



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2026 – 08:54 PM UTC

PDB ID : 6E29 / pdb_00006e29
Title : Crystal structure of Myceliophtheria_thermophila Cps50 (Swd1) beta-propeller domain
Authors : Joshi, M.; Yang, Y.; Brunzelle, J.S.; Couture, J.F.
Deposited on : 2018-07-10
Resolution : 1.82 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

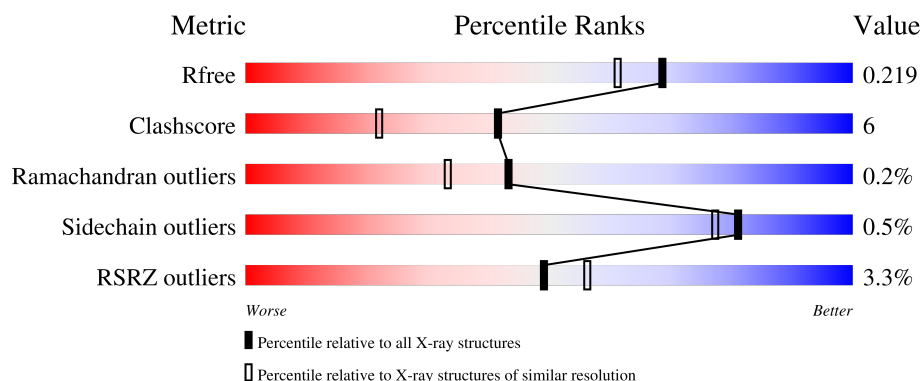
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.82 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	1112 (1.82-1.82)
Clashscore	190562	1148 (1.82-1.82)
Ramachandran outliers	187476	1140 (1.82-1.82)
Sidechain outliers	187428	1140 (1.82-1.82)
RSRZ outliers	180081	1112 (1.82-1.82)

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	345	4% 85% 9% • 5%
1	B	345	3% 75% 16% 9%
1	C	345	3% 77% 14% • 9%
1	D	345	2% 78% 12% • 10%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10930 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 SWD1-like protein.

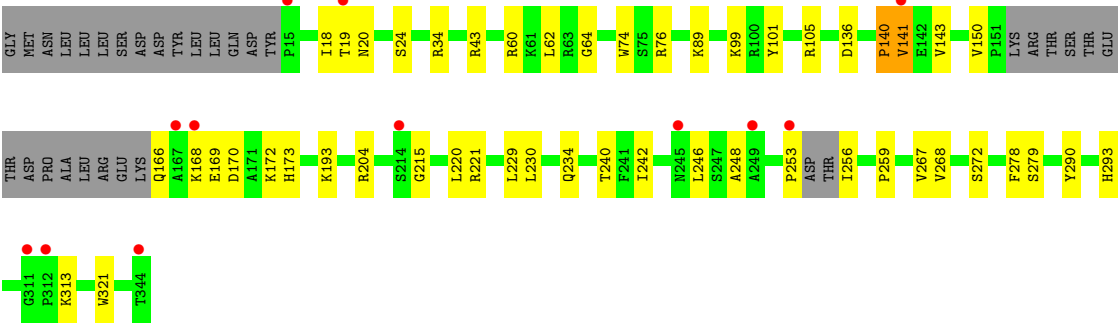
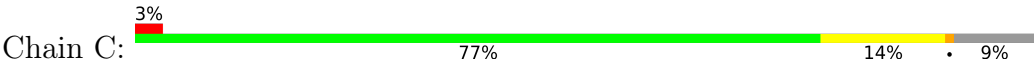
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	312	Total	C	N	O	S	3	7	0
			2514	1594	457	454	9			
1	A	327	Total	C	N	O	S	4	8	0
			2646	1672	481	484	9			
1	B	315	Total	C	N	O	S	0	8	0
			2547	1616	461	461	9			
1	C	314	Total	C	N	O	S	0	10	0
			2547	1613	463	462	9			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	0	GLY	-	expression tag	UNP G2Q1Q9
A	0	GLY	-	expression tag	UNP G2Q1Q9
B	0	GLY	-	expression tag	UNP G2Q1Q9
C	0	GLY	-	expression tag	UNP G2Q1Q9

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	213	Total	O	0	0
			213	213		
2	A	126	Total	O	0	0
			126	126		
2	B	164	Total	O	0	0
			164	164		
2	C	173	Total	O	0	0
			173	173		



