



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 07:07 AM UTC

PDB ID : 3GHP / pdb_00003ghp
Title : Structure of the second type II cohesin module from the adaptor ScaA scaffoldin of *Acetivibrio cellulolyticus* (including long C-terminal linker)
Authors : Noach, I.; Frolow, F.; Bayer, E.A.
Deposited on : 2009-03-04
Resolution : 2.49 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

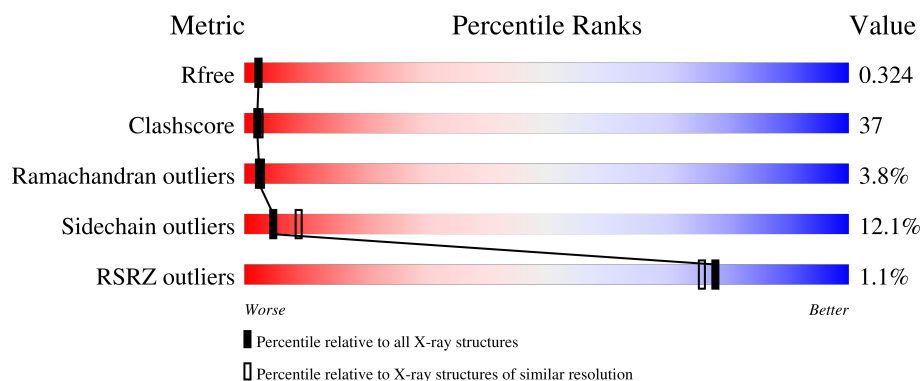
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.49 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	7589 (2.50-2.46)
Clashscore	190562	8295 (2.50-2.46)
Ramachandran outliers	187476	8164 (2.50-2.46)
Sidechain outliers	187428	8166 (2.50-2.46)
RSRZ outliers	180081	7593 (2.50-2.46)

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	227	
1	B	227	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2975 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 Cellulosomal scaffoldin adaptor protein B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	190	Total	C	N	O	S	4	2	0
			1457	929	235	290	3			
1	B	183	Total	C	N	O	S	14	0	0
			1389	885	224	277	3			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	220	LEU	-	expression tag	UNP Q7WYN3
A	221	GLU	-	expression tag	UNP Q7WYN3
A	222	HIS	-	expression tag	UNP Q7WYN3
A	223	HIS	-	expression tag	UNP Q7WYN3
A	224	HIS	-	expression tag	UNP Q7WYN3
A	225	HIS	-	expression tag	UNP Q7WYN3
A	226	HIS	-	expression tag	UNP Q7WYN3
A	227	HIS	-	expression tag	UNP Q7WYN3
B	220	LEU	-	expression tag	UNP Q7WYN3
B	221	GLU	-	expression tag	UNP Q7WYN3
B	222	HIS	-	expression tag	UNP Q7WYN3
B	223	HIS	-	expression tag	UNP Q7WYN3
B	224	HIS	-	expression tag	UNP Q7WYN3
B	225	HIS	-	expression tag	UNP Q7WYN3
B	226	HIS	-	expression tag	UNP Q7WYN3
B	227	HIS	-	expression tag	UNP Q7WYN3

- Molecule 2 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	43	Total 43	O 43	0	0
3	B	30	Total 30	O 30	0	0

- Molecule 1: Cellulosomal scaffoldin adaptor protein B



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	68.76Å 159.22Å 44.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.60 – 2.49 42.60 – 2.49	Depositor EDS
% Data completeness (in resolution range)	65.7 (42.60-2.49) 71.8 (42.60-2.49)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 2.48Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.212 , 0.317 0.222 , 0.324	Depositor DCC
R_{free} test set	662 reflections (5.18%)	wwPDB-VP
Wilson B-factor (Å ²)	52.8	Xtriage
Anisotropy	0.664	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 80.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2975	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.98% 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: EDO

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	0.64	0/1488	1.12	7/2026 (0.3%)
1	B	0.63	0/1416	1.04	8/1930 (0.4%)
All	All	0.64	0/2904	1.08	15/3956 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	2

