



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 17, 2024 – 10:07 AM JST

PDB ID : 8K4D
Title : Structure of the SA2/Sccl/CENP_U complex
Authors : Liu, M.J.; He, X.J.
Deposited on : 2023-07-18
Resolution : 3.52 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.37.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.37.1

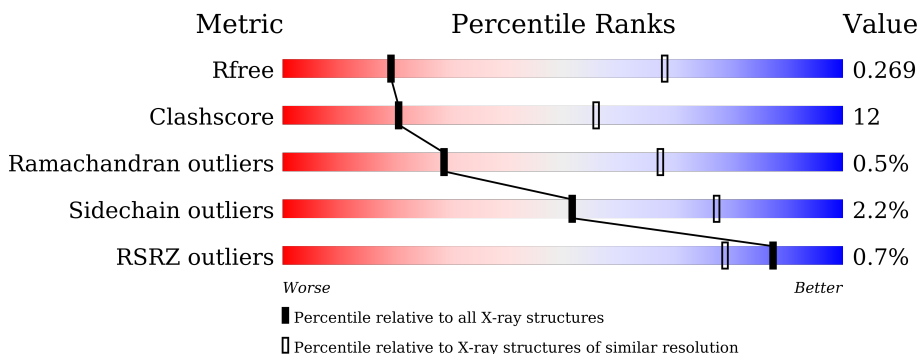
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.52 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}	130704	1161 (3.60-3.44)
Clashscore	141614	1244 (3.60-3.44)
Ramachandran outliers	138981	1206 (3.60-3.44)
Sidechain outliers	138945	1207 (3.60-3.44)
RSRZ outliers	127900	1080 (3.60-3.44)

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	980	 11% 70% 23% 6%
2	B	141	 40% 12% 48%
3	C	9	 11% 33% 67%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8141 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 Cohesin subunit SA-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	920	7478	4771	1245	1408	54	0	0	0

- Molecule 2 is a protein called 64-kDa C-terminal product.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	73	587	379	99	106	3	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	280	MET	-	initiating methionine	UNP O60216

