



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 07:56 am GMT

PDB ID : 2OAJ
Title : Crystal structure of Sro7 from *S. cerevisiae*
Authors : Hattendorf, D.A.; Weis, W.I.
Deposited on : 2006-12-15
Resolution : 2.40 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

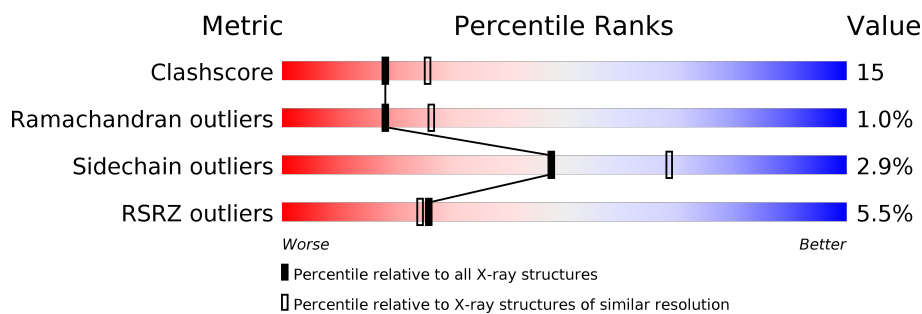
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	112137	3674 (2.40-2.40)
Ramachandran outliers	110173	3616 (2.40-2.40)
Sidechain outliers	110143	3617 (2.40-2.40)
RSRZ outliers	101464	3195 (2.40-2.40)

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	902	<div> <div>5%</div> <div>70%</div> <div>25%</div> <div>..</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NA	A	964	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7118 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 Protein SNI1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	875	Total	C	N	O	S	0	0	0
			6830	4367	1142	1292	29			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Na	0	0
			1	1		

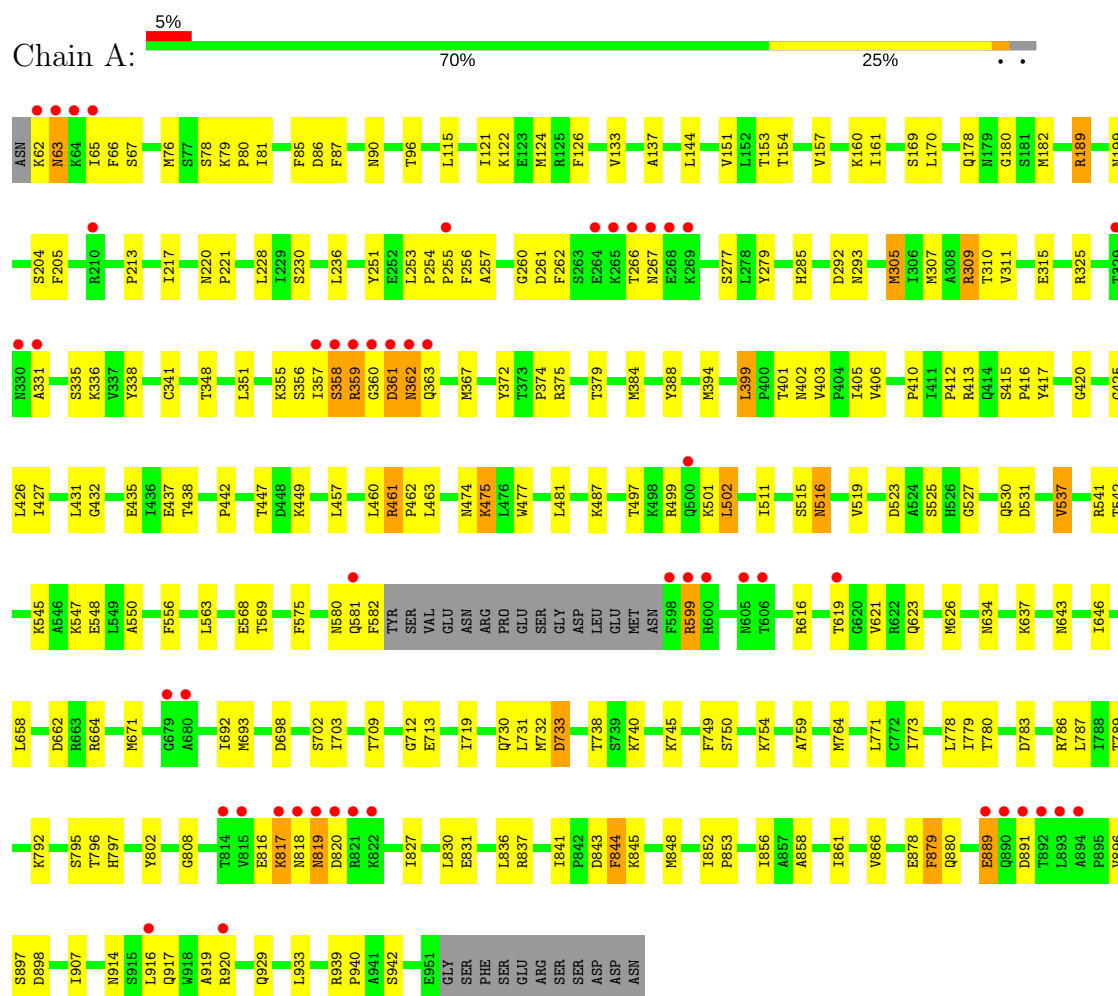
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	286	Total	O	0	0
			286	286		