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	168	GLN	CA-C-N	10.31	130.62	119.28
1	A	168	GLN	C-N-CA	10.31	130.62	119.28
1	B	168	GLN	N-CA-C	7.29	125.93	109.81
1	B	118	GLY	N-CA-C	6.05	119.83	110.88
1	A	180	ALA	N-CA-C	5.87	113.60	108.78
1	A	150	THR	N-CA-C	5.85	118.65	108.76
1	B	143	MET	CA-C-N	5.55	125.21	119.05
1	B	143	MET	C-N-CA	5.55	125.21	119.05
1	B	49	ASP	CA-C-N	5.46	125.07	119.56
1	B	49	ASP	C-N-CA	5.46	125.07	119.56
1	B	127	VAL	CB-CA-C	-5.42	106.05	111.80
1	A	11	TYR	N-CA-C	5.21	116.12	107.73
1	A	143	MET	CA-C-N	5.16	125.20	119.32
1	A	143	MET	C-N-CA	5.16	125.20	119.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	10	SER	N-CA-C	5.14	117.04	107.99

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	167	ILE	Peptide
1	B	174	SER	Peptide

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	1457	0	1468	99	0
1	B	1389	0	1396	120	0
2	A	24	0	36	3	0
2	B	32	0	48	8	0
3	A	43	0	0	2	0
3	B	30	0	0	3	0
All	All	2975	0	2948	216	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 37.

