



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 07:53 AM UTC

PDB ID : 3HUF / pdb_00003huf
Title : Structure of the *S. pombe* Nbs1-Ctp1 complex
Authors : Williams, R.S.; Guenther, G.; Tainer, J.A.
Deposited on : 2009-06-13
Resolution : 2.15 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

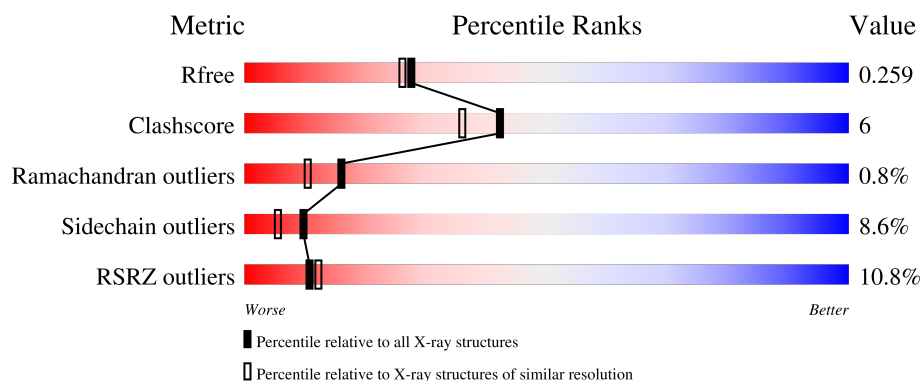
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	325	14% 75% 15% • 7%
1	B	325	7% 82% 12% • •
1	C	325	9% 79% 15% • •
2	E	13	23% 38% 62%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7688 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 DNA repair and telomere maintenance protein nbs1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	301	Total	C	N	O	S	0	0	0
			2308	1467	381	445	15			
1	B	315	Total	C	N	O	S	0	0	0
			2461	1565	407	474	15			
1	C	313	Total	C	N	O	S	0	0	0
			2445	1552	406	472	15			

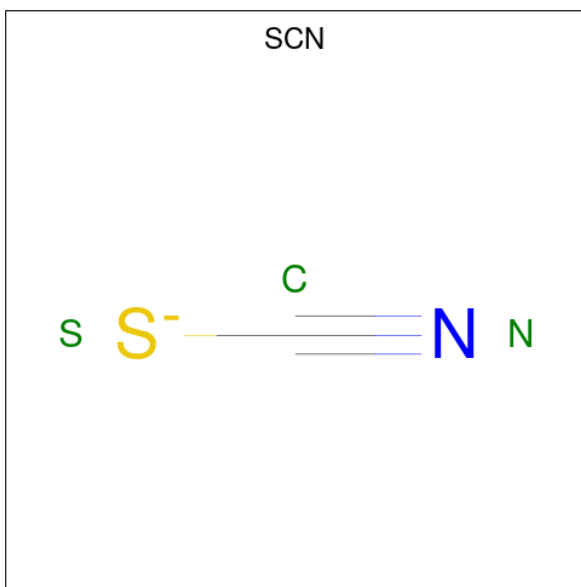
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	322	LEU	-	expression tag	UNP O43070
A	323	VAL	-	expression tag	UNP O43070
A	324	PRO	-	expression tag	UNP O43070
A	325	ARG	-	expression tag	UNP O43070
B	322	LEU	-	expression tag	UNP O43070
B	323	VAL	-	expression tag	UNP O43070
B	324	PRO	-	expression tag	UNP O43070
B	325	ARG	-	expression tag	UNP O43070
C	322	LEU	-	expression tag	UNP O43070
C	323	VAL	-	expression tag	UNP O43070
C	324	PRO	-	expression tag	UNP O43070
C	325	ARG	-	expression tag	UNP O43070

- Molecule 2 is a protein called Double-strand break repair protein ctp1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	5	Total	C	N	O	P	0	0	0
			40	20	5	14	1			

- Molecule 3 is THIOCYANATE ION (CCD ID: SCN) (formula: CNS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	S	0	0
			3	1	1	1		
3	B	1	Total	C	N	S	0	0
			3	1	1	1		

