



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 02:22 am GMT

PDB ID : 4XFV
Title : Crystal Structure of Elp2
Authors : Lin, Z.; Dong, C.; Long, J.; Shen, Y.
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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

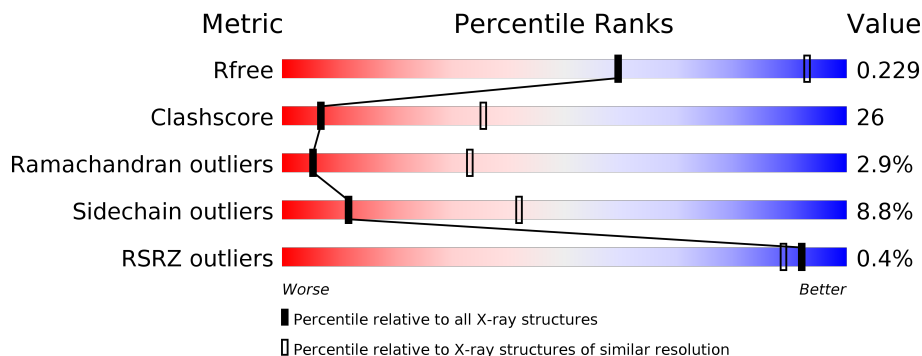
1 Overall quality at a glance

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}	100719	1015 (3.22-3.18)
Clashscore	112137	1009 (3.20-3.20)
Ramachandran outliers	110173	1118 (3.22-3.18)
Sidechain outliers	110143	1117 (3.22-3.18)
RSRZ outliers	101464	1020 (3.22-3.18)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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

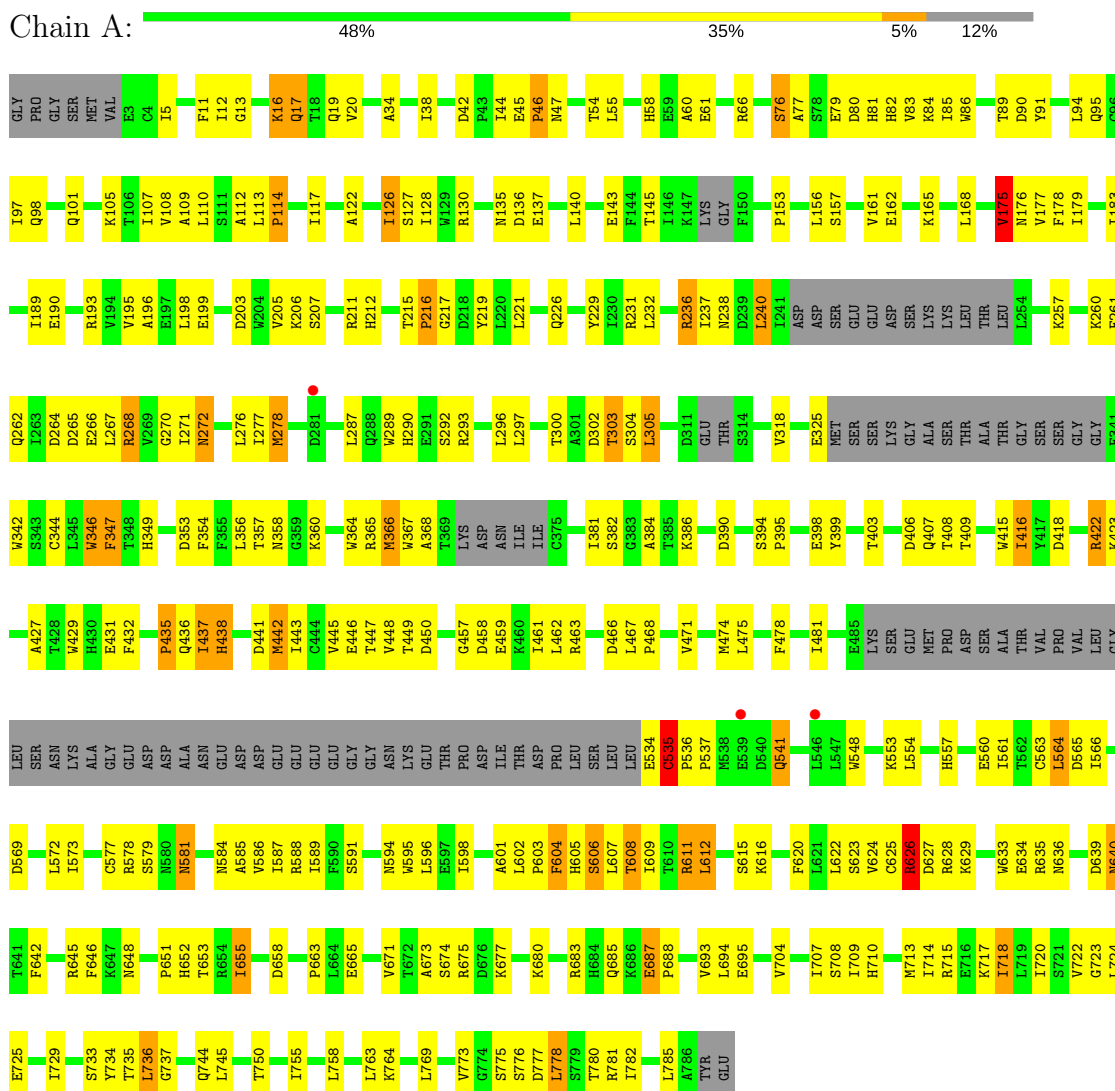
- Molecule 2 is water.

