



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 5, 2026 – 01:06 AM UTC

PDB ID : 9D9Q / pdb_00009d9q
Title : Crystal structure of IgG1 FC M252R at pH 8.0
Authors : Reddem, E.R.; Shapiro, L.
Deposited on : 2024-08-21
Resolution : 2.82 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

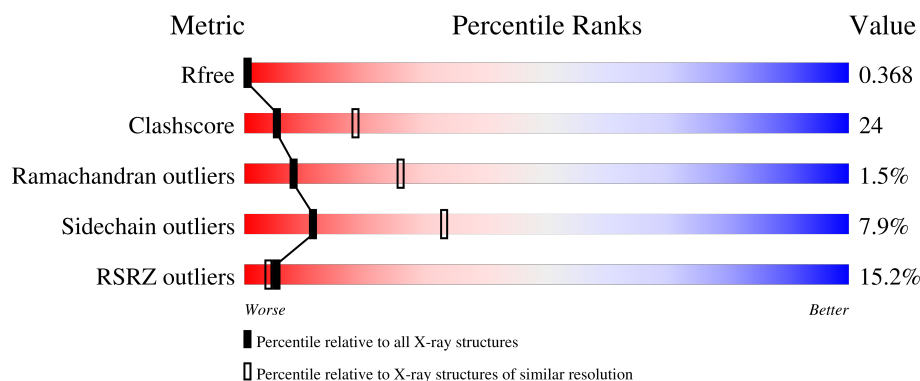
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.82 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	4591 (2.84-2.80)
Clashscore	190562	5010 (2.84-2.80)
Ramachandran outliers	187476	4916 (2.84-2.80)
Sidechain outliers	187428	4918 (2.84-2.80)
RSRZ outliers	180081	4594 (2.84-2.80)

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	231	<div> <div>8%</div> <div>42%</div> <div>42%</div> <div>5%</div> <div>11%</div> </div>
1	B	231	<div> <div>19%</div> <div>44%</div> <div>41%</div> <div>•</div> <div>12%</div> </div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3282 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 Immunoglobulin gamma-1 heavy chain Fc fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	203	Total	C	N	O	S	0	0	0
			1632	1036	278	313	5			
1	A	205	Total	C	N	O	S	0	0	0
			1650	1050	280	315	5			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	217	MET	-	initiating methionine	UNP P0DOX5
B	218	GLY	-	expression tag	UNP P0DOX5
B	219	TRP	-	expression tag	UNP P0DOX5
B	220	SER	-	expression tag	UNP P0DOX5
B	221	CYS	-	expression tag	UNP P0DOX5
B	222	ILE	-	expression tag	UNP P0DOX5
B	223	ILE	-	expression tag	UNP P0DOX5
B	224	LEU	-	expression tag	UNP P0DOX5
B	225	PHE	-	expression tag	UNP P0DOX5
B	226	LEU	-	expression tag	UNP P0DOX5
B	227	VAL	-	expression tag	UNP P0DOX5
B	228	ALA	-	expression tag	UNP P0DOX5
B	229	THR	-	expression tag	UNP P0DOX5
B	230	ALA	-	expression tag	UNP P0DOX5
B	231	THR	-	expression tag	UNP P0DOX5
B	232	GLY	-	expression tag	UNP P0DOX5
B	233	VAL	-	expression tag	UNP P0DOX5
B	234	HIS	-	expression tag	UNP P0DOX5
B	235	SER	-	expression tag	UNP P0DOX5
B	252	ARG	MET	engineered mutation	UNP P0DOX5
A	217	MET	-	initiating methionine	UNP P0DOX5
A	218	GLY	-	expression tag	UNP P0DOX5
A	219	TRP	-	expression tag	UNP P0DOX5
A	220	SER	-	expression tag	UNP P0DOX5
A	221	CYS	-	expression tag	UNP P0DOX5