3 Residue-property plots [i](#)

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: Protein SNI1



4 Data and refinement statistics

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α , β , γ	130.52Å 130.52Å 116.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.12 – 2.40 56.88 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.4 (49.12-2.40) 99.5 (56.88-2.40)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.18 (at 2.40Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.211 , 0.262 0.214 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	24.3	Xtriage
Anisotropy	0.333	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 39.0	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	0.93	EDS
Total number of atoms	7118	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.28% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NA, 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	0.35	0/6972	0.63	0/9462

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	6830	0	6897	206	0
2	A	1	0	0	0	0
3	A	1	0	0	0	0
4	A	286	0	0	11	0
All	All	7118	0	6897	206	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 15.

All (206) 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:502:LEU:HD23	1:A:502:LEU:H	1.30	0.96
1:A:401:THR:HG22	1:A:403:VAL:H	1.29	0.96
1:A:830:LEU:HD21	1:A:861:ILE:HD11	1.48	0.93
1:A:581:GLN:O	1:A:582:PHE:HB2	1.78	0.81
1:A:435:GLU:HG2	1:A:463:LEU:HG	1.63	0.81
1:A:545:LYS:HZ1	1:A:569:THR:HG21	1.48	0.79
1:A:122:LYS:HE3	1:A:137:ALA:HA	1.66	0.77
1:A:447:THR:HG22	1:A:449:LYS:H	1.50	0.75
1:A:796:THR:HG22	1:A:797:HIS:H	1.51	0.74
1:A:599:ARG:HD2	1:A:599:ARG:H	1.53	0.73
1:A:66:PHE:HB3	1:A:848:MET:HE3	1.73	0.71
1:A:182:MET:HE2	1:A:217:ILE:HD13	1.72	0.70
1:A:309:ARG:HG3	1:A:309:ARG:HH11	1.57	0.70
1:A:475:LYS:HE3	1:A:475:LYS:HA	1.75	0.69
1:A:460:LEU:HD21	1:A:879:PHE:HB3	1.73	0.68
1:A:384:MET:HE2	1:A:384:MET:HA	1.75	0.68
1:A:292:ASP:O	1:A:293:ASN:HB2	1.91	0.68
1:A:462:PRO:HG3	1:A:516:ASN:HB2	1.76	0.67
1:A:76:MET:SD	1:A:80:PRO:HG3	2.34	0.67
1:A:384:MET:HA	1:A:384:MET:CE	2.25	0.67
1:A:357:ILE:HG12	1:A:362:ASN:ND2	2.10	0.67
1:A:180:GLY:HA3	1:A:199:ASN:ND2	2.10	0.66
1:A:157:VAL:HG21	1:A:161:ILE:HD11	1.77	0.66
1:A:358:SER:O	1:A:359:ARG:HB2	1.94	0.66
1:A:356:SER:HA	1:A:362:ASN:HD22	1.60	0.66
1:A:261:ASP:HA	1:A:325:ARG:NH1	2.11	0.65
1:A:374:PRO:HB2	1:A:384:MET:HE3	1.77	0.65
1:A:740:LYS:HD3	1:A:783:ASP:OD2	1.97	0.65
1:A:580:ASN:OD1	1:A:581:GLN:N	2.30	0.65
1:A:749:PHE:CE1	1:A:841:ILE:HD12	2.32	0.65
1:A:548:GLU:HA	1:A:568:GLU:OE1	1.97	0.64
