



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 17, 2018 – 12:10 pm GMT

PDB ID : 1HWG  
Title : 1:2 COMPLEX OF HUMAN GROWTH HORMONE WITH ITS SOLUBLE  
BINDING PROTEIN  
Authors : Sundstrom, S.M.; Lundqvist, T.  
Deposited on : 1996-11-13  
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk30686

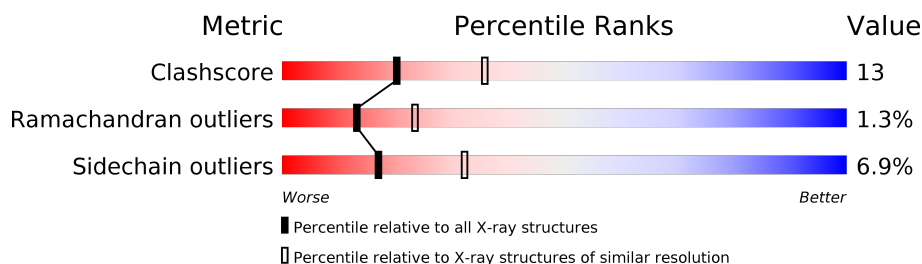
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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	122078	4826 (2.50-2.50)
Ramachandran outliers	120005	4734 (2.50-2.50)
Sidechain outliers	119972	4736 (2.50-2.50)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	191	
2	B	237	
2	C	237	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4718 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 GROWTH HORMONE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	184	Total	C	N	O	S	0	0	0
			1457	933	241	276	7			

- Molecule 2 is a protein called GROWTH HORMONE BINDING PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	194	Total	C	N	O	S	0	0	0
			1559	1001	254	295	9			
2	C	191	Total	C	N	O	S	0	0	0
			1527	979	250	288	10			

- Molecule 3 is water.

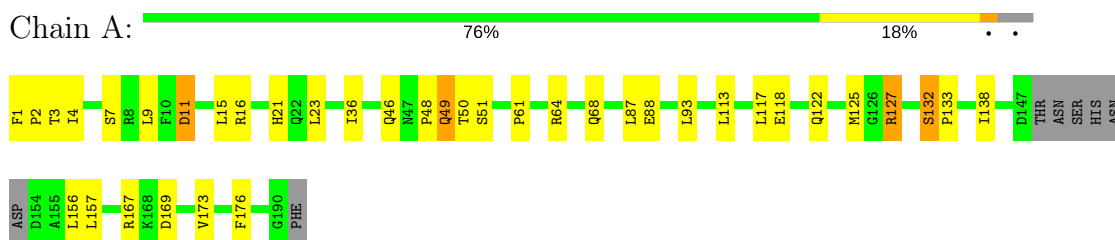
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	50	Total	O	0	0
			50	50		
3	B	57	Total	O	0	0
			57	57		
3	C	68	Total	O	0	0
			68	68		

### 3 Residue-property plots

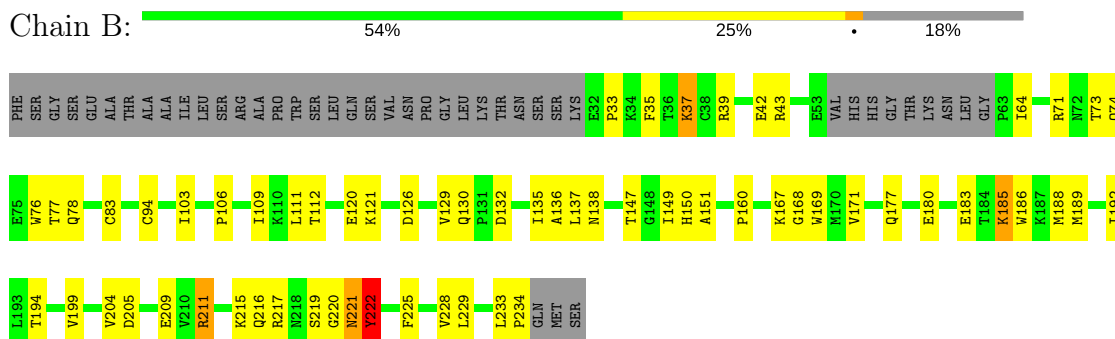
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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.

