



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:59 AM UTC

PDB ID : 3H5W / pdb_00003h5w
Title : Crystal structure of the GluR2-ATD in space group P212121 without solvent
Authors : Jin, R.; Singh, S.K.; Gu, S.; Furukawa, H.; Sobolevsky, A.; Zhou, J.; Jin, Y.; Gouaux, E.
Deposited on : 2009-04-22
Resolution : 2.69 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

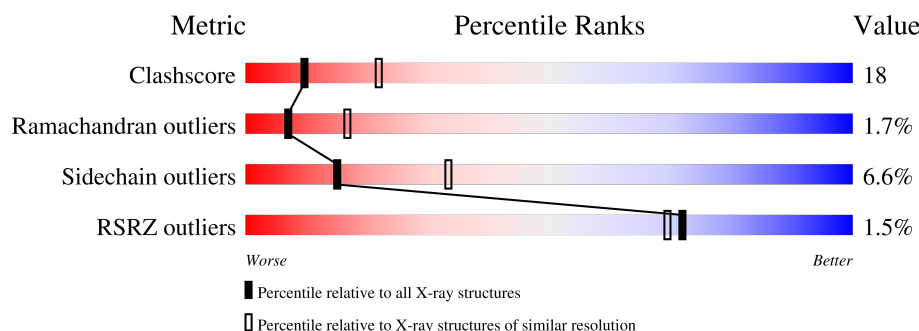
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.69 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))
Clashscore	190562	5409 (2.70-2.66)
Ramachandran outliers	187476	5324 (2.70-2.66)
Sidechain outliers	187428	5324 (2.70-2.66)
RSRZ outliers	180081	5070 (2.70-2.66)

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	394	2% 62% 28% 6%
1	B	394	2% 65% 25% 7%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 5404 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 Glutamate receptor 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	371	Total	C	N	O	S	1	0	0
			2805	1784	472	540	9			
1	B	366	Total	C	N	O	S	0	0	0
			2599	1645	438	508	8			

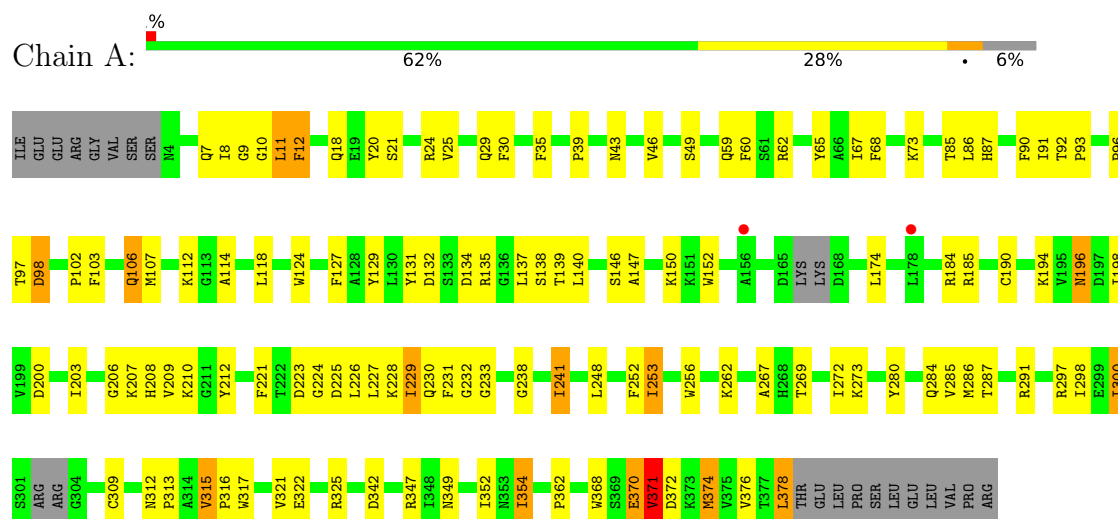
There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	ILE	-	expression tag	UNP P19491
A	-3	GLU	-	expression tag	UNP P19491
A	-2	GLU	-	expression tag	UNP P19491
A	-1	ARG	-	expression tag	UNP P19491
A	384	LEU	-	expression tag	UNP P19491
A	385	GLU	-	expression tag	UNP P19491
A	386	LEU	-	expression tag	UNP P19491
A	387	VAL	-	expression tag	UNP P19491
A	388	PRO	-	expression tag	UNP P19491
A	389	ARG	-	expression tag	UNP P19491
B	-4	ILE	-	expression tag	UNP P19491
B	-3	GLU	-	expression tag	UNP P19491
B	-2	GLU	-	expression tag	UNP P19491
B	-1	ARG	-	expression tag	UNP P19491
B	384	LEU	-	expression tag	UNP P19491
B	385	GLU	-	expression tag	UNP P19491
B	386	LEU	-	expression tag	UNP P19491
B	387	VAL	-	expression tag	UNP P19491
B	388	PRO	-	expression tag	UNP P19491
B	389	ARG	-	expression tag	UNP P19491

