



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 15, 2026 – 12:47 AM UTC

PDB ID : 9XHN / pdb_00009xhn
Title : Crystal structure of the C-terminal domain of AcvB from *Agrobacterium tumefaciens*
Authors : Hoshi, M.; Watanabe, Y.
Deposited on : 2025-11-01
Resolution : 1.83 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

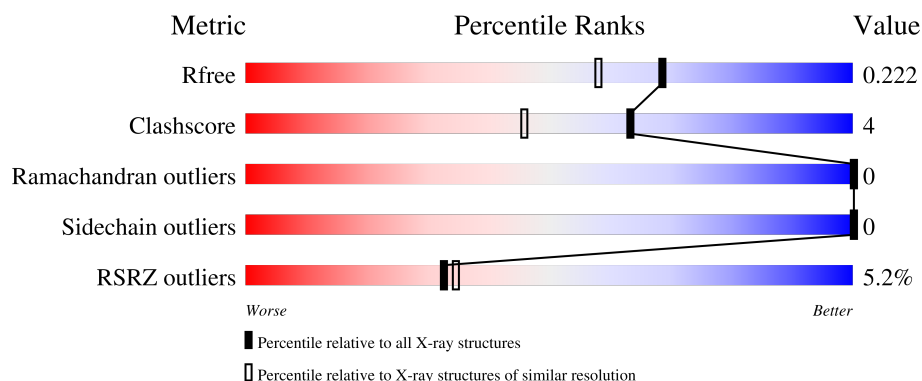
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.83 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	1296 (1.84-1.84)
Clashscore	190562	1329 (1.84-1.84)
Ramachandran outliers	187476	1318 (1.84-1.84)
Sidechain outliers	187428	1318 (1.84-1.84)
RSRZ outliers	180081	1296 (1.84-1.84)

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	212	<div> <div>6%</div> <div> <div></div> <div>91%</div> <div>8%</div> <div>.</div> </div> </div>
1	B	212	<div> <div>5%</div> <div> <div></div> <div>92%</div> <div>6%</div> <div>.</div> </div> </div>
1	C	212	<div> <div>2%</div> <div> <div></div> <div>88%</div> <div>7%</div> <div>5%</div> </div> </div>
1	D	212	<div> <div>7%</div> <div> <div></div> <div>83%</div> <div>10%</div> <div>7%</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
2	PEG	A	501	-	-	X	-
2	PEG	D	501	-	-	X	-

2 Entry composition [i](#)

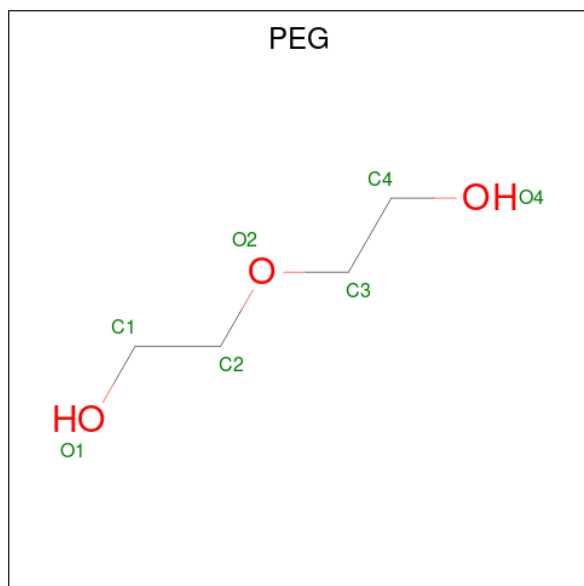
There are 3 unique types of molecules in this entry. The entry contains 6618 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 AcvB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	208	Total	C	N	O	S	0	0	0
			1593	1015	262	312	4			
1	B	206	Total	C	N	O	S	0	0	0
			1563	997	257	305	4			
1	C	202	Total	C	N	O	S	0	0	0
			1551	988	253	306	4			
1	D	198	Total	C	N	O	S	0	0	0
			1498	956	248	290	4			

- Molecule 2 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			7	4	3		
2	B	1	Total	C	O	0	0
			7	4	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			7	4	3		
2	D	1	Total	C	O	0	0
			7	4	3		

