



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:35 PM UTC

PDB ID : 4YWS / pdb_00004yws
Title : Thermostable enolase from *Chloroflexus aurantiacus*
Authors : Zadvornyy, O.A.; Peters, J.W.
Deposited on : 2015-03-20
Resolution : 2.45 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

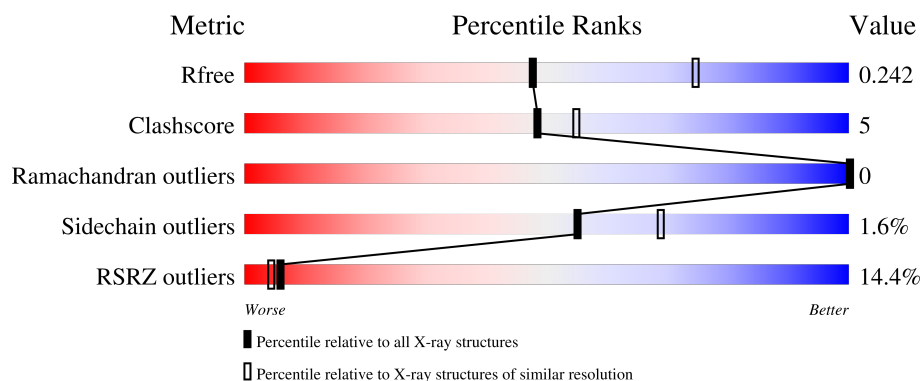
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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}	180053	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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 $\leq 5\%$. 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	426	
1	B	426	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6154 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 Enolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	405	Total	C	N	O	S	0	0	0
			3071	1935	537	591	8			
1	B	393	Total	C	N	O	S	0	0	0
			2986	1884	521	573	8			

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
2	B	1	Total	Mg	0	0
			1	1		

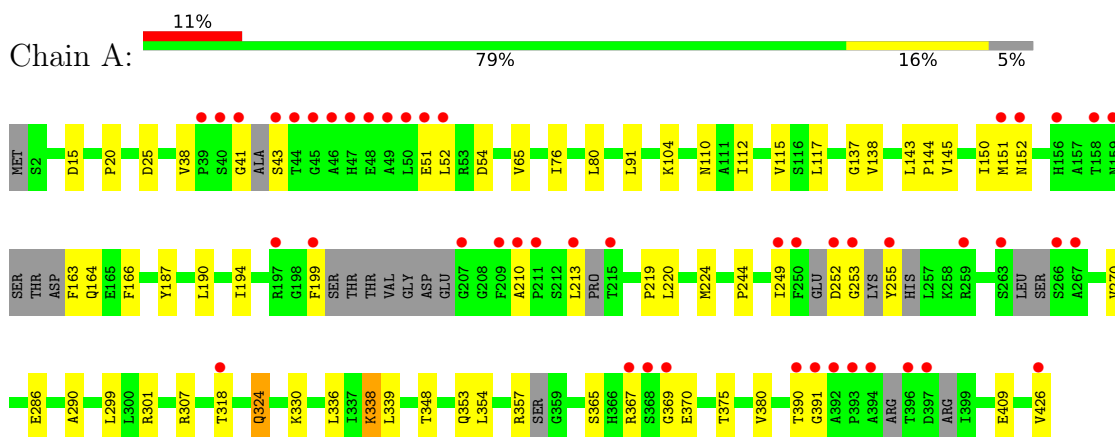
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	57	Total	O	0	0
			57	57		
3	B	38	Total	O	0	0
			38	38		