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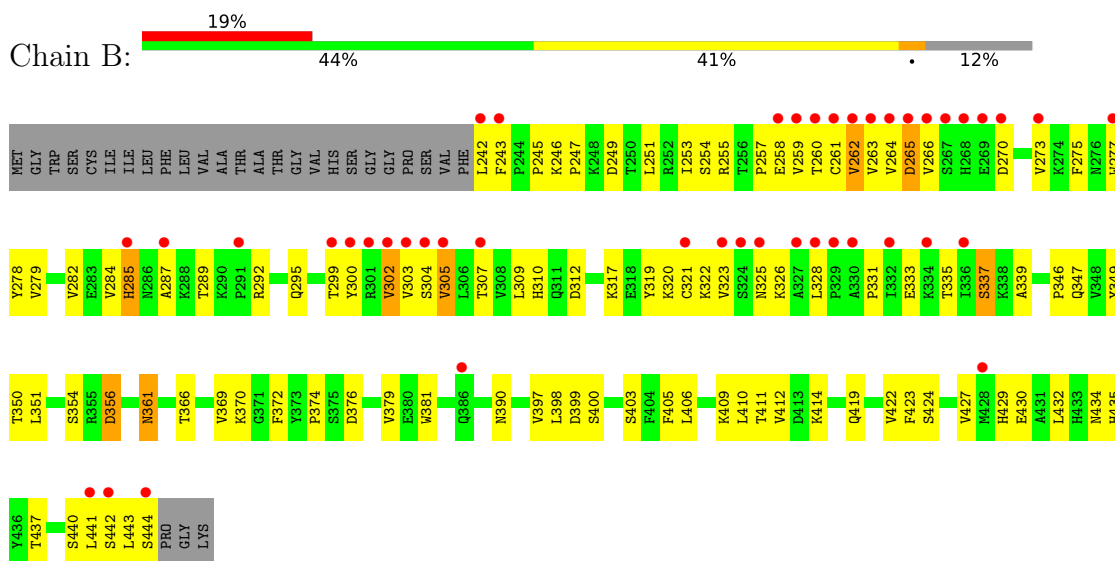
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Chain	Residue	Modelled	Actual	Comment	Reference
A	222	ILE	-	expression tag	UNP P0DOX5
A	223	ILE	-	expression tag	UNP P0DOX5
A	224	LEU	-	expression tag	UNP P0DOX5
A	225	PHE	-	expression tag	UNP P0DOX5
A	226	LEU	-	expression tag	UNP P0DOX5
A	227	VAL	-	expression tag	UNP P0DOX5
A	228	ALA	-	expression tag	UNP P0DOX5
A	229	THR	-	expression tag	UNP P0DOX5
A	230	ALA	-	expression tag	UNP P0DOX5
A	231	THR	-	expression tag	UNP P0DOX5
A	232	GLY	-	expression tag	UNP P0DOX5
A	233	VAL	-	expression tag	UNP P0DOX5
A	234	HIS	-	expression tag	UNP P0DOX5
A	235	SER	-	expression tag	UNP P0DOX5
A	252	ARG	MET	engineered mutation	UNP P0DOX5

