



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:46 AM UTC

PDB ID : 2I7V / pdb_00002i7v
Title : Structure of Human CPSF-73
Authors : Mandel, C.R.; Zhang, H.; Tong, L.
Deposited on : 2006-08-31
Resolution : 2.10 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
Mogul : 2022.3.0, CSD as543be (2022)
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

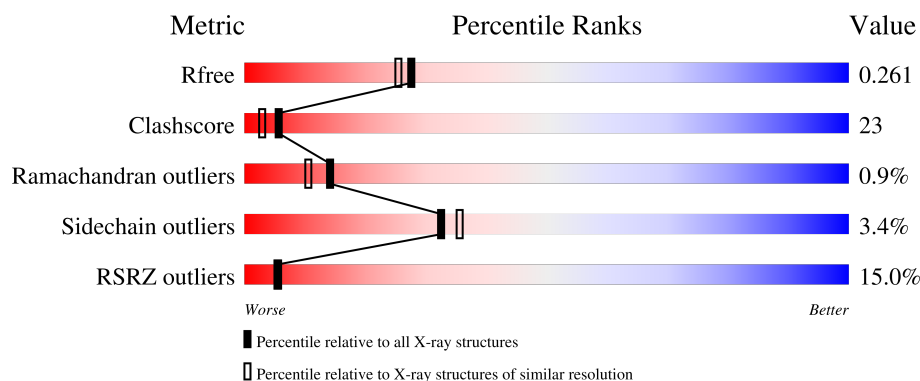
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	459	14% 56% 35% 6%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3727 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 Cleavage and polyadenylation specificity factor 73 kDa subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	433	Total	C	N	O	S	0	0	0
			3471	2214	591	642	24			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	4	Total	Zn	0	0
			4	4		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		

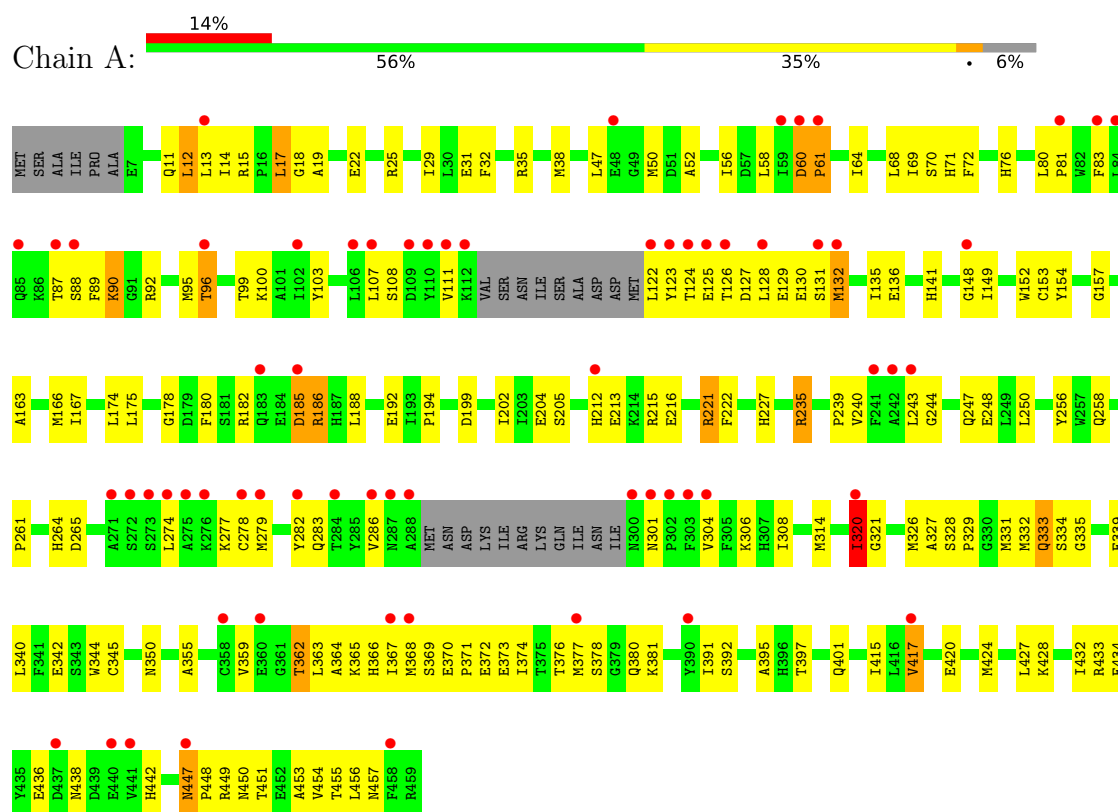
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	247	Total 247	O 247	0	0