All (216) 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:169:PRO:HB2	1:B:170:GLY:CA	1.79	1.12
1:B:7:ILE:HB	1:B:8:LYS:HA	1.11	1.09
1:B:169:PRO:CB	1:B:170:GLY:HA3	1.86	1.05
1:A:188:ALA:CB	1:A:189:PHE:HA	1.92	0.99
1:B:7:ILE:CB	1:B:8:LYS:HA	1.94	0.96
1:B:174:SER:HA	1:B:175:ALA:HB3	1.44	0.96
1:B:175:ALA:N	1:B:176:PRO:HD3	1.81	0.95
1:A:7:ILE:HD12	1:A:8:LYS:H	1.31	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:ILE:HG22	1:A:39:PHE:HB2	1.51	0.93
1:B:7:ILE:HB	1:B:8:LYS:CA	1.99	0.91
1:B:174:SER:HA	1:B:175:ALA:CB	2.01	0.90
1:B:73:GLY:HA3	1:B:119:THR:HG22	1.54	0.90
1:B:60:THR:OG1	1:B:62:VAL:HG12	1.72	0.88
1:B:169:PRO:HB2	1:B:170:GLY:HA3	0.91	0.87
1:A:56:VAL:HG23	1:A:57:ASN:O	1.81	0.81
1:A:188:ALA:HB1	1:A:189:PHE:HA	1.60	0.80
1:B:175:ALA:H	1:B:176:PRO:HD3	1.45	0.80
1:B:176:PRO:O	1:B:177:ILE:HG22	1.82	0.80
1:B:19:ASN:O	1:B:172:ILE:HG22	1.82	0.79
1:A:188:ALA:HB3	1:A:189:PHE:HA	1.65	0.78
1:B:6:MET:HB3	1:B:7:ILE:HD13	1.66	0.77
1:B:57:ASN:HB2	2:B:228:EDO:H12	1.67	0.77
1:A:134:THR:HG23	1:A:170:GLY:O	1.86	0.75
1:B:174:SER:CA	1:B:175:ALA:HB3	2.16	0.75
1:B:173:ASN:O	1:B:174:SER:C	2.30	0.75
1:B:165:PHE:HD2	1:B:166:VAL:H	1.33	0.74
1:B:133:THR:N	1:B:172:ILE:O	2.20	0.74
1:B:175:ALA:N	1:B:176:PRO:CD	2.52	0.73
1:B:83:PRO:HG3	1:B:102:TYR:HE2	1.52	0.73
1:B:161:GLN:O	1:B:162:SER:HB2	1.89	0.72
1:B:56:VAL:HG12	1:B:63:ALA:HA	1.70	0.72
1:B:73:GLY:CA	1:B:119:THR:HG22	2.20	0.72
1:A:188:ALA:CB	1:A:189:PHE:CA	2.69	0.71
1:B:183:THR:HG23	1:B:183:THR:O	1.90	0.71
1:A:22:GLU:O	1:A:25:GLU:HG2	1.92	0.70
1:A:50:PRO:HG3	1:A:94:GLY:CA	2.21	0.70
1:B:68:SER:HA	2:B:229:EDO:H21	1.74	0.70
1:A:175:ALA:HB2	3:A:269:HOH:O	1.94	0.67
1:B:81:TYR:HA	1:B:104:ALA:HB3	1.78	0.66
1:B:17:ASP:CG	1:B:18:LYS:H	2.05	0.65
1:B:81:TYR:CE2	1:B:107:VAL:HG12	2.31	0.65
1:B:8:LYS:CG	1:B:9:ALA:H	2.10	0.65
1:A:50:PRO:HG3	1:A:94:GLY:HA3	1.78	0.65
1:B:8:LYS:HG3	1:B:9:ALA:H	1.62	0.64
1:B:148:THR:HG22	1:B:148:THR:O	1.97	0.64
1:B:20:ALA:HB1	1:B:173:ASN:HB3	1.80	0.63
1:A:65:THR:HG22	1:A:67:SER:H	1.64	0.62
1:B:183:THR:O	1:B:184:THR:C	2.42	0.62
1:B:183:THR:O	1:B:183:THR:CG2	2.46	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:166:VAL:O	2:B:232:EDO:H12	1.99	0.62
1:A:36:ILE:CG2	1:A:39:PHE:HB2	2.26	0.61
1:A:66:ASN:HB2	1:A:89:HIS:O	2.01	0.61
1:A:109:ARG:HD3	1:A:155:TRP:O	2.01	0.61
1:B:75:LEU:HD12	1:B:120:LEU:HD21	1.83	0.60
1:B:38:ASN:O	1:B:154:ASN:HA	2.02	0.60
1:A:89:HIS:HB2	1:A:91:ILE:HD12	1.84	0.60
1:B:174:SER:CA	1:B:175:ALA:CB	2.75	0.60
1:B:165:PHE:HE2	2:B:232:EDO:HO2	1.50	0.60
1:A:21:ALA:HB3	1:A:127:VAL:HG21	1.84	0.59
2:A:232:EDO:O2	2:B:229:EDO:H11	2.02	0.59
1:B:17:ASP:CG	1:B:18:LYS:N	2.61	0.59
1:B:177:ILE:HG23	1:B:178:ALA:O	2.03	0.58
1:B:57:ASN:HB2	2:B:228:EDO:C1	2.33	0.58
1:A:158:ASN:OD1	1:B:106:GLU:HB2	2.04	0.58
1:A:188:ALA:HB3	1:A:189:PHE:CA	2.32	0.58
1:A:159:ARG:HD2	1:A:161:GLN:NE2	2.19	0.58
1:B:83:PRO:HG3	1:B:102:TYR:CE2	2.37	0.57
1:A:7:ILE:HD12	1:A:8:LYS:N	2.12	0.57
1:A:148:THR:HG23	2:A:233:EDO:H22	1.84	0.57
1:A:115:GLU:HG3	2:A:228:EDO:H11	1.87	0.57
1:A:66:ASN:HA	1:A:89:HIS:HD2	1.70	0.57
1:B:47:LYS:HD2	1:B:94:GLY:O	2.05	0.56
