



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 23, 2023 – 08:53 PM EDT

PDB ID : 3B3L  
Title : Crystal structures of alternatively-spliced isoforms of human ketohexokinase  
Authors : Trinh, C.H.; Asipu, A.; Bonthron, D.T.; Phillips, S.E.V.  
Deposited on : 2007-10-22  
Resolution : 2.90 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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) : 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

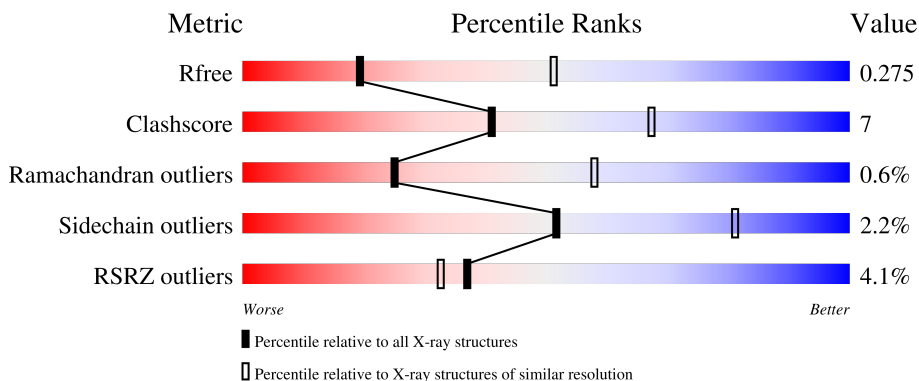
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*


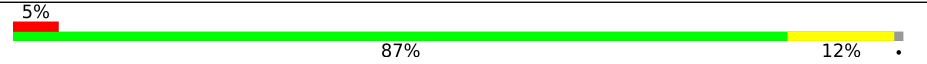
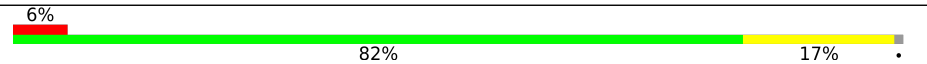

The reported resolution of this entry is 2.90 Å.

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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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  $\geq 3$ , 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  $\leq 5\%$ . 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.

Mol	Chain	Length	Quality of chain
1	A	298	 85% 13% ..
1	B	298	 87% 12% .
1	C	298	 82% 17% .
1	D	298	 79% 19% ..

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8916 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 Ketohexokinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	296	2232	1399	399	422	12	0	0	0
1	B	296	2201	1382	392	415	12	0	0	0
1	C	296	2230	1399	399	420	12	0	0	0
1	D	296	2235	1401	398	424	12	0	0	0

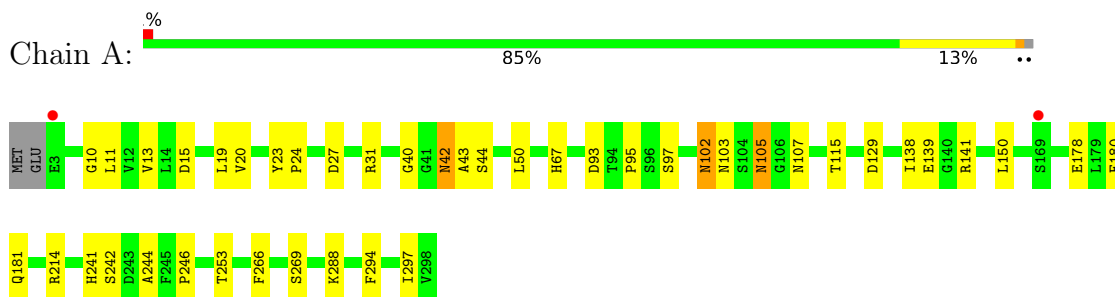
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	7	Total 7	O 7	0	0
2	B	2	Total 2	O 2	0	0
2	C	6	Total 6	O 6	0	0
2	D	3	Total 3	O 3	0	0

