



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 03:41 PM GMT

PDB ID : 1BHT
Title : NK1 FRAGMENT OF HUMAN HEPATOCYTE GROWTH FACTOR
Authors : Ultsch, M.H.; Lokker, N.A.; Godowski, P.J.; De Vos, A.M.
Deposited on : 1998-06-10
Resolution : 2.00 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

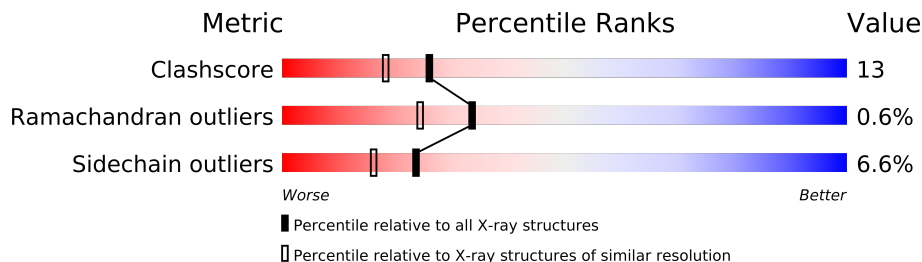
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	176	
1	B	176	

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 3133 atoms, of which 0 are hydrogen and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called HEPATOCYTE GROWTH FACTOR.

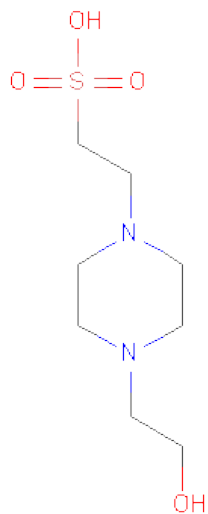
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	176	Total	C	N	O	S	22	0	0
			1421	895	255	259	12			
1	B	174	Total	C	N	O	S	28	0	0
			1401	884	250	255	12			

- Molecule 2 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 4-(2-HYDROXYETHYL)-1-PIPERAZINEETHANESULFONIC ACID (three-letter code: EPE) (formula: C₈H₁₈N₂O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
3	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 4 is water.

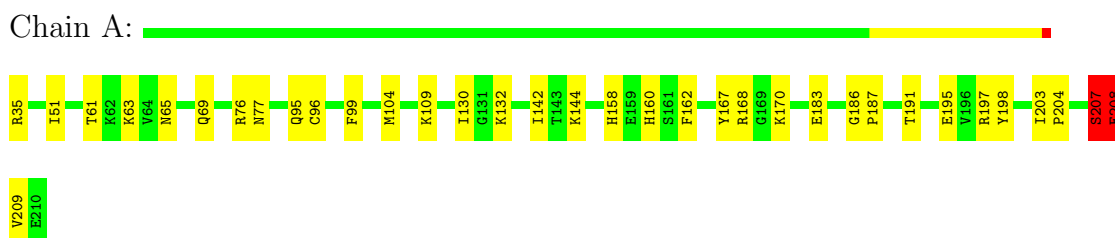
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	139	Total	O	0	0
			139	139		
4	B	127	Total	O	0	0
			127	127		

3 Residue-property plots

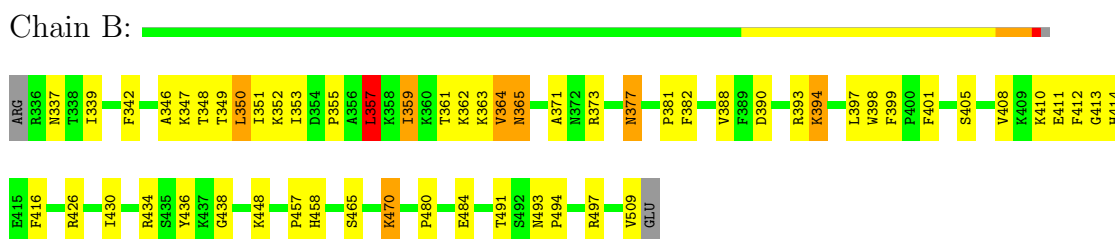
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

• Molecule 1: HEPATOCYTE GROWTH FACTOR



• Molecule 1: HEPATOCYTE GROWTH FACTOR



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	88.32Å 88.32Å 117.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.00	Depositor
% Data completeness (in resolution range)	94.0 (10.00-2.00)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	0.06	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.196 , 0.247	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3133	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: EPE, SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.69	1/1456 (0.1%)	0.94	8/1956 (0.4%)
1	B	0.60	0/1436	0.77	1/1930 (0.1%)
All	All	0.65	1/2892 (0.0%)	0.86	9/3886 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	207	SER	N-CA	5.06	1.56	1.46

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	207	SER	N-CA-C	-10.65	82.23	111.00
1	A	208	GLU	CB-CA-C	-8.71	92.98	110.40
1	A	207	SER	CA-C-N	-7.00	101.80	117.20
1	A	208	GLU	CG-CD-OE2	-6.75	104.81	118.30
1	A	208	GLU	CA-CB-CG	6.27	127.20	113.40
1	A	207	SER	C-N-CA	6.16	137.10	121.70
1	A	208	GLU	CG-CD-OE1	5.77	129.84	118.30
1	B	357	LEU	CA-CB-CG	5.51	127.97	115.30
1	A	99	PHE	CB-CA-C	-5.09	100.21	110.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	207	SER	Mainchain
1	A	208	GLU	Mainchain

5.2 Close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1421	0	1396	26	0
1	B	1401	0	1377	49	0
2	A	10	0	0	0	0
2	B	5	0	0	0	0
3	A	15	0	18	2	0
3	B	15	0	18	0	0
4	A	139	0	0	7	0
4	B	127	0	0	3	0
All	All	3133	0	2809	74	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 13.