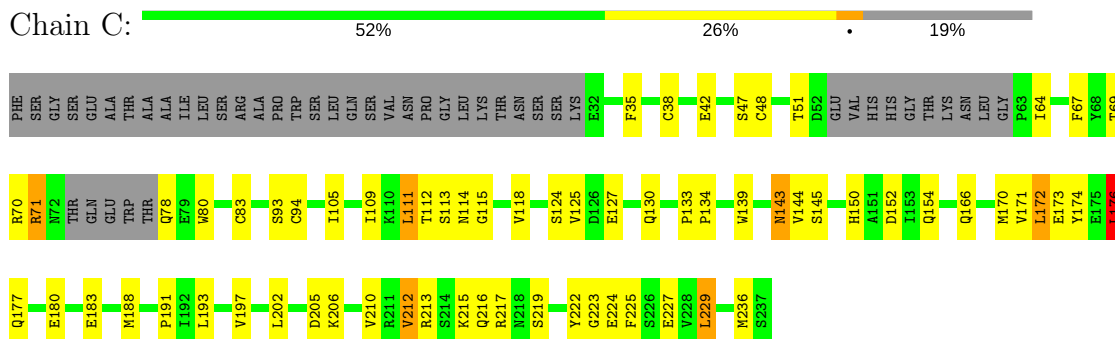
#### • Molecule 1: GROWTH HORMONE



#### • Molecule 2: GROWTH HORMONE BINDING PROTEIN



#### • Molecule 2: GROWTH HORMONE BINDING PROTEIN



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	145.61Å 69.02Å 76.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50	Depositor
% Data completeness (in resolution range)	94.3 (20.00-2.50)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.10	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, $R_{free}$	0.199 , 0.287	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4718	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

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	0.44	0/1486	0.64	1/2010 (0.0%)
2	B	0.48	0/1603	0.76	2/2186 (0.1%)
2	C	0.51	0/1569	0.72	3/2139 (0.1%)
All	All	0.48	0/4658	0.71	6/6335 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	48	PRO	N-CA-CB	6.01	110.51	103.30
2	C	176	LEU	CA-CB-CG	5.88	128.81	115.30
2	B	234	PRO	N-CA-CB	5.32	109.68	103.30
2	C	229	LEU	CA-CB-CG	5.12	127.08	115.30
2	C	173	GLU	N-CA-C	-5.12	97.18	111.00
2	B	222	TYR	N-CA-C	-5.00	97.49	111.00

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-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 hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1457	0	1396	19	0
2	B	1559	0	1463	53	0
2	C	1527	0	1425	47	0
3	A	50	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	57	0	0	4	0
3	C	68	0	0	2	0
All	All	4718	0	4284	113	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 13.