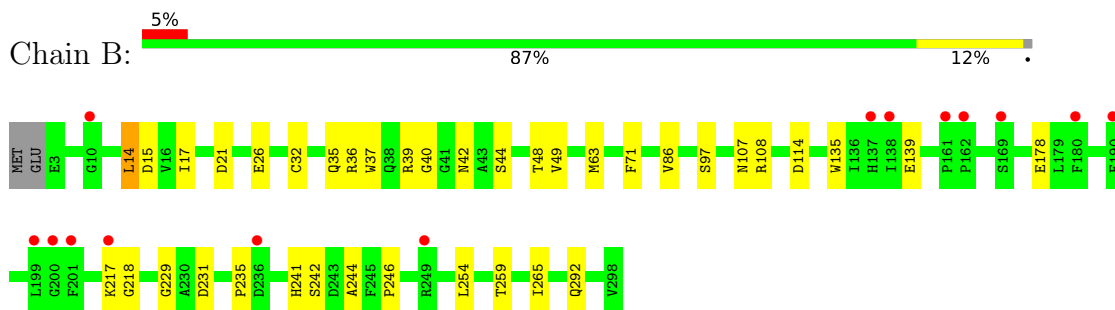
### 3 Residue-property plots [i](#)

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.

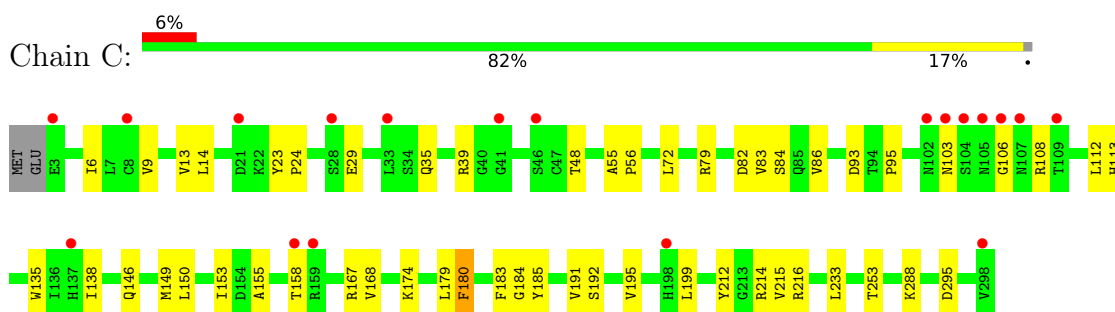
- Molecule 1: Ketohexokinase



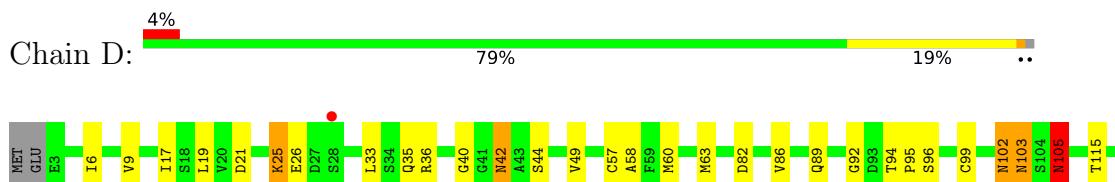
- Molecule 1: Ketohexokinase

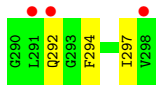
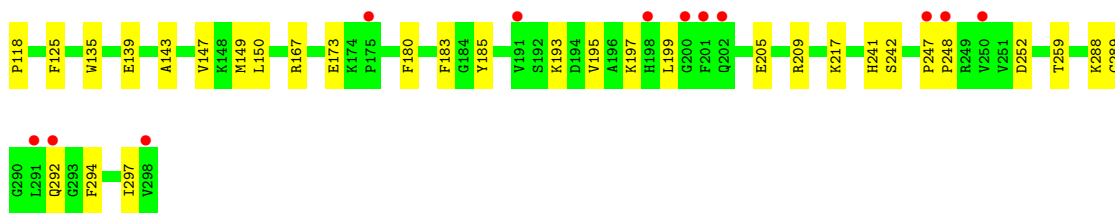


