Full wwPDB X-ray Structure Validation Report (i)

Dec 14, 2023 – 03:43 am GMT

PDB ID	:	2WS2
Title	:	The 2 Angstrom structure of a Nu-class GST from Haemonchus contortus
Authors	:	Line, K.; Isupov, M.N.; vanRossum, A.J.; Brophy, P.M.; Littlechild, J.A.
Deposited on	:	2009-09-03
Resolution	:	2.01 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:

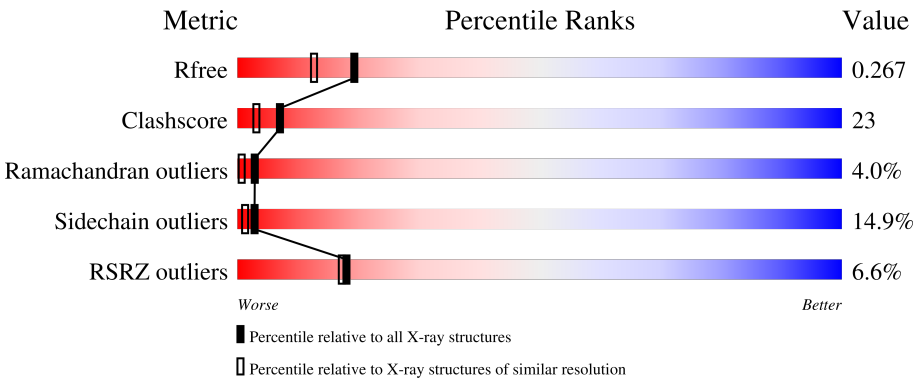
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation 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 2.01 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (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 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% 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.

Mol	Chain	Length	Quality of chain
1	A	204	4% 58% 33% 8%
1	B	204	9% 61% 29% 7% •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3520 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 GLUTATHIONE S-TRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	204	1676	1094	277	301	4	0	2	0
1	B	204	1666	1088	277	297	4	0	1	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	164	GLY	THR	SEE REMARK 999	UNP Q9NAW7
B	164	GLY	THR	SEE REMARK 999	UNP Q9NAW7

- Molecule 2 is water.

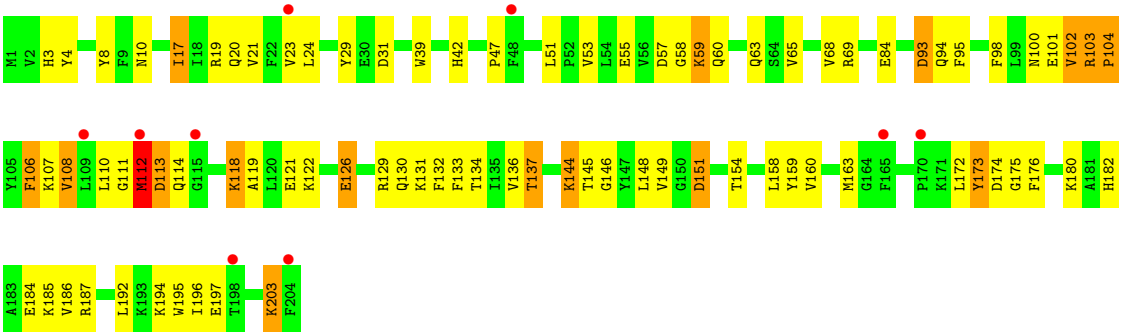
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	78	78	78	0	0
2	B	100	100	100	0	0

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 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.