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: Cleavage and polyadenylation specificity factor 73 kDa subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.50Å 83.20Å 104.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.25 – 2.10 29.25 – 2.10	Depositor EDS
% Data completeness (in resolution range)	96.7 (29.25-2.10) 96.9 (29.25-2.10)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.59 (at 2.10Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.224 , 0.252 0.233 , 0.261	Depositor DCC
R_{free} test set	4265 reflections (7.41%)	wwPDB-VP
Wilson B-factor (Å ²)	35.7	Xtriage
Anisotropy	0.152	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 44.5	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	0.94	EDS
Total number of atoms	3727	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.60% 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: SO4, ZN

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	0.41	0/3552	0.93	17/4802 (0.4%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	320	ILE	N-CA-C	8.88	119.71	111.45
1	A	448	PRO	N-CA-C	6.86	122.43	111.26
1	A	178	GLY	N-CA-C	-6.57	100.07	111.50
1	A	417	VAL	CB-CA-C	-6.49	100.65	111.29
1	A	17	LEU	N-CA-C	-6.39	105.42	113.72
1	A	149	ILE	N-CA-C	6.23	116.84	107.75
1	A	395	ALA	N-CA-C	6.20	118.17	110.91
1	A	417	VAL	N-CA-C	6.17	122.16	109.34
1	A	392	SER	N-CA-C	5.89	118.80	109.96
1	A	72	PHE	N-CA-C	5.67	120.04	113.18
1	A	186	ARG	N-CA-C	-5.55	106.34	113.23
1	A	345	CYS	N-CA-C	5.54	119.14	112.38
1	A	157	GLY	N-CA-C	-5.46	101.87	112.02
1	A	60	ASP	CA-C-N	5.34	126.52	119.84
1	A	60	ASP	C-N-CA	5.34	126.52	119.84
1	A	199	ASP	N-CA-C	-5.02	106.33	112.90
1	A	70	SER	N-CA-C	5.01	116.56	111.14

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	3471	0	3415	161	0
2	A	4	0	0	0	0
3	A	5	0	0	0	0
4	A	247	0	0	10	0
All	All	3727	0	3415	161	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 23.