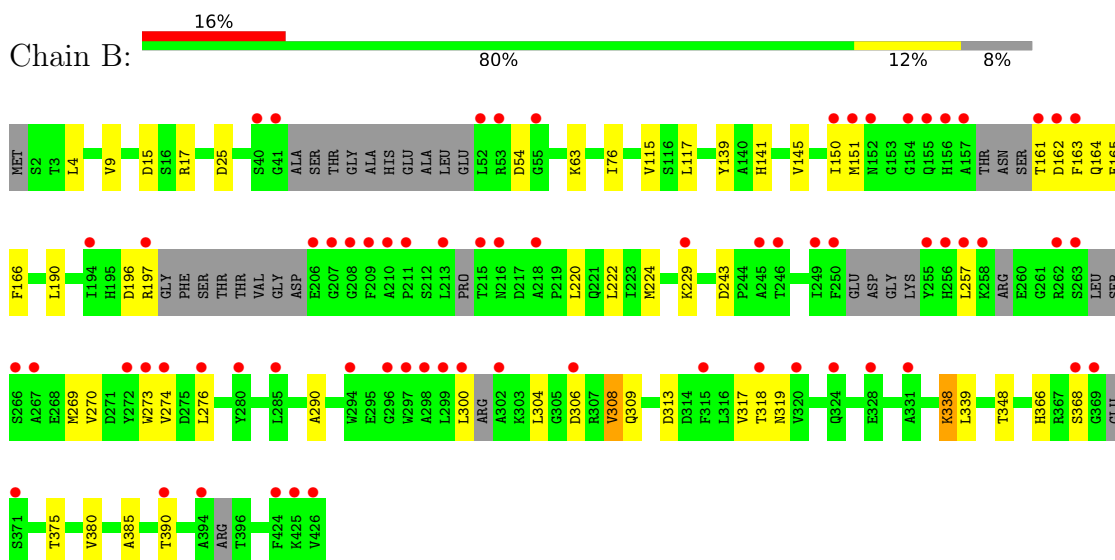
3 Residue-property plots [i](#)

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: Enolase



• Molecule 1: Enolase



4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α , β , γ	146.32Å 146.32Å 102.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.76 – 2.45 37.76 – 2.45	Depositor EDS
% Data completeness (in resolution range)	99.7 (37.76-2.45) 99.7 (37.76-2.45)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.95 (at 2.45Å)	Xtriage
Refinement program	BUSTER 2.10.1	Depositor
R, R_{free}	0.199 , 0.240 (Not available) , 0.242	Depositor DCC
R_{free} test set	2015 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	46.0	Xtriage
Anisotropy	0.478	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 47.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.026 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6154	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.93% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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

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

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	0.88	0/3109	1.36	14/4196 (0.3%)
1	B	0.87	1/3024 (0.0%)	1.34	15/4084 (0.4%)
All	All	0.87	1/6133 (0.0%)	1.35	29/8280 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	317	VAL	CA-C	5.23	1.59	1.52

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	290	ALA	CA-C-N	7.89	132.65	120.82
1	A	290	ALA	C-N-CA	7.89	132.65	120.82
1	B	15	ASP	CA-CB-CG	7.18	119.78	112.60
1	A	15	ASP	CA-CB-CG	6.93	119.53	112.60
1	A	338	LYS	CA-C-N	6.69	129.57	120.54
1	A	338	LYS	C-N-CA	6.69	129.57	120.54
1	A	137	GLY	CA-C-N	6.53	130.79	120.47
1	A	137	GLY	C-N-CA	6.53	130.79	120.47
1	A	145	VAL	N-CA-C	-6.36	101.34	107.76
1	A	164	GLN	N-CA-C	6.30	117.84	110.97
1	B	25	ASP	CA-CB-CG	6.29	118.89	112.60
1	B	164	GLN	N-CA-C	5.89	117.37	111.07
1	B	290	ALA	CA-C-N	5.70	128.19	120.38
1	B	290	ALA	C-N-CA	5.70	128.19	120.38
1	B	54	ASP	N-CA-C	5.57	117.15	111.14
1	B	306	ASP	N-CA-C	5.53	119.75	112.89
1	B	313	ASP	CA-CB-CG	5.52	118.12	112.60
1	B	63	LYS	CA-C-N	5.34	125.17	120.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	63	LYS	C-N-CA	5.34	125.17	120.10
1	A	54	ASP	N-CA-C	5.33	117.17	111.36
1	B	338	LYS	CA-C-N	5.28	127.67	120.54
1	B	338	LYS	C-N-CA	5.28	127.67	120.54
1	A	163	PHE	CA-C-N	5.16	127.46	120.65
1	A	163	PHE	C-N-CA	5.16	127.46	120.65
1	A	290	ALA	N-CA-C	5.14	117.16	110.43
1	B	145	VAL	N-CA-C	-5.12	102.59	107.76
1	A	25	ASP	CA-CB-CG	5.04	117.64	112.60
1	B	319	ASN	CA-C-N	5.03	127.35	120.46
1	B	319	ASN	C-N-CA	5.03	127.35	120.46

