



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:18 PM UTC

PDB ID : 3HQI / pdb_00003hqi
Title : Structures of SPOP-Substrate Complexes: Insights into Molecular Architectures of BTB-Cul3 Ubiquitin Ligases: SPOPMATHx/BTB/3-box-PucSBC 1
Authors : Zhuang, M.; Schulman, B.A.
Deposited on : 2009-06-07
Resolution : 2.62 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
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

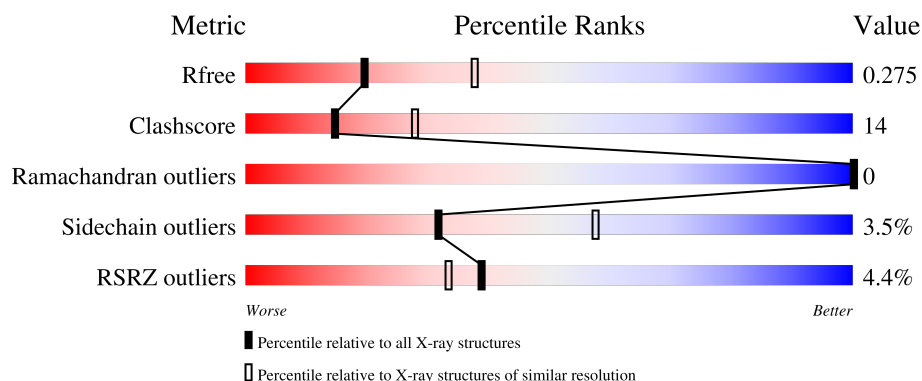
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.62 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	4951 (2.64-2.60)
Clashscore	190562	5303 (2.64-2.60)
Ramachandran outliers	187476	5217 (2.64-2.60)
Sidechain outliers	187428	5217 (2.64-2.60)
RSRZ outliers	180081	4950 (2.64-2.60)

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	312	4% 62% 29% 6%
1	B	312	4% 70% 20% 9%
2	C	7	86% 14%
2	D	7	57% 14% 14% 14%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4668 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 Speckle-type POZ protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	294	Total	C	N	O	S	0	0	0
			2286	1454	387	424	21			
1	B	284	Total	C	N	O	S	0	0	0
			2228	1424	371	413	20			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	26	GLY	-	expression tag	UNP O43791
A	27	SER	-	expression tag	UNP O43791
A	140	GLY	ASP	engineered mutation	UNP O43791
A	330	THR	-	linker	UNP O43791
A	331	ASP	-	linker	UNP O43791
A	332	VAL	-	linker	UNP O43791
A	333	LEU	-	linker	UNP O43791
A	334	GLU	-	linker	UNP O43791
A	335	THR	-	linker	UNP O43791
A	336	SER	-	linker	UNP O43791
A	337	GLY	-	linker	UNP O43791
B	26	GLY	-	expression tag	UNP O43791
B	27	SER	-	expression tag	UNP O43791
B	140	GLY	ASP	engineered mutation	UNP O43791
B	330	THR	-	linker	UNP O43791
B	331	ASP	-	linker	UNP O43791
B	332	VAL	-	linker	UNP O43791
B	333	LEU	-	linker	UNP O43791
B	334	GLU	-	linker	UNP O43791
B	335	THR	-	linker	UNP O43791
B	336	SER	-	linker	UNP O43791
B	337	GLY	-	linker	UNP O43791

- Molecule 2 is a protein called PucSBC1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	6	Total 40	C 23	N 6	O 11	0	0	0
2	D	6	Total 40	C 23	N 6	O 11	0	0	0

