



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 08:14 PM UTC

PDB ID : 4XFV / pdb_00004xfv
Title : Crystal Structure of Elp2
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Deposited on : 2014-12-29
Resolution : 3.20 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

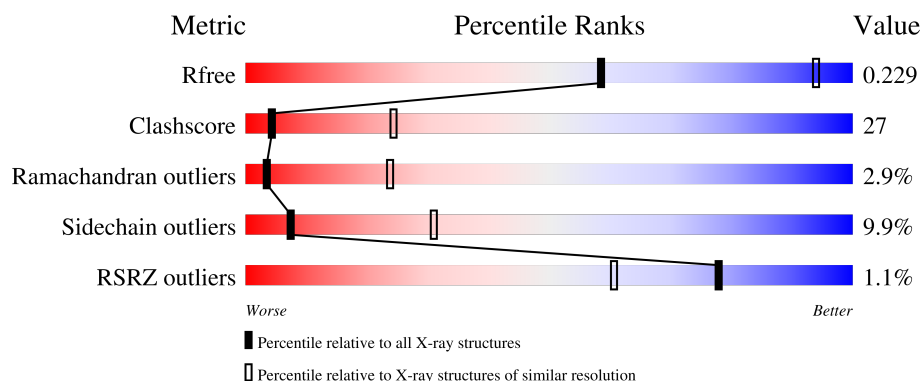
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	792	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5379 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 Elongator complex protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	700	Total	C	N	O	S	0	0	0
			5374	3458	928	964	24			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP P42935
A	-2	PRO	-	expression tag	UNP P42935
A	-1	GLY	-	expression tag	UNP P42935
A	0	SER	-	expression tag	UNP P42935

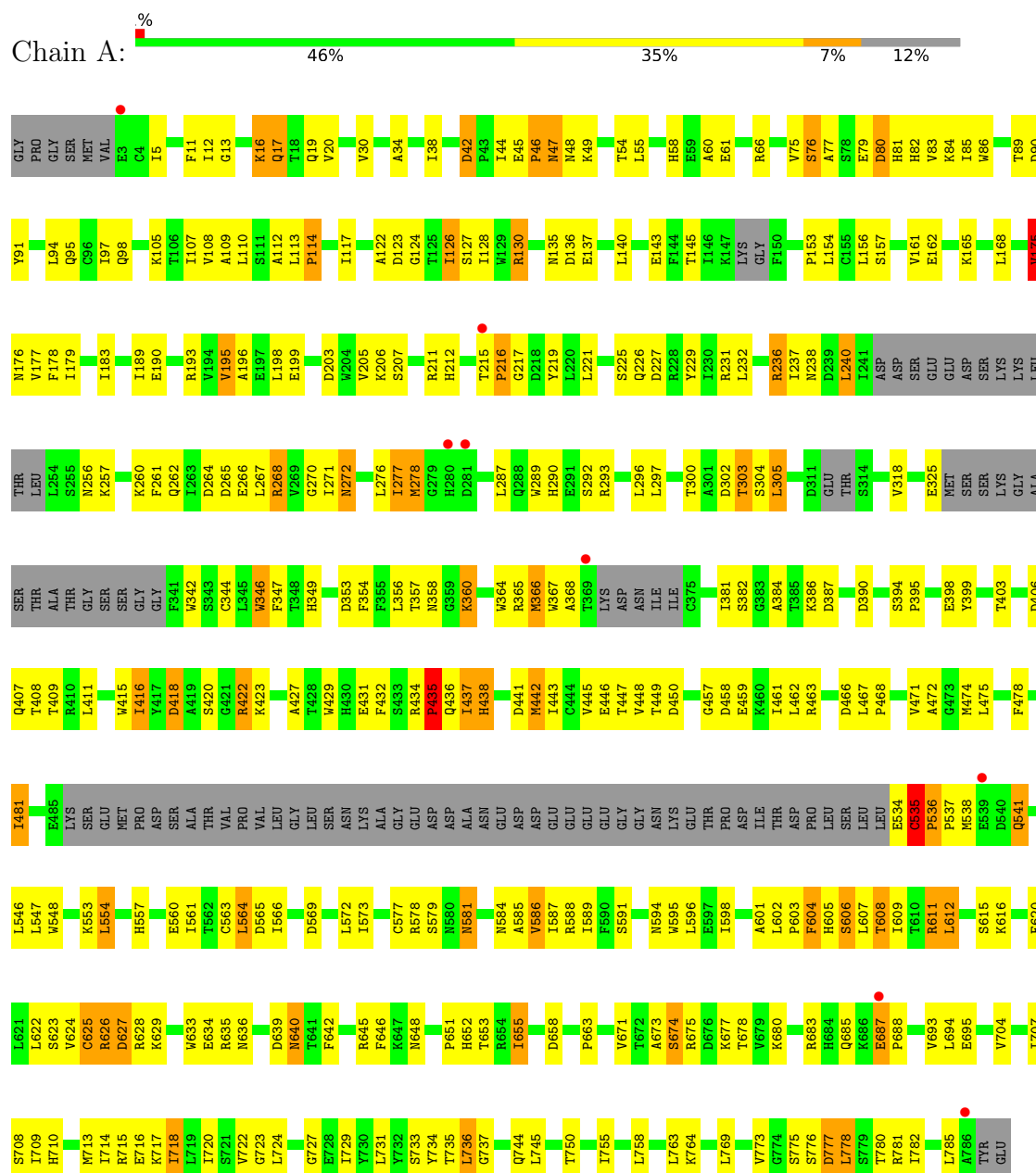
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	5	Total	O	0	0
			5	5		

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: Elongator complex protein 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	80.64Å 80.64Å 535.79Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.00 – 3.20 48.00 – 3.20	Depositor EDS
% Data completeness (in resolution range)	87.8 (48.00-3.20) 87.8 (48.00-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.80 (at 3.19Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, R_{free}	0.199 , 0.241 0.200 , 0.229	Depositor DCC