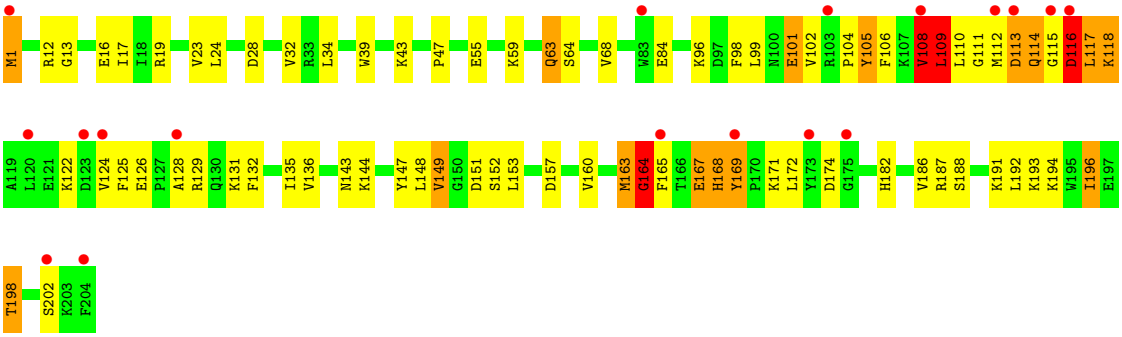
• Molecule 1: GLUTATHIONE S-TRANSFERASE

Chain A:



• Molecule 1: GLUTATHIONE S-TRANSFERASE

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	47.22Å 101.38Å 47.17Å 90.00° 100.46° 90.00°	Depositor
Resolution (Å)	50.71 – 2.01 13.06 – 2.01	Depositor EDS
% Data completeness (in resolution range)	96.8 (50.71-2.01) 97.3 (13.06-2.01)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.24 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.5.0088	Depositor
R, R_{free}	0.233 , 0.290 0.248 , 0.267	Depositor DCC
R_{free} test set	1433 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å^2)	26.0	Xtriage
Anisotropy	0.667	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 20.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.45$, $< L^2 > = 0.27$	Xtriage
Estimated twinning fraction	0.407 for l,-k,h	Xtriage
Reported twinning fraction	0.599 for H,K,L 0.401 for L,-K,H	Depositor
Outliers	0 of 28324 reflections	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	3520	wwPDB-VP
Average B, all atoms (Å^2)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.22% 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	0.82	0/1726	0.88	3/2325 (0.1%)
1	B	0.80	0/1713	0.92	2/2312 (0.1%)
All	All	0.81	0/3439	0.90	5/4637 (0.1%)

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	B	0	2

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	93	ASP	CB-CG-OD1	6.12	123.81	118.30
1	A	69	ARG	NE-CZ-NH2	-5.93	117.33	120.30
1	B	109	LEU	CA-CB-CG	5.63	128.24	115.30
1	B	168	HIS	N-CA-C	-5.44	96.31	111.00
1	A	93	ASP	CB-CG-OD2	-5.20	113.62	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	116	ASP	Peptide
1	B	164	GLY	Peptide

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	1676	0	1676	87	0
1	B	1666	0	1674	67	0
2	A	78	0	0	14	0
2	B	100	0	0	5	1
All	All	3520	0	3350	153	1

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 23.