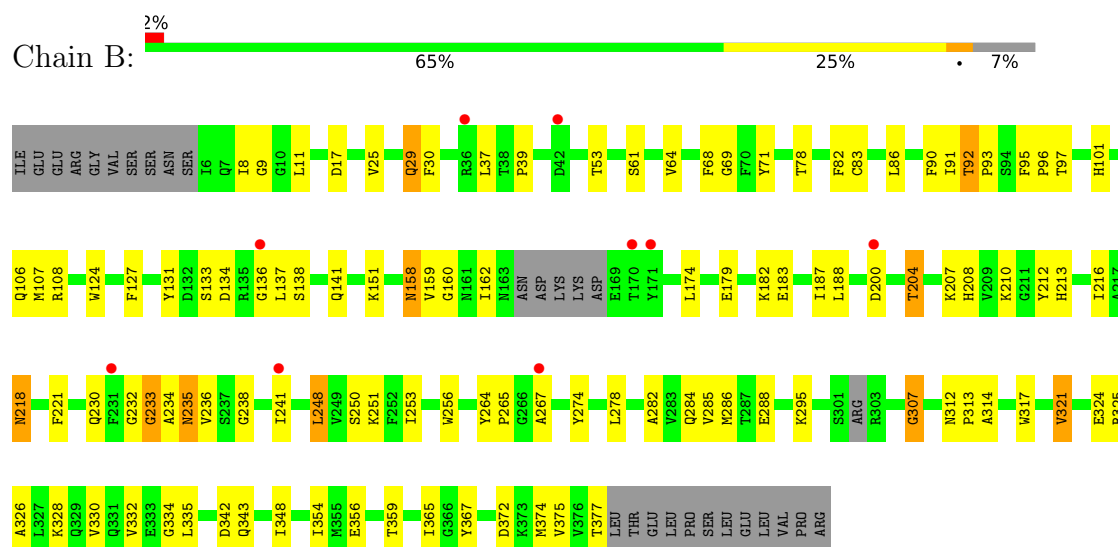
3 Residue-property plots [i](#)

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: Glutamate receptor 2



• Molecule 1: Glutamate receptor 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	85.87Å 96.36Å 108.48Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.96 – 2.69 19.96 – 2.69	Depositor EDS
% Data completeness (in resolution range)	98.6 (19.96-2.69) 93.1 (19.96-2.69)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.46 (at 2.69Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.215 , 0.291 0.220 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	71.3	Xtriage
Anisotropy	0.187	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 91.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5404	wwPDB-VP
Average B, all atoms (Å ²)	95.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.07% 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