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.78Å 88.90Å 124.32Å 90.00° 94.30° 90.00°	Depositor
Resolution (Å)	45.03 – 1.82 45.03 – 1.82	Depositor EDS
% Data completeness (in resolution range)	91.0 (45.03-1.82) 90.9 (45.03-1.82)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.18 (at 1.82Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.168 , 0.216 0.172 , 0.219	Depositor DCC
R_{free} test set	5199 reflections (4.46%)	wwPDB-VP
Wilson B-factor (Å ²)	27.5	Xtriage
Anisotropy	0.166	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	10930	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.03% 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

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	0.61	0/2727	0.83	3/3701 (0.1%)
1	B	0.63	0/2630	0.80	5/3571 (0.1%)
1	C	0.72	1/2633 (0.0%)	0.89	4/3575 (0.1%)
1	D	0.77	2/2592 (0.1%)	0.91	5/3518 (0.1%)
All	All	0.68	3/10582 (0.0%)	0.86	17/14365 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	252	ASP	N-CA	-14.17	1.31	1.46
1	C	64	GLY	C-O	-6.12	1.15	1.23
1	D	251	LEU	C-N	-6.10	1.27	1.33

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	141	VAL	N-CA-C	-17.94	82.99	108.11
1	C	141	VAL	N-CA-C	-17.47	81.75	107.78
1	A	140	PRO	N-CA-C	15.32	144.04	112.47
1	A	141	VAL	N-CA-C	-15.16	77.81	109.34
1	C	140	PRO	N-CA-C	12.78	138.79	112.47
1	D	140	PRO	N-CA-C	10.66	134.43	112.47
1	B	310	GLU	N-CA-C	9.82	122.90	111.11
1	D	251	LEU	O-C-N	9.47	134.69	122.95
1	B	310	GLU	CB-CA-C	-8.36	97.44	110.81
1	D	141	VAL	N-CA-CB	-7.59	100.94	111.41
1	D	252	ASP	N-CA-C	-6.71	94.23	108.73
1	B	305	LEU	CA-C-N	-6.02	114.68	121.85
1	B	305	LEU	C-N-CA	-6.02	114.68	121.85
1	A	55	THR	N-CA-CB	-5.96	101.77	110.53
1	C	18	ILE	CA-C-N	5.79	128.04	120.28
1	C	18	ILE	C-N-CA	5.79	128.04	120.28
1	B	63	ARG	N-CA-C	5.46	117.55	107.73

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	2646	0	2639	33	0
1	B	2547	0	2552	38	0
1	C	2547	0	2542	33	1
1	D	2514	0	2518	42	1
2	A	126	0	0	1	0
2	B	164	0	0	6	0
2	C	173	0	0	11	0
2	D	213	0	0	6	0
All	All	10930	0	10251	131	1

