Full wwPDB X-ray Structure Validation Report (i)

Oct 23, 2023 – 12:00 PM EDT

PDB ID	:	3EMO
Title	:	Crystal structure of transmembrane Hia 973-1098
Authors	:	Meng, G.; Waksman, G.
Deposited on	:	2008-09-24
Resolution	:	3.00 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*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

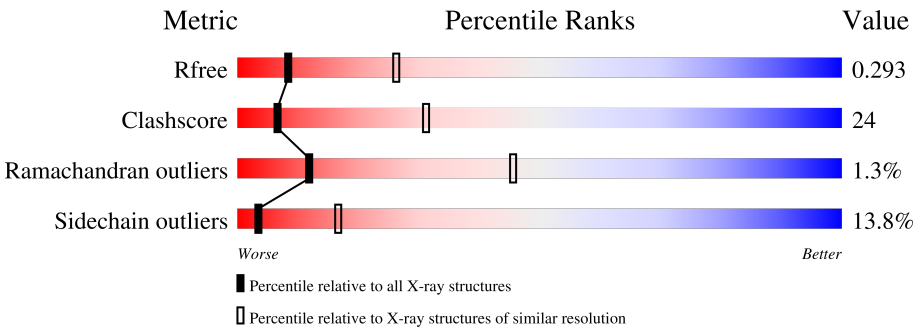
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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))
R_{free}	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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 <=5%

Mol	Chain	Length	Quality of chain
1	A	162	50% 22% 6% 22%
1	B	162	52% 22% • 22%
1	C	162	45% 26% 7% 22%

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 2605 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 Hia (Adhesin).

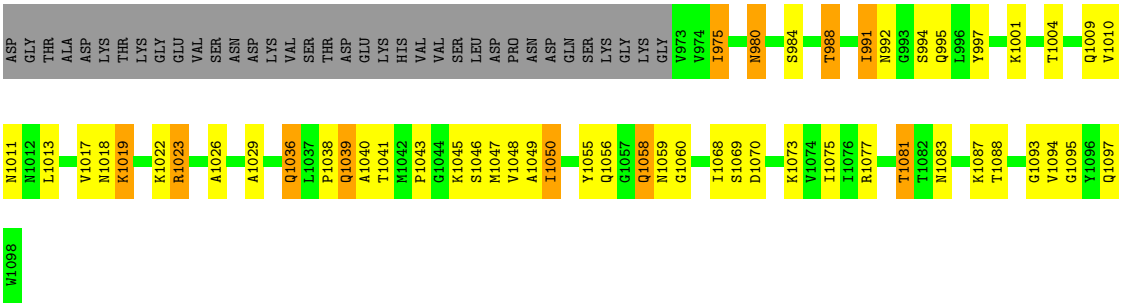
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	C	126	875	535	160	178	2	0	0	0
			Total	C	N	O	S			
1	A	126	861	525	155	179	2	0	0	0
			Total	C	N	O	S			
1	B	126	869	532	157	178	2	0	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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. 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. 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.

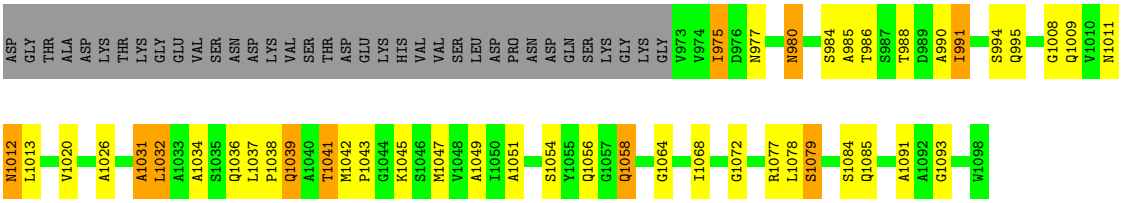
- Molecule 1: Hia (Adhesin)

Chain C: 45% 26% 7% 22%



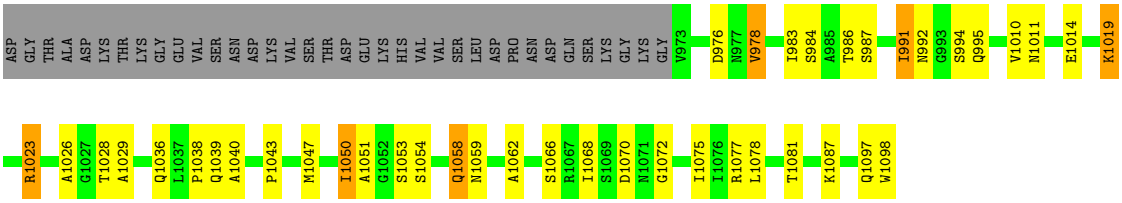
- Molecule 1: Hia (Adhesin)