1:A:184:THR:C	1:A:186:PRO:HD3	2.30	0.56
1:A:43:GLN:HG3	1:A:99:SER:OG	2.05	0.56
1:A:143:MET:N	1:A:144:PRO:HD3	2.22	0.55
1:A:66:ASN:HB2	1:A:89:HIS:C	2.31	0.55
1:A:99:SER:CB	1:B:84:ILE:HD11	2.36	0.55
1:B:45:ASN:ND2	1:B:140:SER:HB2	2.21	0.55
1:A:148:THR:O	1:A:150:THR:HG23	2.07	0.55
1:B:109:ARG:HG3	1:B:155:TRP:CD2	2.42	0.55
1:B:168:GLN:CB	1:B:169:PRO:O	2.54	0.55
1:A:33:ILE:HG23	1:A:120:LEU:HD11	1.89	0.54
1:B:30:THR:HB	2:B:230:EDO:H22	1.88	0.54
1:B:75:LEU:O	1:B:77:VAL:HG13	2.07	0.54
1:B:57:ASN:HB3	1:B:60:THR:OG1	2.07	0.54
1:A:76:LEU:HA	1:A:102:TYR:CZ	2.43	0.54
1:B:33:ILE:HG23	1:B:120:LEU:HD11	1.90	0.54
1:A:7:ILE:CD1	1:A:8:LYS:H	2.11	0.54
1:A:42:TYR:CD1	1:A:42:TYR:C	2.86	0.54
1:B:47:LYS:NZ	1:B:48:TYR:O	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:9:ALA:N	3:B:251:HOH:O	2.41	0.53
1:A:66:ASN:HA	1:A:89:HIS:CD2	2.43	0.53
1:B:81:TYR:HE2	1:B:107:VAL:HG12	1.72	0.53
1:B:33:ILE:CG2	1:B:120:LEU:HD11	2.40	0.52
1:A:175:ALA:N	1:A:176:PRO:HD3	2.24	0.52
1:B:69:LEU:N	1:B:69:LEU:HD12	2.25	0.52
1:A:151:THR:O	1:A:152:LEU:HD23	2.10	0.52
1:B:11:TYR:CZ	1:B:34:ASN:HB2	2.45	0.52
1:A:142:THR:C	1:A:144:PRO:HD3	2.34	0.52
1:A:109:ARG:HG3	1:A:155:TRP:CG	2.45	0.52
1:A:21:ALA:HB1	1:A:25:GLU:HG3	1.92	0.51
1:B:48:TYR:HB3	1:B:135:VAL:HG22	1.93	0.51
1:A:89:HIS:HB3	1:A:96:LEU:HD23	1.92	0.51
1:B:177:ILE:HG23	1:B:178:ALA:N	2.26	0.51
1:B:45:ASN:ND2	1:B:138:GLU:O	2.40	0.51
1:B:36:ILE:O	1:B:116:GLU:HG2	2.11	0.50
1:A:74:GLU:HG3	1:A:119:THR:OG1	2.11	0.50
1:A:85:VAL:HG11	1:A:100:ARG:NH1	2.26	0.50
1:A:173:ASN:O	1:A:174:SER:C	2.54	0.50
1:B:8:LYS:O	1:B:9:ALA:HB3	2.12	0.50
1:A:174:SER:C	1:A:176:PRO:HD3	2.37	0.50
1:B:81:TYR:CD2	1:B:107:VAL:HG12	2.46	0.50
1:B:20:ALA:CB	1:B:173:ASN:HB3	2.41	0.50
1:B:7:ILE:HG12	3:B:254:HOH:O	2.12	0.50
1:B:11:TYR:CD1	1:B:11:TYR:C	2.90	0.50
1:A:18[A]:LYS:NZ	1:A:20:ALA:HB3	2.27	0.49
1:A:14:MET:HB3	1:A:169:PRO:HB3	1.94	0.49
1:B:109:ARG:HG3	1:B:155:TRP:CG	2.47	0.49
1:B:82:GLY:O	1:B:103:THR:OG1	2.30	0.49
1:B:122:VAL:HG12	1:B:123:VAL:N	2.27	0.49
1:B:59:LYS:HG2	1:B:60:THR:N	2.27	0.49
1:A:99:SER:HB3	1:B:84:ILE:HD11	1.95	0.48
1:A:133:THR:O	1:A:171:GLU:HG2	2.13	0.48
1:B:14:MET:HE3	1:B:168:GLN:O	2.12	0.48
1:B:74:GLU:HG3	1:B:119:THR:HB	1.95	0.48
1:A:7:ILE:HG23	1:A:9:ALA:H	1.79	0.48
1:A:85:VAL:HG11	1:A:100:ARG:HH11	1.77	0.48
1:A:139:ASP:OD1	1:A:148:THR:HA	2.14	0.48
1:A:52:VAL:O	1:A:53:LEU:HD23	2.13	0.48
1:A:99:SER:HB2	1:B:84:ILE:HD11	1.95	0.48
1:A:85:VAL:HG22	1:A:86:GLN:N	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:47:LYS:HE2	1:B:93:GLU:O	2.13	0.47
1:A:176:PRO:C	1:A:178:ALA:H	2.23	0.47
1:B:22:GLU:OE2	1:B:130:LYS:HE3	2.14	0.47
1:B:69:LEU:N	1:B:69:LEU:CD1	2.78	0.47
1:A:57:ASN:OD1	1:A:58:PRO:HD2	2.14	0.47
1:B:105:LEU:O	1:B:108:TYR:HB3	2.14	0.47
1:A:53:LEU:HD23	1:A:53:LEU:HA	1.72	0.47
1:A:89:HIS:HB3	1:A:95:ILE:O	2.15	0.47
1:B:132:ALA:HB2	1:B:175:ALA:HB2	1.97	0.47
1:A:7:ILE:HG23	1:A:8:LYS:N	2.29	0.47
1:A:189:PHE:O	1:A:190:ALA:HB2	2.14	0.47
1:B:108:TYR:CD2	1:B:155:TRP:HB3	2.50	0.46
1:A:46:ILE:HG22	1:A:47:LYS:N	2.30	0.46
1:A:176:PRO:O	1:A:178:ALA:N	2.48	0.46
1:B:181:THR:N	1:B:182:PRO:CD	2.78	0.46