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	3071	0	3066	37	0
1	B	2986	0	2990	29	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	57	0	0	0	0
3	B	38	0	0	0	0
All	All	6154	0	6056	66	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 5.

All (66) 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:196:ASP:O	1:B:197:ARG:HG3	1.74	0.87
1:B:318:THR:HG21	1:B:338:LYS:H	1.40	0.87
1:A:80:LEU:HD13	1:A:91:LEU:HD11	1.57	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:318:THR:HG21	1:A:338:LYS:H	1.43	0.83
1:A:143:LEU:CD2	1:A:409:GLU:HB2	2.09	0.83
1:B:161:THR:HG23	1:B:162:ASP:N	2.03	0.72
1:B:163:PHE:CZ	1:B:273:TRP:CH2	2.78	0.71
1:B:163:PHE:CE2	1:B:273:TRP:HH2	2.09	0.70
1:A:253:GLY:O	1:A:255:TYR:N	2.25	0.69
1:A:143:LEU:HD22	1:A:409:GLU:HB2	1.72	0.69
1:A:143:LEU:HD21	1:A:409:GLU:HB2	1.74	0.69
1:B:163:PHE:CE2	1:B:273:TRP:CH2	2.80	0.69
1:B:161:THR:HG23	1:B:163:PHE:H	1.61	0.66
1:B:151:MET:HB3	1:B:166:PHE:HB2	1.78	0.65
1:A:318:THR:HG22	1:A:336:LEU:O	1.98	0.64
1:B:163:PHE:CZ	1:B:273:TRP:HH2	2.14	0.62
1:A:80:LEU:CD1	1:A:91:LEU:HD11	2.30	0.62
1:B:366:HIS:ND1	1:B:390:THR:HA	2.20	0.56
1:A:213:LEU:HD13	1:A:219:PRO:HG3	1.87	0.56
1:A:270:VAL:HG21	1:A:299:LEU:HD22	1.88	0.56
1:A:152:ASN:HB2	1:A:210:ALA:HB2	1.89	0.55
1:B:196:ASP:O	1:B:197:ARG:CG	2.51	0.54
1:A:80:LEU:HD13	1:A:91:LEU:CD1	2.33	0.54
1:B:165:GLU:HB2	1:B:243:ASP:HB3	1.88	0.54
1:A:301:ARG:HD3	1:A:330:LYS:HG2	1.91	0.53
1:B:380:VAL:HG23	1:B:385:ALA:HB3	1.89	0.53
1:B:139:TYR:O	1:B:141:HIS:CD2	2.61	0.53
1:A:318:THR:HG21	1:A:338:LYS:N	2.20	0.51
1:A:324:GLN:HG2	1:A:354:LEU:HD21	1.93	0.50