All (153) 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:A:42:HIS:HB2	2:A:2017:HOH:O	1.21	1.30
1:B:108:VAL:HG23	1:B:109:LEU:N	1.75	1.01
1:B:101:GLU:OE1	1:B:128:ALA:HB2	1.62	0.99
1:A:133:PHE:O	1:A:137:THR:HG22	1.64	0.95
1:B:108:VAL:HG23	1:B:109:LEU:H	1.36	0.90
1:A:29:TYR:CE1	1:A:31[B]:ASP:HB3	2.05	0.90
1:B:104:PRO:O	2:B:2067:HOH:O	1.90	0.89
1:B:39:TRP:CZ2	1:B:43:LYS:HG3	2.08	0.89
1:A:145:THR:HG21	1:A:151:ASP:HB3	1.54	0.88
1:B:163:MET:O	1:B:165:PHE:N	2.07	0.87
1:A:65:VAL:HA	1:A:68:VAL:HG22	1.58	0.86
1:A:194:LYS:HE3	1:A:197:GLU:OE2	1.74	0.86
1:A:8:TYR:HB3	2:A:2009:HOH:O	1.76	0.85
1:A:19:ARG:O	1:A:23:VAL:HG23	1.77	0.85
1:A:42:HIS:NE2	2:A:2019:HOH:O	2.10	0.84
1:A:65:VAL:HA	1:A:68:VAL:CG2	2.12	0.80
1:A:42:HIS:CD2	2:A:2019:HOH:O	2.35	0.79
1:A:17:ILE:HG23	1:A:68:VAL:CG1	2.13	0.79
1:A:154:THR:HG22	2:A:2065:HOH:O	1.82	0.78
1:B:109:LEU:HD12	1:B:111:GLY:O	1.84	0.77
1:A:130:GLN:O	1:A:134:THR:HG23	1.84	0.77
1:A:42:HIS:CB	2:A:2017:HOH:O	1.96	0.77
1:A:17:ILE:HG23	1:A:68:VAL:HG11	1.65	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:GLU:HG3	1:A:149:VAL:CG1	2.15	0.76
1:A:84:GLU:HG3	1:A:149:VAL:HG12	1.68	0.75
1:A:17:ILE:CG2	1:A:68:VAL:HG11	2.17	0.74
1:B:101:GLU:OE1	1:B:128:ALA:CB	2.35	0.74
1:A:119:ALA:HA	1:A:122:LYS:HG2	1.72	0.71
1:B:169:TYR:HD1	1:B:169:TYR:O	1.74	0.70
1:A:17:ILE:HG21	1:A:159:TYR:CZ	2.26	0.70
1:B:157:ASP:OD2	2:B:2086:HOH:O	2.09	0.69
1:A:173:TYR:HD1	1:A:180:LYS:N	1.92	0.68
1:A:194:LYS:HA	1:A:197:GLU:OE2	1.93	0.68
1:A:29:TYR:HE1	1:A:31[B]:ASP:HB3	1.53	0.67
1:B:193:LYS:O	1:B:196:ILE:HG23	1.95	0.67
1:A:59:LYS:N	1:A:59:LYS:HD3	2.07	0.67
1:B:163:MET:C	1:B:165:PHE:H	1.97	0.67
1:A:132:PHE:O	1:A:136:VAL:HG23	1.94	0.66
1:A:187:ARG:HA	1:A:192:LEU:HD23	1.76	0.66
1:B:108:VAL:CG2	1:B:109:LEU:N	2.49	0.66
1:A:118:LYS:O	1:A:122:LYS:HE3	1.97	0.65
1:A:29:TYR:HA	2:A:2012:HOH:O	1.97	0.64
1:A:173:TYR:CD1	1:A:180:LYS:HA	2.33	0.63
1:A:100:ASN:HA	1:A:103:ARG:HD2	1.81	0.62
1:A:111:GLY:O	1:A:112:MET:HB2	1.99	0.61
1:B:117:LEU:O	1:B:118:LYS:HB2	1.99	0.61
1:A:4:TYR:HE1	2:A:2002:HOH:O	1.81	0.61
1:A:65:VAL:CA	1:A:68:VAL:HG22	2.30	0.60
1:B:174:ASP:HB3	2:B:2091:HOH:O	2.01	0.60
1:B:125:PHE:HD2	1:B:126:GLU:HG2	1.68	0.59
1:A:173:TYR:N	1:A:173:TYR:CD2	2.69	0.59