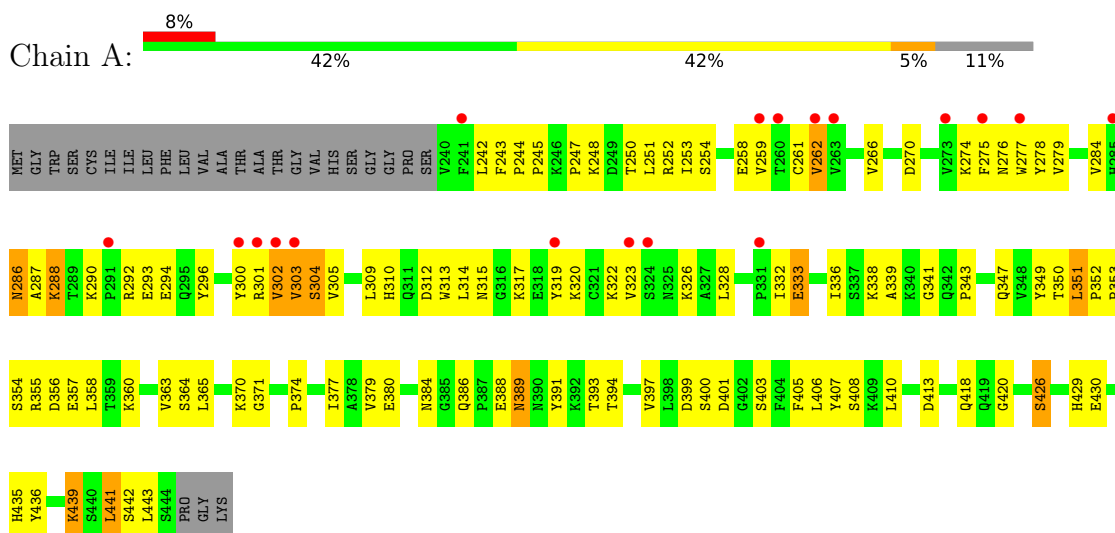
3 Residue-property plots

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: Immunoglobulin gamma-1 heavy chain Fc fragment



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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	48.67Å 78.93Å 137.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.40 – 2.82 45.40 – 2.82	Depositor EDS
% Data completeness (in resolution range)	99.5 (45.40-2.82) 99.4 (45.40-2.82)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.10 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.21.1_5286	Depositor
R, R_{free}	0.284 , 0.367 0.284 , 0.368	Depositor DCC
R_{free} test set	619 reflections (4.63%)	wwPDB-VP
Wilson B-factor (Å ²)	73.7	Xtriage
Anisotropy	0.481	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 55.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	3282	wwPDB-VP
Average B, all atoms (Å ²)	92.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.42% 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 [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	0.51	0/1695	0.67	0/2309
1	B	0.43	0/1676	0.68	0/2283
All	All	0.47	0/3371	0.68	0/4592

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	1650	0	1620	85	2
1	B	1632	0	1602	79	2
All	All	3282	0	3222	157	2