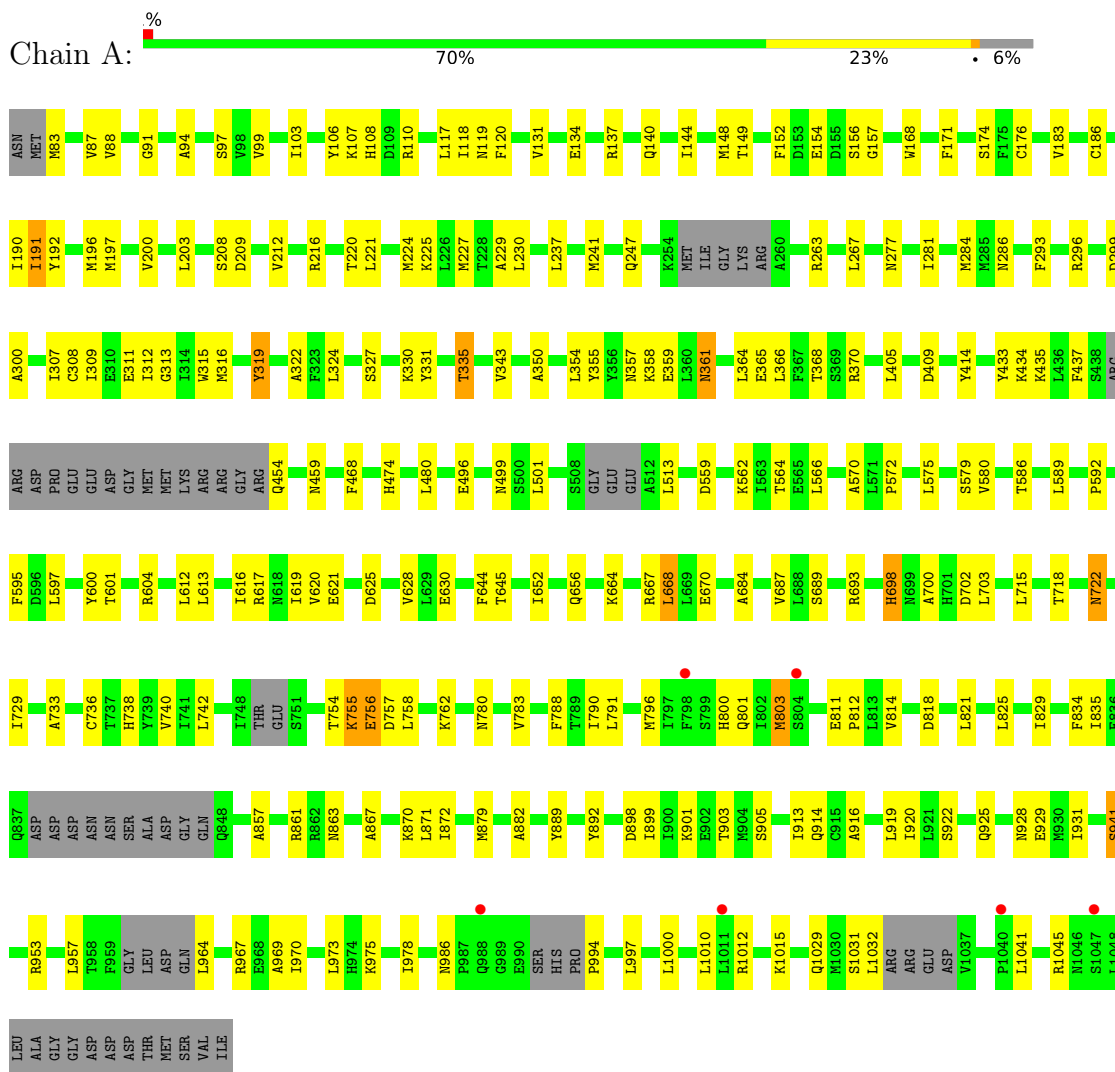
- Molecule 3 is a protein called CENP-U.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	C	9	76	50	10	16	0	0	0

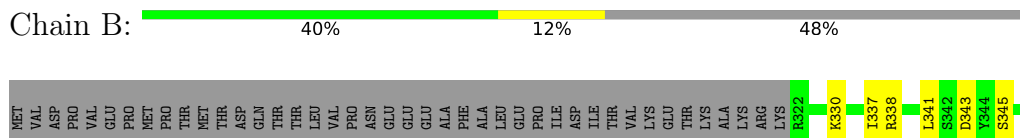
3 Residue-property plots

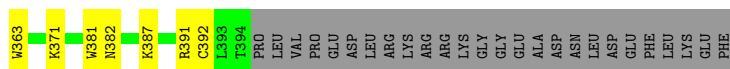
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Cohesin subunit SA-2

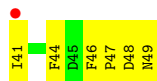


- Molecule 2: 64-kDa C-terminal product





- Molecule 3: CENP-U



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	79.32Å 132.62Å 148.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.43 – 3.52 49.38 – 3.52	Depositor EDS
% Data completeness (in resolution range)	99.2 (49.43-3.52) 99.2 (49.38-3.52)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.07 (at 3.48Å)	Xtrriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.220 , 0.269 0.221 , 0.269	Depositor DCC
R_{free} test set	1036 reflections (5.24%)	wwPDB-VP
Wilson B-factor (Å ²)	124.9	Xtrriage
Anisotropy	0.072	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 98.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8141	wwPDB-VP
Average B, all atoms (Å ²)	150.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% 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	0.59	0/7606	0.65	0/10248
2	B	0.60	0/597	0.62	0/807
3	C	0.85	0/78	0.79	0/106
All	All	0.59	0/8281	0.65	0/11161

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	7478	0	7484	186	0
2	B	587	0	626	15	0
3	C	76	0	62	9	0
All	All	8141	0	8172	201	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 12.