- Molecule 1: Ketohexokinase



- Molecule 1: Ketohexokinase





## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.56Å 140.70Å 179.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.38 – 2.90 25.38 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.6 (25.38-2.90) 99.8 (25.38-2.90)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.91 (at 2.89Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.236 , 0.281 0.230 , 0.275	Depositor DCC
$R_{free}$ test set	2608 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	80.4	Xtrriage
Anisotropy	0.144	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 63.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	8916	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	63.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.77% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 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	0.69	0/2274	0.64	0/3080
1	B	0.53	0/2243	0.58	0/3043
1	C	0.60	0/2272	0.60	0/3077
1	D	0.56	0/2277	0.58	0/3083
All	All	0.60	0/9066	0.60	0/12283

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

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	2232	0	2175	34	0
1	B	2201	0	2131	22	0
1	C	2230	0	2180	37	0
1	D	2235	0	2183	41	0
2	A	7	0	0	0	0
2	B	2	0	0	0	0
2	C	6	0	0	0	0
2	D	3	0	0	0	0
All	All	8916	0	8669	125	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 7.

All (125) 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:D:21:ASP:HA	1:D:103:ASN:HB2	1.52	0.88
1:C:138:ILE:HD12	1:C:150:LEU:HD21	1.62	0.81
1:D:36:ARG:NH1	1:D:292:GLN:HE21	1.80	0.78
1:D:19:LEU:HD22	1:D:33:LEU:HD12	1.72	0.70
1:A:31:ARG:HD3	1:B:114:ASP:OD1	1.92	0.70
1:A:42:ASN:H	1:A:42:ASN:HD22	1.44	0.66
1:D:17:ILE:O	1:D:35:GLN:HA	1.98	0.63
1:D:42:ASN:HB2	1:D:139:GLU:CD	2.18	0.63
1:A:13:VAL:HG22	1:A:95:PRO:HB2	1.80	0.63
1:D:195:VAL:O	1:D:199:LEU:HG	1.99	0.63
1:A:42:ASN:H	1:A:42:ASN:ND2	1.97	0.62
1:C:106:GLY:O	1:C:253:THR:HG23	1.99	0.62
1:B:244:ALA:O	1:B:246:PRO:HD3	1.99	0.62
1:A:107:ASN:ND2	1:B:26:GLU:OE2	2.33	0.61
1:C:35:GLN:NE2	1:D:96:SER:OG	2.34	0.60
1:D:63:MET:HE2	1:D:86:VAL:HG21	1.84	0.59
1:D:92:GLY:HA2	1:D:115:THR:HG21	1.85	0.58
1:A:102:ASN:HD22	1:A:103:ASN:N	2.01	0.58
1:A:105:ASN:N	1:A:105:ASN:OD1	2.32	0.58
1:D:288:LYS:HD2	1:D:288:LYS:O	2.04	0.57
1:A:115:THR:HG22	1:A:115:THR:O	2.05	0.57
1:C:29:GLU:N	1:C:29:GLU:OE2	2.36	0.57