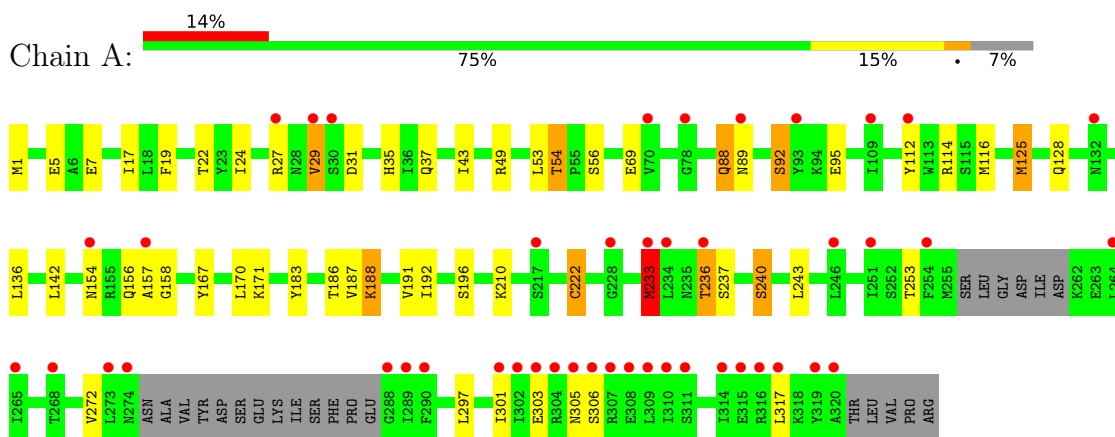
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	143	Total	O	0	0
			143	143		
4	B	155	Total	O	0	0
			155	155		
4	C	129	Total	O	0	0
			129	129		
4	E	1	Total	O	0	0
			1	1		

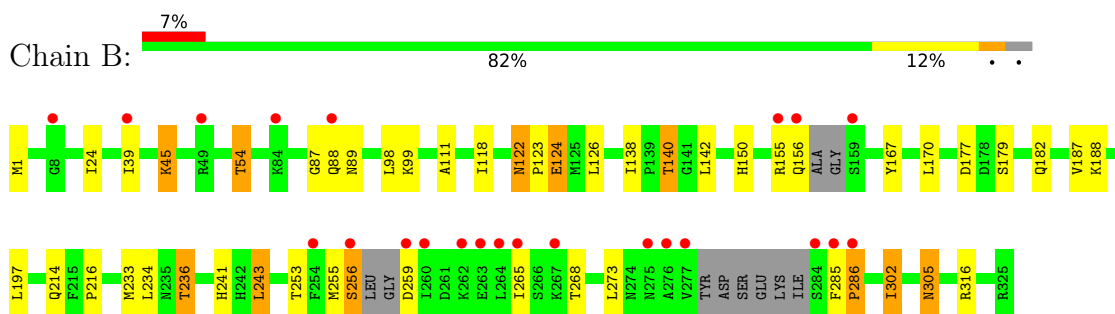
3 Residue-property plots

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.

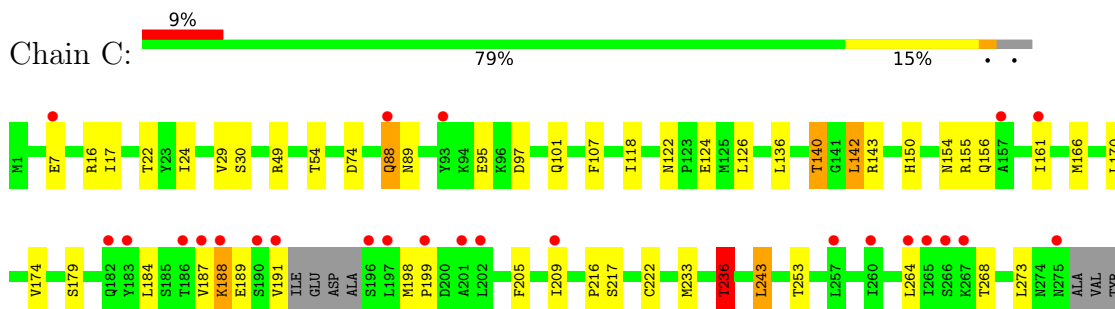
- Molecule 1: DNA repair and telomere maintenance protein nbs1



- Molecule 1: DNA repair and telomere maintenance protein nbs1

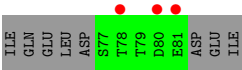


- Molecule 1: DNA repair and telomere maintenance protein nbs1





● Molecule 2: Double-strand break repair protein ctp1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	97.43Å 244.63Å 51.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.15 50.00 – 2.15	Depositor EDS
% Data completeness (in resolution range)	96.9 (50.00-2.15) 96.9 (50.00-2.15)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 2.16Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.215 , 0.262 0.214 , 0.259	Depositor DCC
R_{free} test set	3374 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	38.7	Xtriage
Anisotropy	0.396	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 36.5	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	7688	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 4.32% 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: TPO, SCN, SEP