All (201) 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:209:ASP:HA	1:A:296:ARG:NH2	1.80	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:601:THR:HG21	1:A:645:THR:HB	1.58	0.84
1:A:327:SER:O	1:A:330:LYS:HE2	1.77	0.84
1:A:986:ASN:HB2	1:A:994:PRO:N	1.95	0.81
1:A:595:PHE:CD2	1:A:600:TYR:OH	2.34	0.81
1:A:800:HIS:CE1	1:A:801:GLN:HG2	2.17	0.79
1:A:811:GLU:N	1:A:812:PRO:HD2	1.98	0.78
1:A:209:ASP:HA	1:A:296:ARG:HH22	1.51	0.76
1:A:986:ASN:HB2	1:A:994:PRO:CA	2.17	0.74
1:A:496:GLU:HG3	1:A:566:LEU:HD11	1.69	0.74
1:A:800:HIS:CE1	1:A:801:GLN:HE21	2.05	0.72
1:A:756:GLU:O	1:A:758:LEU:N	2.23	0.72
1:A:818:ASP:HB2	1:A:821:LEU:H	1.53	0.72
3:C:46:PHE:CD1	3:C:47:PRO:HD2	2.25	0.72
1:A:191:ILE:HD12	1:A:191:ILE:O	1.89	0.71
1:A:191:ILE:HD13	1:A:281:ILE:HD11	1.73	0.69
1:A:756:GLU:C	1:A:758:LEU:H	1.95	0.69
1:A:414:TYR:OH	1:A:435:LYS:HG2	1.92	0.69
1:A:762:LYS:NZ	1:A:814:VAL:O	2.23	0.69
1:A:221:LEU:HD13	1:A:307:ILE:HD13	1.75	0.68
1:A:986:ASN:H	1:A:994:PRO:HA	1.59	0.68
1:A:216:ARG:HD2	1:A:296:ARG:CZ	2.24	0.68
1:A:227:MET:HG2	1:A:315:TRP:CZ2	2.28	0.67
1:A:702:ASP:O	1:A:703:LEU:HD23	1.95	0.65
1:A:800:HIS:CE1	1:A:801:GLN:NE2	2.65	0.64
1:A:196:MET:O	1:A:200:VAL:HG23	1.98	0.64
1:A:118:ILE:HG23	1:A:203:LEU:CD2	2.27	0.64
1:A:208:SER:O	1:A:296:ARG:NH2	2.31	0.63
1:A:592:PRO:HA	1:A:595:PHE:CD2	2.34	0.63
1:A:595:PHE:CE2	1:A:600:TYR:OH	2.52	0.62
1:A:216:ARG:HD2	1:A:296:ARG:NH2	2.14	0.62
1:A:106:TYR:CE2	1:A:110:ARG:HG2	2.34	0.62
1:A:601:THR:HG21	1:A:645:THR:CB	2.30	0.62
1:A:803:MET:SD	1:A:811:GLU:HG2	2.38	0.62
1:A:1031:SER:HB2	1:A:1045:ARG:HH22	1.66	0.61
1:A:796:MET:HE2	1:A:871:LEU:HD23	1.83	0.60
1:A:220:THR:O	1:A:224:MET:HG2	2.01	0.60
3:C:41:ILE:HG22	3:C:41:ILE:O	2.00	0.60
1:A:216:ARG:CG	1:A:296:ARG:HE	2.16	0.59
1:A:186:CYS:CB	1:A:191:ILE:HG23	2.33	0.59
1:A:118:ILE:HG23	1:A:203:LEU:HD21	1.83	0.58
1:A:313:GLY:HA3	1:A:350:ALA:HB1	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:811:GLU:N	1:A:812:PRO:CD	2.65	0.58
1:A:1031:SER:HB2	1:A:1045:ARG:HH12	1.68	0.58
1:A:559:ASP:HA	1:A:562:LYS:HB2	1.86	0.58
1:A:1012:ARG:HA	1:A:1015:LYS:HE3	1.85	0.58
1:A:197:MET:HG2	1:A:284:MET:SD	2.44	0.58
1:A:796:MET:HE3	1:A:870:LYS:HB3	1.86	0.57
1:A:586:THR:HG22	1:A:628:VAL:HA	1.86	0.57
1:A:144:ILE:O	1:A:148:MET:HG2	2.03	0.57