R_{free} test set	844 reflections (4.47%)	wwPDB-VP
Wilson B-factor (Å ²)	49.2	Xtriage
Anisotropy	0.243	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 55.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	5379	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% 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	0.38	0/5503	0.90	20/7483 (0.3%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	277	ILE	N-CA-C	10.07	120.94	111.48
1	A	627	ASP	N-CA-C	-8.85	100.90	112.41
1	A	625	CYS	N-CA-C	8.02	121.14	110.88
1	A	716	GLU	N-CA-C	-7.54	101.79	113.02
1	A	80	ASP	N-CA-C	-7.42	98.46	109.15
1	A	727	GLY	N-CA-C	-7.29	104.51	115.00
1	A	256	ASN	N-CA-C	-6.43	105.73	112.93
1	A	418	ASP	N-CA-C	-6.20	102.79	110.41
1	A	538	MET	N-CA-C	-6.18	102.38	110.53
1	A	124	GLY	N-CA-C	-6.03	106.99	115.32
1	A	123	ASP	N-CA-C	-5.79	106.17	113.18
1	A	154	LEU	N-CA-C	-5.65	107.63	114.75
1	A	195	VAL	N-CA-C	5.44	118.21	112.83
1	A	42	ASP	CA-C-N	5.23	124.41	118.97
1	A	42	ASP	C-N-CA	5.23	124.41	118.97
1	A	434	ARG	CA-C-N	5.20	126.34	119.84
1	A	434	ARG	C-N-CA	5.20	126.34	119.84
1	A	777	ASP	N-CA-C	-5.09	107.23	112.93
1	A	678	THR	N-CA-C	5.08	118.29	110.42
1	A	674	SER	N-CA-C	5.04	116.85	109.24

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	5374	0	5141	283	0
2	A	5	0	0	0	0
All	All	5379	0	5141	283	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 27.