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	0.54	0/2860	0.91	6/3878 (0.2%)
1	B	0.48	0/2651	0.90	6/3612 (0.2%)
All	All	0.52	0/5511	0.91	12/7490 (0.2%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	101	HIS	CA-C-N	8.90	128.93	119.76
1	B	101	HIS	C-N-CA	8.90	128.93	119.76
1	A	224	GLY	N-CA-C	6.25	121.59	113.27
1	A	223	ASP	N-CA-C	-6.00	101.53	110.23
1	A	315	VAL	CA-C-N	5.99	126.19	119.90
1	A	315	VAL	C-N-CA	5.99	126.19	119.90
1	B	78	THR	N-CA-C	-5.44	105.25	111.07
1	B	92	THR	CA-C-N	5.26	126.29	120.45
1	B	92	THR	C-N-CA	5.26	126.29	120.45
1	B	374	MET	N-CA-C	5.11	118.04	110.48
1	A	12	PHE	CA-C-N	5.07	124.83	119.76
1	A	12	PHE	C-N-CA	5.07	124.83	119.76

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	2805	0	2639	97	0
1	B	2599	0	2254	85	0
All	All	5404	0	4893	182	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 18.

All (182) 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:230:GLN:HE21	1:A:232:GLY:H	1.01	0.95
1:A:97:THR:HG23	1:A:106:GLN:HE22	1.31	0.93
1:A:241:ILE:CD1	1:A:352:ILE:HG12	2.02	0.89
1:A:35:PHE:HE2	1:A:291:ARG:HG2	1.38	0.89
1:B:282:ALA:O	1:B:286:MET:HG3	1.73	0.88
1:A:35:PHE:CE2	1:A:291:ARG:HG2	2.14	0.82
1:A:97:THR:HG23	1:A:106:GLN:NE2	1.93	0.82
1:A:371:VAL:O	1:A:371:VAL:HG23	1.81	0.80
1:A:210:LYS:HG2	1:A:233:GLY:HA3	1.66	0.77
1:A:230:GLN:NE2	1:A:232:GLY:H	1.80	0.77
1:B:37:LEU:O	1:B:39:PRO:HD3	1.84	0.77
1:A:85:THR:HG22	1:A:86:LEU:HD23	1.65	0.76
1:A:349:ASN:H	1:A:370:GLU:HG2	1.49	0.76
1:B:317:TRP:O	1:B:321:VAL:HG23	1.86	0.76
1:B:83:CYS:SG	1:B:90:PHE:HB2	2.26	0.75
1:A:378:LEU:HD13	1:A:378:LEU:H	1.52	0.74
1:A:8:ILE:HD12	1:A:286:MET:HE1	1.70	0.74
1:B:210:LYS:HA	1:B:233:GLY:HA2	1.68	0.74
1:A:18:GLN:HE22	1:A:273:LYS:H	1.35	0.74
1:A:241:ILE:HD11	1:A:352:ILE:HG12	1.71	0.73
1:B:9:GLY:H	1:B:64:VAL:HG11	1.53	0.72
1:A:97:THR:H	1:A:106:GLN:NE2	1.87	0.72
1:B:356:GLU:HG3	1:B:365:ILE:HG12	1.71	0.72
1:B:221:PHE:CD1	1:B:238:GLY:HA3	2.25	0.71
1:A:262:LYS:HA	1:A:262:LYS:HE2	1.72	0.71
1:A:96:PRO:HA	1:A:106:GLN:HE21	1.56	0.70
1:B:307:GLY:H	1:B:312:ASN:HD22	1.40	0.70
1:A:73:LYS:HD3	1:A:134:ASP:HA	1.73	0.70
1:B:218:ASN:C	1:B:218:ASN:HD22	2.00	0.69
1:B:91:ILE:HD13	1:B:286:MET:HE3	1.75	0.67
1:A:65:TYR:CE1	1:A:300:ILE:HB	2.30	0.67
