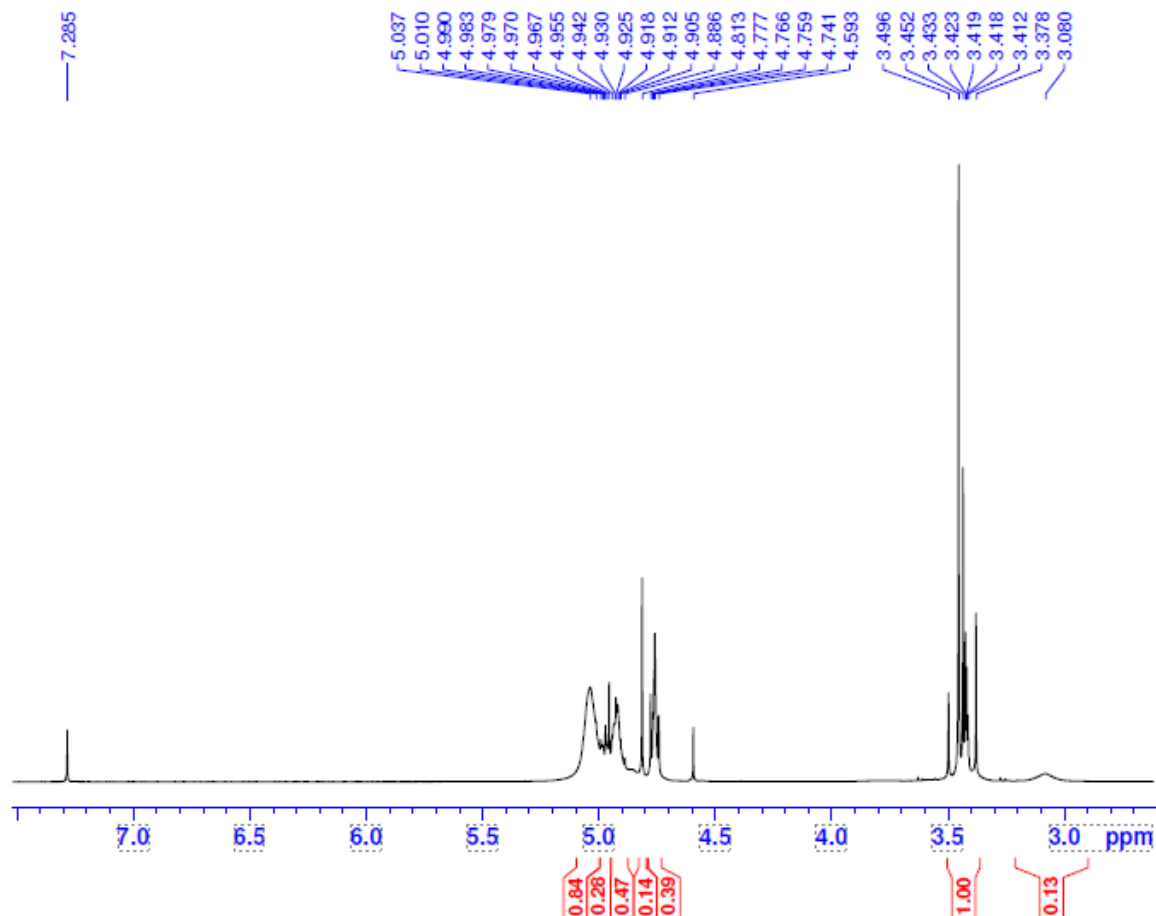
Date_ 20260402
Time 18.03 M
INSTRUM Avance
PROBHD Z166552_0024 (zq30)
PULPROG zg30
TD 65536
SOLVENT CDC13
NS 24
DS 0
SWH 10000.000 Hz
FIDRES 0.305176 Hz
AQ 3.2767999 sec
RG 12
DW 50.000 usec
DE 11.14 usec
TE 296.6 K
D1 1.00000000 sec
ID0 1
SF01 400.1336012 MHz
NUC1 1H
P0 2.67 usec
P1 8.00 usec
PIW1 22.37700081 W

F2 - Processing parameters

SI 65536
SF 400.1300000 MHz
WDW EM
SSE 0
LB 0.30 Hz
GB 0
PC 1.00



0.84
0.28
0.47
0.14
0.39
1.00
0.13

NKCF23M067
 CDC13 PROTON

 Current Data Parameters
 NAME 117-K.M.PHARMA
 EXPNO 1
 PROCNO 1

 F2 - Acquisition Parameters
 Date 20260402
 Time 18.03 H
 INSTRUM Avance
 PROBD 2166552_0024
 PULPROG zg30
 TD 65536
 SOLVENT CDC13
 NS 24
 DS 0
 SWH 10000.000 Hz
 FIDRES 0.305176 Hz
 AQ 1.2767999 sec
 RG 127
 DW 50.000 usec
 DE 11.14 usec
 IE 296.6 K
 DJ 1.00000000 sec
 TD0 1
 SFO1 400.1336012 MHz
 NUC1 1H
 PO 2.67 usec
 P1 8.00 usec
 PLW1 22.37700081 W

 F2 - Processing parameters
 SI 65536
 SF 400.1300000 MHz
 WDW EM
 SSB 0
 LB 0.30 Hz
 GB 0
 PC 1.00