1:A:22:GLU:O	1:A:23:VAL:C	2.58	0.46
1:A:85:VAL:CG2	1:A:86:GLN:N	2.79	0.46
1:A:65:THR:N	1:A:68:SER:OG	2.38	0.46
1:A:65:THR:HG22	1:A:67:SER:N	2.31	0.46
1:A:143:MET:HE1	1:A:149:GLY:O	2.16	0.46
1:A:181:THR:N	1:A:182:PRO:HD3	2.31	0.45
1:A:65:THR:C	1:A:67:SER:H	2.23	0.45
1:B:8:LYS:CG	1:B:9:ALA:N	2.78	0.45
1:B:78:SER:N	1:B:102:TYR:OH	2.38	0.45
1:B:140:SER:OG	1:B:142:THR:HB	2.15	0.45
1:A:48:TYR:O	1:A:50:PRO:HD3	2.16	0.45
1:A:28:LYS:HE3	1:A:28:LYS:HB2	1.73	0.44
1:A:11:TYR:CD1	1:A:11:TYR:C	2.94	0.44
1:A:49:ASP:HB3	1:A:52:VAL:HB	1.99	0.44
1:A:184:THR:O	1:A:186:PRO:HD3	2.17	0.44
1:B:168:GLN:HB3	1:B:169:PRO:O	2.17	0.44
1:A:82:GLY:HA3	1:A:103:THR:HG22	1.98	0.44
1:B:50:PRO:HG3	1:B:94:GLY:CA	2.48	0.44
1:B:173:ASN:OD1	1:B:173:ASN:N	2.44	0.44
1:A:7:ILE:HD13	1:A:7:ILE:HA	1.84	0.44
1:A:72:SER:HB3	1:A:77:VAL:HG11	2.00	0.44
1:B:114:PRO:HB2	1:B:116:GLU:HG3	2.00	0.44
1:B:186:PRO:C	1:B:188:ALA:N	2.75	0.44
1:B:168:GLN:HB2	1:B:169:PRO:O	2.18	0.43
1:B:171:GLU:O	1:B:172:ILE:HG23	2.17	0.43
1:A:57:ASN:OD1	1:A:58:PRO:CD	2.67	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:57:ASN:HA	1:A:58:PRO:HD3	1.82	0.43
1:A:89:HIS:CG	1:A:96:LEU:HD23	2.54	0.43
1:B:59:LYS:H	1:B:59:LYS:HD2	1.82	0.43
1:B:84:ILE:O	1:B:100:ARG:HA	2.18	0.43
1:A:173:ASN:HB3	1:A:174:SER:H	1.60	0.43
1:B:155:TRP:CD1	1:B:156:TYR:CE1	3.07	0.43
1:B:114:PRO:HG3	1:B:155:TRP:CH2	2.54	0.42
1:B:155:TRP:HD1	1:B:156:TYR:CD1	2.36	0.42
1:A:89:HIS:CG	1:A:96:LEU:CD2	3.03	0.42
1:A:114:PRO:HD2	3:A:248:HOH:O	2.20	0.42
1:B:165:PHE:HD2	1:B:166:VAL:N	2.11	0.42
1:B:79:GLU:OE2	1:B:79:GLU:HA	2.19	0.42
1:A:53:LEU:CD2	1:A:127:VAL:HA	2.50	0.42
1:A:90:LYS:HD3	1:A:90:LYS:HA	1.89	0.42
1:B:155:TRP:NE1	1:B:156:TYR:CE1	2.88	0.42
1:A:65:THR:O	1:A:67:SER:N	2.53	0.42
1:A:189:PHE:CG	1:A:190:ALA:N	2.88	0.42
1:A:21:ALA:CB	1:A:127:VAL:HG21	2.49	0.42
1:A:78:SER:O	1:A:83:PRO:HD3	2.19	0.41
1:B:8:LYS:HB3	1:B:8:LYS:HE3	1.65	0.41
1:B:49:ASP:HA	1:B:50:PRO:HD3	1.83	0.41
1:A:75:LEU:HG	1:A:119:THR:O	2.20	0.41
1:A:175:ALA:N	1:A:176:PRO:CD	2.83	0.41
1:A:191:SER:OG	1:A:192:THR:N	2.52	0.41
1:B:81:TYR:CE2	1:B:108:TYR:HA	2.55	0.41
1:B:74:GLU:N	1:B:119:THR:O	2.54	0.41
1:B:88:VAL:O	1:B:97:ASN:HB3	2.21	0.41
1:B:138:GLU:HG2	1:B:139:ASP:N	2.35	0.41
1:A:22:GLU:H	1:A:25:GLU:CD	2.28	0.41
1:A:36:ILE:HA	1:A:36:ILE:HD12	1.84	0.41
1:A:46:ILE:CG2	1:A:47:LYS:N	2.83	0.41
1:A:165:PHE:N	1:A:165:PHE:CD2	2.88	0.41
1:B:43:GLN:HA	1:B:98:LEU:O	2.21	0.41
1:B:122:VAL:C	1:B:123:VAL:HG23	2.46	0.41
1:B:147:ILE:O	1:B:148:THR:HB	2.21	0.41
1:B:165:PHE:HE2	2:B:232:EDO:O2	2.03	0.41
1:B:176:PRO:O	1:B:177:ILE:CG2	2.61	0.41
1:B:50:PRO:HG3	1:B:94:GLY:N	2.37	0.40
1:B:56:VAL:HG12	1:B:63:ALA:CA	2.45	0.40
1:B:9:ALA:HA	3:B:251:HOH:O	2.20	0.40
1:B:47:LYS:HA	1:B:94:GLY:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:42:TYR:CE2	1:B:100:ARG:HB3	2.56	0.40
1:A:84:ILE:O	1:A:84:ILE:HG23	2.21	0.40
1:B:172:ILE:HA	1:B:173:ASN:HA	1.67	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	190/227 (84%)	160 (84%)	22 (12%)	8 (4%)	2	2
1	B	181/227 (80%)	149 (82%)	26 (14%)	6 (3%)	3	3
All	All	371/454 (82%)	309 (83%)	48 (13%)	14 (4%)	2	3