All (283) 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:290:HIS:HD2	1:A:293:ARG:H	1.03	0.99
1:A:658:ASP:HB3	1:A:707:ILE:HG22	1.49	0.94
1:A:240:LEU:H	1:A:240:LEU:HD23	1.32	0.93
1:A:416:ILE:HG23	1:A:427:ALA:HB2	1.54	0.88
1:A:262:GLN:HB2	1:A:267:LEU:O	1.75	0.87
1:A:415:TRP:HB3	1:A:432:PHE:HZ	1.41	0.85
1:A:687:GLU:HB2	1:A:688:PRO:HD3	1.57	0.84
1:A:418:ASP:OD2	1:A:422:ARG:HD2	1.75	0.84
1:A:290:HIS:CD2	1:A:293:ARG:H	1.94	0.82
1:A:76:SER:HB2	1:A:86:TRP:HZ3	1.45	0.79
1:A:90:ASP:OD1	1:A:91:TYR:N	2.15	0.79
1:A:60:ALA:HB1	1:A:79:GLU:HB2	1.66	0.77
1:A:403:THR:HB	1:A:442:MET:HB3	1.67	0.77
1:A:416:ILE:HD12	1:A:416:ILE:H	1.51	0.75
1:A:305:LEU:HD21	1:A:344:CYS:SG	2.27	0.75
1:A:297:LEU:HD21	1:A:305:LEU:HD23	1.67	0.75
1:A:240:LEU:HD23	1:A:240:LEU:N	2.01	0.74
1:A:611:ARG:HD2	1:A:658:ASP:OD1	1.87	0.74
1:A:17:GLN:HB3	1:A:20:VAL:HG13	1.70	0.74
1:A:215:THR:HA	1:A:217:GLY:N	2.02	0.74
1:A:468:PRO:HG2	1:A:471:VAL:HG23	1.68	0.73
1:A:408:THR:HG22	1:A:437:ILE:HG23	1.71	0.72
1:A:89:THR:CG2	1:A:95:GLN:HB2	2.20	0.72
1:A:777:ASP:O	1:A:778:LEU:HG	1.89	0.72
1:A:671:VAL:HG23	1:A:680:LYS:O	1.88	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:717:LYS:HD3	1:A:733:SER:OG	1.90	0.71
1:A:19:GLN:HE22	1:A:66:ARG:HH21	1.39	0.71
1:A:709:ILE:HG12	1:A:720:ILE:HG12	1.72	0.71
1:A:17:GLN:HB3	1:A:20:VAL:CG1	2.21	0.70
1:A:229:TYR:HB3	1:A:276:LEU:HD11	1.72	0.70
1:A:713:MET:HG2	1:A:718:ILE:HD12	1.74	0.69
1:A:236:ARG:HG2	1:A:236:ARG:HH11	1.58	0.68
1:A:76:SER:CB	1:A:86:TRP:HZ3	2.07	0.68
1:A:587:ILE:HB	1:A:602:LEU:HB2	1.74	0.67
1:A:406:ASP:O	1:A:407:GLN:HB2	1.94	0.67
1:A:305:LEU:HD11	1:A:357:THR:HG21	1.76	0.67
1:A:135:ASN:ND2	1:A:137:GLU:HB2	2.10	0.66
1:A:395:PRO:HG2	1:A:449:THR:HA	1.77	0.66
1:A:240:LEU:HD12	1:A:270:GLY:HA3	1.77	0.66
1:A:395:PRO:HD3	1:A:447:THR:HB	1.78	0.66
1:A:616:LYS:HE2	1:A:663:PRO:HA	1.78	0.66
1:A:626:ARG:HB2	1:A:655:ILE:HG12	1.79	0.65
1:A:264:ASP:C	1:A:266:GLU:H	2.03	0.64
1:A:79:GLU:OE2	1:A:79:GLU:HA	1.98	0.64
1:A:457:GLY:C	1:A:459:GLU:H	2.04	0.64
1:A:775:SER:HB3	1:A:777:ASP:OD1	1.97	0.64
1:A:113:LEU:HB3	1:A:114:PRO:HD2	1.79	0.63
1:A:755:ILE:HG21	1:A:758:LEU:HD13	1.80	0.63
1:A:713:MET:HG2	1:A:718:ILE:CD1	2.28	0.63
1:A:76:SER:HB2	1:A:86:TRP:CZ3	2.32	0.62
1:A:287:LEU:HD12	1:A:287:LEU:O	2.00	0.62
1:A:627:ASP:OD2	1:A:629:LYS:HG2	2.00	0.62
1:A:80:ASP:O	1:A:82:HIS:N	2.32	0.62
1:A:625:CYS:O	1:A:625:CYS:SG	2.57	0.61