1:A:545:LYS:HE2	1:A:547:LYS:O	1.98	0.64
1:A:502:LEU:CD2	1:A:502:LEU:H	2.08	0.64
1:A:412:PRO:HG3	1:A:420:GLY:O	1.98	0.64
1:A:221:PRO:HG3	4:A:1226:HOH:O	1.98	0.63
1:A:361:ASP:HA	1:A:402:ASN:HA	1.80	0.63
1:A:435:GLU:CG	1:A:463:LEU:HG	2.28	0.62
1:A:417:TYR:HB3	4:A:1002:HOH:O	1.98	0.62
1:A:361:ASP:H	1:A:402:ASN:HB2	1.64	0.62
1:A:787:LEU:HD11	1:A:844:PHE:HE1	1.64	0.61
1:A:856:ILE:HG21	1:A:861:ILE:HD13	1.82	0.61
1:A:356:SER:HB3	1:A:363:GLN:HE21	1.63	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:356:SER:CB	1:A:363:GLN:HE21	2.14	0.60
1:A:357:ILE:O	1:A:358:SER:HB2	2.01	0.60
1:A:257:ALA:O	1:A:266:THR:HG23	2.02	0.60
1:A:796:THR:HG22	1:A:797:HIS:N	2.16	0.60
1:A:599:ARG:HD2	1:A:599:ARG:N	2.16	0.60
1:A:230:SER:OG	1:A:277:SER:HB2	2.01	0.60
1:A:76:MET:O	1:A:880:GLN:HB2	2.02	0.59
1:A:361:ASP:H	1:A:402:ASN:CB	2.14	0.59
1:A:778:LEU:HD13	1:A:827:ILE:HD12	1.83	0.59
1:A:311:VAL:HG23	4:A:1212:HOH:O	2.02	0.59
1:A:255:PRO:HB3	1:A:267:ASN:O	2.02	0.59
1:A:66:PHE:HB3	1:A:848:MET:CE	2.33	0.58
1:A:153:THR:HG22	1:A:154:THR:N	2.18	0.58
1:A:86:ASP:HB2	1:A:126:PHE:CG	2.37	0.58
1:A:309:ARG:NH1	1:A:309:ARG:HG3	2.20	0.57
1:A:523:ASP:OD1	1:A:525:SER:HB3	2.05	0.56
1:A:497:THR:HG21	1:A:896:VAL:HG11	1.88	0.55
1:A:786:ARG:HD3	1:A:797:HIS:HB3	1.87	0.55
1:A:817:LYS:HB2	1:A:820:ASP:HB3	1.88	0.55
1:A:501:LYS:HE2	4:A:1196:HOH:O	2.07	0.55
1:A:457:LEU:O	1:A:461:ARG:HG2	2.08	0.54
1:A:401:THR:HG22	1:A:402:ASN:N	2.23	0.54
1:A:375:ARG:N	1:A:375:ARG:HD2	2.23	0.54
1:A:401:THR:HG22	1:A:403:VAL:N	2.11	0.54
1:A:305:MET:HE3	1:A:307:MET:N	2.23	0.54
1:A:836:LEU:HD23	1:A:837:ARG:N	2.23	0.53
1:A:375:ARG:HH21	1:A:917:GLN:NE2	2.06	0.53
1:A:745:LYS:HB3	1:A:780:THR:HB	1.91	0.53
1:A:81:ILE:HD12	1:A:406:VAL:HG11	1.90	0.53
1:A:260:GLY:HA3	1:A:292:ASP:O	2.10	0.52
1:A:401:THR:CG2	1:A:403:VAL:HG23	2.38	0.52
1:A:698:ASP:HB2	4:A:1179:HOH:O	2.09	0.52
1:A:160:LYS:HB3	1:A:178:GLN:HB2	1.92	0.52
1:A:646:ILE:HB	1:A:764:MET:HE1	1.92	0.52
1:A:361:ASP:C	1:A:363:GLN:H	2.13	0.51
1:A:787:LEU:O	1:A:795:SER:HB2	2.10	0.51
1:A:730:GLN:HB3	1:A:732:MET:HE3	1.92	0.51
1:A:356:SER:HB2	1:A:362:ASN:HB3	1.93	0.51
1:A:379:THR:HG21	1:A:384:MET:HE2	1.91	0.51
1:A:80:PRO:HG2	1:A:460:LEU:HD13	1.93	0.51
1:A:712:GLY:HA3	1:A:738:THR:O	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:581:GLN:O	1:A:582:PHE:CB	2.51	0.51
1:A:523:ASP:HB3	1:A:530:GLN:HB3	1.93	0.51
1:A:285:HIS:HD2	1:A:388:TYR:OH	1.94	0.50
1:A:341:CYS:HA	1:A:348:THR:HA	1.93	0.50
1:A:914:ASN:HB2	1:A:917:GLN:HB2	1.93	0.50
1:A:178:GLN:NE2	1:A:213:PRO:HB3	2.26	0.50