All (74) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:208:GLU:HG3	4:A:861:HOH:O	1.30	1.27
1:A:208:GLU:CG	4:A:861:HOH:O	1.79	1.24
1:A:208:GLU:OE2	4:A:861:HOH:O	1.61	1.14
1:B:401:PHE:CE1	1:B:408:VAL:HG11	2.21	0.75
1:B:362:LYS:HB3	1:B:398:TRP:HZ3	1.55	0.71
1:B:362:LYS:HB3	1:B:398:TRP:CZ3	2.26	0.71
1:A:183:GLU:OE1	3:A:1:EPE:H51	1.91	0.71
1:A:63:LYS:HE2	1:A:95:GLN:NE2	2.07	0.70
1:A:160:HIS:HD2	1:A:162:PHE:H	1.40	0.68
1:B:352:LYS:HD2	1:B:357:LEU:O	1.95	0.67
1:A:167:TYR:O	1:A:170:LYS:HB2	1.96	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:352:LYS:NZ	1:B:355:PRO:HA	2.14	0.63
1:B:388:VAL:HG23	1:B:399:PHE:HE2	1.63	0.61
1:B:470:LYS:HG2	1:B:480:PRO:HA	1.81	0.61
1:B:352:LYS:HB2	1:B:359:ILE:HD11	1.82	0.60
1:B:401:PHE:CD1	1:B:408:VAL:HG11	2.38	0.58
1:A:35:ARG:HH22	1:A:69:GLN:HG2	1.70	0.56
1:B:348:THR:HG21	1:B:410:LYS:HE2	1.88	0.56
1:B:347:LYS:N	1:B:347:LYS:HD2	2.19	0.56
1:B:353:ILE:HG23	4:B:870:HOH:O	2.05	0.56
1:B:426:ARG:HB2	4:B:692:HOH:O	2.05	0.55
1:B:351:ILE:HD11	1:B:411:GLU:HG2	1.90	0.54
1:A:35:ARG:NH2	1:A:69:GLN:HG2	2.22	0.54
1:B:339:ILE:O	1:B:342:PHE:HB2	2.09	0.53
1:A:187:PRO:HB2	1:A:203:ILE:HD12	1.89	0.53
1:A:35:ARG:HH22	1:A:69:GLN:HA	1.75	0.52
1:A:158:HIS:CD2	1:A:197:ARG:HA	2.45	0.52
1:B:352:LYS:HZ2	1:B:355:PRO:HA	1.75	0.52
1:B:347:LYS:HG3	1:B:414:HIS:HA	1.92	0.52
1:B:365:ASN:HD22	1:B:394:LYS:HE3	1.75	0.51
1:A:158:HIS:NE2	1:A:197:ARG:HA	2.25	0.51
1:A:130:ILE:HD12	4:A:852:HOH:O	2.10	0.50
1:B:348:THR:HG22	1:B:349:THR:H	1.77	0.49
1:A:160:HIS:CD2	1:A:162:PHE:H	2.26	0.49
1:B:388:VAL:HG23	1:B:399:PHE:CE2	2.47	0.49
1:B:339:ILE:HG23	1:B:371:ALA:HB2	1.93	0.49
1:B:349:THR:O	1:B:410:LYS:HA	2.13	0.48
1:B:412:PHE:CD1	1:B:413:GLY:N	2.82	0.48
1:B:436:TYR:CZ	1:B:438:GLY:HA3	2.49	0.47
1:B:346:ALA:HB1	1:B:347:LYS:HD2	1.96	0.47
1:B:457:PRO:HG2	1:B:458:HIS:CD2	2.51	0.46
1:B:426:ARG:HH11	1:B:430:ILE:HG13	1.80	0.45
1:B:363:LYS:O	1:B:364:VAL:HG23	2.16	0.45
1:B:458:HIS:CE1	1:B:497:ARG:HA	2.51	0.45
1:B:350:LEU:C	1:B:351:ILE:HD12	2.37	0.45
1:B:448:LYS:HE3	4:B:660:HOH:O	2.16	0.44
1:A:208:GLU:CD	4:A:861:HOH:O	2.02	0.44
1:B:401:PHE:HB2	1:B:405:SER:OG	2.19	0.43
1:B:348:THR:HG22	1:B:349:THR:N	2.33	0.43
1:A:132:LYS:O	1:A:186:GLY:HA2	2.18	0.42
1:A:198:TYR:CD2	3:A:1:EPE:H22	2.54	0.42
1:B:348:THR:HA	1:B:411:GLU:O	2.19	0.42
1:B:426:ARG:NH1	1:B:430:ILE:HG13	2.35	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:493:ASN:HA	1:B:494:PRO:HD3	1.95	0.42
1:B:361:THR:HG22	1:B:397:LEU:HG	2.02	0.41
1:B:352:LYS:HZ3	1:B:355:PRO:HA	1.85	0.41
1:B:381:PRO:HG2	1:B:382:PHE:CD2	2.55	0.41
1:A:203:ILE:HA	1:A:204:PRO:HD2	1.89	0.41
1:B:373:ARG:HD2	1:B:398:TRP:CZ2	2.55	0.41
1:B:352:LYS:HG3	1:B:357:LEU:CD2	2.51	0.41
1:A:65:ASN:HB2	1:A:69:GLN:OE1	2.20	0.41
1:A:158:HIS:HD2	4:A:645:HOH:O	2.04	0.41
1:A:142:ILE:HD13	1:B:377:ASN:HB2	2.01	0.41
1:B:390:ASP:OD2	1:B:393:ARG:HB2	2.21	0.41
1:A:35:ARG:NH2	1:A:69:GLN:HA	2.34	0.41
1:B:347:LYS:O	1:B:412:PHE:HA	2.21	0.41
1:B:373:ARG:HD2	1:B:398:TRP:CH2	2.56	0.40
1:A:144:LYS:HE2	4:A:893:HOH:O	2.21	0.40
1:B:348:THR:O	1:B:416:PHE:HB3	2.21	0.40
1:B:414:HIS:CD2	1:B:414:HIS:H	2.37	0.40
1:A:51:ILE:HG12	1:A:109:LYS:O	2.21	0.40
1:A:61:THR:HA	1:A:96:CYS:O	2.22	0.40
1:B:347:LYS:CG	1:B:414:HIS:HA	2.50	0.40
1:B:394:LYS:HB3	1:B:394:LYS:HE2	1.78	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	174/176 (99%)	163 (94%)	10 (6%)	1 (1%)	33	24
1	B	172/176 (98%)	153 (89%)	18 (10%)	1 (1%)	33	24
All	All	346/352 (98%)	316 (91%)	28 (8%)	2 (1%)	33	24