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 (157) 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:292:ARG:H	1:A:292:ARG:HD2	1.33	0.94
1:B:273:VAL:HG11	1:B:302:VAL:HG21	1.47	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:365:LEU:HD12	1:A:410:LEU:HD23	1.56	0.88
1:B:243:PHE:HB3	1:B:260:THR:H	1.42	0.85
1:A:292:ARG:HG3	1:A:302:VAL:HB	1.57	0.85
1:B:249:ASP:HA	1:B:255:ARG:HD2	1.67	0.77
1:A:351:LEU:HD13	1:A:352:PRO:HD2	1.69	0.74
1:A:320:LYS:NZ	1:A:333:GLU:OE2	2.18	0.74
1:A:338:LYS:NZ	1:A:430:GLU:OE2	2.16	0.73
1:B:270:ASP:HB3	1:B:326:LYS:HB3	1.71	0.72
1:A:312:ASP:HB3	1:A:319:TYR:OH	1.91	0.70
1:B:275:PHE:HA	1:B:323:VAL:HG12	1.76	0.67
1:B:379:VAL:HG13	1:B:427:VAL:HG22	1.76	0.67
1:B:266:VAL:HB	1:B:300:TYR:HB2	1.77	0.67
1:A:393:THR:HG22	1:A:408:SER:OG	1.94	0.66
1:B:309:LEU:HB2	1:B:312:ASP:HB2	1.77	0.65
1:A:253:ILE:HA	1:A:310:HIS:CD2	2.32	0.65
1:B:277:TRP:HZ2	1:B:304:SER:HB3	1.62	0.64
1:B:325:ASN:H	1:B:328:LEU:HD12	1.62	0.64
1:A:290:LYS:HB3	1:A:303:VAL:HG22	1.79	0.64
1:B:279:VAL:HG23	1:B:284:VAL:HG21	1.80	0.62
1:A:397:VAL:HB	1:A:405:PHE:CE1	2.36	0.61
1:A:379:VAL:HG21	1:A:406:LEU:HD11	1.83	0.61
1:A:258:GLU:HB3	1:A:305:VAL:HG13	1.82	0.60
1:B:292:ARG:HG2	1:B:300:TYR:HB3	1.82	0.60
1:A:270:ASP:OD1	1:A:326:LYS:NZ	2.34	0.60
1:A:275:PHE:CD1	1:A:304:SER:HB3	2.36	0.60
1:B:409:LYS:NZ	1:A:399:ASP:OD2	2.25	0.60
1:B:399:ASP:OD1	1:B:400:SER:N	2.32	0.59
1:A:399:ASP:OD1	1:A:400:SER:N	2.34	0.59
1:A:286:ASN:OD1	1:A:286:ASN:N	2.32	0.59
1:B:422:VAL:HA	1:B:442:SER:HB2	1.85	0.59
1:B:351:LEU:HB2	1:B:366:THR:HB	1.84	0.59
1:B:249:ASP:HB3	1:B:257:PRO:HA	1.86	0.58
1:B:292:ARG:HG3	1:B:302:VAL:HG13	1.84	0.58
1:B:285:HIS:C	1:B:287:ALA:H	2.10	0.57
1:B:397:VAL:HB	1:B:405:PHE:CE1	2.38	0.57
1:A:245:PRO:HD3	1:A:259:VAL:HG22	1.87	0.57
1:A:270:ASP:OD2	1:A:326:LYS:HB2	2.05	0.56
1:A:377:ILE:HG13	1:A:429:HIS:HB2	1.86	0.56
1:B:249:ASP:HB3	1:B:257:PRO:HB3	1.87	0.56
1:A:350:THR:HG23	1:A:439:LYS:HB3	1.88	0.56
1:A:278:TYR:HB2	1:A:320:LYS:HB3	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:275:PHE:HB3	1:B:277:TRP:CD1	2.40	0.56
1:A:389:ASN:N	1:A:389:ASN:OD1	2.38	0.56
1:A:247:PRO:O	1:A:251:LEU:HD13	2.08	0.54
1:B:273:VAL:HG23	1:B:325:ASN:HB2	1.90	0.54
1:B:249:ASP:HB3	1:B:257:PRO:CA	2.37	0.53
1:A:261:CYS:HB2	1:A:277:TRP:CZ2	2.44	0.53
1:B:412:VAL:HG21	1:B:423:PHE:CE2	2.44	0.53