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	0.71	2/2349 (0.1%)	0.95	4/3177 (0.1%)
1	B	0.68	0/2503	0.96	4/3379 (0.1%)
1	C	0.69	0/2488	0.93	3/3360 (0.1%)
2	E	0.58	0/22	0.60	0/27
All	All	0.69	2/7362 (0.0%)	0.95	11/9943 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	317	LEU	C-O	6.19	1.31	1.24
1	A	317	LEU	C-N	6.15	1.42	1.33

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	187	VAL	N-CA-C	9.69	122.42	112.83
1	B	286	PRO	N-CA-CB	7.09	110.69	103.25
1	B	122	ASN	CA-C-N	5.69	125.37	119.05
1	B	122	ASN	C-N-CA	5.69	125.37	119.05
1	C	236	THR	CB-CA-C	5.62	118.28	109.84
1	A	233	MET	N-CA-C	5.30	117.24	109.24
1	C	122	ASN	CA-C-N	5.06	125.09	119.32
1	C	122	ASN	C-N-CA	5.06	125.09	119.32
1	A	54	THR	N-CA-CB	-5.04	102.75	110.11
1	A	196	SER	N-CA-C	5.02	117.48	111.71
1	B	39	ILE	N-CA-C	5.02	114.80	107.37

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	2308	0	2241	28	0
1	B	2461	0	2435	33	0
1	C	2445	0	2417	30	0
2	E	40	0	26	0	0
3	A	3	0	0	0	0
3	B	3	0	0	0	0
4	A	143	0	0	3	0
4	B	155	0	0	6	0
4	C	129	0	0	5	0
4	E	1	0	0	0	0
All	All	7688	0	7119	91	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 6.