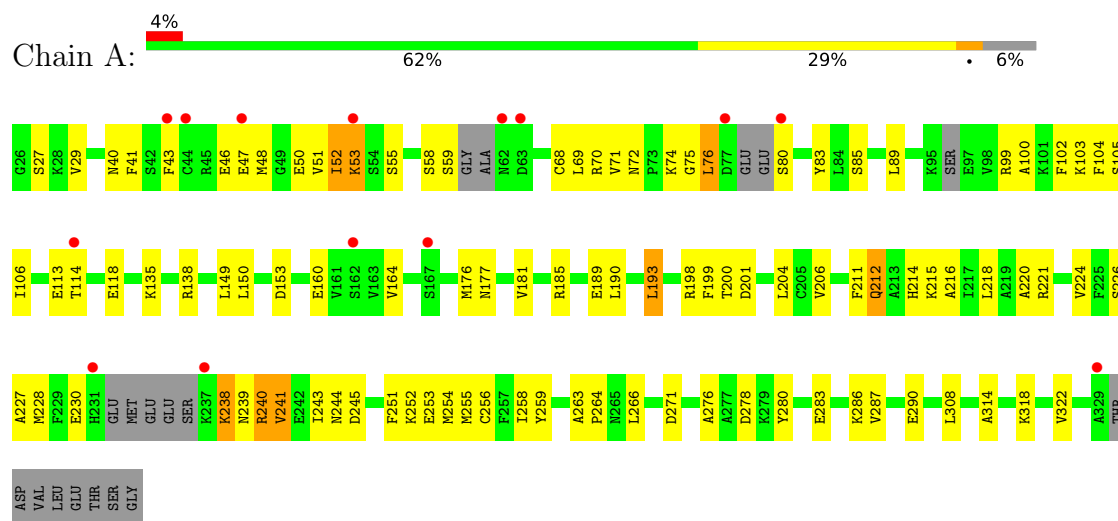
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	54	Total 54	O 54	0	0
3	B	18	Total 18	O 18	0	0
3	C	2	Total 2	O 2	0	0

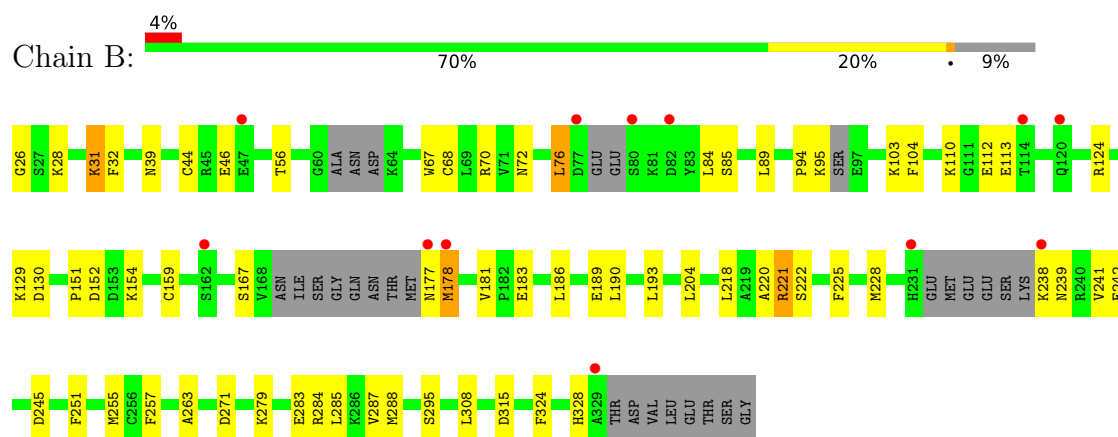
3 Residue-property plots [i](#)

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.