Chain A: 50% 22% 6% 22%



- Molecule 1: Hia (Adhesin)

Chain B: 52% 22% • 22%



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	194.30Å 45.81Å 56.42Å 90.00° 95.12° 90.00°	Depositor
Resolution (Å)	29.57 – 3.00 29.57 – 2.99	Depositor EDS
% Data completeness (in resolution range)	97.2 (29.57-3.00) 83.0 (29.57-2.99)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	5.20 (at 3.00Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.221 , 0.289 0.298 , 0.293	Depositor DCC
R_{free} test set	846 reflections (9.93%)	wwPDB-VP
Wilson B-factor (Å^2)	35.8	Xtriage
Anisotropy	0.517	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.26 , 26.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.41$, $< L^2 > = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.84	EDS
Total number of atoms	2605	wwPDB-VP
Average B, all atoms (Å^2)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.69% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

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	0.32	0/867	0.54	0/1179
1	B	0.34	0/875	0.51	0/1186
1	C	0.33	0/881	0.49	0/1193
All	All	0.33	0/2623	0.52	0/3558

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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	861	0	842	44	0
1	B	869	0	875	43	0
1	C	875	0	886	72	0
All	All	2605	0	2603	123	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 24.

All (123) 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)
1:B:1019:LYS:HD2	1:B:1023:ARG:HD3	1.53	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:980:ASN:H	1:C:980:ASN:HD22	1.27	0.83
1:C:1056:GLN:NE2	1:B:1087:LYS:HE3	1.96	0.81
1:A:1012:ASN:HD22	1:A:1012:ASN:N	1.81	0.79
1:C:980:ASN:ND2	1:C:992:ASN:HD21	1.79	0.78
1:C:980:ASN:HD22	1:C:980:ASN:N	1.82	0.73
1:C:1039:GLN:HE22	1:A:1039:GLN:H	1.34	0.72
1:C:975:ILE:O	1:C:975:ILE:HG13	1.89	0.70
1:A:1026:ALA:HB2	1:A:1058:GLN:HG2	1.73	0.69
1:B:992:ASN:H	1:B:995:GLN:HE21	1.40	0.69
1:C:1045:LYS:HG2	1:B:1098:TRP:O	1.93	0.68
1:C:991:ILE:HD11	1:B:991:ILE:HD11	1.76	0.66
1:C:1010:VAL:HG12	1:C:1010:VAL:O	1.94	0.66
1:C:992:ASN:OD1	1:C:994:SER:HB2	1.96	0.65
1:C:980:ASN:HA	1:C:995:GLN:NE2	2.11	0.65
1:C:1013:LEU:HD21	1:A:1013:LEU:HD23	1.80	0.64
1:C:994:SER:HA	1:B:984:SER:O	1.99	0.62
1:C:1056:GLN:NE2	1:B:1087:LYS:HB3	2.14	0.62
1:C:1081:THR:O	1:C:1088:THR:HA	2.00	0.61
1:C:1097:GLN:HB3	1:A:1047:MET:HB2	1.83	0.61
1:C:1087:LYS:HD2	1:A:1056:GLN:NE2	2.16	0.60
1:A:1058:GLN:HG3	1:A:1084:SER:OG	2.00	0.60
1:C:992:ASN:H	1:C:995:GLN:HE21	1.48	0.60
1:A:1038:PRO:HD2	1:A:1049:ALA:HB1	1.84	0.60
1:C:980:ASN:H	1:C:980:ASN:ND2	1.99	0.60
1:B:1036:GLN:O	1:B:1038:PRO:HD3	2.01	0.60
1:A:984:SER:C	1:A:986:THR:H	2.04	0.60
1:A:1064:GLY:HA2	1:A:1079:SER:HA	1.84	0.60
1:B:983:ILE:HG13	1:B:983:ILE:O	2.02	0.60
1:A:991:ILE:HD11	1:B:991:ILE:HD11	1.83	0.59
1:C:1056:GLN:HE22	1:B:1087:LYS:HE3	1.68	0.58
1:C:975:ILE:O	1:C:975:ILE:CG1	2.51	0.58
1:C:1039:GLN:HE22	1:A:1039:GLN:N	1.99	0.58
1:A:1012:ASN:HD22	1:A:1012:ASN:H	1.49	0.58
1:C:1040:ALA:HA	1:C:1047:MET:HE3	1.86	0.57
1:C:1039:GLN:H	1:B:1039:GLN:HE22	1.53	0.57
1:A:1093:GLY:HA3	1:B:1036:GLN:O	2.04	0.56
1:C:1019:LYS:HE3	1:C:1023:ARG:NH1	2.20	0.56
1:B:1043:PRO:HB3	1:B:1072:GLY:CA	2.36	0.56
1:B:1066:SER:HB3	1:B:1077:ARG:HB3	1.89	0.55
1:C:1047:MET:HE1	1:B:1075:ILE:HD12	1.88	0.55
1:B:992:ASN:N	1:B:995:GLN:HE21	2.05	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1026:ALA:HB1	1:C:1083:ASN:HB2	1.89	0.54
1:C:1068:ILE:HG22	1:C:1069:SER:O	2.07	0.54
1:A:1068:ILE:CG2	1:A:1072:GLY:HA2	2.38	0.54
1:B:1026:ALA:HB2	1:B:1058:GLN:HG2	1.90	0.53
1:C:1050:ILE:O	1:C:1050:ILE:HG23	2.08	0.53