All (91) 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:B:1:MET:HE1	1:B:197:LEU:HD12	1.56	0.87
1:B:243:LEU:C	1:B:243:LEU:HD12	2.04	0.83
1:B:126:LEU:HD11	1:B:140:THR:HG23	1.62	0.82
1:A:233:MET:HG3	1:A:236:THR:HG22	1.64	0.79
1:C:198:MET:HB2	1:C:199:PRO:HD2	1.66	0.76
1:C:17:ILE:H	1:C:191:VAL:HG11	1.51	0.75
1:B:256:SER:HG	1:B:259:ASP:N	1.89	0.69
1:A:29:VAL:HG13	1:A:35:HIS:HB3	1.72	0.69
1:C:295:GLU:O	1:C:299:LYS:HD3	1.94	0.69
1:A:69:GLU:HG2	1:A:92:SER:HB3	1.75	0.67
1:C:7:GLU:HG2	1:C:107:PHE:HD2	1.60	0.67
1:B:45:LYS:HB3	1:B:45:LYS:NZ	2.11	0.66
1:C:7:GLU:HG2	1:C:107:PHE:CD2	2.31	0.66
1:B:140:THR:HG22	4:B:376:HOH:O	1.98	0.63
1:C:170:LEU:HD12	1:C:243:LEU:HD13	1.81	0.62
1:C:88:GLN:O	1:C:89:ASN:HB2	1.99	0.61
1:B:236:THR:HG23	1:B:253:THR:HG23	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:233:MET:HG3	1:A:236:THR:CG2	2.31	0.61
1:C:299:LYS:HD2	4:C:446:HOH:O	2.00	0.61
1:B:305:ASN:H	1:B:305:ASN:HD22	1.50	0.59
1:B:305:ASN:H	1:B:305:ASN:ND2	1.99	0.59
1:C:22:THR:HG21	1:C:49:ARG:NH1	2.17	0.59
1:A:156:GLN:O	1:A:158:GLY:N	2.35	0.58
1:B:243:LEU:HD12	1:B:243:LEU:O	2.03	0.58
1:B:45:LYS:HB3	1:B:45:LYS:HZ2	1.69	0.57
1:A:237:SER:HG	1:A:240:SER:HG	1.52	0.57
1:C:7:GLU:CD	4:C:366:HOH:O	2.47	0.57
1:B:243:LEU:C	1:B:243:LEU:CD1	2.78	0.57
1:B:54:THR:HG23	4:B:412:HOH:O	2.05	0.57
1:C:136:LEU:HD21	1:C:188:LYS:HD2	1.89	0.54
1:A:27:ARG:NH1	1:A:43:ILE:O	2.40	0.54
1:A:183:TYR:O	1:A:186:THR:HB	2.09	0.53
1:A:170:LEU:HD21	1:A:222:CYS:HB3	1.90	0.52
1:B:170:LEU:CD1	1:B:243:LEU:HD13	2.40	0.52
1:B:98:LEU:HB2	1:B:111:ALA:HB3	1.92	0.51
1:C:17:ILE:H	1:C:191:VAL:CG1	2.23	0.51
1:C:126:LEU:HD22	1:C:142:LEU:HD23	1.92	0.50
1:A:29:VAL:O	1:A:37:GLN:HB2	2.12	0.50
1:C:304:ARG:NH2	4:C:454:HOH:O	2.20	0.49
1:A:136:LEU:HD23	1:A:188:LYS:HG2	1.94	0.49
1:B:233:MET:SD	1:B:236:THR:HB	2.52	0.49
1:B:177:ASP:OD2	1:B:179:SER:OG	2.26	0.49
1:C:236:THR:HG23	1:C:253:THR:OG1	2.13	0.48
1:C:101:GLN:NE2	4:C:390:HOH:O	2.46	0.48
1:C:205:PHE:CE1	1:C:209:ILE:HD11	2.48	0.48
1:A:22:THR:HG21	1:A:49:ARG:HD3	1.96	0.47
1:A:1:MET:HE1	1:A:17:ILE:CG2	2.44	0.47
1:A:116:MET:N	4:A:386:HOH:O	2.47	0.47
1:A:233:MET:HB3	1:A:272:VAL:O	2.15	0.47
1:A:125:MET:HE3	4:A:340:HOH:O	2.14	0.47
1:A:125:MET:HE2	1:A:154:ASN:HA	1.97	0.47
1:C:118:ILE:HG12	1:C:150:HIS:HB2	1.97	0.47
1:C:216:PRO:HG3	1:C:222:CYS:SG	2.55	0.47
1:A:1:MET:CE	1:A:17:ILE:CG2	2.92	0.46
1:B:87:GLY:O	4:B:428:HOH:O	2.20	0.46
1:C:233:MET:SD	1:C:236:THR:HB	2.56	0.46
1:A:53:LEU:HG	4:A:388:HOH:O	2.15	0.46
1:A:236:THR:HG23	1:A:253:THR:HG23	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:MET:HE1	1:A:19:PHE:CE2	2.50	0.45
1:A:5:GLU:HG2	1:A:112:TYR:HE2	1.81	0.45
1:A:1:MET:HB3	1:A:1:MET:HE2	1.29	0.45
1:B:1:MET:HE1	1:B:197:LEU:CD1	2.39	0.45
1:A:17:ILE:HD12	1:A:192:ILE:HG13	1.98	0.44
1:A:167:TYR:O	1:A:171:LYS:HG2	2.17	0.44
1:C:236:THR:HG23	1:C:253:THR:HG23	2.00	0.44
1:C:205:PHE:CZ	1:C:209:ILE:HD11	2.53	0.44
1:A:88:GLN:O	1:A:89:ASN:HB2	2.18	0.43
1:B:122:ASN:OD1	1:B:124:GLU:HG2	2.18	0.43
1:B:54:THR:HG22	4:B:348:HOH:O	2.17	0.43
1:C:17:ILE:N	1:C:191:VAL:HG11	2.28	0.43
1:A:1:MET:N	1:A:114:ARG:O	2.52	0.43
1:B:236:THR:HG23	1:B:253:THR:CG2	2.46	0.42
1:B:241:HIS:HE1	4:B:451:HOH:O	2.01	0.42
1:B:122:ASN:HA	1:B:123:PRO:HD3	1.83	0.42
1:C:161:ILE:HG23	1:C:166:MET:HE2	2.01	0.42
1:B:123:PRO:HG3	1:B:142:LEU:HD21	2.01	0.42
1:B:155:ARG:O	1:B:156:GLN:C	2.62	0.42
1:A:88:GLN:H	1:A:88:GLN:HG2	1.67	0.42
1:C:154:ASN:C	1:C:156:GLN:H	2.27	0.42
1:B:140:THR:CG2	4:B:376:HOH:O	2.63	0.42
1:B:167:TYR:CD1	1:B:243:LEU:HB2	2.55	0.42
1:C:304:ARG:NE	4:C:454:HOH:O	2.47	0.42
1:C:236:THR:HG23	1:C:253:THR:CG2	2.50	0.42
1:B:216:PRO:HD3	1:B:302:ILE:HD11	2.03	0.41
1:C:198:MET:HE3	1:C:199:PRO:HG2	2.02	0.41
1:B:118:ILE:HG12	1:B:150:HIS:HB2	2.02	0.41
1:B:118:ILE:HD12	1:B:138:ILE:HG21	2.02	0.40
1:B:305:ASN:HD22	1:B:305:ASN:N	2.19	0.40
1:C:126:LEU:HD11	1:C:140:THR:HG23	2.03	0.40
1:C:174:VAL:HG13	1:C:205:PHE:HB2	2.03	0.40
1:B:265:ILE:HD13	1:B:265:ILE:HA	1.92	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	295/325 (91%)	277 (94%)	13 (4%)	5 (2%)	7	2
1	B	307/325 (94%)	300 (98%)	5 (2%)	2 (1%)	18	13
1	C	307/325 (94%)	297 (97%)	10 (3%)	0	100	100
2	E	2/13 (15%)	1 (50%)	1 (50%)	0	100	100
All	All	911/988 (92%)	875 (96%)	29 (3%)	7 (1%)	16	10