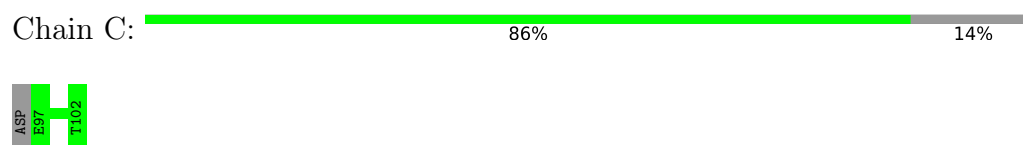
• Molecule 1: Speckle-type POZ protein



• Molecule 1: Speckle-type POZ protein

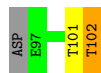


• Molecule 2: PucSBC1



● Molecule 2: PucSBC1

Chain D: 57% 14% 14% 14%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	57.23Å 106.87Å 130.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.62 50.00 – 2.62	Depositor EDS
% Data completeness (in resolution range)	98.2 (50.00-2.62) 98.1 (50.00-2.62)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.77 (at 2.61Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.226 , 0.275 0.226 , 0.275	Depositor DCC
R_{free} test set	1236 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	35.0	Xtriage
Anisotropy	1.127	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 35.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4668	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.53% 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	0.44	0/2325	0.91	9/3130 (0.3%)
1	B	0.44	0/2266	0.89	4/3047 (0.1%)
2	C	0.63	0/39	0.97	0/52
2	D	0.61	0/39	1.03	0/52
All	All	0.44	0/4669	0.90	13/6281 (0.2%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	263	ALA	CA-C-N	6.76	127.03	119.32
1	A	263	ALA	C-N-CA	6.76	127.03	119.32
1	A	181	VAL	N-CA-C	5.74	113.56	107.76
1	B	263	ALA	CA-C-N	5.62	125.46	119.28
1	B	263	ALA	C-N-CA	5.62	125.46	119.28
1	B	28	LYS	N-CA-C	-5.62	99.75	108.90
1	A	100	ALA	N-CA-C	5.56	118.25	108.75
1	A	283	GLU	N-CA-C	5.49	117.69	111.11
1	A	238	LYS	N-CA-C	-5.39	106.28	112.92
1	A	244	ASN	N-CA-C	5.39	119.17	112.59
1	B	315	ASP	N-CA-C	5.38	117.57	111.11
1	A	55	SER	N-CA-C	-5.06	103.72	110.55
1	A	51	VAL	N-CA-C	5.04	115.37	108.12

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	2286	0	2227	83	0
1	B	2228	0	2189	58	0
2	C	40	0	36	0	0
2	D	40	0	36	3	0
3	A	54	0	0	2	0
3	B	18	0	0	0	0
3	C	2	0	0	0	0
All	All	4668	0	4488	126	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 14.