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 and DNA 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 48.34 – 3.17	Depositor EDS
% Data completeness (in resolution range)	87.8 (48.00-3.20) 86.6 (48.34-3.17)	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	828 reflections (5.14%)	DCC
Wilson B-factor (Å ²)	49.2	Xtriage
Anisotropy	0.243	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 49.4	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 [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	0.32	0/5503	0.59	0/7483

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	5374	0	5141	270	0
2	A	5	0	0	0	0
All	All	5379	0	5141	270	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 26.

All (270) 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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:403:THR:HB	1:A:442:MET:HB3	1.67	0.77
1:A:60:ALA:HB1	1:A:79:GLU:HB2	1.66	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:N	1:A:240:LEU:HD23	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
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:79:GLU:OE2	1:A:79:GLU:HA	1.98	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:264:ASP:C	1:A:266:GLU:H	2.03	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:457:GLY:C	1:A:459:GLU:H	2.04	0.59
1:A:212:HIS:HB2	1:A:219:TYR:CE1	2.37	0.59
1:A:5:ILE:HD12	1:A:5:ILE:N	2.18	0.59
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:290:HIS:CD2	1:A:292:SER:H	2.21	0.58
1:A:13:GLY:HA3	1:A:384:ALA:O	2.03	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:429:TRP:CZ3	1:A:782:ILE:HD13	2.40	0.57
1:A:5:ILE:H	1:A:5:ILE:HD12	1.68	0.57
1:A:107:ILE:HD12	1:A:107:ILE:N	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: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:297:LEU:HD21	1:A:305:LEU:CD2	2.34	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:561:ILE:HD12	1:A:561:ILE:N	2.23	0.54
1:A:581:ASN:H	1:A:581:ASN:HD22	1.55	0.54
1:A:19:GLN:NE2	1:A:66:ARG:HH21	2.03	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: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
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: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:215:THR:HA	1:A:216:PRO:C	2.30	0.52
1:A:665:GLU:HB3	1:A:713:MET:HE2	1.92	0.51
1:A:675:ARG:C	1:A:677:LYS:H	2.13	0.51
1:A:777:ASP:C	1:A:778:LEU:HG	2.31	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:723:GLY:C	1:A:724:LEU:HD12	2.32	0.51
1:A:176:ASN:ND2	1:A:199:GLU:HA	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:H	1:A:360:LYS:NZ	2.09	0.50
1:A:156:LEU:HD12	1:A:156:LEU:O	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:CG	1:A:236:ARG:HH11	2.22	0.49
1:A:398:GLU:HG2	1:A:481:ILE:CG2	2.43	0.49
1:A:54:THR:HG21	1:A:781:ARG:NH2	2.28	0.49
1:A:457:GLY:C	1:A:459:GLU:N	2.66	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
1:A:237:ILE:HG22	1:A:238:ASN:N	2.27	0.48
1:A:177:VAL:HB	1:A:198:LEU:HB2	1.94	0.48
1:A:687:GLU:HB2	1:A:688:PRO:CD	2.37	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:85:ILE:O	1:A:97:ILE:HG12	2.12	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:N	1:A:724:LEU:HD12	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:694:LEU:HD23	1:A:694:LEU:C	2.35	0.48
1:A:744:GLN:O	1:A:745:LEU:C	2.52	0.48
1:A:110:LEU:C	1:A:110:LEU:HD23	2.34	0.47
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:394:SER:HB3	1:A:399:TYR:CE2	2.50	0.47
1:A:342:TRP:CE2	1:A:360:LYS:HA	2.50	0.47
1:A:609:ILE:HG21	1:A:612:LEU:CD1	2.45	0.47
1:A:109:ALA:HB1	1:A:156:LEU:HG	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:ILE:C	1:A:183:ILE:HD12	2.35	0.47
1:A:360:LYS:H	1:A:360:LYS:CE	2.28	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:461:ILE:HD11	1:A:463:ARG:HE	1.79	0.47
1:A:38:ILE:HB	1:A:55:LEU:HB2	1.97	0.47
1:A:271:ILE:HG22	1:A:272:ASN:N	2.30	0.47
1:A:386:LYS:HB3	1:A:386:LYS:NZ	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:264:ASP:C	1:A:266:GLU:N	2.68	0.46
1:A:652:HIS:ND1	1:A:674:SER:HB3	2.31	0.46
1:A:429:TRP:CE3	1:A:782:ILE:HD13	2.51	0.46
1:A:763:LEU:HD12	1:A:764:LYS:H	1.80	0.46
1:A:462:LEU:HD23	1:A:462:LEU:H	1.80	0.46
1:A:58:HIS:CE1	1:A:84:LYS:HD2	2.50	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
1:A:135:ASN:HD22	1:A:137:GLU:HB2	1.78	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:189:ILE:CG2	1:A:190:GLU:N	2.79	0.45
1:A:177:VAL:HG22	1:A:205:VAL:HG21	1.99	0.45
1:A:395:PRO:HD2	1:A:450:ASP:OD2	2.17	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:HG3	1:A:645:ARG:HH11	1.82	0.45
1:A:745:LEU:HB3	1:A:750:THR:CG2	2.47	0.45
1:A:195:VAL:O	1:A:195:VAL:HG22	2.18	0.44
1:A:611:ARG:HB3	1:A:624:VAL:HG22	1.99	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:602:LEU:HB3	1:A:633:TRP:CZ3	2.52	0.44
1:A:80:ASP:OD1	1:A:82:HIS:CG	2.71	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:297:LEU:HB2	1:A:346:TRP:CD1	2.53	0.44
1:A:591:SER:HB2	1:A:598:ILE:HD11	2.00	0.44
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: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:459:GLU:O	1:A:560:GLU:HA	2.18	0.43
1:A:61:GLU:HB3	1:A:79:GLU:HG2	2.00	0.43
1:A:128:ILE:CG2	1:A:189:ILE:HD11	2.48	0.42
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:553:LYS:C	1:A:554:LEU:HD23	2.39	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:687:GLU:N	1:A:687:GLU:CD	2.73	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:126:ILE:O	1:A:126:ILE:HG12	2.20	0.42
1:A:11:PHE:HB2	1:A:780:THR:HB	2.02	0.42
1:A:211:ARG:HB2	1:A:289:TRP:CE2	2.55	0.41
1:A:541:GLN:HB3	1:A:541:GLN:HE21	1.62	0.41
1:A:620:PHE:CD2	1:A:634:GLU:HA	2.55	0.41
1:A:94:LEU:N	1:A:94:LEU:HD12	2.35	0.41
1:A:409:THR:HB	1:A:436:GLN:HG2	2.01	0.41
1:A:231:ARG:C	1:A:232:LEU:HD12	2.41	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:277:ILE:CD1	1:A:296:LEU:HD23	2.51	0.41
1:A:12:ILE:HD13	1:A:34:ALA:HB3	2.03	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:584:ASN:C	1:A:586:VAL:H	2.24	0.41
1:A:625:CYS:O	1:A:626:ARG:C	2.59	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
1:A:84:LYS:C	1:A:85:ILE:HD12	2.41	0.41
1:A:277:ILE:O	1:A:278:MET:C	2.58	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:429:TRP:CE3	1:A:782:ILE:CD1	3.04	0.41
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:195:VAL:CG2	1:A:261:PHE:HB2	2.52	0.40
1:A:360:LYS:HE2	1:A:360:LYS:H	1.86	0.40
1:A:80:ASP:OD1	1:A:80:ASP:O	2.39	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%)	5 33