1:B:229:GLY:HA3	1:B:242:SER:O	2.05	0.56
1:D:241:HIS:CG	1:D:242:SER:N	2.74	0.55
1:D:17:ILE:HG13	1:D:99:CYS:HB3	1.89	0.55
1:D:40:GLY:O	1:D:44:SER:HB3	2.05	0.55
1:A:107:ASN:HD22	1:B:26:GLU:CD	2.10	0.55
1:A:11:LEU:HD21	1:A:141:ARG:CD	2.37	0.54
1:D:25:LYS:O	1:D:25:LYS:HD3	2.07	0.54
1:C:214:ARG:HB3	1:C:214:ARG:HH11	1.73	0.54
1:A:20:VAL:O	1:A:103:ASN:HB2	2.07	0.54
1:D:6:ILE:HG23	1:D:135:TRP:HB3	1.89	0.54
1:A:138:ILE:HD12	1:A:150:LEU:HD21	1.90	0.53
1:C:135:TRP:HA	1:C:167:ARG:O	2.08	0.53
1:B:15:ASP:OD1	1:B:97:SER:OG	2.24	0.53
1:A:178:GLU:O	1:A:181:GLN:HG2	2.09	0.52
1:B:63:MET:CE	1:B:86:VAL:HG21	2.40	0.52
1:C:138:ILE:HD12	1:C:150:LEU:CD2	2.36	0.52
1:D:42:ASN:HD21	1:D:173:GLU:HG3	1.74	0.52
1:D:294:PHE:O	1:D:297:ILE:HG12	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:42:ASN:HB2	1:A:139:GLU:OE1	2.10	0.51
1:C:180:PHE:CE2	1:C:199:LEU:HD21	2.46	0.51
1:A:10:GLY:HA3	1:A:43:ALA:HB2	1.93	0.50
1:C:192:SER:OG	1:C:195:VAL:HG23	2.11	0.50
1:A:266:PHE:O	1:A:269:SER:HB3	2.11	0.50
1:B:42:ASN:HB2	1:B:139:GLU:CD	2.32	0.50
1:C:112:LEU:HD12	1:C:113:HIS:H	1.77	0.50
1:A:67:HIS:HB3	1:B:37:TRP:CZ3	2.47	0.50
1:C:48:THR:HG23	1:C:79:ARG:HD2	1.93	0.49
1:C:6:ILE:HG23	1:C:135:TRP:HB3	1.93	0.49
1:D:58:ALA:HB2	1:D:82:ASP:HB3	1.93	0.49
1:A:23:TYR:HB2	1:B:26:GLU:OE1	2.13	0.49
1:C:13:VAL:HG22	1:C:95:PRO:HB2	1.93	0.49
1:C:180:PHE:HD1	1:C:180:PHE:N	2.11	0.49
1:C:153:ILE:HG21	1:C:168:VAL:HG21	1.94	0.48
1:C:83:VAL:O	1:C:86:VAL:HG23	2.13	0.48
1:A:27:ASP:HB2	1:B:107:ASN:HD22	1.79	0.48
1:B:135:TRP:CE2	1:B:265:ILE:HG23	2.49	0.47
1:C:180:PHE:N	1:C:180:PHE:CD1	2.80	0.47
1:B:17:ILE:HG21	1:B:36:ARG:NH1	2.29	0.47
1:C:150:LEU:HD13	1:C:185:TYR:HB2	1.96	0.47
1:D:63:MET:CE	1:D:86:VAL:HG21	2.44	0.47
1:C:9:VAL:HB	1:C:138:ILE:HG12	1.97	0.47
1:B:14:LEU:C	1:B:14:LEU:HD12	2.35	0.47
1:A:15:ASP:OD1	1:A:97:SER:OG	2.33	0.47
1:B:218:GLY:HA2	1:B:235:PRO:HB3	1.96	0.46
1:C:23:TYR:HA	1:C:24:PRO:HD3	1.74	0.46
1:C:146:GLN:HG2	1:C:149:MET:HE3	1.97	0.46
1:D:49:VAL:HG21	1:D:259:THR:HA	1.97	0.46
1:D:150:LEU:HD13	1:D:185:TYR:CB	2.46	0.46
1:C:39:ARG:HG3	1:C:72:LEU:HD22	1.98	0.46
1:A:102:ASN:HD22	1:A:103:ASN:H	1.64	0.46
1:D:25:LYS:HD3	1:D:25:LYS:C	2.36	0.46
1:C:48:THR:HG23	1:C:79:ARG:CD	2.46	0.45
1:C:184:GLY:O	1:C:216:ARG:HB3	2.17	0.45
1:A:294:PHE:O	1:A:297:ILE:HG23	2.16	0.45
1:B:36:ARG:HH12	1:B:292:GLN:HE21	1.65	0.45