1:A:667:ARG:NH2	1:A:670:GLU:OE1	2.37	0.57
1:A:755:LYS:HE2	1:A:758:LEU:HD13	1.86	0.57
1:A:361:ASN:HB2	1:A:364:LEU:HB2	1.87	0.57
1:A:916:ALA:HB1	1:A:973:LEU:HD12	1.86	0.57
1:A:811:GLU:H	1:A:812:PRO:HD2	1.66	0.57
1:A:791:LEU:HD23	1:A:825:LEU:HD21	1.87	0.56
1:A:863:ASN:HB2	2:B:363:TRP:HZ3	1.69	0.56
1:A:617:ARG:O	1:A:621:GLU:HG2	2.04	0.56
1:A:928:ASN:HA	1:A:931:ILE:HG12	1.86	0.56
2:B:354:ALA:O	2:B:355:PRO:C	2.44	0.56
1:A:433:TYR:HA	1:A:437:PHE:HD2	1.70	0.56
1:A:216:ARG:HD2	1:A:296:ARG:NE	2.22	0.55
1:A:756:GLU:C	1:A:758:LEU:N	2.58	0.55
1:A:964:LEU:HB3	1:A:967:ARG:HH12	1.72	0.55
1:A:405:LEU:O	1:A:409:ASP:HB2	2.08	0.54
2:B:387:LYS:O	2:B:391:ARG:N	2.41	0.53
1:A:879:MET:HE1	1:A:914:GLN:C	2.28	0.53
1:A:355:TYR:OH	1:A:368:THR:CG2	2.58	0.52
1:A:572:PRO:HG3	1:A:612:LEU:HB2	1.90	0.52
1:A:925:GLN:O	1:A:929:GLU:HB2	2.08	0.52
1:A:186:CYS:HB2	1:A:191:ILE:CG2	2.40	0.52
1:A:209:ASP:HA	1:A:296:ARG:HH21	1.66	0.52
1:A:87:VAL:O	1:A:91:GLY:HA3	2.10	0.52
1:A:134:GLU:OE1	1:A:137:ARG:NH1	2.43	0.52
1:A:355:TYR:CZ	1:A:364:LEU:HD13	2.45	0.52
1:A:149:THR:HG22	1:A:212:VAL:HG21	1.91	0.52
1:A:192:TYR:OH	1:A:237:LEU:HD11	2.10	0.51
2:B:343:ASP:OD1	2:B:345:SER:OG	2.23	0.51
1:A:99:VAL:O	1:A:103:ILE:HG13	2.10	0.51
1:A:119:ASN:HD21	1:A:131:VAL:HB	1.75	0.50
1:A:366:LEU:HD11	3:C:41:ILE:HG13	1.93	0.50
1:A:698:HIS:NE2	1:A:740:VAL:HG22	2.26	0.50
1:A:913:ILE:HG13	1:A:969:ALA:HA	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:474:HIS:CD2	2:B:356:PRO:HA	2.46	0.50
1:A:625:ASP:HB3	1:A:628:VAL:HG23	1.93	0.50
1:A:800:HIS:CE1	1:A:801:GLN:CG	2.93	0.50
1:A:149:THR:CG2	1:A:212:VAL:HG21	2.42	0.50
1:A:152:PHE:CD1	1:A:154:GLU:HB3	2.47	0.50
3:C:41:ILE:HG23	3:C:44:PHE:HD2	1.76	0.50
3:C:47:PRO:HG2	3:C:49:ASN:HD21	1.75	0.50
1:A:94:ALA:O	1:A:97:SER:OG	2.27	0.49
1:A:192:TYR:OH	1:A:277:ASN:HB3	2.12	0.49
2:B:387:LYS:HE3	2:B:391:ARG:HG2	1.95	0.49
1:A:601:THR:HG21	1:A:645:THR:CG2	2.42	0.49
1:A:316:MET:HB3	1:A:324:LEU:HD13	1.94	0.49
1:A:575:LEU:O	1:A:579:SER:OG	2.23	0.49
1:A:800:HIS:ND1	1:A:801:GLN:HG2	2.28	0.49
1:A:118:ILE:HG23	1:A:203:LEU:HD22	1.94	0.49
2:B:338:ARG:HH21	3:C:48:ASP:HA	1.78	0.49
1:A:496:GLU:CG	1:A:566:LEU:HD11	2.40	0.48
1:A:293:PHE:CE1	1:A:308:CYS:HB3	2.49	0.48
1:A:237:LEU:O	1:A:241:MET:HB2	2.14	0.48