All (126) 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:257:PHE:HE1	1:B:288:MET:HE1	1.30	0.97
1:B:285:LEU:HA	1:B:288:MET:HE3	1.50	0.94
1:B:257:PHE:CE1	1:B:288:MET:HE1	2.02	0.94
1:A:218:LEU:HD21	1:A:255:MET:HE1	1.53	0.90
1:B:72:ASN:HB2	1:B:85:SER:HB2	1.59	0.84
1:A:176:MET:HG2	1:B:39:ASN:ND2	1.91	0.83
1:B:284:ARG:HG2	1:B:288:MET:HE2	1.62	0.82
1:B:228:MET:HE1	1:B:241:VAL:HG11	1.63	0.80
1:A:72:ASN:HB2	1:A:85:SER:HB2	1.64	0.78
1:A:53:LYS:HD2	1:A:53:LYS:C	2.10	0.77
1:A:254:MET:HE1	1:A:280:TYR:CD1	2.21	0.75
1:A:176:MET:HG2	1:B:39:ASN:HD21	1.49	0.74
1:B:26:GLY:N	1:B:167:SER:HG	1.86	0.74
1:B:178:MET:SD	1:B:178:MET:N	2.63	0.71
1:A:240:ARG:HG2	1:A:240:ARG:HH11	1.56	0.71
1:A:214:HIS:HD2	1:A:259:TYR:OH	1.77	0.67
1:B:76:LEU:HD11	2:D:101:THR:HG22	1.77	0.66
1:A:204:LEU:HD13	1:A:251:PHE:HZ	1.60	0.66
1:A:220:ALA:HB2	1:B:193:LEU:HD13	1.79	0.65
1:B:245:ASP:OD2	1:B:279:LYS:HD3	1.96	0.65
1:A:254:MET:HE3	1:A:276:ALA:HB1	1.80	0.64
1:A:271:ASP:HB2	1:A:308:LEU:HD21	1.80	0.64
1:B:183:GLU:H	1:B:183:GLU:CD	2.06	0.64
1:A:254:MET:HE1	1:A:280:TYR:HD1	1.58	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:110:LYS:HE3	1:B:112:GLU:OE1	1.98	0.63
1:B:44:CYS:HB3	1:B:46:GLU:OE1	1.99	0.62
1:A:243:ILE:HD13	1:A:251:PHE:CE1	2.34	0.62
1:A:99:ARG:HH11	1:A:99:ARG:HG2	1.64	0.61
1:A:76:LEU:N	1:A:76:LEU:HD12	2.15	0.61
1:A:198:ARG:O	1:A:214:HIS:HE1	1.83	0.61
1:A:48:MET:HA	1:A:74:LYS:HB2	1.84	0.60
1:A:216:ALA:HB1	1:B:193:LEU:HD21	1.83	0.60
1:B:104:PHE:HE1	1:B:159:CYS:HG	1.50	0.60
1:B:178:MET:H	1:B:178:MET:CE	2.15	0.59
1:A:53:LYS:HD2	1:A:53:LYS:O	2.02	0.59
1:A:215:LYS:HE2	1:A:239:ASN:OD1	2.02	0.59
1:A:204:LEU:HD23	1:A:241:VAL:HG13	1.83	0.59
1:A:71:VAL:HA	1:A:85:SER:O	2.03	0.59
1:A:221:ARG:HD2	1:B:186:LEU:HD12	1.85	0.58
1:B:39:ASN:OD1	1:B:154:LYS:HE2	2.03	0.58
1:B:94:PRO:O	1:B:95:LYS:C	2.46	0.57
1:A:190:LEU:HD13	1:B:190:LEU:HD11	1.87	0.57
1:A:83:TYR:CE1	1:A:138:ARG:HG2	2.39	0.57
1:A:149:LEU:O	1:A:150:LEU:HD12	2.03	0.57
1:A:103:LYS:HD3	1:A:160:GLU:OE2	2.04	0.57
1:A:46:GLU:HA	1:A:50:GLU:OE1	2.03	0.57
1:A:198:ARG:HH21	1:A:199:PHE:HE1	1.51	0.57
1:B:183:GLU:CD	1:B:183:GLU:N	2.63	0.57
1:A:228:MET:C	1:A:230:GLU:H	2.14	0.56
1:A:105:SER:OG	1:A:113:GLU:HG2	2.06	0.56
1:B:70:ARG:HH11	1:B:70:ARG:HB2	1.70	0.56
1:B:283:GLU:O	1:B:287:VAL:HG23	2.06	0.55
1:A:138:ARG:HB3	1:A:138:ARG:NH1	2.21	0.55
1:A:75:GLY:C	1:A:76:LEU:HD12	2.32	0.54
1:B:324:PHE:O	1:B:328:HIS:HD2	1.91	0.54
1:A:106:ILE:O	1:A:114:THR:HG22	2.09	0.53
1:A:318:LYS:O	1:A:322:VAL:HG23	2.08	0.53
1:B:284:ARG:HG2	1:B:288:MET:CE	2.37	0.53
1:A:198:ARG:NH2	1:A:199:PHE:HE1	2.07	0.52
1:A:200:THR:HG21	1:A:212:GLN:HB3	1.92	0.51
1:A:256:CYS:HB2	3:A:368:HOH:O	2.09	0.51
1:B:31:LYS:O	1:B:31:LYS:HG3	2.08	0.51
1:B:225:PHE:HD2	1:B:228:MET:HE2	1.75	0.51
1:A:266:LEU:O	1:A:266:LEU:HD23	2.11	0.50
1:A:40:ASN:HA	1:A:153:ASP:OD2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:LEU:HD22	1:A:104:PHE:HE1	1.77	0.50
1:A:206:VAL:HG22	1:A:243:ILE:HB	1.93	0.50
1:A:221:ARG:HG2	1:B:189:GLU:OE1	2.12	0.49