1:A:63:ASN:C	1:A:65:ILE:H	2.13	0.50
1:A:307:MET:SD	1:A:309:ARG:HD2	2.51	0.50
1:A:789:THR:HG21	1:A:792:LYS:HD2	1.94	0.50
1:A:817:LYS:HD2	1:A:820:ASP:CB	2.41	0.50
1:A:462:PRO:CG	1:A:516:ASN:HB2	2.42	0.49
1:A:487:LYS:O	1:A:487:LYS:HG2	2.12	0.49
1:A:399:LEU:HD21	1:A:405:ILE:HD11	1.93	0.49
1:A:351:LEU:CD2	1:A:427:ILE:HD13	2.42	0.49
1:A:236:LEU:C	1:A:236:LEU:HD12	2.32	0.49
1:A:580:ASN:HB2	1:A:626:MET:SD	2.53	0.49
1:A:750:SER:O	1:A:754:LYS:HA	2.11	0.49
1:A:463:LEU:HD13	1:A:878:GLU:CD	2.33	0.49
1:A:730:GLN:O	1:A:732:MET:HE2	2.12	0.49
1:A:787:LEU:HD11	1:A:844:PHE:CE1	2.46	0.49
1:A:537:VAL:HG23	4:A:1186:HOH:O	2.11	0.49
1:A:374:PRO:HG3	4:A:1054:HOH:O	2.12	0.48
1:A:361:ASP:OD1	1:A:361:ASP:O	2.30	0.48
1:A:375:ARG:NH2	1:A:917:GLN:NE2	2.61	0.48
1:A:568:GLU:O	1:A:637:LYS:HE2	2.14	0.48
1:A:62:LYS:HD2	1:A:845:LYS:HE2	1.95	0.48
1:A:502:LEU:HD11	1:A:891:ASP:CB	2.43	0.48
1:A:379:THR:HG21	1:A:384:MET:CE	2.44	0.48
1:A:942:SER:HA	4:A:1043:HOH:O	2.13	0.48
1:A:236:LEU:O	1:A:236:LEU:HD12	2.12	0.47
1:A:541:ARG:NH1	1:A:616:ARG:O	2.47	0.47
1:A:255:PRO:O	1:A:256:PHE:HB2	2.14	0.47
1:A:384:MET:HE2	1:A:384:MET:CA	2.44	0.47
1:A:426:LEU:HD12	1:A:427:ILE:H	1.79	0.47
1:A:502:LEU:N	1:A:502:LEU:HD23	2.12	0.47
1:A:732:MET:O	1:A:733:ASP:HB3	2.14	0.47
1:A:351:LEU:HD22	1:A:427:ILE:HD13	1.97	0.47
1:A:204:SER:HB2	1:A:205:PHE:CE1	2.50	0.47
1:A:410:PRO:O	1:A:412:PRO:HD3	2.14	0.47
1:A:189:ARG:NH1	1:A:940:PRO:O	2.47	0.47
1:A:519:VAL:HB	1:A:537:VAL:HG12	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:929:GLN:O	1:A:933:LEU:HG	2.14	0.47
1:A:499:ARG:NH2	1:A:898:ASP:OD1	2.48	0.47
1:A:889:GLU:OE1	1:A:889:GLU:HA	2.14	0.47
1:A:919:ALA:O	1:A:920:ARG:HB2	2.15	0.47
1:A:703:ILE:HD13	1:A:764:MET:HE2	1.97	0.46
1:A:144:LEU:CD2	1:A:151:VAL:HG22	2.45	0.46
1:A:262:PHE:CE1	1:A:331:ALA:HB1	2.50	0.46
1:A:81:ILE:CD1	1:A:406:VAL:HG11	2.46	0.46
1:A:852:ILE:HG21	1:A:856:ILE:HD11	1.97	0.46
1:A:124:MET:HA	1:A:133:VAL:O	2.15	0.45
1:A:693:MET:HE2	4:A:1229:HOH:O	2.17	0.45
1:A:786:ARG:HD3	1:A:797:HIS:CB	2.46	0.45
1:A:817:LYS:HB2	1:A:820:ASP:CB	2.46	0.45
1:A:545:LYS:NZ	1:A:569:THR:HG21	2.26	0.45
1:A:537:VAL:HG21	1:A:575:PHE:CE2	2.52	0.45
1:A:527:GLY:O	1:A:531:ASP:HB2	2.17	0.45
1:A:285:HIS:HE1	4:A:1180:HOH:O	2.00	0.45
1:A:153:THR:CG2	1:A:154:THR:N	2.80	0.45
1:A:251:TYR:OH	1:A:253:LEU:HD21	2.17	0.44
1:A:621:VAL:HG12	1:A:623:GLN:O	2.18	0.44
1:A:87:PHE:HB3	1:A:416:PRO:HB3	2.00	0.44
1:A:426:LEU:HD11	1:A:438:THR:HB	1.99	0.44
1:A:779:ILE:HB	1:A:786:ARG:HB2	1.99	0.44
1:A:220:ASN:HA	1:A:279:TYR:CE2	2.52	0.44
1:A:818:ASN:O	1:A:819:ASN:HB2	2.18	0.44