1:A:253:THR:O	1:A:253:THR:CG2	2.64	0.45
1:D:252:ASP:O	1:D:289:CYS:HA	2.17	0.45
1:D:125:PHE:CB	1:D:149:MET:HG2	2.48	0.44
1:B:241:HIS:CG	1:B:242:SER:N	2.85	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:42:ASN:ND2	1:D:173:GLU:HG3	2.32	0.44
1:B:40:GLY:O	1:B:44:SER:HB3	2.18	0.43
1:C:23:TYR:HB2	1:D:26:GLU:OE1	2.19	0.43
1:A:42:ASN:ND2	1:A:42:ASN:N	2.66	0.43
1:B:39:ARG:NH2	1:B:48:THR:OG1	2.50	0.43
1:D:89:GLN:OE1	1:D:118:PRO:HG2	2.18	0.43
1:A:10:GLY:HA3	1:A:43:ALA:CB	2.48	0.43
1:A:67:HIS:CE1	1:B:71:PHE:HE1	2.35	0.43
1:D:205:GLU:HB3	1:D:209:ARG:HH12	1.84	0.43
1:B:49:VAL:HG21	1:B:259:THR:HA	2.00	0.43
1:D:9:VAL:HA	1:D:60:MET:O	2.19	0.43
1:C:174:LYS:HB2	1:C:179:LEU:HD11	2.00	0.43
1:D:6:ILE:O	1:D:57:CYS:HB2	2.19	0.43
1:D:150:LEU:HD13	1:D:185:TYR:HB2	2.00	0.43
1:D:193:LYS:HE2	1:D:197:LYS:HD2	1.99	0.43
1:A:23:TYR:HA	1:A:24:PRO:HD3	1.91	0.42
1:A:50:LEU:HD23	1:A:50:LEU:HA	1.84	0.42
1:D:94:THR:HG22	1:D:95:PRO:HD2	2.01	0.42
1:C:183:PHE:O	1:C:215:VAL:HA	2.19	0.42
1:C:55:ALA:HA	1:C:56:PRO:HD2	1.94	0.42
1:A:129:ASP:OD1	1:A:129:ASP:C	2.58	0.42
1:D:115:THR:HG22	1:D:115:THR:O	2.20	0.42
1:A:244:ALA:C	1:A:246:PRO:HD3	2.39	0.42
1:C:82:ASP:OD1	1:C:84:SER:HB3	2.19	0.42
1:C:155:ALA:O	1:C:158:THR:HB	2.19	0.42
1:C:214:ARG:HB3	1:C:214:ARG:NH1	2.33	0.42
1:D:217:LYS:HE2	1:D:217:LYS:HB3	1.83	0.42
1:A:40:GLY:O	1:A:44:SER:HB3	2.19	0.42
1:D:102:ASN:HD22	1:D:103:ASN:H	1.68	0.42
1:A:180:PHE:CD2	1:A:214:ARG:NH1	2.88	0.42
1:C:212:TYR:HB2	1:C:233:LEU:HD23	2.01	0.42
1:D:247:PRO:O	1:D:248:PRO:C	2.58	0.41
1:C:82:ASP:OD1	1:C:84:SER:CB	2.68	0.41
1:D:143:ALA:O	1:D:147:VAL:HG23	2.20	0.41
1:C:146:GLN:HA	1:C:149:MET:CE	2.51	0.41
1:C:191:VAL:O	1:C:191:VAL:HG13	2.21	0.41
1:D:102:ASN:HB3	1:D:105:ASN:HB2	2.03	0.41
1:B:217:LYS:HE2	1:B:217:LYS:HB3	1.87	0.40
1:D:180:PHE:HD1	1:D:183:PHE:HE1	1.69	0.40
1:C:146:GLN:O	1:C:149:MET:HB2	2.21	0.40
1:D:105:ASN:OD1	1:D:105:ASN:N	2.52	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:153:ILE:HG21	1:C:168:VAL:CG2	2.51	0.40
1:A:11:LEU:HD21	1:A:141:ARG:NE	2.36	0.40
1:A:241:HIS:CG	1:A:242:SER:N	2.90	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	294/298 (99%)	283 (96%)	11 (4%)	0	100	100
1	B	294/298 (99%)	276 (94%)	15 (5%)	3 (1%)	15	45
1	C	294/298 (99%)	274 (93%)	19 (6%)	1 (0%)	41	71
1	D	294/298 (99%)	276 (94%)	15 (5%)	3 (1%)	15	45
All	All	1176/1192 (99%)	1109 (94%)	60 (5%)	7 (1%)	25	58