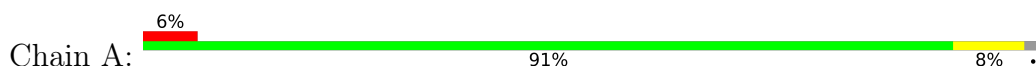
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	103	Total	O	0	0
			103	103		
3	B	96	Total	O	0	0
			96	96		
3	C	123	Total	O	0	0
			123	123		
3	D	63	Total	O	0	0
			63	63		

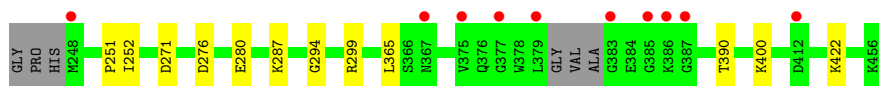
3 Residue-property plots [i](#)

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.

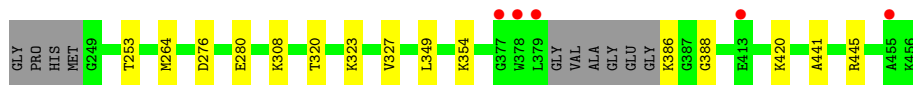
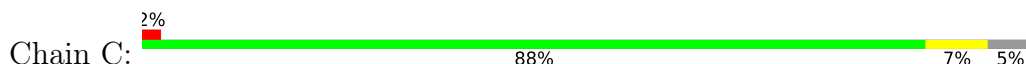
- Molecule 1: AcvB



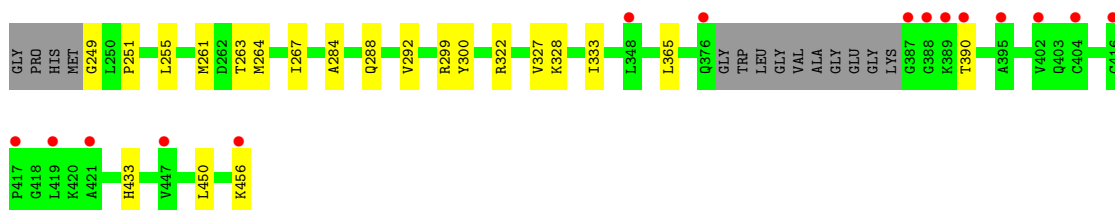
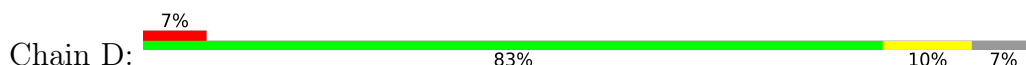
- Molecule 1: AcvB



- Molecule 1: AcvB



- Molecule 1: AcvB



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	47.16Å 64.15Å 68.82Å 85.83° 81.44° 87.70°	Depositor
Resolution (Å)	46.61 – 1.83 46.61 – 1.83	Depositor EDS
% Data completeness (in resolution range)	97.6 (46.61-1.83) 97.6 (46.61-1.83)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.69 (at 1.83Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.197 , 0.222 0.198 , 0.222	Depositor DCC
R_{free} test set	3319 reflections (4.74%)	wwPDB-VP
Wilson B-factor (Å ²)	30.9	Xtriage
Anisotropy	0.276	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 33.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6618	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.75% 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: PEG

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	0.28	0/1622	0.48	0/2196
1	B	0.32	0/1591	0.53	0/2156
1	C	0.35	0/1579	0.53	0/2141
1	D	0.30	0/1524	0.50	0/2067
All	All	0.31	0/6316	0.51	0/8560

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	1593	0	1593	16	0
1	B	1563	0	1543	9	0
1	C	1551	0	1532	12	0
1	D	1498	0	1486	16	0
2	A	7	0	10	4	0
2	B	7	0	10	2	0
2	C	7	0	10	0	0
2	D	7	0	10	4	0
3	A	103	0	0	3	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	96	0	0	1	0
3	C	123	0	0	2	0
3	D	63	0	0	2	1
All	All	6618	0	6194	50	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 4.