1:B:992:ASN:HB3	1:B:994:SER:H	1.74	0.52
1:B:1050:ILE:HD13	1:B:1051:ALA:N	2.25	0.52
1:A:1068:ILE:HG22	1:A:1072:GLY:HA2	1.92	0.52
1:B:1068:ILE:HG22	1:B:1072:GLY:HA2	1.92	0.52
1:C:1068:ILE:HG12	1:C:1075:ILE:HG23	1.92	0.52
1:C:1039:GLN:NE2	1:A:1039:GLN:H	2.05	0.51
1:C:1019:LYS:HD2	1:C:1023:ARG:NH1	2.25	0.51
1:C:1038:PRO:HD2	1:C:1049:ALA:HB1	1.93	0.51
1:C:1018:ASN:O	1:C:1022:LYS:HG3	2.11	0.51
1:B:1040:ALA:HA	1:B:1047:MET:HE3	1.93	0.51
1:A:975:ILE:HG13	1:A:975:ILE:O	2.11	0.50
1:C:1013:LEU:O	1:C:1017:VAL:HG23	2.11	0.49
1:A:1037:LEU:HD23	1:A:1051:ALA:HB2	1.94	0.49
1:C:1029:ALA:O	1:C:1060:GLY:HA3	2.13	0.48
1:C:1039:GLN:O	1:C:1047:MET:HE1	2.14	0.48
1:C:1010:VAL:HG11	1:B:1010:VAL:HG22	1.96	0.48
1:A:990:ALA:HB2	1:B:978:VAL:HG21	1.95	0.47
1:C:1058:GLN:C	1:C:1059:ASN:HD22	2.18	0.47
1:C:1043:PRO:HA	1:C:1068:ILE:HB	1.97	0.47
1:B:1026:ALA:O	1:B:1029:ALA:N	2.47	0.47
1:C:980:ASN:ND2	1:C:980:ASN:N	2.54	0.47
1:A:984:SER:C	1:A:986:THR:N	2.68	0.47
1:C:980:ASN:HD22	1:C:992:ASN:HD21	1.62	0.47
1:C:1038:PRO:HD2	1:C:1049:ALA:CB	2.44	0.46
1:C:984:SER:O	1:A:994:SER:HA	2.15	0.46
1:B:1068:ILE:HG23	1:B:1075:ILE:HG12	1.97	0.46
1:C:1094:VAL:HG22	1:C:1095:GLY:N	2.31	0.45
1:A:1012:ASN:H	1:A:1012:ASN:ND2	2.14	0.45
1:C:1048:VAL:HG23	1:C:1048:VAL:O	2.16	0.45
1:C:1026:ALA:HB2	1:C:1058:GLN:CG	2.47	0.45
1:C:1013:LEU:CD2	1:A:1013:LEU:HD23	2.47	0.45
1:C:991:ILE:HB	1:C:995:GLN:NE2	2.32	0.45
1:C:1038:PRO:HA	1:B:1039:GLN:NE2	2.32	0.44
1:C:1010:VAL:O	1:C:1010:VAL:CG1	2.63	0.44
1:C:1026:ALA:HB1	1:C:1083:ASN:CB	2.47	0.44
1:C:1094:VAL:C	1:A:1038:PRO:HG2	2.38	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1078:LEU:O	1:A:1079:SER:HB3	2.16	0.44
1:A:1084:SER:O	1:A:1085:GLN:HG2	2.17	0.44
1:C:1055:TYR:CD2	1:C:1056:GLN:HG2	2.52	0.44
1:C:1010:VAL:CG1	1:B:1010:VAL:HG22	2.49	0.43
1:A:1008:GLY:O	1:A:1011:ASN:HB3	2.18	0.43
1:C:1056:GLN:HE21	1:B:1087:LYS:HE3	1.78	0.43
1:C:1093:GLY:HA3	1:A:1036:GLN:O	2.18	0.43
1:A:988:THR:O	1:B:978:VAL:HG23	2.19	0.43
1:B:1039:GLN:O	1:B:1047:MET:HE1	2.18	0.43
1:C:1019:LYS:CD	1:C:1023:ARG:NH1	2.82	0.43
1:A:1012:ASN:N	1:A:1012:ASN:ND2	2.53	0.43
1:A:990:ALA:HB2	1:B:978:VAL:HG11	2.02	0.42
1:C:988:THR:O	1:C:988:THR:HG23	2.19	0.42
1:C:992:ASN:H	1:C:995:GLN:NE2	2.16	0.42
1:B:1062:ALA:HB2	1:B:1081:THR:HG22	2.02	0.42
1:C:1019:LYS:HB3	1:C:1019:LYS:HE2	1.65	0.42
1:C:1026:ALA:HB2	1:C:1058:GLN:HG2	2.01	0.42
1:C:1045:LYS:HG2	1:B:1098:TRP:C	2.40	0.42
1:C:1036:GLN:OE1	1:C:1036:GLN:HA	2.20	0.42
1:C:1046:SER:O	1:B:1097:GLN:HA	2.20	0.41
1:C:997:TYR:O	1:C:1001:LYS:HB2	2.20	0.41
1:C:1010:VAL:HG11	1:B:1010:VAL:CG2	2.50	0.41
1:A:1031:ALA:O	1:A:1032:LEU:C	2.59	0.41
1:C:1073:LYS:HB2	1:C:1073:LYS:HE2	1.88	0.41
1:A:1091:ALA:O	1:B:1036:GLN:HG3	2.20	0.41
1:A:975:ILE:HD12	1:A:975:ILE:C	2.40	0.41
1:B:1053:SER:O	1:B:1059:ASN:HA	2.20	0.41
1:A:1042:MET:HA	1:A:1043:PRO:HD3	1.96	0.41
1:C:988:THR:HG23	1:A:977:ASN:H	1.86	0.41
1:A:1042:MET:O	1:A:1045:LYS:HB2	2.21	0.41
1:C:1038:PRO:HA	1:B:1039:GLN:HE22	1.84	0.41
1:A:1031:ALA:O	1:A:1034:ALA:N	2.54	0.41
1:B:1077:ARG:C	1:B:1078:LEU:HD12	2.40	0.41
1:A:1009:GLN:HE21	1:B:1010:VAL:HG12	1.86	0.40
1:C:1009:GLN:C	1:C:1011:ASN:H	2.24	0.40
1:C:1041:THR:HA	1:A:1041:THR:HG21	2.04	0.40
1:A:980:ASN:CA	1:A:995:GLN:OE1	2.70	0.40
1:B:1026:ALA:CA	1:B:1058:GLN:HG2	2.51	0.40
1:A:1077:ARG:HG3	1:B:1036:GLN:HE21	1.86	0.40
1:A:1078:LEU:N	1:A:1078:LEU:CD1	2.84	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	124/162 (76%)	103 (83%)	16 (13%)	5 (4%)	3 17
1	B	124/162 (76%)	113 (91%)	11 (9%)	0	100 100
1	C	124/162 (76%)	115 (93%)	9 (7%)	0	100 100
All	All	372/486 (76%)	331 (89%)	36 (10%)	5 (1%)	12 45