1:A:758:LEU:HD11	1:A:812:PRO:HG2	1.95	0.48
1:A:791:LEU:HD21	1:A:821:LEU:HD11	1.96	0.47
2:B:337:ILE:O	2:B:341:LEU:HG	2.13	0.47
1:A:630:GLU:HB2	2:B:381:TRP:CH2	2.50	0.47
1:A:156:SER:OG	1:A:157:GLY:N	2.48	0.47
1:A:309:ILE:HG13	1:A:335:THR:HG21	1.97	0.47
1:A:867:ALA:HA	1:A:870:LYS:HE2	1.97	0.47
1:A:330:LYS:HE3	1:A:331:TYR:CZ	2.49	0.47
1:A:857:ALA:O	1:A:861:ARG:HG3	2.14	0.47
1:A:186:CYS:HB2	1:A:191:ILE:HG23	1.97	0.47
1:A:186:CYS:HB3	1:A:191:ILE:HG23	1.96	0.47
1:A:370:ARG:HB2	3:C:41:ILE:HD13	1.97	0.47
1:A:899:ILE:O	1:A:903:THR:HG22	2.14	0.47
1:A:358:LYS:HG3	1:A:358:LYS:O	2.14	0.47
1:A:994:PRO:CG	1:A:997:LEU:HD11	2.45	0.47
1:A:83:MET:O	1:A:87:VAL:HG23	2.15	0.46
1:A:191:ILE:CD1	1:A:281:ILE:HD11	2.43	0.46
1:A:176:CYS:SG	1:A:225:LYS:HE3	2.54	0.46
1:A:589:LEU:HD11	1:A:619:ILE:HD11	1.98	0.46
1:A:1031:SER:CB	1:A:1045:ARG:HH12	2.29	0.46
2:B:371:LYS:HA	2:B:371:LYS:HD3	1.82	0.46
3:C:47:PRO:HG2	3:C:49:ASN:ND2	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:738:HIS:CD2	1:A:790:ILE:HG12	2.50	0.46
1:A:834:PHE:HD1	1:A:892:TYR:CE1	2.34	0.46
1:A:997:LEU:HD22	1:A:1041:LEU:HD21	1.98	0.46
1:A:88:VAL:HG12	1:A:174:SER:HB3	1.97	0.46
1:A:192:TYR:CZ	1:A:277:ASN:HB3	2.51	0.45
1:A:293:PHE:CD1	1:A:308:CYS:HB3	2.50	0.45
1:A:190:ILE:C	1:A:192:TYR:H	2.19	0.45
1:A:689:SER:O	1:A:693:ARG:HG3	2.16	0.45
1:A:700:ALA:HA	2:B:382:ASN:HB2	1.98	0.45
1:A:645:THR:O	1:A:645:THR:HG22	2.16	0.45
1:A:168:TRP:O	1:A:171:PHE:HB3	2.17	0.45
1:A:616:ILE:O	1:A:619:ILE:HG12	2.17	0.45
1:A:119:ASN:ND2	1:A:131:VAL:HB	2.32	0.45
1:A:617:ARG:HA	1:A:620:VAL:HG22	1.99	0.45
1:A:117:LEU:O	1:A:120:PHE:HB3	2.17	0.44
1:A:801:GLN:HB2	2:B:392:CYS:HB3	1.99	0.44
1:A:872:ILE:HD13	1:A:882:ALA:HB2	1.98	0.44
2:B:362:MET:CE	2:B:363:TRP:NE1	2.80	0.44
1:A:986:ASN:N	1:A:994:PRO:HA	2.30	0.44
1:A:281:ILE:HD13	1:A:281:ILE:HA	1.82	0.44
1:A:729:ILE:O	1:A:733:ALA:N	2.49	0.44
1:A:872:ILE:HG21	1:A:882:ALA:HB2	1.99	0.44
1:A:1000:LEU:HD12	1:A:1041:LEU:HD22	1.99	0.44
1:A:916:ALA:HB1	1:A:973:LEU:HB2	1.99	0.44
1:A:88:VAL:CG1	1:A:174:SER:HB3	2.48	0.44
1:A:613:LEU:O	1:A:617:ARG:HG2	2.17	0.44
1:A:953:ARG:O	1:A:957:LEU:HG	2.18	0.44
1:A:1029:GLN:O	1:A:1032:LEU:HB2	2.17	0.44
1:A:1032:LEU:HD23	1:A:1032:LEU:HA	1.81	0.43
1:A:564:THR:HA	1:A:595:PHE:CD1	2.53	0.43