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	177	ILE
1	A	188	ALA
1	B	177	ILE
1	A	174	SER
1	A	176	PRO
1	A	182	PRO
1	A	191	SER
1	B	9	ALA
1	B	162	SER
1	B	184	THR
1	A	190	ALA
1	B	175	ALA
1	A	186	PRO
1	B	169	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	162/195 (83%)	143 (88%)	19 (12%)	5	10
1	B	154/195 (79%)	134 (87%)	20 (13%)	4	7
All	All	316/390 (81%)	277 (88%)	39 (12%)	5	8

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LEU
1	A	6	MET
1	A	7	ILE
1	A	8	LYS
1	A	18[A]	LYS
1	A	18[B]	LYS
1	A	36	ILE
1	A	56	VAL
1	A	65	THR
1	A	85	VAL
1	A	91	ILE
1	A	92	SER
1	A	97	ASN
1	A	103	THR
1	A	105	LEU
1	A	113	SER
1	A	133	THR
1	A	162	SER
1	A	177	ILE
1	B	7	ILE
1	B	8	LYS
1	B	30	THR
1	B	35	LYS
1	B	59	LYS
1	B	69	LEU
1	B	84	ILE
1	B	91	ILE

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Mol	Chain	Res	Type
1	B	101	SER
1	B	103	THR
1	B	117	THR
1	B	119	THR
1	B	142	THR
1	B	165	PHE
1	B	167	ILE
1	B	171	GLU
1	B	173	ASN
1	B	177	ILE
1	B	183	THR
1	B	184	THR

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	3	ASN
1	A	43	GLN
1	A	45	ASN
1	A	97	ASN
1	B	38	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