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	303	GLU
1	B	285	PHE
1	B	286	PRO
1	A	31	ASP
1	A	306	SER
1	A	157	ALA
1	A	305	ASN

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	252/295 (85%)	231 (92%)	21 (8%)	10	6
1	B	276/295 (94%)	254 (92%)	22 (8%)	11	7
1	C	275/295 (93%)	249 (90%)	26 (10%)	8	4

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	E	3/11 (27%)	3 (100%)	0	100	100
All	All	806/896 (90%)	737 (91%)	69 (9%)	10	5

All (69) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	GLU
1	A	24	ILE
1	A	29	VAL
1	A	54	THR
1	A	56	SER
1	A	88	GLN
1	A	92	SER
1	A	95	GLU
1	A	125	MET
1	A	128	GLN
1	A	142	LEU
1	A	188	LYS
1	A	191	VAL
1	A	210	LYS
1	A	222	CYS
1	A	233	MET
1	A	236	THR
1	A	240	SER
1	A	243	LEU
1	A	297	LEU
1	A	301	ILE
1	B	24	ILE
1	B	45	LYS
1	B	54	THR
1	B	88	GLN
1	B	89	ASN
1	B	99	LYS
1	B	124	GLU
1	B	140	THR
1	B	182	GLN
1	B	187	VAL
1	B	188	LYS
1	B	214	GLN
1	B	234	LEU
1	B	236	THR
1	B	243	LEU

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Mol	Chain	Res	Type
1	B	255	MET
1	B	256	SER
1	B	268	THR
1	B	273	LEU
1	B	302	ILE
1	B	305	ASN
1	B	316	ARG
1	C	16	ARG
1	C	24	ILE
1	C	29	VAL
1	C	30	SER
1	C	54	THR
1	C	74	ASP
1	C	88	GLN
1	C	95	GLU
1	C	97	ASP
1	C	124	GLU
1	C	140	THR
1	C	142	LEU
1	C	143	ARG
1	C	155	ARG
1	C	179	SER
1	C	184	LEU
1	C	187	VAL
1	C	188	LYS
1	C	189	GLU
1	C	217	SER
1	C	236	THR
1	C	243	LEU
1	C	264	LEU
1	C	268	THR
1	C	273	LEU
1	C	325	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	242	HIS
1	B	28	ASN
1	B	154	ASN
1	B	224	ASN
1	B	305	ASN

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Mol	Chain	Res	Type
1	C	132	ASN
1	C	134	ASN
1	C	156	GLN
1	C	275	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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	$\# Z > 2$
2	SEP	E	77	2	3,4,10	0.66	0	2,4,14	0.82	0
2	TPO	E	79	2	8,10,11	0.79	0	10,14,16	1.03	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SEP	E	77	2	-	0/1/2/10	-
2	TPO	E	79	2	-	0/9/11/13	-

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.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

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	SCN	A	326	-	1,2,2	1.02	0	0,1,1	-	-
3	SCN	B	326	-	1,2,2	0.68	0	0,1,1	-	-