1:A:791:LEU:HD12	1:A:791:LEU:HA	1.88	0.43
1:A:589:LEU:HD11	1:A:619:ILE:CD1	2.48	0.43
1:A:788:PHE:CZ	1:A:829:ILE:HD11	2.53	0.43
1:A:319:TYR:CE2	1:A:322:ALA:HB2	2.54	0.43
1:A:738:HIS:CE1	1:A:742:LEU:HD11	2.52	0.43
3:C:41:ILE:O	3:C:41:ILE:CG2	2.67	0.43
1:A:652:ILE:O	1:A:656:GLN:HB3	2.18	0.43
1:A:919:LEU:O	1:A:922:SER:OG	2.34	0.43
1:A:186:CYS:O	1:A:191:ILE:HG23	2.19	0.43
1:A:357:ASN:HB3	1:A:359:GLU:HG2	1.99	0.43
1:A:758:LEU:HD21	1:A:812:PRO:HG2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:780:ASN:OD1	1:A:783:VAL:N	2.45	0.42
1:A:499:ASN:HD21	1:A:570:ALA:CB	2.32	0.42
1:A:898:ASP:HA	1:A:901:LYS:HE2	2.02	0.42
1:A:140:GLN:O	1:A:144:ILE:HG13	2.19	0.42
1:A:736:CYS:O	1:A:740:VAL:HG23	2.19	0.42
1:A:183:VAL:HB	1:A:229:ALA:HB1	2.01	0.42
1:A:309:ILE:HA	1:A:312:ILE:HG22	2.01	0.42
1:A:722:ASN:N	1:A:722:ASN:OD1	2.53	0.42
1:A:224:MET:HB3	1:A:311:GLU:HG3	2.02	0.42
1:A:324:LEU:HD11	1:A:354:LEU:HD13	2.00	0.42
1:A:468:PHE:HD1	1:A:480:LEU:HD23	1.84	0.42
1:A:811:GLU:H	1:A:812:PRO:CD	2.31	0.42
1:A:247:GLN:HG3	1:A:267:LEU:HD11	2.01	0.42
1:A:513:LEU:HD23	1:A:513:LEU:HA	1.88	0.42
1:A:1031:SER:HB2	1:A:1045:ARG:NH2	2.33	0.42
1:A:664:LYS:O	1:A:668:LEU:HD12	2.20	0.41
1:A:684:ALA:HA	1:A:687:VAL:HG12	2.02	0.41
1:A:216:ARG:HG2	1:A:296:ARG:HE	1.84	0.41
1:A:454:GLN:CD	1:A:501:LEU:HD23	2.40	0.41
1:A:964:LEU:HB3	1:A:967:ARG:NH1	2.33	0.41
1:A:355:TYR:O	1:A:358:LYS:HB3	2.20	0.41
1:A:889:TYR:CE2	1:A:901:LYS:HB3	2.56	0.41
1:A:103:ILE:HG22	1:A:107:LYS:HE3	2.02	0.41
1:A:300:ALA:N	2:B:330:LYS:O	2.54	0.41
1:A:597:LEU:HB2	1:A:644:PHE:CD1	2.56	0.41
1:A:227:MET:HA	1:A:230:LEU:HD12	2.02	0.40
1:A:335:THR:HG22	1:A:343:VAL:CG1	2.52	0.40
1:A:296:ARG:CD	1:A:299:ASP:OD2	2.70	0.40
1:A:361:ASN:O	1:A:361:ASN:ND2	2.54	0.40
1:A:970:ILE:HG12	1:A:1010:LEU:HD13	2.02	0.40
1:A:975:LYS:HD2	1:A:978:ILE:HD11	2.03	0.40
1:A:986:ASN:HB2	1:A:994:PRO:HA	1.99	0.40
1:A:118:ILE:HG12	1:A:203:LEU:HD22	2.03	0.40
1:A:920:ILE:HD13	1:A:920:ILE:HA	1.93	0.40
1:A:434:LYS:HE2	1:A:434:LYS:HB2	1.90	0.40
2:B:387:LYS:HA	2:B:387:LYS:HD2	1.91	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	902/980 (92%)	847 (94%)	50 (6%)	5 (1%)	25	64
2	B	71/141 (50%)	66 (93%)	5 (7%)	0	100	100
3	C	7/9 (78%)	7 (100%)	0	0	100	100
All	All	980/1130 (87%)	920 (94%)	55 (6%)	5 (0%)	29	67