1:A:229:ILE:HG12	1:A:231:PHE:CE2	2.29	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:GLN:HE21	1:A:232:GLY:N	1.85	0.67
1:B:8:ILE:HA	1:B:64:VAL:HG13	1.77	0.67
1:B:8:ILE:O	1:B:8:ILE:HG13	1.96	0.65
1:B:234:ALA:O	1:B:236:VAL:HG23	1.98	0.63
1:B:354:ILE:O	1:B:365:ILE:HG13	1.99	0.63
1:A:207:LYS:HB3	1:A:208:HIS:HA	1.81	0.60
1:B:317:TRP:CZ3	1:B:321:VAL:HG21	2.37	0.60
1:B:288:GLU:HG3	1:B:332:VAL:HG11	1.84	0.59
1:A:112:LYS:HE2	1:A:138:SER:OG	2.03	0.59
1:B:367:TYR:CZ	1:B:375:VAL:HG21	2.40	0.57
1:B:218:ASN:C	1:B:218:ASN:ND2	2.59	0.57
1:A:206:GLY:C	1:A:207:LYS:HG2	2.30	0.57
1:B:8:ILE:O	1:B:39:PRO:HA	2.05	0.56
1:A:241:ILE:HD12	1:A:352:ILE:HG12	1.87	0.56
1:A:371:VAL:O	1:A:371:VAL:CG2	2.53	0.56
1:A:107:MET:HE1	1:A:285:VAL:HG11	1.87	0.55
1:A:196:ASN:C	1:A:196:ASN:HD22	2.14	0.55
1:A:349:ASN:N	1:A:370:GLU:HG2	2.21	0.55
1:B:9:GLY:H	1:B:64:VAL:CG1	2.19	0.55
1:A:11:LEU:HD22	1:A:67:ILE:HG21	1.89	0.55
1:A:368:TRP:CZ2	1:A:372:ASP:HA	2.41	0.55
1:A:8:ILE:O	1:A:39:PRO:HA	2.06	0.55
1:A:118:LEU:HA	1:A:374:MET:HE1	1.89	0.54
1:A:370:GLU:C	1:A:372:ASP:H	2.16	0.53
1:B:107:MET:HE1	1:B:285:VAL:HG11	1.91	0.53
1:A:184:ARG:HA	1:A:212:TYR:CD2	2.43	0.53
1:A:9:GLY:HA3	1:A:59:GLN:OE1	2.09	0.53
1:A:98:ASP:CG	1:A:138:SER:HB2	2.34	0.53
1:A:91:ILE:HG21	1:A:107:MET:HE3	1.90	0.53
1:A:229:ILE:HG12	1:A:231:PHE:CZ	2.44	0.53
1:A:12:PHE:O	1:A:43:ASN:HA	2.09	0.53
1:A:226:LEU:O	1:A:229:ILE:HG23	2.09	0.53
1:B:210:LYS:C	1:B:212:TYR:H	2.16	0.53
1:A:91:ILE:CG2	1:A:107:MET:HE3	2.39	0.52
1:B:107:MET:HE1	1:B:285:VAL:CB	2.39	0.52
1:A:112:LYS:CE	1:A:138:SER:HB3	2.39	0.52
1:B:131:TYR:CG	1:B:158:ASN:HA	2.43	0.52
1:B:307:GLY:HA3	1:B:312:ASN:ND2	2.25	0.52
1:B:317:TRP:CE3	1:B:321:VAL:HG21	2.44	0.52
1:A:86:LEU:O	1:A:87:HIS:HB2	2.09	0.52
1:B:158:ASN:C	1:B:160:GLY:H	2.18	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:GLN:O	1:A:65:TYR:HD2	1.92	0.52
1:A:312:ASN:HA	1:A:313:PRO:C	2.33	0.52
1:B:210:LYS:HA	1:B:233:GLY:CA	2.39	0.52
1:A:97:THR:H	1:A:106:GLN:HE21	1.57	0.52
1:A:132:ASP:HB2	1:A:190:CYS:HA	1.91	0.52
1:A:221:PHE:CD1	1:A:238:GLY:HA3	2.46	0.51
1:B:96:PRO:HD3	1:B:108:ARG:HD2	1.92	0.51
1:B:68:PHE:C	1:B:68:PHE:CD2	2.88	0.51
1:B:86:LEU:HA	1:B:314:ALA:CB	2.41	0.51