1:A:293:GLU:O	1:A:300:TYR:HA	2.09	0.53
1:A:347:GLN:HB3	1:A:349:TYR:CE1	2.44	0.53
1:A:353:PRO:HD3	1:A:365:LEU:HD23	1.92	0.52
1:A:355:ARG:O	1:A:358:LEU:HG	2.09	0.52
1:A:323:VAL:HG22	1:A:332:ILE:HG23	1.92	0.51
1:A:418:GLN:C	1:A:420:GLY:H	2.17	0.51
1:B:279:VAL:O	1:B:282:VAL:HG22	2.10	0.51
1:A:279:VAL:HG23	1:A:284:VAL:HG21	1.91	0.51
1:A:394:THR:HG23	1:A:407:TYR:O	2.11	0.51
1:B:289:THR:HA	1:B:304:SER:HA	1.93	0.51
1:A:266:VAL:HG23	1:A:300:TYR:HB2	1.94	0.50
1:B:251:LEU:HG	1:B:435:HIS:ND1	2.27	0.49
1:B:278:TYR:CE2	1:B:322:LYS:HD2	2.48	0.49
1:B:253:ILE:HA	1:B:310:HIS:CD2	2.47	0.49
1:A:339:ALA:HB3	1:A:374:PRO:CB	2.42	0.48
1:B:242:LEU:HB3	1:B:243:PHE:H	1.46	0.48
1:A:326:LYS:C	1:A:328:LEU:H	2.21	0.48
1:B:423:PHE:O	1:B:441:LEU:N	2.46	0.48
1:B:349:TYR:HB3	1:A:354:SER:HB3	1.94	0.48
1:B:243:PHE:O	1:B:259:VAL:HA	2.13	0.48
1:A:341:GLY:O	1:A:343:PRO:HD3	2.13	0.48
1:A:242:LEU:HD23	1:A:243:PHE:N	2.28	0.48
1:B:289:THR:HG23	1:B:304:SER:HB2	1.94	0.48
1:A:262:VAL:HG21	1:A:301:ARG:HE	1.79	0.48
1:B:424:SER:HA	1:B:440:SER:HA	1.96	0.47
1:B:432:LEU:HD22	1:B:437:THR:HB	1.94	0.47
1:B:285:HIS:C	1:B:287:ALA:N	2.72	0.47
1:A:287:ALA:C	1:A:288:LYS:HG3	2.39	0.47
1:B:263:VAL:HB	1:B:302:VAL:O	2.15	0.47
1:B:369:VAL:O	1:B:372:PHE:HE1	1.98	0.47
1:A:244:PRO:HB3	1:A:336:ILE:HD11	1.97	0.47
1:A:248:LYS:NZ	1:A:380:GLU:OE2	2.30	0.47
1:A:338:LYS:HE3	1:A:374:PRO:HG3	1.96	0.47
1:B:243:PHE:HB3	1:B:260:THR:N	2.18	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:279:VAL:HG22	1:B:319:TYR:CD1	2.49	0.47
1:B:260:THR:CG2	1:B:303:VAL:HG13	2.45	0.46
1:B:264:VAL:O	1:B:265:ASP:HB2	2.16	0.46
1:B:249:ASP:HB3	1:B:257:PRO:CB	2.45	0.46
1:A:371:GLY:HA2	1:A:403:SER:OG	2.16	0.46
1:B:278:TYR:HE2	1:B:322:LYS:HD2	1.80	0.46
1:B:361:ASN:HA	1:B:414:LYS:HE2	1.96	0.46
1:A:288:LYS:HB2	1:A:288:LYS:HE2	1.59	0.46
1:B:398:LEU:HD12	1:B:403:SER:O	2.16	0.46
1:A:275:PHE:CE1	1:A:304:SER:HB3	2.50	0.46
1:A:292:ARG:HD2	1:A:292:ARG:N	2.15	0.46
1:A:350:THR:HB	1:A:441:LEU:HD13	1.98	0.46
1:A:339:ALA:HB3	1:A:374:PRO:HB3	1.97	0.46
1:A:242:LEU:HD22	1:A:336:ILE:HD12	1.98	0.45
1:B:260:THR:HG21	1:B:303:VAL:HG13	1.99	0.45
1:B:253:ILE:HA	1:B:310:HIS:NE2	2.31	0.45
1:A:353:PRO:HG3	1:A:363:VAL:CG2	2.45	0.45
1:B:245:PRO:HD3	1:B:259:VAL:HG22	1.97	0.45
1:B:321:CYS:SG	1:B:323:VAL:HG13	2.57	0.45
1:B:242:LEU:HG	1:B:261:CYS:HA	1.98	0.44