1:A:29:VAL:HG22	1:A:164:VAL:HG12	1.95	0.49
1:A:185:ARG:O	1:A:189:GLU:HG3	2.12	0.49
1:A:52:ILE:HG13	1:A:71:VAL:CG1	2.42	0.49
1:A:201:ASP:OD1	1:A:215:LYS:HG3	2.12	0.49
1:A:70:ARG:HH21	1:A:72:ASN:HD21	1.60	0.48
1:A:177:ASN:HD21	1:B:295:SER:HB3	1.79	0.48
1:B:76:LEU:CD1	2:D:101:THR:HG22	2.43	0.48
1:B:238:LYS:O	1:B:239:ASN:HB2	2.13	0.48
1:B:103:LYS:NZ	1:B:113:GLU:OE1	2.47	0.47
1:A:83:TYR:CE2	1:A:135:LYS:HE3	2.49	0.47
1:A:240:ARG:HH11	1:A:240:ARG:CG	2.27	0.47
1:B:204:LEU:HD13	1:B:251:PHE:HZ	1.79	0.47
1:A:190:LEU:HD11	1:B:190:LEU:CD1	2.44	0.47
1:A:211:PHE:HE2	1:A:252:LYS:HE2	1.80	0.47
1:A:27:SER:HB3	1:B:31:LYS:HG2	1.97	0.46
1:A:68:CYS:SG	1:A:89:LEU:HD23	2.56	0.46
1:A:266:LEU:HD23	1:A:266:LEU:C	2.40	0.46
1:A:238:LYS:HA	3:A:11:HOH:O	2.16	0.45
1:B:218:LEU:HD21	1:B:255:MET:HE1	1.98	0.45
1:A:193:LEU:HD13	1:B:220:ALA:HB2	1.97	0.45
1:A:190:LEU:CD1	1:B:190:LEU:CD1	2.95	0.45
1:B:70:ARG:HD2	1:B:89:LEU:HD22	1.98	0.45
1:A:245:ASP:N	1:A:245:ASP:OD1	2.49	0.44
1:B:130:ASP:OD1	1:B:130:ASP:C	2.60	0.44
1:A:103:LYS:HG3	1:A:118:GLU:HB3	1.98	0.44
1:A:224:VAL:HG12	1:A:228:MET:CE	2.48	0.44
1:A:287:VAL:HG12	1:B:181:VAL:HG22	1.99	0.44
1:B:284:ARG:CG	1:B:288:MET:HE2	2.39	0.44
1:A:224:VAL:HG12	1:A:228:MET:HE3	1.99	0.44
1:B:70:ARG:HB2	1:B:70:ARG:NH1	2.33	0.43
1:A:190:LEU:CD1	1:B:190:LEU:HD11	2.48	0.43
1:B:177:ASN:CG	1:B:178:MET:SD	3.01	0.43
1:A:227:ALA:O	1:A:230:GLU:HB3	2.19	0.43
1:B:178:MET:H	1:B:178:MET:HE2	1.82	0.43
1:A:198:ARG:O	1:A:214:HIS:CE1	2.69	0.43
1:A:200:THR:HG21	1:A:212:GLN:CB	2.49	0.42
1:A:278:ASP:OD1	1:A:286:LYS:HE3	2.19	0.42
1:B:31:LYS:HE2	1:B:32:PHE:N	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:324:PHE:O	1:B:328:HIS:CD2	2.72	0.42
1:A:290:GLU:HG2	1:A:314:ALA:HB1	2.02	0.42
1:B:222:SER:HB3	1:B:225:PHE:HB2	2.01	0.42
1:B:271:ASP:HB2	1:B:308:LEU:HD21	2.01	0.42
1:A:218:LEU:HD23	1:A:258:ILE:HG21	2.02	0.42
1:B:151:PRO:O	1:B:152:ASP:HB2	2.19	0.42
1:A:189:GLU:OE1	1:B:221:ARG:NH1	2.49	0.41
1:A:221:ARG:HG2	1:A:221:ARG:HH11	1.86	0.41
1:A:240:ARG:HG2	1:A:240:ARG:NH1	2.30	0.41
1:B:70:ARG:NH1	1:B:70:ARG:CB	2.83	0.41
1:A:204:LEU:HD23	1:A:241:VAL:CG1	2.48	0.41
1:B:56:THR:HA	1:B:67:TRP:O	2.20	0.41
1:A:58:SER:OG	1:A:59:SER:N	2.52	0.41
1:A:41:PHE:C	1:A:43:PHE:H	2.29	0.41
1:A:198:ARG:HB3	1:A:199:PHE:HD1	1.86	0.41
1:A:253:GLU:OE2	1:A:264:PRO:HD2	2.21	0.41
1:B:129:LYS:HD2	2:D:102:THR:HG23	2.02	0.40
1:A:102:PHE:HA	1:A:160:GLU:O	2.22	0.40
1:A:226:SER:O	1:A:230:GLU:HB2	2.22	0.40
1:A:47:GLU:N	1:A:50:GLU:OE1	2.48	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	284/312 (91%)	272 (96%)	12 (4%)	0	100	100
1	B	272/312 (87%)	262 (96%)	10 (4%)	0	100	100
2	C	4/7 (57%)	4 (100%)	0	0	100	100
2	D	4/7 (57%)	4 (100%)	0	0	100	100
All	All	564/638 (88%)	542 (96%)	22 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	242/271 (89%)	234 (97%)	8 (3%)	33	59
1	B	238/271 (88%)	230 (97%)	8 (3%)	32	58
2	C	5/7 (71%)	5 (100%)	0	100	100
2	D	5/7 (71%)	4 (80%)	1 (20%)	1	2
All	All	490/556 (88%)	473 (96%)	17 (4%)	32	57