1:A:252:PHE:HD2	1:A:253:ILE:HD12	1.76	0.51
1:A:21:SER:O	1:A:25:VAL:HG23	2.11	0.50
1:A:112:LYS:HE2	1:A:138:SER:CB	2.42	0.50
1:A:124:TRP:HA	1:A:185:ARG:NH1	2.26	0.50
1:B:107:MET:HE1	1:B:285:VAL:HB	1.94	0.50
1:A:147:ALA:HA	1:A:152:TRP:HB2	1.94	0.50
1:B:29:GLN:O	1:B:29:GLN:OE1	2.29	0.50
1:B:321:VAL:O	1:B:325:ARG:HB2	2.12	0.50
1:A:87:HIS:CG	1:A:316:PRO:HB3	2.47	0.49
1:B:179:GLU:HA	1:B:179:GLU:OE2	2.12	0.49
1:B:183:GLU:O	1:B:183:GLU:HG3	2.11	0.49
1:B:187:ILE:C	1:B:188:LEU:HD12	2.38	0.49
1:A:90:PHE:CE2	1:A:92:THR:HB	2.47	0.49
1:B:86:LEU:HA	1:B:314:ALA:HB2	1.96	0.48
1:B:97:THR:H	1:B:106:GLN:NE2	2.12	0.48
1:A:368:TRP:CZ3	1:A:370:GLU:HA	2.48	0.48
1:B:342:ASP:HB3	1:B:348:ILE:HD13	1.95	0.48
1:B:274:TYR:O	1:B:278:LEU:HG	2.13	0.48
1:B:179:GLU:OE2	1:B:182:LYS:HA	2.13	0.47
1:A:267:ALA:O	1:A:269:THR:HG23	2.14	0.47
1:B:354:ILE:CG2	1:B:365:ILE:HD11	2.44	0.47
1:B:107:MET:HE1	1:B:285:VAL:CG1	2.45	0.47
1:A:114:ALA:HA	1:A:368:TRP:CD1	2.49	0.47
1:B:200:ASP:O	1:B:204:THR:HG22	2.14	0.47
1:B:90:PHE:CE2	1:B:92:THR:HB	2.50	0.47
1:A:30:PHE:CZ	1:A:284:GLN:HB2	2.50	0.46
1:A:129:TYR:CE1	1:A:131:TYR:HB3	2.50	0.46
1:A:10:GLY:HA2	1:A:68:PHE:O	2.15	0.46
1:A:209:VAL:HG23	1:A:210:LYS:N	2.30	0.46
1:A:98:ASP:OD2	1:A:138:SER:HB2	2.15	0.46
1:A:67:ILE:O	1:A:90:PHE:HA	2.16	0.46
1:A:207:LYS:CB	1:A:208:HIS:HA	2.44	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:354:ILE:N	1:B:354:ILE:HD12	2.31	0.45
1:A:225:ASP:O	1:A:228:LYS:HB2	2.16	0.45
1:A:253:ILE:HA	1:A:256:TRP:HB3	1.98	0.45
1:B:90:PHE:HE2	1:B:92:THR:HB	1.81	0.45
1:A:20:TYR:CE2	1:A:24:ARG:HD2	2.52	0.45
1:B:158:ASN:O	1:B:160:GLY:N	2.42	0.45
1:A:8:ILE:HD12	1:A:286:MET:CE	2.44	0.45
1:A:298:ILE:HD13	1:A:322:GLU:HG2	1.97	0.45
1:B:71:TYR:C	1:B:71:TYR:CD1	2.94	0.45
1:B:241:ILE:HG13	1:B:241:ILE:O	2.17	0.45
1:B:232:GLY:O	1:B:233:GLY:C	2.60	0.44
1:B:208:HIS:O	1:B:212:TYR:CE2	2.70	0.44
1:B:317:TRP:O	1:B:321:VAL:CG2	2.63	0.44
1:A:106:GLN:O	1:A:347:ARG:HG2	2.18	0.44
1:B:95:PHE:HA	1:B:96:PRO:HD3	1.85	0.44
1:B:230:GLN:C	1:B:232:GLY:H	2.26	0.44
1:A:127:PHE:HB2	1:A:185:ARG:O	2.18	0.44
1:A:227:LEU:CD1	1:A:362:PRO:HG3	2.47	0.44
1:B:216:ILE:HD12	1:B:216:ILE:N	2.33	0.44
1:B:248:LEU:HD13	1:B:334:GLY:O	2.17	0.44
1:B:17:ASP:HB3	1:B:265:PRO:HB2	2.00	0.43