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	980	ASN
1	A	985	ALA
1	A	1031	ALA
1	A	1032	LEU
1	A	1079	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	84/121 (69%)	76 (90%)	8 (10%)	8 32
1	B	88/121 (73%)	74 (84%)	14 (16%)	2 12
1	C	89/121 (74%)	75 (84%)	14 (16%)	2 13
All	All	261/363 (72%)	225 (86%)	36 (14%)	3 17

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	975	ILE
1	C	980	ASN
1	C	988	THR
1	C	991	ILE
1	C	1004	THR
1	C	1019	LYS
1	C	1023	ARG
1	C	1036	GLN
1	C	1039	GLN
1	C	1050	ILE
1	C	1058	GLN
1	C	1070	ASP
1	C	1077	ARG
1	C	1081	THR
1	A	975	ILE
1	A	991	ILE
1	A	1012	ASN
1	A	1020	VAL
1	A	1039	GLN
1	A	1041	THR
1	A	1054	SER
1	A	1058	GLN
1	B	976	ASP
1	B	978	VAL
1	B	986	THR
1	B	987	SER
1	B	991	ILE
1	B	1011	ASN
1	B	1014	GLU
1	B	1019	LYS
1	B	1023	ARG
1	B	1028	THR
1	B	1050	ILE
1	B	1054	SER
1	B	1058	GLN
1	B	1070	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	980	ASN
1	C	995	GLN
1	C	1012	ASN

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Mol	Chain	Res	Type
1	C	1039	GLN
1	C	1059	ASN
1	A	977	ASN
1	A	980	ASN
1	A	1012	ASN
1	B	995	GLN
1	B	1036	GLN
1	B	1039	GLN
1	B	1059	ASN

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 monosaccharides 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 (i)

6.1 Protein, DNA and RNA chains (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers (i)

Unable to reproduce the depositors R factor - this section is therefore empty.