1:A:375:ARG:HE	1:A:917:GLN:NE2	2.15	0.44
1:A:515:SER:HA	1:A:550:ALA:HA	2.00	0.44
1:A:802:TYR:HB2	1:A:831:GLU:OE1	2.18	0.44
1:A:475:LYS:CA	1:A:475:LYS:HE3	2.46	0.43
1:A:545:LYS:HE2	1:A:548:GLU:HB2	2.00	0.43
1:A:646:ILE:HD13	1:A:759:ALA:HB1	2.00	0.43
1:A:413:ARG:CZ	1:A:425:GLY:HA3	2.48	0.43
1:A:771:LEU:HD13	1:A:773:ILE:HG13	2.01	0.43
1:A:67:SER:N	1:A:848:MET:HE2	2.34	0.43
1:A:309:ARG:NH1	1:A:310:THR:O	2.51	0.43
1:A:474:ASN:HA	1:A:477:TRP:HB3	2.01	0.43
1:A:169:SER:HB3	1:A:907:ILE:HD11	1.99	0.43
1:A:115:LEU:HD12	1:A:121:ILE:HD11	2.01	0.43
1:A:375:ARG:NE	1:A:917:GLN:HE22	2.17	0.42
1:A:401:THR:CG2	1:A:402:ASN:N	2.81	0.42
1:A:477:TRP:CE2	1:A:481:LEU:HD11	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:740:LYS:HA	1:A:740:LYS:HE2	2.00	0.42
1:A:939:ARG:HA	1:A:940:PRO:HD3	1.82	0.42
1:A:662:ASP:OD1	1:A:664:ARG:HB2	2.19	0.42
1:A:305:MET:HE1	1:A:307:MET:HA	2.02	0.42
1:A:335:SER:HB2	1:A:355:LYS:HG2	2.02	0.42
1:A:406:VAL:HG21	1:A:432:GLY:HA2	2.01	0.42
1:A:375:ARG:HE	1:A:917:GLN:HE22	1.68	0.42
1:A:78:SER:OG	1:A:96:THR:HB	2.19	0.42
1:A:363:GLN:O	1:A:399:LEU:HB2	2.19	0.42
1:A:709:THR:OG1	1:A:713:GLU:HB2	2.19	0.42
1:A:808:GLY:HA3	1:A:866:VAL:HB	2.02	0.42
1:A:220:ASN:HA	1:A:221:PRO:HD3	1.89	0.42
1:A:858:ALA:O	1:A:861:ILE:HG22	2.18	0.42
1:A:896:VAL:HG12	1:A:897:SER:N	2.35	0.42
1:A:415:SER:HB2	1:A:416:PRO:HD2	2.02	0.42
1:A:599:ARG:HH11	1:A:599:ARG:HG2	1.84	0.42
1:A:749:PHE:CZ	1:A:841:ILE:HD12	2.53	0.42
1:A:394:MET:HE1	4:A:1012:HOH:O	2.18	0.42
1:A:836:LEU:HD23	1:A:836:LEU:C	2.40	0.42
1:A:357:ILE:H	1:A:362:ASN:ND2	2.18	0.42
1:A:356:SER:HB3	1:A:363:GLN:HG3	2.02	0.41
1:A:658:LEU:O	1:A:671:MET:HA	2.20	0.41
1:A:542:THR:O	1:A:616:ARG:HD3	2.21	0.41
1:A:63:ASN:C	1:A:65:ILE:N	2.73	0.41
1:A:228:LEU:HB2	1:A:279:TYR:CE2	2.56	0.41
1:A:372:TYR:OH	1:A:375:ARG:HD3	2.19	0.41
1:A:375:ARG:NE	1:A:917:GLN:NE2	2.69	0.41
1:A:692:ILE:HA	1:A:702:SER:O	2.21	0.41
1:A:914:ASN:C	1:A:916:LEU:N	2.74	0.41
1:A:254:PRO:HD2	1:A:257:ALA:HB2	2.01	0.41
1:A:336:LYS:HD3	1:A:338:TYR:OH	2.21	0.41
1:A:360:GLY:O	1:A:362:ASN:N	2.54	0.41
1:A:497:THR:HG21	1:A:896:VAL:CG1	2.51	0.41
1:A:367:MET:HE2	1:A:442:PRO:CB	2.51	0.41
1:A:85:PHE:CD2	1:A:412:PRO:HD2	2.56	0.41
1:A:703:ILE:HG22	1:A:719:ILE:HD12	2.03	0.41
1:A:79:LYS:HA	1:A:80:PRO:HD3	1.81	0.41
1:A:852:ILE:HA	1:A:853:PRO:HD3	1.87	0.41
1:A:817:LYS:HD2	1:A:820:ASP:HB2	2.04	0.40
1:A:511:ILE:HD11	1:A:556:PHE:CG	2.57	0.40
1:A:431:LEU:HD11	1:A:437:GLU:OE2	2.20	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	871/902 (97%)	817 (94%)	45 (5%)	9 (1%)	18 26