All (113) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:130:GLN:HA	2:B:221:ASN:HB3	1.51	0.91
2:B:216:GLN:HB3	2:B:221:ASN:HD21	1.45	0.82
2:C:134:PRO:HG2	2:C:212:VAL:HG22	1.70	0.73
2:C:144:VAL:HG11	2:C:236:MET:HG3	1.72	0.71
2:B:71:ARG:HB3	2:B:106:PRO:HD2	1.73	0.70
2:C:133:PRO:HB3	2:C:224:GLU:O	1.92	0.69
1:A:64:ARG:O	1:A:68:GLN:HG2	1.93	0.68
2:B:219:SER:HB2	2:B:221:ASN:OD1	1.95	0.67
2:B:221:ASN:O	2:B:222:TYR:HB2	1.96	0.64
2:B:74:GLN:O	2:B:76:TRP:HD1	1.82	0.63
2:B:220:GLY:HA2	3:B:250:HOH:O	1.99	0.62
2:B:186:TRP:CZ3	2:B:211:ARG:HG2	2.37	0.60
2:C:35:PHE:CZ	2:C:109:ILE:HD12	2.36	0.60
2:C:176:LEU:HD12	2:C:177:GLN:N	2.16	0.60
2:C:83:CYS:CB	2:C:94:CYS:SG	2.90	0.60
2:C:171:VAL:HG11	2:C:217:ARG:HG3	1.82	0.60
2:C:193:LEU:HD21	2:C:217:ARG:HD3	1.82	0.60
1:A:4:ILE:O	1:A:127:ARG:NH2	2.35	0.60
2:C:71:ARG:HE	2:C:78:GLN:HE22	1.48	0.60
1:A:11:ASP:O	1:A:15:LEU:HD13	2.02	0.59
2:C:205:ASP:O	2:C:206:LYS:HG3	2.02	0.59
1:A:36:ILE:HD11	1:A:156:LEU:HD21	1.86	0.58
2:B:215:LYS:HB2	2:B:222:TYR:CD1	2.39	0.57
2:B:211:ARG:HD3	2:B:225:PHE:CD1	2.41	0.56
2:B:130:GLN:HA	2:B:221:ASN:CB	2.30	0.56
2:C:193:LEU:CD2	2:C:217:ARG:HD3	2.35	0.56
2:B:168:GLY:HA2	3:B:245:HOH:O	2.05	0.56
2:B:215:LYS:HG3	2:B:221:ASN:OD1	2.06	0.55
2:B:189:MET:HG2	3:B:272:HOH:O	2.06	0.55
2:B:103:ILE:CG2	2:B:126:ASP:HB3	2.36	0.55
2:B:37:LYS:C	2:B:37:LYS:HD2	2.27	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:150:HIS:HD2	2:C:152:ASP:OD2	1.91	0.53
2:C:171:VAL:O	2:C:172:LEU:HB2	2.09	0.53
2:B:64:ILE:HA	2:B:112:THR:O	2.08	0.53
1:A:68:GLN:NE2	3:A:214:HOH:O	2.41	0.53
2:C:64:ILE:CG2	2:C:111:LEU:HD22	2.39	0.52
1:A:1:PHE:HB2	1:A:2:PRO:HD2	1.91	0.52
2:B:171:VAL:HG12	2:B:217:ARG:HG3	1.92	0.52
2:B:83:CYS:CB	2:B:94:CYS:SG	2.99	0.51
2:B:177:GLN:HG2	2:B:188:MET:HB3	1.92	0.50
2:B:130:GLN:CA	2:B:221:ASN:HB3	2.34	0.50
2:C:176:LEU:HD23	2:C:197:VAL:CG1	2.42	0.50
2:C:177:GLN:HG2	2:C:188:MET:HG2	1.94	0.50
1:A:118:GLU:O	1:A:122:GLN:HG3	2.12	0.50
2:B:209:GLU:HB2	2:B:228:VAL:CG1	2.42	0.49
2:C:180:GLU:HB2	2:C:183:GLU:HB2	1.93	0.49
2:C:83:CYS:SG	2:C:94:CYS:CB	3.00	0.49
1:A:64:ARG:HH22	2:B:167:LYS:HD2	1.78	0.49
2:B:37:LYS:O	2:B:37:LYS:HD2	2.12	0.49
2:B:185:LYS:H	2:B:185:LYS:HD3	1.78	0.49
2:C:172:LEU:HA	2:C:215:LYS:O	2.12	0.49
1:A:156:LEU:HD23	1:A:157:LEU:N	2.27	0.49
2:C:48:CYS:O	2:C:93:SER:HA	2.12	0.49
2:C:130:GLN:HG2	2:C:223:GLY:HA2	1.93	0.49
2:B:233:LEU:H	2:B:233:LEU:HD12	1.77	0.48
2:C:176:LEU:HD13	2:C:210:VAL:HG13	1.96	0.48
2:B:76:TRP:O	2:B:78:GLN:N	2.47	0.48
1:A:49:GLN:O	1:A:51:SER:N	2.47	0.48
1:A:88:GLU:HB2	3:A:201:HOH:O	2.13	0.48
1:A:64:ARG:NH2	2:B:167:LYS:HD2	2.28	0.47
2:C:176:LEU:CD1	2:C:210:VAL:HG13	2.45	0.47
2:C:139:TRP:HA	2:C:154:GLN:O	2.15	0.47
2:B:211:ARG:HD3	2:B:225:PHE:CE1	2.50	0.47
2:B:71:ARG:CB	2:B:106:PRO:HD2	2.42	0.47
1:A:125:MET:HG2	3:A:240:HOH:O	2.15	0.47
1:A:169:ASP:O	1:A:173:VAL:HG23	2.15	0.47
2:B:204:VAL:HA	2:B:233:LEU:HD13	1.97	0.46
2:B:151:ALA:HB3	2:B:233:LEU:HD23	1.95	0.46
2:B:73:THR:O	2:B:76:TRP:HB3	2.15	0.46
2:C:113:SER:HB3	2:C:118:VAL:HG11	1.98	0.46
1:A:132:SER:HA	1:A:133:PRO:HD2	1.86	0.45
2:B:120:GLU:HG2	2:B:121:LYS:N	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:151:ALA:HB3	2:B:233:LEU:CD2	2.46	0.45
2:C:150:HIS:HA	2:C:202:LEU:O	2.16	0.45
2:B:39:ARG:HG3	2:B:130:GLN:HB3	1.99	0.45
2:C:67:PHE:HB3	2:C:80:TRP:HB3	1.99	0.45
2:C:176:LEU:HD23	2:C:197:VAL:HG11	2.00	0.44
2:B:180:GLU:HB2	2:B:183:GLU:HG3	1.98	0.44
2:C:112:THR:HG22	2:C:112:THR:O	2.18	0.44
2:B:216:GLN:HB3	2:B:221:ASN:ND2	2.22	0.44
2:C:113:SER:HB3	2:C:118:VAL:CG1	2.47	0.44
2:B:135:ILE:HG22	2:B:136:ALA:N	2.32	0.43
2:C:38:CYS:HA	2:C:47:SER:O	2.18	0.43
2:C:174:TYR:O	2:C:191:PRO:HA	2.19	0.43
2:C:150:HIS:HD2	3:C:241:HOH:O	2.01	0.43
2:C:216:GLN:CG	2:C:219:SER:HB3	2.49	0.43
2:B:149:ILE:HD12	2:C:143:ASN:ND2	2.33	0.43
1:A:61:PRO:HD2	1:A:176:PHE:CZ	2.53	0.43
2:C:215:LYS:HB2	2:C:222:TYR:CD1	2.54	0.42
2:B:209:GLU:HA	2:B:229:LEU:O	2.19	0.42
2:C:133:PRO:HA	2:C:134:PRO:HD3	1.86	0.42
2:C:70:ARG:HE	2:C:105:ILE:HD11	1.85	0.42
2:B:35:PHE:CZ	2:B:111:LEU:HB2	2.54	0.42
2:C:124:SER:OG	2:C:127:GLU:HG3	2.20	0.42
2:C:70:ARG:HB3	2:C:105:ILE:HD12	2.00	0.42
1:A:46:GLN:HA	2:B:76:TRP:HZ2	1.85	0.42
2:B:129:VAL:O	2:B:221:ASN:HB3	2.20	0.42
2:B:132:ASP:HB2	2:B:160:PRO:HB3	2.02	0.42
1:A:138:ILE:HA	1:A:138:ILE:HD12	1.83	0.41
2:B:192:ILE:HG22	2:B:194:THR:H	1.84	0.41
2:C:112:THR:HG23	2:C:115:GLY:C	2.40	0.41
2:B:33:PRO:HG2	2:B:64:ILE:CD1	2.50	0.41
2:C:213:ARG:HD3	2:C:225:PHE:CE1	2.55	0.41
2:B:43:ARG:HB3	2:B:169:TRP:HE1	1.86	0.41
2:B:209:GLU:HB2	2:B:228:VAL:HG11	2.03	0.41
2:B:147:THR:HG22	2:C:145:SER:HB3	2.01	0.41
2:B:222:TYR:HA	2:B:222:TYR:HD1	1.66	0.41
2:C:213:ARG:HD3	2:C:225:PHE:CZ	2.56	0.41
1:A:16:ARG:HB2	1:A:117:LEU:HD13	2.02	0.40
2:C:125:VAL:HG23	3:C:275:HOH:O	2.21	0.40
2:C:176:LEU:C	2:C:176:LEU:HD12	2.42	0.40
2:B:135:ILE:HD11	3:B:265:HOH:O	2.21	0.40
2:C:171:VAL:CG1	2:C:217:ARG:HG3	2.49	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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	180/191 (94%)	166 (92%)	11 (6%)	3 (2%)	10	17
2	B	190/237 (80%)	167 (88%)	21 (11%)	2 (1%)	16	28
2	C	185/237 (78%)	165 (89%)	18 (10%)	2 (1%)	16	28
All	All	555/665 (84%)	498 (90%)	50 (9%)	7 (1%)	13	23