All (50) 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:365:LEU:HD22	1:A:390:THR:HG21	1.78	0.66
1:B:400:LYS:HE3	1:B:400:LYS:HA	1.78	0.66
1:C:386:LYS:HG3	1:C:388:GLY:H	1.61	0.64
1:B:251:PRO:HD3	2:B:501:PEG:H21	1.80	0.64
1:B:276:ASP:O	1:B:280:GLU:HG3	1.99	0.62
1:C:276:ASP:O	1:C:280:GLU:HG3	1.99	0.60
1:B:365:LEU:HD22	1:B:390:THR:HG21	1.83	0.59
1:D:300:TYR:HD1	2:D:501:PEG:H41	1.67	0.59
1:C:253:THR:HG21	3:C:608:HOH:O	2.02	0.58
1:D:365:LEU:HD22	1:D:390:THR:HG21	1.85	0.58
1:D:299:ARG:NH2	3:D:603:HOH:O	2.36	0.57
1:A:299:ARG:NH2	2:A:501:PEG:O4	2.37	0.57
1:C:264:MET:HB3	1:C:327:VAL:HG21	1.86	0.56
1:D:264:MET:HB3	1:D:327:VAL:HG21	1.88	0.55
1:D:249:GLY:N	3:D:604:HOH:O	2.40	0.54
1:A:271:ASP:HB3	1:C:320:THR:HG22	1.90	0.53
1:C:420:LYS:NZ	3:C:602:HOH:O	2.43	0.50
1:D:322:ARG:HH21	1:D:328:LYS:HA	1.77	0.49
1:A:309:GLU:OE2	3:A:601:HOH:O	2.20	0.48
1:D:251:PRO:HD3	2:D:501:PEG:H22	1.94	0.48
1:C:308:LYS:HD2	1:C:308:LYS:C	2.38	0.48
1:C:441:ALA:O	1:C:445:ARG:HG3	2.12	0.48
3:B:616:HOH:O	1:D:433:HIS:HE1	1.95	0.47
1:D:267:ILE:HG12	1:D:333:ILE:HB	1.96	0.47
1:C:349:LEU:HB2	1:C:354:LYS:HG3	1.96	0.47
1:B:299:ARG:NH2	2:B:501:PEG:H41	2.31	0.46
1:D:300:TYR:CD1	2:D:501:PEG:H41	2.50	0.46
1:A:258:LYS:NZ	3:A:610:HOH:O	2.49	0.45
3:A:623:HOH:O	1:C:320:THR:HG23	2.17	0.45
1:A:271:ASP:OD1	1:C:323:LYS:HE2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:255:LEU:HB2	1:D:292:VAL:HB	1.98	0.45
1:D:263:THR:HG21	1:D:450:LEU:HD21	1.99	0.44
1:A:251:PRO:HD3	2:A:501:PEG:H32	2.00	0.44
1:A:271:ASP:HB3	1:C:320:THR:CG2	2.48	0.43
1:D:456:LYS:N	1:D:456:LYS:HD3	2.33	0.43
1:A:258:LYS:HD3	1:A:259:PRO:HD2	2.00	0.43
1:D:284:ALA:O	1:D:288:GLN:HG3	2.19	0.42
1:A:320:THR:HG23	1:B:271:ASP:HB3	2.00	0.42
1:B:422:LYS:HA	1:B:422:LYS:HE3	2.01	0.42
1:A:346:TYR:O	1:A:354:LYS:HE3	2.19	0.42
1:A:323:LYS:HB2	1:A:323:LYS:HE2	1.72	0.41
1:D:261:MET:HE3	1:D:261:MET:HB2	1.77	0.41
1:B:287:LYS:HB3	1:B:287:LYS:HE2	1.54	0.41
2:A:501:PEG:H31	2:A:501:PEG:H12	1.67	0.41
1:D:251:PRO:HD3	2:D:501:PEG:H42	2.02	0.41
1:A:258:LYS:HD3	1:A:258:LYS:HA	1.83	0.40
1:A:306:ASP:O	1:A:309:GLU:HG2	2.21	0.40
1:A:406:TYR:CZ	1:A:417:PRO:HG3	2.56	0.40
1:A:300:TYR:HD1	2:A:501:PEG:H21	1.86	0.40
1:B:252:ILE:HA	1:B:294:GLY:O	2.22	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)
3:A:700:HOH:O	3:D:650:HOH:O[1_655]	2.19	0.01