All (161) 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:13:LEU:HD11	1:A:453:ALA:HB1	1.31	1.12
1:A:274:LEU:HD21	1:A:277:LYS:HD2	1.43	0.97
1:A:103:TYR:HE2	1:A:132:MET:HE1	1.37	0.89
1:A:376:THR:HG22	1:A:378:SER:H	1.35	0.89
1:A:61:PRO:HG3	1:A:87:THR:HG22	1.54	0.89
1:A:103:TYR:CE2	1:A:132:MET:HE1	2.12	0.83
1:A:279:MET:O	1:A:283:GLN:HG3	1.83	0.79
1:A:424:MET:HG2	1:A:447:ASN:HD21	1.45	0.79
1:A:244:GLY:H	1:A:247:GLN:NE2	1.82	0.77
1:A:342:GLU:OE1	1:A:376:THR:HG23	1.86	0.75
1:A:17:LEU:HD11	1:A:29:ILE:HD11	1.68	0.74
1:A:61:PRO:HG2	1:A:88:SER:OG	1.87	0.74
1:A:92:ARG:HD2	1:A:136:GLU:OE2	1.87	0.73
1:A:61:PRO:HG3	1:A:87:THR:CG2	2.17	0.72
1:A:154:TYR:CZ	1:A:194:PRO:HB3	2.26	0.70
1:A:174:LEU:HD11	1:A:202:ILE:HG13	1.74	0.70
1:A:424:MET:HG2	1:A:447:ASN:ND2	2.06	0.69
1:A:15:ARG:HB2	4:A:631:HOH:O	1.92	0.69
1:A:424:MET:HE2	1:A:447:ASN:ND2	2.07	0.69
1:A:186:ARG:HG2	1:A:215:ARG:HH21	1.56	0.68
1:A:243:LEU:HD21	1:A:278:CYS:HB3	1.74	0.68
1:A:96:THR:HG22	1:A:99:THR:H	1.59	0.67
1:A:424:MET:CE	1:A:447:ASN:ND2	2.58	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:449:ARG:HD3	4:A:644:HOH:O	1.95	0.66
1:A:212:HIS:CD2	1:A:213:GLU:H	2.13	0.65
1:A:35:ARG:HH21	1:A:148:GLY:HA3	1.61	0.65
1:A:240:VAL:HG21	1:A:326:MET:HE3	1.78	0.65
1:A:216:GLU:H	1:A:216:GLU:CD	2.05	0.64
1:A:125:GLU:O	1:A:129:GLU:HG2	1.97	0.64
1:A:274:LEU:HG	1:A:277:LYS:HB2	1.78	0.64
1:A:278:CYS:O	1:A:282:TYR:HB2	1.98	0.63
1:A:128:LEU:O	1:A:132:MET:HE3	1.99	0.63
1:A:132:MET:N	1:A:132:MET:HE2	2.13	0.63
1:A:258:GLN:CD	1:A:306:LYS:HD2	2.23	0.63
1:A:265:ASP:HB2	4:A:541:HOH:O	1.98	0.63
1:A:35:ARG:HE	1:A:148:GLY:H	1.44	0.62
1:A:283:GLN:O	1:A:286:VAL:HG13	2.00	0.62
1:A:397:THR:HG22	1:A:401:GLN:NE2	2.15	0.62
1:A:35:ARG:NH2	1:A:148:GLY:HA3	2.16	0.61
1:A:397:THR:HG22	1:A:401:GLN:HE21	1.65	0.61
1:A:152:TRP:CZ2	1:A:166:MET:HG2	2.36	0.61
1:A:227:HIS:HD2	1:A:256:TYR:OH	1.84	0.60
1:A:12:LEU:HD12	1:A:13:LEU:N	2.17	0.60
1:A:174:LEU:HD12	1:A:175:LEU:N	2.17	0.60
1:A:247:GLN:OE1	1:A:279:MET:HE1	2.02	0.60
1:A:100:LYS:O	1:A:103:TYR:HB3	2.02	0.59
1:A:12:LEU:HD13	1:A:32:PHE:HB2	1.83	0.59
1:A:424:MET:HE1	4:A:513:HOH:O	2.02	0.59
1:A:180:PHE:CE1	1:A:397:THR:HG21	2.37	0.59
1:A:359:VAL:O	1:A:362:THR:HB	2.04	0.57
1:A:61:PRO:CG	1:A:87:THR:HG22	2.31	0.57
1:A:279:MET:HE2	1:A:279:MET:HA	1.87	0.56
1:A:35:ARG:HH11	1:A:35:ARG:HG2	1.70	0.56