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	21	ASP
1	C	295	ASP
1	D	42	ASN
1	D	103	ASN
1	B	178	GLU
1	B	254	LEU
1	D	105	ASN

### 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	236/248 (95%)	230 (98%)	6 (2%)	47	78
1	B	230/248 (93%)	225 (98%)	5 (2%)	52	81
1	C	236/248 (95%)	230 (98%)	6 (2%)	47	78
1	D	238/248 (96%)	234 (98%)	4 (2%)	60	86
All	All	940/992 (95%)	919 (98%)	21 (2%)	52	81

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

Mol	Chain	Res	Type
1	A	19	LEU
1	A	42	ASN
1	A	93	ASP
1	A	102	ASN
1	A	105	ASN
1	A	288	LYS
1	B	14	LEU
1	B	32	CYS
1	B	35	GLN
1	B	108	ARG
1	B	231	ASP
1	C	14	LEU
1	C	93	ASP
1	C	103	ASN
1	C	108	ARG
1	C	180	PHE
1	C	288	LYS
1	D	25	LYS
1	D	102	ASN
1	D	105	ASN
1	D	167	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	GLN
1	A	42	ASN

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Mol	Chain	Res	Type
1	A	102	ASN
1	A	107	ASN
1	B	89	GLN
1	B	102	ASN
1	B	103	ASN
1	B	283	GLN
1	B	292	GLN
1	C	5	GLN
1	C	35	GLN
1	C	113	HIS
1	C	283	GLN
1	D	292	GLN

### 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

### 6.1 Protein, DNA and RNA chains

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	296/298 (99%)	-0.33	2 (0%) 87 87	39, 54, 77, 97	0
1	B	296/298 (99%)	-0.03	14 (4%) 31 28	48, 66, 84, 94	0
1	C	296/298 (99%)	0.13	19 (6%) 19 15	44, 63, 104, 130	0
1	D	296/298 (99%)	0.02	13 (4%) 34 30	37, 59, 82, 97	0
All	All	1184/1192 (99%)	-0.05	48 (4%) 37 32	37, 61, 86, 130	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	107	ASN	5.5
1	D	248	PRO	5.3
1	C	159	ARG	4.6
1	B	137	HIS	3.8
1	D	202	GLN	3.7
1	C	8	CYS	3.7
1	B	200	GLY	3.5
1	C	104	SER	3.4
1	D	175	PRO	3.3
1	C	158	THR	3.2
1	C	137	HIS	3.2
1	B	180	PHE	3.2
1	D	292	GLN	3.2
1	C	3	GLU	3.1
1	D	201	PHE	3.1
1	D	291	LEU	3.0
1	C	103	ASN	3.0
1	C	106	GLY	3.0
1	B	162	PRO	2.9
1	C	28	SER	2.8
1	D	247	PRO	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	169	SER	2.8
1	D	28	SER	2.7
1	C	105	ASN	2.6
1	D	198	HIS	2.6
1	B	217	LYS	2.6
1	D	250	VAL	2.6
1	C	102	ASN	2.4
1	C	33	LEU	2.4
1	C	198	HIS	2.4
1	C	46	SER	2.3
1	D	200	GLY	2.3
1	B	236	ASP	2.3
1	C	41	GLY	2.3
1	D	298	VAL	2.2
1	B	161	PRO	2.2
1	B	10	GLY	2.2
1	B	169	SER	2.2
1	C	298	VAL	2.2
1	D	191	VAL	2.2
1	B	201	PHE	2.1
1	A	3	GLU	2.1
1	C	109	THR	2.1
1	B	249	ARG	2.1
1	B	138	ILE	2.0
1	B	190	PHE	2.0
1	C	21	ASP	2.0
1	B	199	LEU	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.