1:A:175:GLY:O	2:A:2070:HOH:O	2.17	0.59
1:B:125:PHE:CD2	1:B:126:GLU:HG2	2.37	0.59
1:B:169:TYR:O	1:B:169:TYR:CD1	2.55	0.58
1:A:4:TYR:CE1	2:A:2002:HOH:O	2.51	0.58
1:B:19:ARG:O	1:B:23:VAL:HG13	2.04	0.58
1:B:153:LEU:HD11	1:B:186:VAL:HG13	1.85	0.57
1:A:144:LYS:HB2	1:A:144:LYS:NZ	2.19	0.57
1:A:17:ILE:HD13	1:A:17:ILE:O	2.05	0.57
1:A:187:ARG:HD3	2:A:2076:HOH:O	2.05	0.56
1:A:23:VAL:HG11	1:A:195:TRP:HB2	1.87	0.56
1:A:17:ILE:HG23	1:A:68:VAL:HG12	1.87	0.55
1:A:172:LEU:HD23	1:A:173:TYR:HE2	1.71	0.55
1:B:169:TYR:CD1	1:B:169:TYR:C	2.79	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:LEU:HD13	1:B:113:ASP:HA	1.88	0.55
1:B:98:PHE:HE2	1:B:163:MET:SD	2.30	0.55
1:B:163:MET:C	1:B:165:PHE:N	2.57	0.54
1:A:100:ASN:HA	1:A:103:ARG:CD	2.38	0.54
1:B:24[B]:LEU:HD12	1:B:191:LYS:HZ2	1.72	0.54
1:A:173:TYR:CE1	1:A:180:LYS:HA	2.43	0.53
1:A:173:TYR:CD1	1:A:180:LYS:CA	2.91	0.53
1:B:109:LEU:HG	1:B:109:LEU:O	2.07	0.53
1:A:173:TYR:N	1:A:173:TYR:HD2	2.05	0.53
1:B:192:LEU:O	1:B:196:ILE:HG22	2.08	0.53
1:B:113:ASP:O	1:B:115:GLY:N	2.41	0.53
1:A:42:HIS:CG	2:A:2017:HOH:O	2.49	0.53
1:A:173:TYR:HD1	1:A:180:LYS:CA	2.21	0.53
1:B:132:PHE:HE2	1:B:163:MET:HE1	1.75	0.52
1:A:55:GLU:HG2	1:A:60:GLN:HG3	1.92	0.52
1:B:167:GLU:C	1:B:169:TYR:H	2.12	0.52
1:A:136:VAL:HG13	1:A:148:LEU:HD21	1.91	0.52
1:B:64:SER:O	1:B:68:VAL:HG23	2.09	0.52
1:A:103:ARG:O	1:A:106:PHE:N	2.38	0.51
1:B:109:LEU:C	1:B:111:GLY:H	2.13	0.51
1:B:32:VAL:CG1	1:B:34:LEU:HD21	2.41	0.51
1:B:112:MET:O	1:B:113:ASP:C	2.49	0.51
1:B:117:LEU:HD12	1:B:117:LEU:N	2.26	0.51
1:A:121:GLU:O	1:A:126:GLU:HG3	2.11	0.50
1:A:58:GLY:C	1:A:59:LYS:HD3	2.32	0.50
1:A:184:GLU:OE2	2:A:2074:HOH:O	2.20	0.50
1:A:84:GLU:CG	1:A:149:VAL:HG12	2.39	0.49
1:B:55:GLU:OE1	2:B:2042:HOH:O	2.19	0.49
1:A:17:ILE:HD13	1:A:17:ILE:C	2.33	0.49
1:A:112:MET:HB3	1:A:113:ASP:OD1	2.12	0.49
1:A:94:GLN:OE1	1:B:63:GLN:NE2	2.44	0.49
1:A:129:ARG:HD2	1:A:172:LEU:O	2.14	0.48
1:A:134:THR:HA	1:A:137:THR:CG2	2.43	0.48
1:A:182:HIS:O	1:A:186:VAL:HG23	2.13	0.48
1:A:103:ARG:CB	1:A:104:PRO:CD	2.92	0.48
1:A:146:GLY:HA3	1:A:185:LYS:HE2	1.97	0.47
1:B:105:TYR:O	1:B:105:TYR:CD1	2.68	0.47
1:B:182:HIS:O	1:B:186:VAL:HG23	2.15	0.47
1:B:1:MET:SD	1:B:1:MET:N	2.83	0.46
1:B:24[A]:LEU:HD11	1:B:192:LEU:HD22	1.96	0.46