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 (131) 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:D:227:ARG:NH1	1:A:118:TRP:CZ3	2.11	1.17
1:D:227:ARG:NH1	1:A:118:TRP:CE3	2.14	1.16
1:D:195:ARG:NH1	1:D:211:LYS:HB2	1.67	1.09
1:D:227:ARG:NH1	1:A:118:TRP:CH2	2.32	0.95
1:D:227:ARG:NH1	1:A:118:TRP:CD2	2.36	0.93
1:D:252:ASP:OD2	2:D:401:HOH:O	1.96	0.82
1:D:195:ARG:HH12	1:D:211:LYS:HB2	1.45	0.81
1:B:270:ARG:NH1	2:B:401:HOH:O	2.11	0.80
1:D:227:ARG:NH1	1:A:118:TRP:CE2	2.51	0.78
1:D:136:ASP:HB3	1:D:143:VAL:HG13	1.68	0.76
1:D:227:ARG:NH1	1:A:118:TRP:CZ2	2.54	0.76
1:D:225:SER:HB3	1:A:224:GLU:HG3	1.68	0.75
1:B:136:ASP:OD1	1:B:138:THR:HG23	1.88	0.73
1:C:166:GLN:N	2:C:401:HOH:O	2.22	0.71
1:B:145:TYR:CD2	1:B:204:ARG:NH1	2.61	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:24:SER:O	1:C:43:ARG:HD2	1.93	0.68
1:D:224:GLU:O	2:D:403:HOH:O	2.12	0.68
1:D:227:ARG:HD3	1:A:118:TRP:CZ2	2.29	0.68
1:D:18:ILE:HG23	1:D:342[B]:SER:HB3	1.74	0.68
1:B:103:GLU:OE2	1:B:105:ARG:NH1	2.27	0.67
1:D:250:ASP:OD2	2:D:402:HOH:O	2.10	0.67
1:C:60:ARG:NH1	2:C:405:HOH:O	2.28	0.66
1:A:154:THR:HB	1:B:290[B]:TYR:OH	1.97	0.65
1:D:166:GLN:N	2:D:406:HOH:O	2.30	0.63
1:D:18:ILE:HG23	1:D:342[A]:SER:HB2	1.80	0.63
1:D:49:VAL:HG13	1:D:58:VAL:HG13	1.81	0.62
1:A:150[A]:VAL:HG22	1:C:268:VAL:HG13	1.81	0.62
1:B:34:ARG:NH2	1:B:74:TRP:O	2.32	0.62
1:C:60:ARG:HD3	1:C:62:LEU:HD21	1.82	0.61
1:D:49:VAL:HG22	1:D:61:LYS:HG2	1.82	0.61
1:B:145:TYR:CE2	1:B:204:ARG:NH1	2.69	0.61
1:D:133:MET:HE2	1:D:144:ARG:HB3	1.85	0.59
1:C:19:THR:HG22	1:C:20:ASN:OD1	2.03	0.59
1:D:136:ASP:HB3	1:D:143:VAL:CG1	2.32	0.59
1:D:227:ARG:HD2	2:D:450:HOH:O	2.02	0.59
1:B:145:TYR:CD2	1:B:204:ARG:CZ	2.86	0.59
1:B:262:HIS:ND1	2:B:404:HOH:O	2.32	0.59
1:A:24:SER:O	1:A:43:ARG:HD2	2.03	0.58
1:B:34:ARG:NH2	2:B:406:HOH:O	2.36	0.58
1:B:100:ARG:HD2	1:B:103:GLU:OE1	2.04	0.57
1:D:24:SER:O	1:D:43:ARG:HD2	2.05	0.57
1:A:16:GLU:OE2	1:A:311:GLY:N	2.35	0.57
1:B:221:ARG:HH22	1:B:275[B]:HIS:CE1	2.23	0.57
1:C:242:ILE:HB	1:C:259:PRO:HG2	1.87	0.57
1:A:67:LYS:HE3	1:A:87:GLY:HA2	1.88	0.56
1:A:103:GLU:CD	1:A:105:ARG:HH12	2.14	0.56
1:C:20:ASN:ND2	2:C:407:HOH:O	2.32	0.56
1:B:132:PRO:HG2	1:B:147:LEU:HB2	1.87	0.56
1:D:227:ARG:HD3	1:A:118:TRP:CH2	2.40	0.55
1:A:60:ARG:HD3	1:A:62:LEU:HD21	1.89	0.55
1:B:253:PRO:HA	1:B:256:ILE:HD12	1.89	0.54
1:A:43:ARG:HG3	1:A:47:THR:O	2.08	0.53
1:A:163:ARG:NH1	1:B:234:GLN:OE1	2.42	0.53
1:B:18:ILE:HD11	1:B:339:ASN:HB3	1.91	0.53
1:C:172:LYS:HB3	1:C:193:LYS:NZ	2.23	0.52
1:B:330:ALA:HB3	1:B:341:TRP:HZ3	1.74	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:136:ASP:HB3	1:C:143:VAL:HG13	1.91	0.51
1:B:253:PRO:HA	1:B:256:ILE:CD1	2.40	0.51
1:C:248:ALA:N	2:C:406:HOH:O	2.28	0.51
1:B:35:LYS:HE3	2:B:464:HOH:O	2.11	0.50
1:A:16:GLU:HG3	1:A:309:LEU:HD22	1.94	0.50
1:C:169:GLU:OE1	1:C:169:GLU:N	2.41	0.50
1:D:34:ARG:NE	1:D:34:ARG:HA	2.25	0.50
1:C:99:LYS:HE3	1:C:101:TYR:HA	1.94	0.49
1:C:168:LYS:NZ	2:C:403:HOH:O	2.26	0.49