1:A:157:SER:O	1:A:168:LEU:HD12	2.00	0.61
1:A:615:SER:HB2	1:A:620:PHE:HB2	1.82	0.61
1:A:409:THR:O	1:A:435:PRO:HD2	2.02	0.60
1:A:685:GLN:HG3	1:A:687:GLU:OE1	2.01	0.60
1:A:342:TRP:HB2	1:A:358:ASN:HD22	1.66	0.60
1:A:443:ILE:HG12	1:A:457:GLY:HA2	1.83	0.60
1:A:603:PRO:O	1:A:604:PHE:HB2	2.01	0.60
1:A:108:VAL:HG21	1:A:153:PRO:O	2.02	0.60
1:A:622:LEU:HD12	1:A:623:SER:H	1.66	0.59
1:A:42:ASP:OD2	1:A:45:GLU:HB2	2.01	0.59
1:A:5:ILE:HD12	1:A:5:ILE:N	2.18	0.59
1:A:212:HIS:HB2	1:A:219:TYR:CE1	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:663:PRO:CG	1:A:718:ILE:HD11	2.33	0.59
1:A:553:LYS:O	1:A:554:LEU:HD23	2.03	0.59
1:A:13:GLY:HA3	1:A:384:ALA:O	2.03	0.58
1:A:290:HIS:CD2	1:A:292:SER:H	2.21	0.58
1:A:44:ILE:HD11	1:A:347:PHE:HE1	1.68	0.58
1:A:45:GLU:HG3	1:A:46:PRO:HD2	1.85	0.58
1:A:467:LEU:HD13	1:A:548:TRP:CZ2	2.39	0.58
1:A:290:HIS:HD2	1:A:293:ARG:N	1.88	0.58
1:A:303:THR:OG1	1:A:325:GLU:HA	2.04	0.57
1:A:474:MET:SD	1:A:537:PRO:HG3	2.44	0.57
1:A:704:VAL:HG13	1:A:722:VAL:HG13	1.85	0.57
1:A:5:ILE:HD12	1:A:5:ILE:H	1.68	0.57
1:A:429:TRP:CZ3	1:A:782:ILE:HD13	2.40	0.57
1:A:107:ILE:N	1:A:107:ILE:HD12	2.20	0.56
1:A:354:PHE:HA	1:A:367:TRP:O	2.04	0.56
1:A:12:ILE:HG12	1:A:34:ALA:HB1	1.87	0.56
1:A:215:THR:HA	1:A:216:PRO:C	2.30	0.56
1:A:240:LEU:N	1:A:240:LEU:CD2	2.68	0.56
1:A:687:GLU:CB	1:A:688:PRO:HD3	2.33	0.56
1:A:175:VAL:HG12	1:A:203:ASP:N	2.20	0.56
1:A:777:ASP:C	1:A:778:LEU:HG	2.31	0.56
1:A:675:ARG:C	1:A:677:LYS:H	2.13	0.56
1:A:723:GLY:C	1:A:724:LEU:HD12	2.32	0.55
1:A:297:LEU:HD21	1:A:305:LEU:CD2	2.34	0.55
1:A:179:ILE:HB	1:A:196:ALA:HB3	1.88	0.55
1:A:395:PRO:HG2	1:A:449:THR:CA	2.36	0.55
1:A:671:VAL:CG1	1:A:709:ILE:HD11	2.38	0.54
1:A:17:GLN:O	1:A:20:VAL:HG22	2.08	0.54
1:A:19:GLN:NE2	1:A:66:ARG:HH21	2.03	0.54
1:A:561:ILE:N	1:A:561:ILE:HD12	2.23	0.54
1:A:581:ASN:H	1:A:581:ASN:HD22	1.55	0.54
1:A:671:VAL:HG12	1:A:709:ILE:HD11	1.90	0.54
1:A:229:TYR:HA	1:A:278:MET:CB	2.37	0.54
1:A:189:ILE:HG22	1:A:190:GLU:N	2.23	0.54
1:A:356:LEU:CD2	1:A:366:MET:HG2	2.38	0.54
1:A:105:LYS:HB3	1:A:122:ALA:HB3	1.90	0.54
1:A:457:GLY:C	1:A:459:GLU:N	2.66	0.54
1:A:579:SER:H	1:A:608:THR:HB	1.73	0.53
1:A:735:THR:O	1:A:737:GLY:N	2.41	0.53
1:A:415:TRP:HB3	1:A:432:PHE:CZ	2.32	0.53
1:A:635:ARG:HB2	1:A:642:PHE:CE1	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:ALA:HB1	1:A:117:ILE:HG12	1.90	0.53