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	208	GLU
1	B	359	ILE

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	159/159 (100%)	151 (95%)	8 (5%)	34	27
1	B	157/159 (99%)	144 (92%)	13 (8%)	16	10
All	All	316/318 (99%)	295 (93%)	21 (7%)	24	16

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	ARG
1	A	77	ASN
1	A	104	MET
1	A	168	ARG
1	A	191	THR
1	A	195	GLU
1	A	207	SER
1	A	209	VAL
1	B	337	ASN
1	B	350	LEU
1	B	357	LEU
1	B	364	VAL
1	B	365	ASN
1	B	377	ASN
1	B	394	LYS
1	B	434	ARG
1	B	465	SER
1	B	470	LYS
1	B	484	GLU
1	B	491	THR
1	B	509	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	ASN
1	A	72	ASN
1	A	158	HIS
1	A	160	HIS
1	B	365	ASN
1	B	414	HIS

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

5 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	EPE	A	1	-	15,15,15	1.88	5 (33%)	20,20,20	1.52	3 (15%)
2	SO4	A	5	-	4,4,4	1.40	1 (25%)	6,6,6	0.74	0
2	SO4	A	7	-	4,4,4	1.36	0	6,6,6	0.49	0
3	EPE	B	2	-	15,15,15	2.02	5 (33%)	20,20,20	1.18	1 (5%)
2	SO4	B	6	-	4,4,4	1.20	0	6,6,6	0.78	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical

component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EPE	A	1	-	-	0/9/19/19	0/1/1/1
2	SO4	A	5	-	-	0/0/0/0	0/0/0/0
2	SO4	A	7	-	-	0/0/0/0	0/0/0/0
3	EPE	B	2	-	-	0/9/19/19	0/1/1/1
2	SO4	B	6	-	-	0/0/0/0	0/0/0/0

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	2	EPE	C7-N4	4.94	1.59	1.47
3	A	1	EPE	C7-N4	3.17	1.54	1.47
3	B	2	EPE	C5-N4	2.84	1.54	1.47
3	A	1	EPE	O8-C8	2.67	1.56	1.42
3	A	1	EPE	C3-N4	2.62	1.54	1.47
3	B	2	EPE	C3-N4	2.47	1.53	1.47
2	A	5	SO4	O1-S	-2.44	1.39	1.47
3	B	2	EPE	C2-N1	2.44	1.53	1.47
3	B	2	EPE	C9-N1	2.21	1.52	1.47
3	A	1	EPE	C2-N1	2.18	1.53	1.47
3	A	1	EPE	C5-N4	2.09	1.52	1.47

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1	EPE	O2S-S-C10	4.67	110.81	106.81
3	B	2	EPE	O3S-S-C10	2.59	109.20	105.93
3	A	1	EPE	O1S-S-C10	2.58	109.02	106.81
3	A	1	EPE	C9-C10-S	-2.32	104.73	113.01

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.