1:A:20:GLN:O	1:A:24:LEU:HD12	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:136:VAL:HG13	1:B:148:LEU:HD21	1.98	0.46
1:B:160:VAL:O	1:B:163:MET:HB2	2.15	0.46
1:A:20:GLN:HG2	1:A:192:LEU:HD13	1.98	0.45
1:B:194:LYS:O	1:B:198:THR:HB	2.17	0.45
1:A:114:GLN:O	1:A:114:GLN:HG2	2.16	0.45
1:A:108:VAL:HG13	1:A:108:VAL:O	2.16	0.45
1:A:203:LYS:HA	1:A:203:LYS:HD2	1.61	0.45
1:B:105:TYR:CD1	1:B:105:TYR:C	2.91	0.45
1:B:163:MET:O	1:B:164:GLY:C	2.55	0.44
1:B:143:ASN:OD1	1:B:147:TYR:O	2.35	0.44
1:B:147:TYR:CE1	1:B:153:LEU:HD13	2.53	0.44
1:A:65:VAL:O	1:A:68:VAL:CG2	2.66	0.44
1:A:129:ARG:NH1	1:A:172:LEU:O	2.49	0.44
1:B:169:TYR:OH	1:B:171:LYS:NZ	2.48	0.44
1:A:4:TYR:HA	1:A:55:GLU:O	2.17	0.44
1:A:106:PHE:C	1:A:107:LYS:HE2	2.38	0.43
1:B:39:TRP:CH2	1:B:43:LYS:HG3	2.52	0.43
1:B:186:VAL:O	1:B:192:LEU:HD23	2.18	0.43
1:A:20:GLN:HB3	1:A:158:LEU:HD13	2.00	0.43
1:A:103:ARG:CB	1:A:104:PRO:HD3	2.48	0.43
1:B:98:PHE:CE2	1:B:163:MET:SD	3.09	0.43
1:A:111:GLY:O	1:A:112:MET:CB	2.67	0.43
1:A:20:GLN:O	1:A:24:LEU:CD1	2.66	0.43
1:A:98:PHE:O	1:A:102:VAL:HB	2.19	0.43
1:B:143:ASN:O	1:B:144:LYS:HG2	2.18	0.43
1:B:84:GLU:HG2	1:B:149:VAL:CG2	2.48	0.43
1:B:55:GLU:CD	2:B:2042:HOH:O	2.58	0.42
1:A:118:LYS:O	1:A:122:LYS:HG2	2.19	0.42
1:B:108:VAL:CG2	1:B:109:LEU:H	2.09	0.42
1:B:12:ARG:HB3	1:B:16:GLU:HB2	2.02	0.42
1:A:95:PHE:O	1:A:98:PHE:HB3	2.20	0.42
1:B:104:PRO:HA	1:B:110:LEU:CD1	2.50	0.42
1:A:60:GLN:HE21	1:A:60:GLN:HB2	1.60	0.42
1:A:103:ARG:HB2	1:A:104:PRO:CD	2.49	0.42
1:B:12:ARG:NH1	1:B:202:SER:OG	2.52	0.42
1:A:3:HIS:HB3	2:A:2003:HOH:O	2.20	0.41
1:B:84:GLU:HG2	1:B:149:VAL:HG23	2.02	0.41
1:A:53:VAL:CG1	1:A:60:GLN:HG2	2.51	0.41
1:A:187:ARG:HA	1:A:192:LEU:CD2	2.46	0.41
1:B:116:ASP:HA	1:B:117:LEU:HD12	2.02	0.41
1:B:117:LEU:HD12	1:B:117:LEU:H	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:19:ARG:HH11	1:B:19:ARG:HD2	1.73	0.41
1:B:135:ILE:HG22	1:B:136:VAL:N	2.36	0.41
1:B:125:PHE:CE2	1:B:129:ARG:HD3	2.56	0.40
1:B:109:LEU:HD13	1:B:113:ASP:N	2.36	0.40
1:A:134:THR:HG22	1:A:176:PHE:HE1	1.86	0.40
1:A:39:TRP:CZ3	1:A:51:LEU:HG	2.57	0.40
1:B:109:LEU:HD13	1:B:113:ASP:CA	2.50	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:2008:HOH:O	2:B:2072:HOH:O[1_655]	2.01	0.19