1:A:277:ILE:HG13	1:A:278:MET:H	1.74	0.53
1:A:415:TRP:HB2	1:A:478:PHE:HB3	1.91	0.53
1:A:708:SER:HB2	1:A:758:LEU:O	2.08	0.53
1:A:628:ARG:O	1:A:651:PRO:HD2	2.09	0.53
1:A:605:HIS:HB3	1:A:625:CYS:SG	2.49	0.53
1:A:260:LYS:HB3	1:A:268:ARG:NH1	2.24	0.53
1:A:457:GLY:O	1:A:459:GLU:N	2.33	0.53
1:A:112:ALA:CB	1:A:117:ILE:HG12	2.39	0.53
1:A:44:ILE:HD11	1:A:347:PHE:CE1	2.44	0.52
1:A:694:LEU:C	1:A:694:LEU:HD23	2.35	0.52
1:A:744:GLN:O	1:A:745:LEU:C	2.52	0.52
1:A:110:LEU:C	1:A:110:LEU:HD23	2.34	0.52
1:A:625:CYS:O	1:A:627:ASP:N	2.41	0.52
1:A:300:THR:OG1	1:A:304:SER:HB3	2.10	0.52
1:A:128:ILE:HG21	1:A:189:ILE:HD11	1.91	0.52
1:A:211:ARG:HG3	1:A:289:TRP:CE3	2.45	0.52
1:A:416:ILE:CG2	1:A:427:ALA:HB2	2.35	0.52
1:A:693:VAL:HG12	1:A:694:LEU:N	2.25	0.52
1:A:183:ILE:C	1:A:183:ILE:HD12	2.35	0.51
1:A:606:SER:O	1:A:607:LEU:HD23	2.11	0.51
1:A:353:ASP:O	1:A:368:ALA:HA	2.10	0.51
1:A:566:ILE:CD1	1:A:573:ILE:HG12	2.41	0.51
1:A:176:ASN:ND2	1:A:199:GLU:HA	2.26	0.51
1:A:566:ILE:HD12	1:A:573:ILE:HG12	1.91	0.50
1:A:585:ALA:HB2	1:A:607:LEU:O	2.11	0.50
1:A:673:ALA:HB1	1:A:704:VAL:HG12	1.92	0.50
1:A:360:LYS:NZ	1:A:360:LYS:H	2.09	0.50
1:A:156:LEU:O	1:A:156:LEU:HD12	2.11	0.50
1:A:718:ILE:HG22	1:A:734:TYR:HB3	1.93	0.50
1:A:586:VAL:CG1	1:A:601:ALA:HB1	2.42	0.50
1:A:113:LEU:HB3	1:A:114:PRO:CD	2.42	0.49
1:A:231:ARG:HH11	1:A:231:ARG:HG2	1.76	0.49
1:A:629:LYS:HB2	1:A:648:ASN:O	2.12	0.49
1:A:603:PRO:O	1:A:633:TRP:HH2	1.95	0.49
1:A:236:ARG:HH11	1:A:236:ARG:CG	2.22	0.49
1:A:54:THR:HG21	1:A:781:ARG:NH2	2.28	0.49
1:A:398:GLU:HG2	1:A:481:ILE:CG2	2.43	0.49
1:A:603:PRO:O	1:A:604:PHE:CB	2.59	0.49
1:A:126:ILE:HD11	1:A:168:LEU:HD21	1.95	0.49
1:A:176:ASN:HD21	1:A:199:GLU:HG2	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:VAL:HB	1:A:198:LEU:HB2	1.94	0.48
1:A:237:ILE:HG22	1:A:238:ASN:N	2.27	0.48
1:A:687:GLU:HB2	1:A:688:PRO:CD	2.37	0.48
1:A:85:ILE:O	1:A:97:ILE:HG12	2.12	0.48
1:A:127:SER:HB2	1:A:140:LEU:HD11	1.95	0.48
1:A:448:VAL:HG12	1:A:449:THR:HG23	1.96	0.48
1:A:475:LEU:HB3	1:A:481:ILE:HD11	1.96	0.48
1:A:77:ALA:HA	1:A:82:HIS:O	2.13	0.48
1:A:635:ARG:HH21	1:A:640:ASN:HD21	1.61	0.48
1:A:724:LEU:HD12	1:A:724:LEU:N	2.29	0.48
1:A:16:LYS:HG2	1:A:342:TRP:CE2	2.49	0.48
1:A:416:ILE:O	1:A:423:LYS:HA	2.13	0.48
1:A:178:PHE:N	1:A:178:PHE:CD2	2.82	0.48
1:A:264:ASP:O	1:A:266:GLU:N	2.47	0.47
1:A:611:ARG:HB3	1:A:624:VAL:CG2	2.44	0.47
1:A:5:ILE:HG13	1:A:785:LEU:CD1	2.45	0.47