1:A:212:HIS:CG	1:A:213:GLU:H	2.24	0.56
1:A:363:LEU:O	1:A:367:ILE:HG13	2.06	0.56
1:A:13:LEU:HB3	1:A:31:GLU:HB2	1.87	0.56
1:A:239:PRO:HG3	1:A:332:MET:HE2	1.87	0.55
1:A:221:ARG:NH1	1:A:391:ILE:HG12	2.22	0.54
1:A:434:GLU:O	1:A:438:ASN:HB3	2.08	0.54
1:A:185:ASP:HB2	1:A:188:LEU:HB2	1.90	0.54
1:A:186:ARG:HG3	1:A:248:GLU:OE2	2.07	0.54
1:A:367:ILE:C	1:A:369:SER:H	2.14	0.53
1:A:362:THR:HG22	1:A:364:ALA:N	2.22	0.53
1:A:58:LEU:HD13	1:A:451:THR:HG23	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:ILE:HD12	1:A:69:ILE:N	2.24	0.52
1:A:314:MET:HE1	1:A:340:LEU:CD2	2.39	0.52
1:A:433:ARG:HD2	4:A:612:HOH:O	2.09	0.52
1:A:334:SER:HB2	1:A:377:MET:HE1	1.92	0.52
1:A:239:PRO:HA	1:A:327:ALA:O	2.09	0.51
1:A:244:GLY:H	1:A:247:GLN:HE22	1.52	0.51
1:A:331:MET:HB3	1:A:362:THR:HG23	1.93	0.51
1:A:362:THR:HG22	1:A:365:LYS:H	1.74	0.51
1:A:212:HIS:CD2	1:A:213:GLU:HG3	2.45	0.51
1:A:261:PRO:HA	1:A:264:HIS:CD2	2.46	0.51
1:A:314:MET:HE1	1:A:340:LEU:HD23	1.93	0.51
1:A:450:ASN:O	1:A:451:THR:HB	2.10	0.51
1:A:424:MET:HE3	1:A:447:ASN:HD22	1.76	0.50
1:A:141:HIS:HE1	1:A:192:GLU:OE1	1.94	0.50
1:A:373:GLU:OE1	1:A:381:LYS:HD2	2.10	0.50
1:A:50:MET:HE1	1:A:122:LEU:HD23	1.93	0.50
1:A:68:LEU:HD12	1:A:167:ILE:HD11	1.93	0.50
1:A:180:PHE:CZ	1:A:397:THR:HG21	2.47	0.49
1:A:107:LEU:O	1:A:111:VAL:HG23	2.12	0.49
1:A:38:MET:HE3	1:A:64:ILE:HD13	1.95	0.49
1:A:108:SER:HA	1:A:125:GLU:OE1	2.13	0.49
1:A:154:TYR:CE2	1:A:194:PRO:HB3	2.47	0.49
1:A:329:PRO:HD2	1:A:334:SER:O	2.13	0.48
1:A:342:GLU:CD	1:A:376:THR:HG23	2.38	0.48
1:A:320:ILE:HG23	1:A:321:GLY:N	2.28	0.48
1:A:274:LEU:HD11	1:A:277:LYS:HG3	1.95	0.48
1:A:127:ASP:HA	1:A:130:GLU:OE1	2.13	0.48
1:A:61:PRO:O	1:A:89:PHE:HA	2.13	0.48
1:A:335:GLY:O	1:A:339:GLU:HG3	2.13	0.48
1:A:222:PHE:HA	1:A:391:ILE:HD13	1.96	0.48
1:A:182:ARG:H	1:A:401:GLN:NE2	2.12	0.47
1:A:227:HIS:HE1	4:A:511:HOH:O	1.96	0.47
1:A:301:ASN:HB3	1:A:304:VAL:HG23	1.95	0.47
1:A:235:ARG:HG2	1:A:350:ASN:OD1	2.14	0.47
1:A:235:ARG:HG3	1:A:344:TRP:CE2	2.49	0.47
1:A:424:MET:HE3	1:A:447:ASN:HA	1.96	0.47
1:A:331:MET:HE1	1:A:359:VAL:HG23	1.97	0.47
1:A:374:ILE:C	1:A:374:ILE:HD12	2.39	0.47
1:A:240:VAL:HG21	1:A:326:MET:CE	2.44	0.47
1:A:320:ILE:HG23	1:A:321:GLY:H	1.80	0.47
1:A:333:GLN:OE1	1:A:333:GLN:HA	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:PRO:HG2	1:A:123:TYR:CD2	2.51	0.46
1:A:244:GLY:N	1:A:247:GLN:HE22	2.13	0.46