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	206/212 (97%)	202 (98%)	4 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	202/212 (95%)	199 (98%)	3 (2%)	0	100	100
1	C	198/212 (93%)	194 (98%)	4 (2%)	0	100	100
1	D	194/212 (92%)	188 (97%)	6 (3%)	0	100	100
All	All	800/848 (94%)	783 (98%)	17 (2%)	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	169/173 (98%)	169 (100%)	0	100	100
1	B	163/173 (94%)	163 (100%)	0	100	100
1	C	165/173 (95%)	165 (100%)	0	100	100
1	D	157/173 (91%)	157 (100%)	0	100	100
All	All	654/692 (94%)	654 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	288	GLN
1	A	367	ASN
1	B	288	GLN
1	D	433	HIS

5.3.3 RNA ⓘ

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](#)

4 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	$\# Z > 2$
2	PEG	D	501	-	6,6,6	0.44	0	5,5,5	0.40	0
2	PEG	A	501	-	6,6,6	0.46	0	5,5,5	0.35	0
2	PEG	C	501	-	6,6,6	0.48	0	5,5,5	0.28	0
2	PEG	B	501	-	6,6,6	0.45	0	5,5,5	0.56	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	PEG	D	501	-	-	4/4/4/4	-
2	PEG	A	501	-	-	3/4/4/4	-
2	PEG	C	501	-	-	2/4/4/4	-
2	PEG	B	501	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	501	PEG	O1-C1-C2-O2
2	D	501	PEG	O2-C3-C4-O4
2	D	501	PEG	C4-C3-O2-C2
2	A	501	PEG	C1-C2-O2-C3
2	A	501	PEG	O2-C3-C4-O4
2	A	501	PEG	O1-C1-C2-O2
2	D	501	PEG	C1-C2-O2-C3
2	C	501	PEG	O1-C1-C2-O2
2	C	501	PEG	C4-C3-O2-C2

There are no ring outliers.

3 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	501	PEG	4	0
2	A	501	PEG	4	0
2	B	501	PEG	2	0

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	208/212 (98%)	0.57	12 (5%) 29 30	21, 37, 54, 68	0
1	B	206/212 (97%)	0.35	10 (4%) 35 36	19, 32, 52, 70	0
1	C	202/212 (95%)	0.20	5 (2%) 58 64	18, 30, 48, 62	0
1	D	198/212 (93%)	0.70	15 (7%) 20 20	22, 37, 53, 62	0
All	All	814/848 (95%)	0.45	42 (5%) 33 34	18, 34, 52, 70	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	421	ALA	4.3
1	A	378	TRP	4.0
1	B	248	MET	3.6
1	C	379	LEU	3.5
1	D	416	CYS	3.4
1	A	249	GLY	3.4
1	A	427	ILE	3.4
1	D	395	ALA	3.3
1	A	421	ALA	3.1
1	B	387	GLY	3.0
1	B	386	LYS	3.0
1	D	447	VAL	2.9
1	C	378	TRP	2.8
1	C	455	ALA	2.7
1	D	388	GLY	2.7
1	B	383	GLY	2.6
1	D	390	THR	2.6
1	D	387	GLY	2.6
1	D	376	GLN	2.5
1	B	385	GLY	2.5
1	A	426	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	404	CYS	2.4
1	D	402	VAL	2.4
1	B	377	GLY	2.4
1	C	377	GLY	2.4
1	D	389	LYS	2.4
1	D	456	LYS	2.4
1	A	381	VAL	2.3
1	D	419	LEU	2.3
1	B	379	LEU	2.2
1	A	412	ASP	2.2
1	A	411	GLU	2.2
1	B	375	VAL	2.2
1	A	456	LYS	2.2
1	D	348	LEU	2.1
1	B	412	ASP	2.1
1	A	405	VAL	2.1
1	A	379	LEU	2.1
1	C	413	GLU	2.0
1	B	367	ASN	2.0
1	D	417	PRO	2.0
1	A	310	VAL	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](#)

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(\AA^2)	Q<0.9
2	PEG	B	501	7/7	0.74	0.14	34,40,50,51	0
2	PEG	D	501	7/7	0.83	0.12	32,40,45,47	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	PEG	C	501	7/7	0.86	0.12	35,41,47,49	0
2	PEG	A	501	7/7	0.86	0.11	37,41,45,52	0

6.5 Other polymers [i](#)

There are no such residues in this entry.