1:A:553:LYS:C	1:A:554:LEU:HD23	2.39	0.47
1:A:394:SER:HB3	1:A:399:TYR:CE2	2.50	0.47
1:A:109:ALA:HB1	1:A:156:LEU:HG	1.97	0.47
1:A:342:TRP:CE2	1:A:360:LYS:HA	2.50	0.47
1:A:360:LYS:H	1:A:360:LYS:CE	2.28	0.47
1:A:609:ILE:HG21	1:A:612:LEU:CD1	2.45	0.47
1:A:655:ILE:HD12	1:A:675:ARG:NH2	2.30	0.47
1:A:237:ILE:O	1:A:268:ARG:O	2.33	0.47
1:A:687:GLU:N	1:A:687:GLU:CD	2.73	0.47
1:A:38:ILE:HB	1:A:55:LEU:HB2	1.97	0.47
1:A:461:ILE:HD11	1:A:463:ARG:HE	1.79	0.47
1:A:271:ILE:HG22	1:A:272:ASN:N	2.30	0.47
1:A:386:LYS:NZ	1:A:386:LYS:HB3	2.29	0.47
1:A:683:ARG:NE	1:A:695:GLU:OE1	2.47	0.47
1:A:572:LEU:HD22	1:A:589:ILE:CG2	2.45	0.46
1:A:162:GLU:H	1:A:165:LYS:HB2	1.80	0.46
1:A:429:TRP:CE3	1:A:782:ILE:HD13	2.51	0.46
1:A:652:HIS:ND1	1:A:674:SER:HB3	2.31	0.46
1:A:763:LEU:HD12	1:A:764:LYS:H	1.80	0.46
1:A:58:HIS:CE1	1:A:84:LYS:HD2	2.50	0.46
1:A:231:ARG:C	1:A:232:LEU:HD12	2.41	0.46
1:A:462:LEU:H	1:A:462:LEU:HD23	1.80	0.46
1:A:305:LEU:HD11	1:A:357:THR:CG2	2.44	0.46
1:A:364:TRP:HB2	1:A:381:ILE:HG12	1.97	0.46
1:A:394:SER:HA	1:A:447:THR:HG21	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:ASN:HD22	1:A:137:GLU:HB2	1.78	0.46
1:A:625:CYS:O	1:A:626:ARG:C	2.59	0.46
1:A:652:HIS:CE1	1:A:674:SER:HB3	2.51	0.46
1:A:193:ARG:NH2	1:A:262:GLN:O	2.47	0.45
1:A:416:ILE:HD12	1:A:416:ILE:N	2.26	0.45
1:A:17:GLN:HE21	1:A:17:GLN:HB2	1.56	0.45
1:A:277:ILE:O	1:A:278:MET:C	2.58	0.45
1:A:177:VAL:HG22	1:A:205:VAL:HG21	1.99	0.45
1:A:189:ILE:CG2	1:A:190:GLU:N	2.79	0.45
1:A:395:PRO:HD2	1:A:450:ASP:OD2	2.17	0.45
1:A:584:ASN:C	1:A:586:VAL:H	2.24	0.45
1:A:84:LYS:C	1:A:85:ILE:HD12	2.41	0.45
1:A:407:GLN:HB3	1:A:438:HIS:O	2.17	0.45
1:A:673:ALA:HB1	1:A:704:VAL:CG1	2.47	0.45
1:A:342:TRP:CZ2	1:A:360:LYS:HA	2.52	0.45
1:A:645:ARG:HH11	1:A:645:ARG:HG3	1.82	0.45
1:A:745:LEU:HB3	1:A:750:THR:CG2	2.47	0.45
1:A:195:VAL:HG22	1:A:195:VAL:O	2.18	0.44
1:A:611:ARG:HB3	1:A:624:VAL:HG22	1.99	0.44
1:A:546:LEU:C	1:A:547:LEU:HD23	2.42	0.44
1:A:578:ARG:HG3	1:A:578:ARG:HH11	1.83	0.44
1:A:588:ARG:HG3	1:A:588:ARG:HH11	1.83	0.44
1:A:608:THR:O	1:A:626:ARG:HG2	2.18	0.44
1:A:127:SER:HB3	1:A:143:GLU:HG3	1.99	0.44
1:A:436:GLN:C	1:A:437:ILE:HG13	2.42	0.44
1:A:80:ASP:OD1	1:A:82:HIS:CG	2.71	0.44
1:A:602:LEU:HB3	1:A:633:TRP:CZ3	2.52	0.44
1:A:297:LEU:HB2	1:A:346:TRP:CD1	2.53	0.44
1:A:418:ASP:C	1:A:420:SER:N	2.75	0.44
1:A:591:SER:HB2	1:A:598:ILE:HD11	2.00	0.44