1:A:314:MET:HG3	4:A:720:HOH:O	2.14	0.46
1:A:61:PRO:HB3	1:A:83:PHE:CE1	2.50	0.46
1:A:126:THR:O	1:A:130:GLU:HG3	2.16	0.45
1:A:331:MET:HB3	1:A:362:THR:CG2	2.46	0.45
1:A:60:ASP:HA	1:A:61:PRO:HD2	1.76	0.45
1:A:212:HIS:CG	1:A:213:GLU:N	2.84	0.45
1:A:19:ALA:HA	1:A:22:GLU:HG2	1.98	0.45
1:A:370:GLU:O	1:A:370:GLU:HG3	2.16	0.45
1:A:35:ARG:HE	1:A:148:GLY:N	2.14	0.44
1:A:380:GLN:NE2	1:A:381:LYS:H	2.15	0.44
1:A:80:LEU:N	1:A:81:PRO:HD2	2.32	0.44
1:A:50:MET:CE	1:A:122:LEU:HD23	2.47	0.44
1:A:95:MET:HE3	1:A:135:ILE:HG21	1.99	0.44
1:A:153:CYS:HB2	1:A:163:ALA:HB1	1.99	0.44
1:A:107:LEU:HD12	1:A:128:LEU:HD13	1.99	0.44
1:A:433:ARG:NH2	4:A:732:HOH:O	2.51	0.44
1:A:13:LEU:HD21	1:A:15:ARG:HH11	1.82	0.44
1:A:250:LEU:HD22	1:A:308:ILE:HD13	2.00	0.44
1:A:455:THR:O	1:A:456:LEU:HD23	2.17	0.44
1:A:247:GLN:NE2	1:A:247:GLN:H	2.16	0.43
1:A:182:ARG:HB2	1:A:401:GLN:HE21	1.83	0.43
1:A:301:ASN:HB3	1:A:304:VAL:CG2	2.49	0.43
1:A:13:LEU:HD12	1:A:454:VAL:O	2.18	0.43
1:A:18:GLY:HA3	1:A:25:ARG:O	2.18	0.43
1:A:38:MET:HE3	1:A:64:ILE:CD1	2.49	0.43
1:A:247:GLN:HG2	1:A:282:TYR:CE2	2.53	0.43
1:A:366:HIS:HD2	4:A:591:HOH:O	2.01	0.43
1:A:131:SER:O	1:A:132:MET:C	2.61	0.43
1:A:87:THR:HG22	1:A:88:SER:H	1.84	0.43
1:A:328:SER:HB2	1:A:329:PRO:HA	2.00	0.43
1:A:369:SER:O	1:A:370:GLU:HB3	2.19	0.43
1:A:17:LEU:HD13	1:A:56:ILE:HA	2.00	0.43
1:A:92:ARG:CD	1:A:136:GLU:OE2	2.63	0.43
1:A:243:LEU:HA	1:A:247:GLN:HE22	1.84	0.42
1:A:17:LEU:HD11	1:A:29:ILE:CD1	2.45	0.42
1:A:90:LYS:HD2	1:A:90:LYS:C	2.44	0.42
1:A:436:GLU:OE2	1:A:442:HIS:HA	2.19	0.42
1:A:107:LEU:HD22	1:A:123:TYR:OH	2.19	0.42
1:A:355:ALA:HA	1:A:391:ILE:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:LEU:HD12	1:A:13:LEU:H	1.83	0.42
1:A:205:SER:HB3	1:A:427:LEU:HD23	2.00	0.42
1:A:124:THR:C	1:A:126:THR:N	2.78	0.41
1:A:47:LEU:HB2	1:A:52:ALA:HA	2.03	0.41
1:A:71:HIS:CE1	1:A:76:HIS:CD2	3.09	0.41
1:A:370:GLU:N	1:A:371:PRO:CD	2.84	0.41
1:A:380:GLN:HE21	1:A:381:LYS:H	1.69	0.41
1:A:14:ILE:CD1	1:A:415:ILE:HD12	2.50	0.41
1:A:81:PRO:HG2	1:A:123:TYR:CE2	2.56	0.41
1:A:331:MET:HB2	4:A:524:HOH:O	2.21	0.40
1:A:329:PRO:CD	1:A:334:SER:O	2.70	0.40
1:A:428:LYS:O	1:A:432:ILE:HG13	2.22	0.40
1:A:451:THR:O	1:A:451:THR:HG22	2.21	0.40
1:A:11:GLN:HG3	1:A:457:ASN:OD1	2.22	0.40
1:A:367:ILE:C	1:A:369:SER:N	2.78	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	427/459 (93%)	394 (92%)	29 (7%)	4 (1%)	14 10