1:B:161:THR:HG23	1:B:162:ASP:H	1.76	0.50
1:A:244:PRO:HD2	1:A:286:GLU:O	2.11	0.50
1:A:336:LEU:HD21	1:A:365:SER:HB3	1.94	0.48
1:A:41:GLY:HA2	1:A:43:SER:HB3	1.94	0.48
1:B:339:LEU:HD12	1:B:348:THR:HG21	1.95	0.48
1:A:20:PRO:HB2	1:A:65:VAL:HG23	1.94	0.48
1:B:274:VAL:HA	1:B:304:LEU:HD11	1.94	0.48
1:B:117:LEU:HD22	1:B:375:THR:HG21	1.94	0.48
1:B:76:ILE:HG21	1:B:115:VAL:HG21	1.96	0.47
1:A:150:ILE:HD12	1:A:190:LEU:HD22	1.97	0.47
1:A:336:LEU:CD2	1:A:338:LYS:HG3	2.45	0.47
1:B:150:ILE:HD12	1:B:190:LEU:HD22	1.96	0.47
1:A:151:MET:HB3	1:A:166:PHE:HB2	1.96	0.46
1:A:367:ARG:HG3	1:A:369:GLY:H	1.81	0.46
1:A:76:ILE:HG21	1:A:115:VAL:HG21	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:353:GLN:O	1:A:357:ARG:HG2	2.17	0.45
1:A:144:PRO:HG3	1:A:380:VAL:HG11	1.98	0.44
1:A:367:ARG:H	1:A:370:GLU:HG2	1.82	0.44
1:A:252:ASP:O	1:A:252:ASP:OD1	2.35	0.44
1:A:38:VAL:HG11	1:A:110:ASN:HA	2.00	0.44
1:A:117:LEU:HD22	1:A:375:THR:HG21	2.00	0.44
1:A:220:LEU:O	1:A:224:MET:HG2	2.18	0.44
1:A:150:ILE:HG22	1:A:187:TYR:HD1	1.82	0.43
1:A:194:ILE:HG23	1:A:199:PHE:HB2	1.99	0.43
1:B:366:HIS:HD1	1:B:390:THR:HA	1.83	0.43
1:B:17:ARG:HD2	1:B:368:SER:HB3	2.00	0.43
1:A:307:ARG:HA	1:A:426:VAL:HG23	2.00	0.42
1:B:257:LEU:HD21	1:B:269:MET:HE1	2.01	0.42
1:A:339:LEU:CD1	1:A:348:THR:HG21	2.50	0.42
1:A:390:THR:HG22	1:A:391:GLY:H	1.84	0.42
1:B:339:LEU:CD1	1:B:348:THR:HG21	2.50	0.42
1:B:196:ASP:C	1:B:197:ARG:HG3	2.43	0.41
1:B:270:VAL:HG22	1:B:300:LEU:HB2	2.02	0.41
1:A:104:LYS:HB3	1:A:112:ILE:HD12	2.01	0.41
1:B:220:LEU:O	1:B:224:MET:HG2	2.21	0.41
1:B:308:VAL:CG1	1:B:309:GLN:N	2.82	0.41
1:B:197:ARG:HD2	1:B:222:LEU:HD21	2.03	0.41