1:A:225:SER:C	1:A:227:ASP:H	2.26	0.43
1:A:212:HIS:HE1	1:A:215:THR:CB	2.31	0.43
1:A:704:VAL:HG13	1:A:722:VAL:CG1	2.48	0.43
1:A:360:LYS:HZ3	1:A:360:LYS:N	2.15	0.43
1:A:418:ASP:C	1:A:420:SER:H	2.27	0.43
1:A:471:VAL:O	1:A:472:ALA:C	2.62	0.43
1:A:563:CYS:O	1:A:564:LEU:HB3	2.19	0.43
1:A:645:ARG:NH2	1:A:646:PHE:CE1	2.87	0.43
1:A:61:GLU:HB3	1:A:79:GLU:HG2	2.00	0.43
1:A:459:GLU:O	1:A:560:GLU:HA	2.18	0.43
1:A:128:ILE:CG2	1:A:189:ILE:HD11	2.48	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:ILE:HG13	1:A:190:GLU:O	2.18	0.42
1:A:462:LEU:HD21	1:A:554:LEU:HB2	2.00	0.42
1:A:534:GLU:O	1:A:535:CYS:CB	2.67	0.42
1:A:303:THR:O	1:A:325:GLU:N	2.52	0.42
1:A:622:LEU:HD12	1:A:623:SER:N	2.33	0.42
1:A:710:HIS:NE2	1:A:769:LEU:HD22	2.34	0.42
1:A:44:ILE:HD13	1:A:349:HIS:CG	2.55	0.42
1:A:636:ASN:ND2	1:A:639:ASP:H	2.18	0.42
1:A:685:GLN:HG2	1:A:688:PRO:HD2	2.01	0.42
1:A:722:VAL:O	1:A:729:ILE:HA	2.18	0.42
1:A:47:ASN:O	1:A:48:ASN:C	2.63	0.42
1:A:76:SER:O	1:A:83:VAL:HA	2.20	0.42
1:A:206:LYS:NZ	1:A:226:GLN:CB	2.83	0.42
1:A:445:VAL:HG22	1:A:446:GLU:N	2.35	0.42
1:A:557:HIS:HD2	1:A:577:CYS:SG	2.42	0.42
1:A:626:ARG:CB	1:A:655:ILE:HG23	2.49	0.42
1:A:11:PHE:HB2	1:A:780:THR:HB	2.02	0.42
1:A:126:ILE:O	1:A:126:ILE:HG12	2.20	0.42
1:A:94:LEU:HD12	1:A:94:LEU:N	2.35	0.41
1:A:211:ARG:HB2	1:A:289:TRP:CE2	2.55	0.41
1:A:347:PHE:CD1	1:A:347:PHE:C	2.98	0.41
1:A:541:GLN:HE21	1:A:541:GLN:HB3	1.62	0.41
1:A:620:PHE:CD2	1:A:634:GLU:HA	2.55	0.41
1:A:386:LYS:O	1:A:387:ASP:C	2.63	0.41
1:A:625:CYS:C	1:A:627:ASP:N	2.78	0.41
1:A:409:THR:HB	1:A:436:GLN:HG2	2.01	0.41
1:A:704:VAL:CG1	1:A:722:VAL:HG13	2.50	0.41
1:A:357:THR:OG1	1:A:367:TRP:HZ3	2.04	0.41
1:A:585:ALA:CB	1:A:607:LEU:O	2.68	0.41
1:A:98:GLN:HB2	1:A:137:GLU:HG2	2.02	0.41
1:A:236:ARG:CG	1:A:236:ARG:NH1	2.82	0.41
1:A:441:ASP:O	1:A:457:GLY:CA	2.69	0.41
1:A:12:ILE:HD13	1:A:34:ALA:HB3	2.03	0.41
1:A:277:ILE:CD1	1:A:296:LEU:HD23	2.51	0.41
1:A:462:LEU:CD2	1:A:554:LEU:HB2	2.51	0.41
1:A:565:ASP:HB2	1:A:612:LEU:O	2.21	0.41
1:A:581:ASN:HD22	1:A:581:ASN:C	2.28	0.41
1:A:635:ARG:HB2	1:A:642:PHE:HE1	1.84	0.41
1:A:277:ILE:HG13	1:A:278:MET:N	2.35	0.41
1:A:758:LEU:HD12	1:A:773:VAL:HG22	2.03	0.41
1:A:573:ILE:HD11	1:A:595:TRP:HZ3	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:581:ASN:C	1:A:581:ASN:ND2	2.78	0.41
1:A:429:TRP:CE3	1:A:782:ILE:CD1	3.04	0.41
1:A:587:ILE:CG2	1:A:602:LEU:HD12	2.50	0.41