1:A:29:GLN:HG2	1:A:280:TYR:OH	2.19	0.43
1:B:264:TYR:O	1:B:267:ALA:HB3	2.19	0.43
1:A:316:PRO:O	1:A:317:TRP:C	2.62	0.43
1:B:295:LYS:HB2	1:B:295:LYS:HE3	1.70	0.43
1:A:146:SER:HB3	1:A:150:LYS:HE3	2.01	0.43
1:A:112:LYS:HD3	1:A:139:THR:HA	2.01	0.43
1:A:194:LYS:O	1:A:198:ILE:HG13	2.19	0.43
1:A:9:GLY:CA	1:A:59:GLN:OE1	2.67	0.42
1:A:62:ARG:HA	1:A:62:ARG:HD3	1.74	0.42
1:B:312:ASN:HB3	1:B:313:PRO:CD	2.48	0.42
1:A:112:LYS:HE3	1:A:138:SER:HB3	2.00	0.42
1:A:196:ASN:O	1:A:200:ASP:HB2	2.18	0.42
1:B:68:PHE:CD2	1:B:69:GLY:N	2.86	0.42
1:B:188:LEU:HD12	1:B:188:LEU:N	2.34	0.42
1:B:96:PRO:HG3	1:B:108:ARG:O	2.20	0.42
1:B:133:SER:HA	1:B:137:LEU:HD21	2.01	0.42
1:B:207:LYS:HA	1:B:208:HIS:HA	1.57	0.42
1:B:210:LYS:CA	1:B:233:GLY:HA2	2.42	0.42
1:A:92:THR:HA	1:A:93:PRO:HD3	1.84	0.42
1:B:250:SER:O	1:B:251:LYS:C	2.62	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:375:VAL:O	1:B:375:VAL:HG23	2.20	0.42
1:B:30:PHE:CZ	1:B:284:GLN:HB2	2.55	0.42
1:A:102:PRO:O	1:A:103:PHE:HB2	2.20	0.42
1:B:326:ALA:O	1:B:330:VAL:HG23	2.20	0.42
1:A:96:PRO:HA	1:A:106:GLN:NE2	2.30	0.41
1:B:133:SER:HB3	1:B:137:LEU:HD11	2.01	0.41
1:B:213:HIS:HA	1:B:235:ASN:HB3	2.01	0.41
1:A:96:PRO:HA	1:A:106:GLN:HG2	2.02	0.41
1:A:60:PHE:CD2	1:A:309:CYS:HB3	2.56	0.41
1:A:256:TRP:CZ2	1:A:267:ALA:HA	2.56	0.41
1:A:354:ILE:HD11	1:A:368:TRP:HB2	2.01	0.41
1:B:158:ASN:C	1:B:160:GLY:N	2.79	0.41
1:B:124:TRP:HB3	1:B:127:PHE:HD2	1.86	0.41
1:B:324:GLU:HG2	1:B:328:LYS:HD2	2.03	0.41
1:A:374:MET:HB2	1:A:374:MET:HE2	1.85	0.41
1:B:53:THR:HG23	1:B:82:PHE:CZ	2.56	0.41
1:A:342:ASP:OD1	1:A:342:ASP:C	2.64	0.41
1:B:92:THR:HA	1:B:93:PRO:HD3	1.77	0.41
1:A:35:PHE:CD1	1:A:35:PHE:C	2.99	0.41
1:B:25:VAL:HG21	1:B:256:TRP:HZ3	1.86	0.40
1:B:138:SER:HA	1:B:141:GLN:HG3	2.03	0.40
1:A:137:LEU:HD13	1:A:140:LEU:HD22	2.02	0.40
1:A:321:VAL:O	1:A:325:ARG:HB2	2.21	0.40
1:B:248:LEU:HD22	1:B:335:LEU:HD23	2.03	0.40
1:A:325:ARG:HH11	1:A:325:ARG:HG3	1.87	0.40
1:B:133:SER:HA	1:B:134:ASP:CB	2.50	0.40
1:B:307:GLY:N	1:B:312:ASN:HD22	2.13	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	365/394 (93%)	335 (92%)	27 (7%)	3 (1%)	16	34
1	B	360/394 (91%)	308 (86%)	43 (12%)	9 (2%)	4	10
All	All	725/788 (92%)	643 (89%)	70 (10%)	12 (2%)	7	17