Proton Interpretation Table

δ (ppm)	Multiplicity	Integration (approx)	Assignment
7.26	s	—	Residual CHCl_3 (solvent)
5.03– 4.75	m	~3–4H	–CH / =CH / O–CH protons (possibly attached to O or unsaturated system)
4.60– 4.55	m	~1–2H	O–CH / CH–X (X = O, N)
3.50– 3.35	strong m	~2–3H	–CH ₂ –O / –CH ₃ near electronegative group
~1.8–1.2	weak	~1–2H	Aliphatic CH ₂ /CH ₃ (minor impurity or tail)

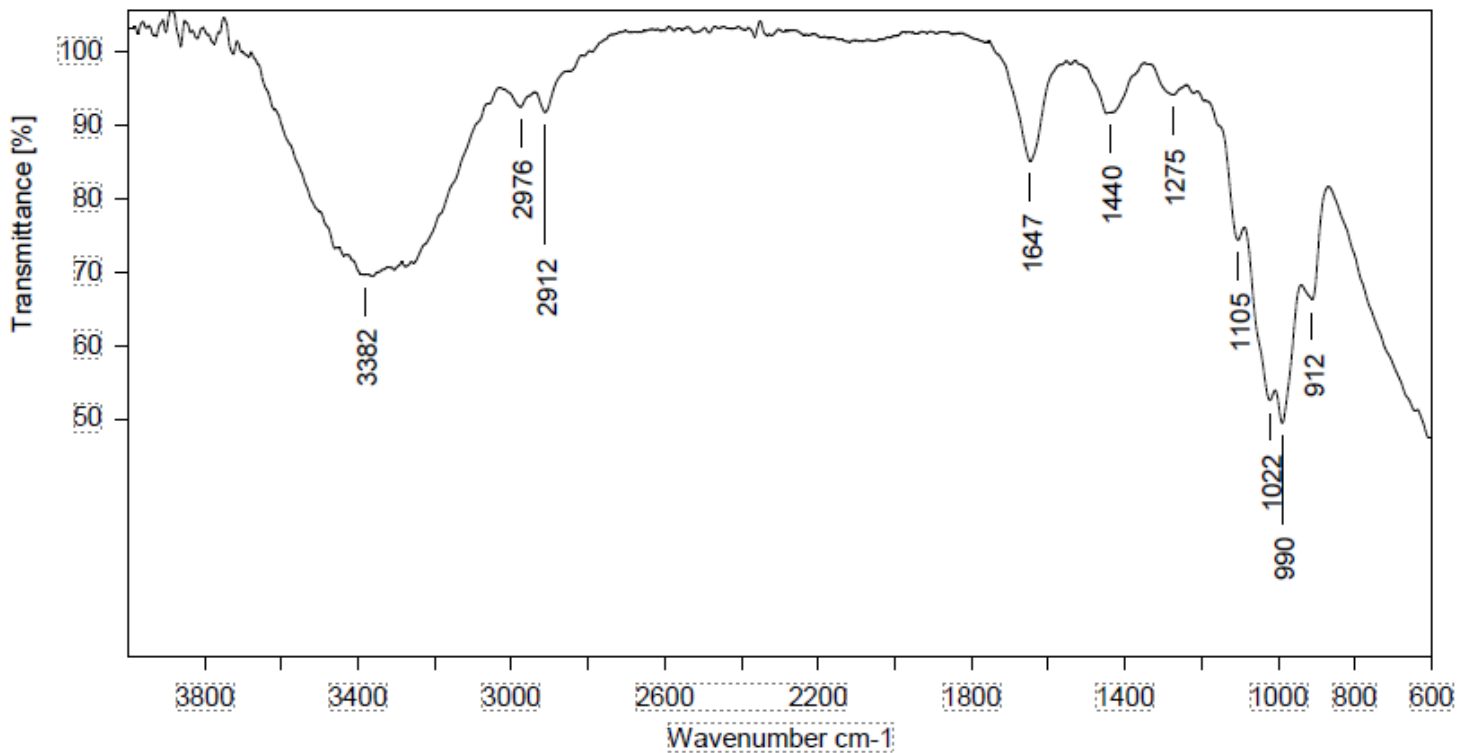
Chemical Formula : HCHO

Conclusion: The structure is confirmed with the signals of spectrum and their interpretation

K.M.Pharma Solution

Batch No.: NKCF23M067

Measurement Date & Time : 03-04-2026 09:41:03



Wavenumber	Abs. intensity	Rel. intensity	Width	Found if threshold <	Shoulder
3382.3493	0.697	0.007	310.8883	2.703315	0
2912.0024	0.918	0.049	689.5941	58.980331	0
2975.6625	0.926	0.023	32.0296	60.567791	0
1646.7835	0.852	0.145	67.2741	92.746864	0
1439.7968	0.918	0.002	22.5778	0.205588	0
1275.1634	0.942	0.023	1759.0049	31.573101	0
989.7926	0.495	0.332	112.3016	171.084076	0
1022.1487	0.526	0.016	15.2539	5.063652	0
911.6151	0.663	0.074	186.0024	6.452155	0
1105.2357	0.744	0.021	18.1022	5.009481	0

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