14 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	EDO	A	232	-	3,3,3	0.44	0	2,2,2	0.37	0
2	EDO	B	233	-	3,3,3	0.50	0	2,2,2	0.26	0
2	EDO	B	234	-	3,3,3	0.46	0	2,2,2	0.30	0
2	EDO	A	229	-	3,3,3	0.47	0	2,2,2	0.40	0
2	EDO	A	231	-	3,3,3	0.44	0	2,2,2	0.36	0
2	EDO	B	229	-	3,3,3	0.47	0	2,2,2	0.28	0
2	EDO	B	230	-	3,3,3	0.48	0	2,2,2	0.24	0
2	EDO	B	235	-	3,3,3	0.46	0	2,2,2	0.34	0
2	EDO	A	230	-	3,3,3	0.38	0	2,2,2	0.62	0
2	EDO	B	231	-	3,3,3	0.72	0	2,2,2	0.27	0
2	EDO	A	228	-	3,3,3	0.57	0	2,2,2	0.18	0
2	EDO	A	233	-	3,3,3	0.52	0	2,2,2	0.23	0
2	EDO	B	232	-	3,3,3	0.56	0	2,2,2	0.17	0
2	EDO	B	228	-	3,3,3	0.50	0	2,2,2	0.37	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	EDO	A	232	-	-	1/1/1/1	-
2	EDO	B	233	-	-	1/1/1/1	-
2	EDO	B	234	-	-	0/1/1/1	-
2	EDO	A	229	-	-	0/1/1/1	-
2	EDO	A	231	-	-	0/1/1/1	-
2	EDO	B	229	-	-	0/1/1/1	-
2	EDO	B	230	-	-	1/1/1/1	-
2	EDO	B	235	-	-	0/1/1/1	-
2	EDO	A	230	-	-	0/1/1/1	-
2	EDO	B	231	-	-	1/1/1/1	-
2	EDO	A	228	-	-	1/1/1/1	-
2	EDO	A	233	-	-	1/1/1/1	-
2	EDO	B	232	-	-	1/1/1/1	-
2	EDO	B	228	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	231	EDO	O1-C1-C2-O2
2	A	228	EDO	O1-C1-C2-O2
2	A	232	EDO	O1-C1-C2-O2
2	B	232	EDO	O1-C1-C2-O2
2	A	233	EDO	O1-C1-C2-O2
2	B	230	EDO	O1-C1-C2-O2
2	B	233	EDO	O1-C1-C2-O2

There are no ring outliers.

7 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	232	EDO	1	0
2	B	229	EDO	2	0
2	B	230	EDO	1	0
2	A	228	EDO	1	0
2	A	233	EDO	1	0
2	B	232	EDO	3	0
2	B	228	EDO	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 [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	190/227 (83%)	-0.07	1 (0%) 87 85	25, 57, 172, 198	6 (3%)
1	B	183/227 (80%)	0.08	3 (1%) 70 68	44, 65, 147, 177	11 (6%)
All	All	373/454 (82%)	0.00	4 (1%) 78 76	25, 61, 169, 198	17 (4%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	39	PHE	3.3
1	B	179	THR	2.8
1	B	157	GLY	2.2
1	A	40	SER	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(Å ²)	Q<0.9
2	EDO	A	229	4/4	0.61	0.11	107,111,112,113	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	EDO	B	231	4/4	0.71	0.22	51,63,75,76	0
2	EDO	B	235	4/4	0.74	0.12	87,88,89,89	0
2	EDO	B	233	4/4	0.75	0.11	87,88,92,92	0
2	EDO	B	229	4/4	0.75	0.13	95,96,101,102	0
2	EDO	A	231	4/4	0.78	0.08	93,96,99,104	0
2	EDO	A	232	4/4	0.78	0.12	108,109,109,110	0
2	EDO	B	230	4/4	0.81	0.18	71,82,93,100	0
2	EDO	A	230	4/4	0.82	0.08	101,104,107,110	0
2	EDO	B	234	4/4	0.83	0.14	102,108,112,115	0
2	EDO	A	228	4/4	0.84	0.19	69,70,77,77	0
2	EDO	B	232	4/4	0.87	0.13	62,66,67,71	0
2	EDO	A	233	4/4	0.92	0.09	65,73,83,84	0
2	EDO	B	228	4/4	0.94	0.10	87,88,91,91	0

6.5 Other polymers [i](#)

There are no such residues in this entry.