1:A:251:LEU:HD23	1:A:435:HIS:HB3	1.98	0.44
1:B:258:GLU:HB3	1:B:305:VAL:HG22	1.98	0.44
1:B:390:ASN:ND2	1:B:411:THR:HB	2.32	0.44
1:B:320:LYS:HE2	1:B:333:GLU:CD	2.43	0.44
1:B:339:ALA:HB3	1:B:374:PRO:HB3	1.97	0.44
1:B:409:LYS:HE2	1:A:405:PHE:CD2	2.53	0.44
1:B:251:LEU:HD23	1:B:435:HIS:HB3	1.99	0.44
1:A:288:LYS:O	1:A:304:SER:HA	2.18	0.44
1:B:379:VAL:HG21	1:B:406:LEU:HD11	1.99	0.43
1:A:347:GLN:OE1	1:A:349:TYR:OH	2.12	0.43
1:A:290:LYS:HD2	1:A:303:VAL:HG13	2.01	0.43
1:B:354:SER:CB	1:A:349:TYR:HB3	2.48	0.43
1:A:401:ASP:OD1	1:A:401:ASP:C	2.61	0.43
1:A:312:ASP:CG	1:A:317:LYS:HD2	2.44	0.43
1:A:250:THR:OG1	1:A:251:LEU:HD12	2.18	0.43
1:B:351:LEU:HD21	1:A:354:SER:HB2	2.00	0.43
1:B:409:LYS:HB3	1:B:409:LYS:HE3	1.89	0.43
1:A:242:LEU:HD21	1:A:336:ILE:HG21	2.01	0.43
1:A:439:LYS:HD3	1:A:439:LYS:HA	1.85	0.43
1:B:429:HIS:H	1:B:432:LEU:HD12	1.83	0.42
1:A:274:LYS:C	1:A:275:PHE:HD2	2.26	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:275:PHE:HZ	1:B:302:VAL:O	2.02	0.42
1:A:313:TRP:C	1:A:315:ASN:H	2.28	0.42
1:A:323:VAL:CG2	1:A:332:ILE:HG23	2.49	0.42
1:B:253:ILE:HA	1:B:253:ILE:HD12	1.90	0.42
1:B:347:GLN:HB3	1:B:349:TYR:CZ	2.55	0.42
1:A:357:GLU:OE1	1:A:364:SER:HB2	2.19	0.42
1:B:443:LEU:HD23	1:B:444:SER:N	2.34	0.42
1:B:370:LYS:HE2	1:A:357:GLU:OE2	2.20	0.42
1:B:247:PRO:HG3	1:B:376:ASP:HB3	2.01	0.41
1:A:312:ASP:O	1:A:317:LYS:HB2	2.20	0.41
1:A:391:TYR:HB3	1:A:410:LEU:CD1	2.49	0.41
1:B:317:LYS:O	1:B:337:SER:HB2	2.21	0.41
1:A:391:TYR:HB3	1:A:410:LEU:HD12	2.02	0.41
1:B:356:ASP:OD1	1:B:356:ASP:N	2.51	0.41
1:B:260:THR:O	1:B:262:VAL:HG23	2.21	0.41
1:A:276:ASN:HB3	1:A:322:LYS:HE2	2.02	0.41
1:A:279:VAL:HG22	1:A:319:TYR:CD2	2.55	0.41
1:B:253:ILE:HD12	1:B:310:HIS:CE1	2.56	0.41
1:B:350:THR:HB	1:B:441:LEU:HD22	2.03	0.41
1:B:351:LEU:CD2	1:A:354:SER:HB2	2.51	0.41
1:B:257:PRO:O	1:B:307:THR:HA	2.20	0.41
1:A:296:TYR:N	1:A:296:TYR:CD1	2.89	0.41
1:A:309:LEU:HA	1:A:309:LEU:HD13	1.88	0.41
1:A:426:SER:HB3	1:A:436:TYR:OH	2.21	0.41
1:A:443:LEU:HD12	1:A:443:LEU:HA	1.86	0.41
1:A:388:GLU:HG3	1:A:410:LEU:HD21	2.03	0.40
1:B:346:PRO:HB3	1:B:372:PHE:HB3	2.02	0.40
1:B:381:TRP:CD2	1:B:410:LEU:HB2	2.56	0.40
1:A:261:CYS:HB2	1:A:277:TRP:CH2	2.56	0.40
1:A:333:GLU:H	1:A:333:GLU:HG2	1.57	0.40
1:A:253:ILE:HA	1:A:310:HIS:NE2	2.37	0.40
1:A:242:LEU:HD23	1:A:242:LEU:C	2.46	0.40