1:D:257:GLN:HG2	1:D:259:PRO:HD3	1.93	0.49
1:A:103:GLU:OE1	1:A:105:ARG:NH1	2.45	0.49
1:B:312:PRO:HB3	1:B:337:ARG:HH22	1.77	0.49
1:A:30:VAL:O	1:A:320:GLU:HG3	2.13	0.49
1:B:168:LYS:HD2	1:B:168:LYS:H	1.77	0.48
1:D:225:SER:HB3	1:A:224:GLU:CG	2.41	0.48
1:D:34:ARG:NH2	2:D:417:HOH:O	2.47	0.48
1:C:173:HIS:ND1	2:C:402:HOH:O	2.25	0.48
1:A:279:SER:HA	1:A:321:TRP:CG	2.49	0.48
1:C:293[B]:HIS:HD2	1:C:313:LYS:HD2	1.80	0.47
1:B:242:ILE:HB	1:B:259:PRO:HG2	1.96	0.47
1:C:34:ARG:HH12	1:C:74:TRP:HB2	1.80	0.47
1:D:147:LEU:HD23	1:D:204:ARG:HG2	1.97	0.47
1:D:170:ASP:OD2	1:D:170:ASP:N	2.48	0.47
1:C:204:ARG:NE	2:C:409:HOH:O	2.35	0.47
1:C:230:LEU:HB2	1:C:278:PHE:CZ	2.50	0.47
1:D:43:ARG:HG3	1:D:47:THR:O	2.15	0.47
1:B:307:ARG:NH1	1:B:344:THR:H	2.13	0.47
1:C:267:VAL:HG23	2:C:454:HOH:O	2.15	0.46
1:C:293[B]:HIS:CD2	1:C:313:LYS:HD2	2.50	0.46
1:B:195:ARG:HG3	2:B:476:HOH:O	2.15	0.46
1:D:245:ASN:HB2	1:A:118:TRP:CE3	2.51	0.46
1:A:92:LEU:HD23	1:A:101:TYR:HD2	1.81	0.46
1:C:279:SER:HA	1:C:321:TRP:CG	2.51	0.46
1:B:230:LEU:HB2	1:B:278:PHE:CZ	2.51	0.45
1:B:260:LEU:HD21	1:B:263:LYS:HG3	1.97	0.45
1:B:97:ASP:OD1	1:B:99:LYS:HE2	2.16	0.45
1:B:133:MET:HE3	1:B:144:ARG:HB3	1.99	0.45
1:B:169:GLU:OE1	1:B:172:LYS:HD2	2.17	0.45
1:C:272[A]:SER:HB3	1:C:290:TYR:HB3	1.97	0.45
1:D:202:ARG:O	1:D:204:ARG:NH2	2.50	0.44
1:D:227:ARG:CZ	1:A:118:TRP:CE2	3.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:279:SER:HA	1:D:321:TRP:CG	2.52	0.44
1:D:195:ARG:HH11	1:D:211:LYS:HB2	1.73	0.44
1:B:145:TYR:CE2	1:B:204:ARG:CZ	3.01	0.44
1:A:242:ILE:HB	1:A:259:PRO:HG2	1.99	0.44
1:C:34:ARG:HD2	1:C:76:ARG:HA	2.00	0.44
1:B:94:ASP:HB3	1:B:99:LYS:HE3	1.99	0.44
1:C:89:LYS:HG2	1:C:105:ARG:HG2	2.00	0.44
1:C:253:PRO:HA	1:C:256:ILE:HG13	1.99	0.43
1:B:274:ASN:OD1	1:B:290[A]:TYR:CE1	2.72	0.43
1:D:126:ALA:C	1:D:127:LEU:HD22	2.44	0.43
1:A:309:LEU:HD13	1:A:341:TRP:CD1	2.54	0.43
1:D:80:TYR:HB3	1:D:92:LEU:HD11	2.01	0.42
1:D:230:LEU:HB2	1:D:278:PHE:CZ	2.54	0.42
1:B:29[B]:CYS:SG	1:B:69:ILE:HG22	2.59	0.42
1:A:53:LEU:HD23	1:A:53:LEU:HA	1.81	0.42
1:B:34:ARG:HG2	2:B:509:HOH:O	2.18	0.42
1:A:92:LEU:HD12	1:A:92:LEU:HA	1.88	0.42
1:B:91:ILE:HG23	1:B:100:ARG:HG3	2.02	0.42
1:C:166:GLN:OE1	1:C:170:ASP:OD2	2.37	0.42
1:C:215:GLY:HA3	1:C:234:GLN:HG2	2.01	0.42
1:A:221:ARG:HG2	2:A:493:HOH:O	2.19	0.41
1:A:230:LEU:HB2	1:A:278:PHE:CZ	2.55	0.41
1:C:34:ARG:HG2	2:C:416:HOH:O	2.19	0.41
1:C:220:LEU:O	1:C:221:ARG:HG2	2.20	0.41
1:C:234:GLN:NE2	2:C:413:HOH:O	2.44	0.41
1:D:228:GLU:HG2	1:D:300:ARG:NH2	2.35	0.41
1:A:229:LEU:O	1:A:240:THR:HA	2.20	0.41
1:B:279:SER:HA	1:B:321:TRP:CG	2.55	0.41
1:C:246:LEU:C	2:C:406:HOH:O	2.63	0.41
1:B:43:ARG:HG3	1:B:47:THR:O	2.21	0.40
1:D:238:ILE:HB	1:D:264:PHE:HB2	2.03	0.40
1:B:136:ASP:HB3	1:B:143:VAL:HB	2.03	0.40
1:C:229:LEU:O	1:C:240:THR:HA	2.22	0.40
1:D:174:MET:HE3	1:D:174:MET:HB2	1.84	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:195:ARG:NH2	1:C:141:VAL:CB[1_545]	2.16	0.04