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	301/325 (92%)	1.02	45 (14%) 5 6	32, 43, 66, 76	0
1	B	315/325 (96%)	0.65	23 (7%) 21 24	28, 42, 62, 76	0
1	C	313/325 (96%)	0.79	30 (9%) 13 15	30, 41, 70, 84	0
2	E	3/13 (23%)	2.68	3 (100%) 0 0	58, 58, 60, 60	0
All	All	932/988 (94%)	0.82	101 (10%) 11 12	28, 42, 66, 84	0

All (101) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	254	PHE	5.7
1	A	264	LEU	5.6
1	A	314	ILE	5.2
1	A	305	ASN	5.1
1	C	186	THR	4.7
1	C	191	VAL	4.7
1	B	259	ASP	4.7
1	B	277	VAL	4.6
1	B	275	ASN	4.6
1	C	202	LEU	4.5
1	B	276	ALA	4.5
1	A	307	ARG	4.4
1	A	306	SER	4.4
1	C	197	LEU	4.4
1	C	201	ALA	4.3
1	A	265	ILE	4.1
1	C	284	SER	4.1
1	A	288	GLY	4.1
1	A	289	ILE	4.1
1	C	257	LEU	4.0
1	A	157	ALA	3.9

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Mol	Chain	Res	Type	RSRZ
1	B	284	SER	3.9
1	C	196	SER	3.9
1	C	266	SER	3.8
1	B	286	PRO	3.8
1	A	302	ILE	3.7
1	B	264	LEU	3.6
1	A	310	ILE	3.4
1	B	254	PHE	3.4
1	A	274	ASN	3.3
1	B	8	GLY	3.3
2	E	81	GLU	3.3
1	B	156	GLN	3.3
1	C	286	PRO	3.2
1	A	303	GLU	3.2
1	C	264	LEU	3.1
1	B	262	LYS	3.1
1	A	93	TYR	3.1
1	A	317	LEU	3.1
1	C	187	VAL	3.0
1	C	305	ASN	2.9
1	B	265	ILE	2.9
1	A	309	LEU	2.9
1	B	260	ILE	2.9
1	C	285	PHE	2.9
1	B	88	GLN	2.8
1	A	236	THR	2.8
1	A	251	ILE	2.8
1	C	260	ILE	2.8
1	C	182	GLN	2.7
1	A	320	ALA	2.7
1	C	209	ILE	2.7
1	A	78	GLY	2.7
1	C	199	PRO	2.7
1	C	183	TYR	2.7
1	B	39	ILE	2.7
1	A	273	LEU	2.6
1	B	84	LYS	2.6
1	A	319	TYR	2.6
1	A	29	VAL	2.6
1	A	315	GLU	2.5
1	A	132	ASN	2.5
1	C	93	TYR	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	267	LYS	2.5
1	A	308	GLU	2.5
1	B	263	GLU	2.5
1	C	157	ALA	2.5
1	A	290	PHE	2.5
1	C	325	ARG	2.5
2	E	80	ASP	2.5
1	C	188	LYS	2.4
1	B	285	PHE	2.4
1	B	256	SER	2.4
1	A	301	ILE	2.4
1	C	275	ASN	2.4
1	A	233	MET	2.4
1	A	311	SER	2.4
1	B	159	SER	2.4
1	A	112	TYR	2.3
1	B	155	ARG	2.3
1	A	89	ASN	2.3
1	B	49	ARG	2.3
2	E	78	THR	2.3
1	C	88	GLN	2.2
1	C	161	ILE	2.2
1	A	304	ARG	2.2
1	A	268	THR	2.2
1	A	246	LEU	2.1
1	C	7	GLU	2.1
1	A	228	GLY	2.1
1	A	70	VAL	2.1
1	B	267	LYS	2.1
1	A	30	SER	2.1
1	A	234	LEU	2.1
1	A	316	ARG	2.1
1	A	109	ILE	2.1
1	A	217	SER	2.1
1	C	190	SER	2.1
1	C	265	ILE	2.0
1	A	154	ASN	2.0
1	A	27	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
2	SEP	E	77	5/11	0.43	0.25	59,59,59,59	0
2	TPO	E	79	11/12	0.75	0.18	58,59,60,61	0

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(Å ²)	Q<0.9
3	SCN	B	326	3/3	0.96	0.16	28,28,30,30	0
3	SCN	A	326	3/3	0.97	0.15	28,28,28,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.