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	359	ARG
1	A	817	LYS
1	A	819	ASN
1	A	63	ASN
1	A	361	ASP
1	A	362	ASN
1	A	358	SER
1	A	816	GLU
1	A	619	THR

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	762/787 (97%)	740 (97%)	22 (3%)	48 68

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	90	ASN
1	A	170	LEU
1	A	189	ARG
1	A	305	MET
1	A	309	ARG
1	A	315	GLU
1	A	399	LEU
1	A	461	ARG
1	A	475	LYS
1	A	502	LEU
1	A	516	ASN
1	A	537	VAL
1	A	563	LEU
1	A	599	ARG
1	A	634	ASN
1	A	643	ASN
1	A	731	LEU
1	A	733	ASP
1	A	843	ASP
1	A	844	PHE
1	A	879	PHE
1	A	889	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	285	HIS
1	A	362	ASN
1	A	363	GLN
1	A	516	ASN
1	A	581	GLN
1	A	819	ASN
1	A	917	GLN

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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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 [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	875/902 (97%)	0.13	48 (5%) 26 24	8, 22, 55, 85	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	892	THR	5.8
1	A	893	LEU	5.5
1	A	600	ARG	5.2
1	A	815	VAL	5.0
1	A	819	ASN	4.6
1	A	63	ASN	4.5
1	A	889	GLU	4.4
1	A	894	ALA	4.3
1	A	357	ILE	4.1
1	A	605	ASN	4.0
1	A	62	LYS	4.0
1	A	891	ASP	3.9
1	A	329	THR	3.8
1	A	822	LYS	3.8
1	A	818	ASN	3.7
1	A	821	ARG	3.6
1	A	331	ALA	3.5
1	A	362	ASN	3.5
1	A	820	ASP	3.2
1	A	267	ASN	3.2
1	A	268	GLU	3.2
1	A	814	THR	3.1
1	A	361	ASP	3.1
1	A	358	SER	3.1
1	A	890	GLN	3.0
1	A	606	THR	3.0
1	A	679	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	599	ARG	3.0
1	A	265	LYS	2.9
1	A	359	ARG	2.8
1	A	363	GLN	2.7
1	A	264	GLU	2.7
1	A	581	GLN	2.5
1	A	330	ASN	2.5
1	A	920	ARG	2.5
1	A	65	ILE	2.4
1	A	619	THR	2.4
1	A	500	GLN	2.3
1	A	266	THR	2.3
1	A	360	GLY	2.3
1	A	817	LYS	2.2
1	A	64	LYS	2.1
1	A	255	PRO	2.1
1	A	269	LYS	2.1
1	A	680	ALA	2.1
1	A	598	PHE	2.0
1	A	210	ARG	2.0
1	A	916	LEU	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	NA	A	964	1/1	0.95	0.20	2.67	33,33,33,33	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	ZN	A	963	1/1	0.99	0.07	-	32,32,32,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.