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	ILE
1	A	53	LYS
1	A	76	LEU
1	A	80	SER
1	A	193	LEU
1	A	212	GLN
1	A	240	ARG
1	A	241	VAL
1	B	31	LYS
1	B	68	CYS
1	B	76	LEU
1	B	84	LEU
1	B	124	ARG
1	B	178	MET
1	B	221	ARG
1	B	242	GLU
2	D	102	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	ASN
1	A	72	ASN
1	A	147	ASN
1	A	177	ASN
1	A	209	GLN
1	A	214	HIS
1	A	265	ASN
1	B	147	ASN
1	B	165	GLN
1	B	231	HIS
1	B	320	GLN
1	B	328	HIS

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 oligosaccharides 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 ⓘ

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	294/312 (94%)	0.18	14 (4%) 35 30	21, 38, 60, 83	6 (2%)
1	B	284/312 (91%)	0.13	12 (4%) 40 35	18, 37, 56, 76	5 (1%)
2	C	6/7 (85%)	0.25	0 100 100	31, 35, 40, 43	0
2	D	6/7 (85%)	0.33	0 100 100	34, 37, 44, 53	0
All	All	590/638 (92%)	0.16	26 (4%) 39 33	18, 37, 57, 83	11 (1%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	80	SER	6.2
1	A	77	ASP	4.1
1	A	237	LYS	4.1
1	A	43	PHE	3.9
1	A	231	HIS	3.2
1	B	114	THR	3.2
1	B	177	ASN	3.2
1	B	238	LYS	3.1
1	A	47	GLU	3.0
1	B	231	HIS	2.9
1	A	114	THR	2.8
1	A	329	ALA	2.8
1	A	162	SER	2.7
1	B	329	ALA	2.6
1	A	63	ASP	2.6
1	A	53	LYS	2.6
1	B	178	MET	2.5
1	B	120	GLN	2.5
1	B	162	SER	2.5
1	A	62	ASN	2.3
1	B	82	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	80	SER	2.1
1	A	167	SER	2.1
1	B	47	GLU	2.1
1	A	44	CYS	2.1
1	B	77	ASP	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 oligosaccharides 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.