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	757	ASP
1	A	604	ARG
1	A	754	THR
1	A	941	SER
1	A	191	ILE

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	831/881 (94%)	811 (98%)	20 (2%)	49	76
2	B	68/130 (52%)	68 (100%)	0	100	100
3	C	9/9 (100%)	9 (100%)	0	100	100
All	All	908/1020 (89%)	888 (98%)	20 (2%)	52	77

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	108	HIS
1	A	263	ARG
1	A	286	ASN
1	A	319	TYR
1	A	335	THR
1	A	361	ASN
1	A	365	GLU
1	A	459	ASN
1	A	580	VAL
1	A	668	LEU
1	A	698	HIS
1	A	715	LEU
1	A	718	THR
1	A	722	ASN
1	A	755	LYS
1	A	756	GLU
1	A	803	MET
1	A	835	ILE
1	A	905	SER
1	A	941	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	123	GLN
1	A	499	ASN
1	A	738	HIS
1	A	801	GLN
3	C	49	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 monosaccharides 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	920/980 (93%)	-0.00	6 (0%) 87 79	79, 151, 212, 273	0
2	B	73/141 (51%)	-0.28	0 100 100	78, 124, 181, 187	0
3	C	9/9 (100%)	0.93	1 (11%) 5 5	98, 111, 153, 159	0
All	All	1002/1130 (88%)	-0.02	7 (0%) 87 79	78, 149, 211, 273	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1040	PRO	2.9
1	A	1011	LEU	2.7
1	A	804	SER	2.5
1	A	798	PHE	2.5
1	A	988	GLN	2.4
1	A	1047	SER	2.4
3	C	41	ILE	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 monosaccharides 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.