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	81	HIS

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Mol	Chain	Res	Type
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
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%)	498 (91%)	48 (9%)	12	42

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	LYS
1	A	17	GLN
1	A	47	ASN
1	A	76	SER
1	A	101	GLN
1	A	114	PRO
1	A	126	ILE

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Mol	Chain	Res	Type
1	A	130	ARG
1	A	136	ASP
1	A	145	THR
1	A	175	VAL
1	A	207	SER
1	A	221	LEU
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	347	PHE
1	A	365	ARG
1	A	366	MET
1	A	390	ASP
1	A	416	ILE
1	A	422	ARG
1	A	437	ILE
1	A	442	MET
1	A	466	ASP
1	A	535	CYS
1	A	541	GLN
1	A	564	LEU
1	A	569	ASP
1	A	581	ASN
1	A	606	SER
1	A	608	THR
1	A	611	ARG
1	A	612	LEU
1	A	626	ARG
1	A	640	ASN
1	A	653	THR
1	A	718	ILE
1	A	725	GLU
1	A	736	LEU
1	A	776	SER
1	A	778	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	19	GLN
1	A	48	ASN
1	A	82	HIS
1	A	101	GLN
1	A	131	GLN
1	A	132	ASN
1	A	135	ASN
1	A	176	ASN
1	A	212	HIS
1	A	258	GLN
1	A	288	GLN
1	A	290	HIS
1	A	358	ASN
1	A	557	HIS
1	A	581	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 carbohydrates 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 ⓘ

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%)	-0.46	3 (0%) 92 89	15, 42, 77, 105	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	539	GLU	3.2
1	A	281	ASP	2.3
1	A	546	LEU	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 carbohydrates 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.