1:A:594:ASN:O	1:A:596:LEU:HG	2.22	0.41
1:A:130:ARG:C	1:A:130:ARG:HD3	2.46	0.40
1:A:382:SER:HB3	1:A:431:GLU:OE2	2.21	0.40
1:A:416:ILE:HG23	1:A:427:ALA:CB	2.38	0.40
1:A:394:SER:O	1:A:395:PRO:C	2.65	0.40
1:A:195:VAL:CG2	1:A:261:PHE:HB2	2.52	0.40
1:A:80:ASP:O	1:A:80:ASP:OD1	2.39	0.40
1:A:360:LYS:H	1:A:360:LYS:HE2	1.86	0.40
1:A:411:LEU:HD23	1:A:411:LEU:C	2.47	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	686/792 (87%)	595 (87%)	71 (10%)	20 (3%)	3	24

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	81	HIS
1	A	535	CYS
1	A	604	PHE
1	A	626	ARG
1	A	714	ILE
1	A	736	LEU
1	A	265	ASP
1	A	438	HIS
1	A	536	PRO

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Mol	Chain	Res	Type
1	A	278	MET
1	A	303	THR
1	A	458	ASP
1	A	655	ILE
1	A	715	ARG
1	A	161	VAL
1	A	46	PRO
1	A	435	PRO
1	A	175	VAL
1	A	687	GLU
1	A	216	PRO

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	546/694 (79%)	492 (90%)	54 (10%)	7	31

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	LYS
1	A	17	GLN
1	A	30	VAL
1	A	47	ASN
1	A	49	LYS
1	A	75	VAL
1	A	76	SER
1	A	114	PRO
1	A	126	ILE
1	A	130	ARG
1	A	136	ASP
1	A	145	THR
1	A	175	VAL
1	A	207	SER
1	A	221	LEU

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Mol	Chain	Res	Type
1	A	236	ARG
1	A	240	LEU
1	A	257	LYS
1	A	268	ARG
1	A	272	ASN
1	A	302	ASP
1	A	305	LEU
1	A	318	VAL
1	A	346	TRP
1	A	360	LYS
1	A	365	ARG
1	A	366	MET
1	A	390	ASP
1	A	416	ILE
1	A	422	ARG
1	A	435	PRO
1	A	437	ILE
1	A	442	MET
1	A	466	ASP
1	A	481	ILE
1	A	535	CYS
1	A	536	PRO
1	A	541	GLN
1	A	554	LEU
1	A	564	LEU
1	A	569	ASP
1	A	581	ASN
1	A	586	VAL
1	A	606	SER
1	A	608	THR
1	A	611	ARG
1	A	612	LEU
1	A	640	ASN
1	A	653	THR
1	A	718	ILE
1	A	731	LEU
1	A	736	LEU
1	A	776	SER
1	A	778	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	19	GLN
1	A	47	ASN
1	A	48	ASN
1	A	82	HIS
1	A	101	GLN
1	A	131	GLN
1	A	135	ASN
1	A	176	ASN
1	A	212	HIS
1	A	258	GLN
1	A	272	ASN
1	A	288	GLN
1	A	290	HIS
1	A	295	GLN
1	A	358	ASN
1	A	557	HIS
1	A	581	ASN
1	A	584	ASN
1	A	636	ASN
1	A	640	ASN
1	A	726	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	700/792 (88%)	-0.24	8 (1%)	78 61	15, 42, 77, 105	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	281	ASP	3.5
1	A	687	GLU	3.2
1	A	369	THR	3.2
1	A	215	THR	2.7
1	A	786	ALA	2.4
1	A	280	HIS	2.3
1	A	539	GLU	2.2
1	A	3	GLU	2.2

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.