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	368	MET
1	A	132	MET
1	A	417	VAL
1	A	61	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	377/400 (94%)	364 (97%)	13 (3%)	32 35

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	LEU
1	A	90	LYS
1	A	96	THR
1	A	185	ASP
1	A	204	GLU
1	A	221	ARG
1	A	235	ARG
1	A	320	ILE
1	A	333	GLN
1	A	362	THR
1	A	372	GLU
1	A	420	GLU
1	A	447	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	GLN
1	A	97	HIS
1	A	141	HIS
1	A	212	HIS
1	A	227	HIS
1	A	247	GLN
1	A	307	HIS
1	A	310	ASN
1	A	316	HIS
1	A	380	GLN
1	A	401	GLN
1	A	413	HIS

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Mol	Chain	Res	Type
1	A	422	ASN
1	A	447	ASN
1	A	457	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	SO4	A	491	2	4,4,4	0.39	0	6,6,6	0.07	0

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

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 ⓘ

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	433/459 (94%)	0.82	65 (15%) 5 5	22, 42, 85, 99	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	273	SER	6.0
1	A	212	HIS	5.4
1	A	107	LEU	4.2
1	A	243	LEU	4.2
1	A	278	CYS	3.9
1	A	110	TYR	3.8
1	A	367	ILE	3.8
1	A	123	TYR	3.8
1	A	111	VAL	3.7
1	A	320	ILE	3.6
1	A	128	LEU	3.6
1	A	288	ALA	3.5
1	A	131	SER	3.5
1	A	122	LEU	3.4
1	A	102	ILE	3.3
1	A	417	VAL	3.3
1	A	441	VAL	3.3
1	A	148	GLY	3.2
1	A	274	LEU	3.2
1	A	124	THR	3.2
1	A	125	GLU	3.2
1	A	360	GLU	3.0
1	A	185	ASP	3.0
1	A	275	ALA	3.0
1	A	61	PRO	3.0
1	A	84	LEU	3.0
1	A	241	PHE	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	282	TYR	2.9
1	A	126	THR	2.8
1	A	242	ALA	2.8
1	A	301	ASN	2.8
1	A	109	ASP	2.7
1	A	271	ALA	2.6
1	A	106	LEU	2.6
1	A	287	ASN	2.6
1	A	60	ASP	2.5
1	A	81	PRO	2.5
1	A	272	SER	2.5
1	A	85	GLN	2.5
1	A	88	SER	2.5
1	A	112	LYS	2.5
1	A	447	ASN	2.5
1	A	358	CYS	2.4
1	A	87	THR	2.4
1	A	132	MET	2.4
1	A	48	GLU	2.3
1	A	390	TYR	2.3
1	A	377	MET	2.3
1	A	286	VAL	2.3
1	A	458	PHE	2.3
1	A	279	MET	2.3
1	A	440	GLU	2.2
1	A	368	MET	2.2
1	A	437	ASP	2.2
1	A	284	THR	2.2
1	A	302	PRO	2.1
1	A	276	LYS	2.1
1	A	96	THR	2.1
1	A	59	ILE	2.1
1	A	13	LEU	2.1
1	A	183	GLN	2.1
1	A	303	PHE	2.1
1	A	83	PHE	2.1
1	A	304	VAL	2.1
1	A	300	ASN	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](#)

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
3	SO4	A	491	5/5	0.89	0.12	68,68,70,71	0
2	ZN	A	484	1/1	0.90	0.12	98,98,98,98	0
2	ZN	A	483	1/1	0.99	0.03	49,49,49,49	0
2	ZN	A	481	1/1	0.99	0.06	38,38,38,38	0
2	ZN	A	482	1/1	0.99	0.02	38,38,38,38	0

6.5 Other polymers [i](#)

There are no such residues in this entry.