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	50	THR
2	C	71	ARG
2	B	222	TYR
2	B	77	THR
1	A	49	GLN
2	C	172	LEU
1	A	132	SER

### 5.3.2 Protein sidechains [i](#)

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	156/176 (89%)	145 (93%)	11 (7%)	16	31
2	B	170/217 (78%)	159 (94%)	11 (6%)	19	35

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	C	166/217 (76%)	154 (93%)	12 (7%)	16	30
All	All	492/610 (81%)	458 (93%)	34 (7%)	17	32

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	THR
1	A	7	SER
1	A	9	LEU
1	A	11	ASP
1	A	21	HIS
1	A	23	LEU
1	A	87	LEU
1	A	93	LEU
1	A	113	LEU
1	A	127	ARG
1	A	167	ARG
2	B	37	LYS
2	B	42	GLU
2	B	109	ILE
2	B	137	LEU
2	B	138	ASN
2	B	185	LYS
2	B	199	VAL
2	B	205	ASP
2	B	211	ARG
2	B	221	ASN
2	B	222	TYR
2	C	42	GLU
2	C	51	THR
2	C	69	THR
2	C	111	LEU
2	C	114	ASN
2	C	143	ASN
2	C	166	GLN
2	C	170	MET
2	C	176	LEU
2	C	212	VAL
2	C	227	GLU
2	C	229	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	GLN
1	A	68	GLN
1	A	181	GLN
2	B	49	HIS
2	C	78	GLN
2	C	114	ASN
2	C	143	ASN
2	C	150	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

### 5.7 Other polymers [i](#)

There are no such residues in this entry.

### 5.8 Polymer linkage issues [i](#)

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 is therefore empty.

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

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

### 6.4 Ligands

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers

EDS was not executed - this section is therefore empty.