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	329/345 (95%)	314 (95%)	14 (4%)	1 (0%)	36	26
1	B	319/345 (92%)	308 (97%)	11 (3%)	0	100	100
1	C	318/345 (92%)	311 (98%)	6 (2%)	1 (0%)	36	26
1	D	313/345 (91%)	304 (97%)	8 (3%)	1 (0%)	36	26
All	All	1279/1380 (93%)	1237 (97%)	39 (3%)	3 (0%)	43	33

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	140	PRO
1	C	140	PRO
1	D	140	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	287/296 (97%)	284 (99%)	3 (1%)	68	60
1	B	276/296 (93%)	274 (99%)	2 (1%)	76	70
1	C	277/296 (94%)	275 (99%)	2 (1%)	76	70
1	D	272/296 (92%)	271 (100%)	1 (0%)	84	81
All	All	1112/1184 (94%)	1104 (99%)	8 (1%)	81	70

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	141	VAL
1	A	141	VAL
1	A	150[A]	VAL
1	A	150[B]	VAL
1	B	150[A]	VAL
1	B	150[B]	VAL
1	C	150[A]	VAL
1	C	150[B]	VAL

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

Mol	Chain	Res	Type
1	D	26	HIS
1	D	173	HIS
1	A	20	ASN
1	A	166	GLN
1	A	265	GLN
1	A	292	ASN
1	B	26	HIS
1	B	122	GLN
1	B	302	HIS
1	C	120	HIS
1	C	122	GLN
1	C	166	GLN

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

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	327/345 (94%)	0.13	13 (3%) 42 48	15, 39, 65, 113	8 (2%)
1	B	315/345 (91%)	-0.10	9 (2%) 53 60	13, 32, 56, 102	8 (2%)
1	C	314/345 (91%)	-0.06	12 (3%) 44 50	11, 31, 62, 152	10 (3%)
1	D	312/345 (90%)	-0.20	8 (2%) 57 62	13, 29, 55, 129	7 (2%)
All	All	1268/1380 (91%)	-0.06	42 (3%) 49 56	11, 32, 61, 152	33 (2%)

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	253	PRO	7.2
1	C	15	PRO	6.0
1	D	251	LEU	4.4
1	B	165	LYS	4.1
1	A	250	ASP	3.9
1	B	18	ILE	3.9
1	C	344	THR	3.8
1	A	253	PRO	3.7
1	B	344	THR	3.7
1	C	249	ALA	3.7
1	C	141	VAL	3.7
1	A	344	THR	3.4
1	C	253	PRO	3.4
1	C	312	PRO	3.4
1	C	19	THR	3.0
1	B	20	ASN	3.0
1	A	249	ALA	2.9
1	C	168	LYS	2.9
1	A	140	PRO	2.8
1	A	252	ASP	2.8
1	D	18	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	247	SER	2.7
1	B	204	ARG	2.7
1	A	141	VAL	2.7
1	B	290[A]	TYR	2.6
1	A	15	PRO	2.6
1	D	249	ALA	2.6
1	C	167	ALA	2.6
1	A	302	HIS	2.5
1	A	256	ILE	2.5
1	B	19	THR	2.5
1	D	195	ARG	2.4
1	D	252	ASP	2.4
1	B	152	LYS	2.4
1	B	169	GLU	2.4
1	A	245	ASN	2.4
1	D	344	THR	2.4
1	D	256	ILE	2.2
1	C	214	SER	2.1
1	A	88	TRP	2.1
1	C	245	ASN	2.0
1	C	311	GLY	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.