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	159	VAL
1	B	151	LYS
1	B	235	ASN
1	A	174	LEU
1	B	158	ASN
1	B	162	ILE
1	B	174	LEU
1	A	371	VAL
1	B	307	GLY
1	B	136	GLY
1	A	203	ILE
1	B	233	GLY

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	284/342 (83%)	262 (92%)	22 (8%)	12	27
1	B	233/342 (68%)	221 (95%)	12 (5%)	21	43
All	All	517/684 (76%)	483 (93%)	34 (7%)	15	33

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	LEU
1	A	46	VAL
1	A	49	SER
1	A	98	ASP

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Mol	Chain	Res	Type
1	A	106	GLN
1	A	135	ARG
1	A	196	ASN
1	A	229	ILE
1	A	241	ILE
1	A	248	LEU
1	A	253	ILE
1	A	272	ILE
1	A	287	THR
1	A	297	ARG
1	A	300	ILE
1	A	315	VAL
1	A	354	ILE
1	A	370	GLU
1	A	371	VAL
1	A	374	MET
1	A	376	VAL
1	A	378	LEU
1	B	11	LEU
1	B	29	GLN
1	B	61	SER
1	B	204	THR
1	B	218	ASN
1	B	248	LEU
1	B	253	ILE
1	B	321	VAL
1	B	343	GLN
1	B	359	THR
1	B	372	ASP
1	B	377	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	101	HIS
1	A	106	GLN
1	A	158	ASN
1	A	196	ASN
1	A	230	GLN
1	A	284	GLN
1	A	292	ASN

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Mol	Chain	Res	Type
1	A	319	GLN
1	B	18	GLN
1	B	54	ASN
1	B	77	ASN
1	B	141	GLN
1	B	208	HIS
1	B	218	ASN
1	B	240	GLN
1	B	292	ASN
1	B	312	ASN
1	B	319	GLN
1	B	344	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 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 [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	371/394 (94%)	-0.23	2 (0%) 87 86	43, 76, 149, 195	1 (0%)
1	B	366/394 (92%)	0.09	9 (2%) 58 54	52, 104, 177, 247	0
All	All	737/788 (93%)	-0.07	11 (1%) 72 69	43, 88, 170, 247	1 (0%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	267	ALA	3.1
1	B	36	ARG	3.0
1	B	171	TYR	2.8
1	B	42	ASP	2.6
1	B	231	PHE	2.4
1	B	170	THR	2.3
1	B	200	ASP	2.3
1	A	178	LEU	2.3
1	B	241	ILE	2.2
1	A	156	ALA	2.0
1	B	136	GLY	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.