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	203/204 (100%)	179 (88%)	18 (9%)	6 (3%)	4 1
1	B	203/204 (100%)	177 (87%)	16 (8%)	10 (5%)	2 0
All	All	406/408 (100%)	356 (88%)	34 (8%)	16 (4%)	3 1

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	112	MET
1	B	106	PHE
1	B	114	GLN
1	B	118	LYS
1	B	164	GLY
1	A	57	ASP

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Mol	Chain	Res	Type
1	A	108	VAL
1	A	110	LEU
1	B	108	VAL
1	B	13	GLY
1	A	63	GLN
1	B	113	ASP
1	B	172	LEU
1	A	174	ASP
1	B	167	GLU
1	B	63	GLN

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	176/174 (101%)	152 (86%)	24 (14%)	3 2
1	B	175/174 (101%)	147 (84%)	28 (16%)	2 1
All	All	351/348 (101%)	299 (85%)	52 (15%)	3 1

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ASN
1	A	17	ILE
1	A	21	VAL
1	A	47	PRO
1	A	59	LYS
1	A	93	ASP
1	A	101	GLU
1	A	102	VAL
1	A	103	ARG
1	A	104	PRO
1	A	106	PHE
1	A	112	MET
1	A	113	ASP

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Mol	Chain	Res	Type
1	A	118	LYS
1	A	126	GLU
1	A	131	LYS
1	A	137	THR
1	A	144	LYS
1	A	151	ASP
1	A	160	VAL
1	A	163	MET
1	A	173	TYR
1	A	196	ILE
1	A	203	LYS
1	B	1	MET
1	B	17	ILE
1	B	28	ASP
1	B	47	PRO
1	B	59	LYS
1	B	96	LYS
1	B	99	LEU
1	B	101	GLU
1	B	102	VAL
1	B	105	TYR
1	B	108	VAL
1	B	109	LEU
1	B	114	GLN
1	B	116	ASP
1	B	117	LEU
1	B	122	LYS
1	B	124	VAL
1	B	131	LYS
1	B	149	VAL
1	B	151	ASP
1	B	152	SER
1	B	163	MET
1	B	168	HIS
1	B	169	TYR
1	B	187	ARG
1	B	188	SER
1	B	196	ILE
1	B	198	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	GLN
1	A	189	ASN
1	B	130	GLN
1	B	143	ASN
1	B	168	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 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)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	204/204 (100%)	0.60	9 (4%)	34	33	20, 29, 41, 57	0
1	B	204/204 (100%)	0.70	18 (8%)	10	9	21, 30, 44, 55	0
All	All	408/408 (100%)	0.65	27 (6%)	18	17	20, 30, 43, 57	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	113	ASP	3.9
1	B	120	LEU	3.6
1	B	165	PHE	3.4
1	A	165	PHE	3.4
1	B	112	MET	3.2
1	B	173	TYR	3.1
1	B	123	ASP	2.9
1	A	204[A]	PHE	2.9
1	A	48	PHE	2.8
1	B	204	PHE	2.8
1	A	115	GLY	2.8
1	B	169	TYR	2.7
1	B	175	GLY	2.7
1	B	124	VAL	2.7
1	B	1	MET	2.4
1	A	23	VAL	2.4
1	B	103	ARG	2.4
1	B	115	GLY	2.3
1	A	170	PRO	2.1
1	A	198	THR	2.1
1	B	108	VAL	2.1
1	B	116	ASP	2.1
1	A	112	MET	2.0
1	A	109	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	128	ALA	2.0
1	B	202	SER	2.0
1	B	83	TRP	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.