All (2) 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)
1:B:434:ASN:O	1:A:252:ARG:NH2[3_445]	2.09	0.11
1:B:246:LYS:NZ	1:A:384:ASN:O[3_445]	2.17	0.03

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	203/231 (88%)	180 (89%)	22 (11%)	1 (0%)	24	53
1	B	201/231 (87%)	166 (83%)	30 (15%)	5 (2%)	4	15
All	All	404/462 (87%)	346 (86%)	52 (13%)	6 (2%)	8	26

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	295	GLN
1	B	265	ASP
1	A	441	LEU
1	B	331	PRO
1	B	419	GLN
1	B	262	VAL

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	192/211 (91%)	172 (90%)	20 (10%)	7	21
1	B	190/211 (90%)	180 (95%)	10 (5%)	20	50
All	All	382/422 (90%)	352 (92%)	30 (8%)	11	33

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	254	SER
1	B	285	HIS
1	B	299	THR
1	B	302	VAL
1	B	305	VAL
1	B	335	THR
1	B	337	SER
1	B	356	ASP
1	B	361	ASN
1	B	430	GLU
1	A	254	SER
1	A	262	VAL
1	A	286	ASN
1	A	288	LYS
1	A	294	GLU
1	A	302	VAL
1	A	303	VAL
1	A	304	SER
1	A	314	LEU
1	A	333	GLU
1	A	351	LEU
1	A	356	ASP
1	A	360	LYS
1	A	370	LYS
1	A	386	GLN
1	A	389	ASN
1	A	413	ASP
1	A	426	SER
1	A	439	LYS
1	A	442	SER

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

Mol	Chain	Res	Type
1	B	384	ASN
1	B	438	GLN
1	A	311	GLN
1	A	342	GLN
1	A	435	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 oligosaccharides 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, 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	205/231 (88%)	0.69	18 (8%) 15 11	45, 76, 123, 139	0
1	B	203/231 (87%)	1.14	44 (21%) 2 2	40, 81, 187, 201	0
All	All	408/462 (88%)	0.91	62 (15%) 5 4	40, 79, 182, 201	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	263	VAL	10.4
1	B	302	VAL	7.3
1	B	264	VAL	7.3
1	B	262	VAL	6.3
1	B	273	VAL	6.3
1	B	265	ASP	5.2
1	B	242	LEU	4.7
1	B	243	PHE	4.4
1	B	266	VAL	4.3
1	B	261	CYS	4.2
1	B	328	LEU	4.0
1	B	269	GLU	3.6
1	B	303	VAL	3.6
1	B	305	VAL	3.5
1	A	300	TYR	3.5
1	B	299	THR	3.4
1	A	277	TRP	3.3
1	A	323	VAL	3.3
1	B	323	VAL	3.3
1	B	285	HIS	3.2
1	A	302	VAL	3.2
1	B	267	SER	3.2
1	B	260	THR	3.1
1	A	303	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	321	CYS	3.0
1	B	301	ARG	2.9
1	B	325	ASN	2.9
1	B	332	ILE	2.9
1	B	259	VAL	2.7
1	B	268	HIS	2.6
1	A	291	PRO	2.6
1	A	241	PHE	2.6
1	B	304	SER	2.6
1	B	291	PRO	2.5
1	B	329	PRO	2.5
1	A	285	HIS	2.5
1	B	270	ASP	2.4
1	A	263	VAL	2.4
1	A	301	ARG	2.4
1	B	300	TYR	2.4
1	B	258	GLU	2.4
1	B	287	ALA	2.4
1	B	428	MET	2.3
1	B	330	ALA	2.3
1	B	277	TRP	2.3
1	A	275	PHE	2.2
1	A	260	THR	2.2
1	B	324	SER	2.2
1	A	331	PRO	2.2
1	B	334	LYS	2.2
1	A	273	VAL	2.2
1	B	307	THR	2.2
1	A	259	VAL	2.1
1	A	319	TYR	2.1
1	B	444	SER	2.1
1	A	324	SER	2.1
1	B	441	LEU	2.1
1	B	442	SER	2.0
1	B	386	GLN	2.0
1	B	336	ILE	2.0
1	B	327	ALA	2.0
1	A	262	VAL	2.0

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

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.