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	382/426 (90%)	367 (96%)	15 (4%)	0	100	100
1	B	371/426 (87%)	357 (96%)	14 (4%)	0	100	100
All	All	753/852 (88%)	724 (96%)	29 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	313/332 (94%)	308 (98%)	5 (2%)	55	69
1	B	306/332 (92%)	301 (98%)	5 (2%)	55	69
All	All	619/664 (93%)	609 (98%)	10 (2%)	55	69

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	GLU
1	A	52	LEU
1	A	138	VAL
1	A	249	ILE
1	A	324	GLN
1	B	4	LEU
1	B	9	VAL
1	B	229	LYS
1	B	276	LEU
1	B	308	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	164	GLN
1	A	221	GLN
1	A	321	GLN
1	A	353	GLN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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, 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	405/426 (95%)	0.47	47 (11%) 9 8	31, 51, 113, 139	0
1	B	393/426 (92%)	0.70	68 (17%) 4 3	31, 59, 121, 139	0
All	All	798/852 (93%)	0.58	115 (14%) 6 4	31, 55, 116, 139	0

All (115) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	426	VAL	6.7
1	A	397	ASP	6.0
1	A	253	GLY	5.3
1	B	157	ALA	5.2
1	B	300	LEU	5.1
1	A	40	SER	5.0
1	B	306	ASP	4.9
1	B	163	PHE	4.8
1	B	426	VAL	4.7
1	A	396	THR	4.6
1	A	209	PHE	4.6
1	B	213	LEU	4.4
1	A	213	LEU	4.4
1	A	45	GLY	4.3
1	A	207	GLY	4.3
1	B	161	THR	4.2
1	B	250	PHE	4.1
1	A	46	ALA	4.1
1	A	44	THR	4.1
1	B	249	ILE	4.0
1	A	199	PHE	4.0
1	A	263	SER	4.0
1	B	52	LEU	3.9
1	B	209	PHE	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	51	GLU	3.8
1	B	41	GLY	3.7
1	B	369	GLY	3.7
1	A	43	SER	3.6
1	B	258	LYS	3.6
1	A	390	THR	3.6
1	B	294	TRP	3.6
1	A	394	ALA	3.5
1	B	150	ILE	3.4
1	A	215	THR	3.4
1	A	255	TYR	3.3
1	A	41	GLY	3.3
1	A	52	LEU	3.3
1	B	263	SER	3.2
1	B	194	ILE	3.2
1	A	39	PRO	3.0
1	A	50	LEU	3.0
1	A	266	SER	3.0
1	A	259	ARG	2.9
1	B	216	ASN	2.9
1	A	368	SER	2.8
1	B	257	LEU	2.8
1	B	155	GLN	2.8
1	B	266	SER	2.8
1	A	250	PHE	2.8
1	B	215	THR	2.7
1	B	255	TYR	2.7
1	B	246	THR	2.7
1	A	391	GLY	2.6
1	B	394	ALA	2.6
1	B	272	TYR	2.6
1	A	369	GLY	2.6
1	B	151	MET	2.6
1	A	47	HIS	2.6
1	B	197	ARG	2.5
1	B	274	VAL	2.5
1	A	392	ALA	2.5
1	B	262	ARG	2.5
1	B	425	LYS	2.5
1	A	318	THR	2.5
1	B	162	ASP	2.5
1	A	159	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	393	PRO	2.4
1	B	154	GLY	2.4
1	A	252	ASP	2.4
1	B	280	TYR	2.4
1	B	315	PHE	2.4
1	B	245	ALA	2.4
1	B	318	THR	2.4
1	A	49	ALA	2.4
1	B	218	ALA	2.4
1	A	367	ARG	2.3
1	B	297	TRP	2.3
1	B	207	GLY	2.3
1	B	276	LEU	2.3
1	B	331	ALA	2.3
1	B	55	GLY	2.3
1	B	285	LEU	2.3
1	A	48	GLU	2.3
1	B	320	VAL	2.3
1	A	249	ILE	2.2
1	B	229	LYS	2.2
1	B	273	TRP	2.2
1	A	267	ALA	2.2
1	A	156	HIS	2.2
1	B	206	GLU	2.2
1	B	256	HIS	2.2
1	A	151	MET	2.2
1	B	210	ALA	2.2
1	B	208	GLY	2.2
1	B	424	PHE	2.2
1	B	267	ALA	2.1
1	B	296	GLY	2.1
1	B	390	THR	2.1
1	A	210	ALA	2.1
1	B	156	HIS	2.1
1	B	328	GLU	2.1
1	B	371	SER	2.1
1	B	298	ALA	2.1
1	B	40	SER	2.1
1	B	368	SER	2.1
1	A	211	PRO	2.1
1	B	211	PRO	2.1
1	B	53	ARG	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	152	ASN	2.0
1	B	152	ASN	2.0
1	B	324	GLN	2.0
1	A	197	ARG	2.0
1	B	299	LEU	2.0
1	B	302	ALA	2.0
1	A	158	THR	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MG	A	501	1/1	0.87	0.09	51,51,51,51	0
2	MG	B	501	1/1	0.91	0.11	71,71,71,71	0

6.5 Other polymers [i](#)

There are no such residues in this entry.