



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:14 AM UTC

PDB ID : 4WIW / pdb_00004wiw
Title : Crystal structure of C-terminal domain of putative chitinase from *Desulfotobacterium hafniense* DCB-2
Authors : Chang, C.; Tesar, C.; Endres, M.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2014-09-26
Resolution : 2.64 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

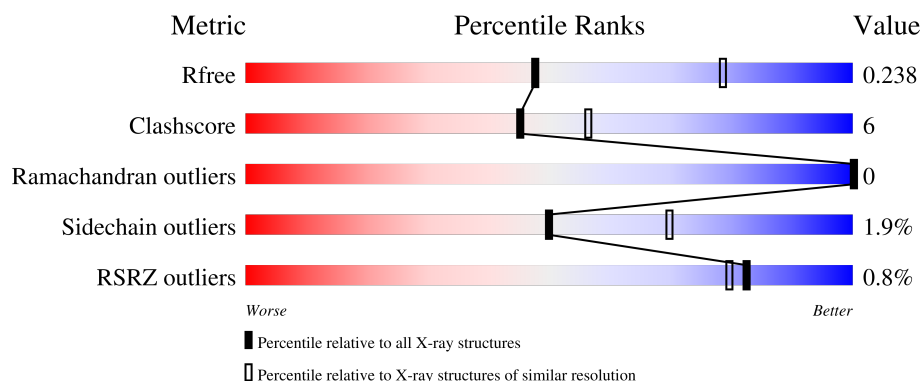
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

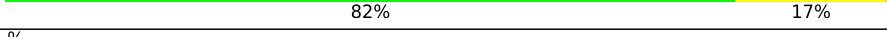
The reported resolution of this entry is 2.64 Å.

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	2053 (2.66-2.62)
Clashscore	190562	2097 (2.66-2.62)
Ramachandran outliers	187476	2066 (2.66-2.62)
Sidechain outliers	187428	2066 (2.66-2.62)
RSRZ outliers	180081	2052 (2.66-2.62)

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	349	
1	B	349	
1	C	349	
1	D	349	
1	E	349	

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Mol	Chain	Length	Quality of chain
1	F	349	 84% 15% .

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 16700 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 Glycoside hydrolase family 18.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	348	Total	C	N	O	Se	0	0	0
			2732	1753	440	528	11			
1	B	348	Total	C	N	O	Se	0	0	0
			2734	1753	442	528	11			
1	C	348	Total	C	N	O	Se	0	0	0
			2720	1746	440	523	11			
1	D	349	Total	C	N	O	Se	0	0	0
			2716	1741	440	524	11			
1	E	348	Total	C	N	O	Se	0	2	0
			2739	1756	440	532	11			
1	F	348	Total	C	N	O	Se	0	1	0
			2745	1759	446	529	11			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	340	SER	-	expression tag	UNP B8FWJ2
A	341	ASN	-	expression tag	UNP B8FWJ2
A	342	ALA	-	expression tag	UNP B8FWJ2
B	340	SER	-	expression tag	UNP B8FWJ2
B	341	ASN	-	expression tag	UNP B8FWJ2
B	342	ALA	-	expression tag	UNP B8FWJ2
C	340	SER	-	expression tag	UNP B8FWJ2
C	341	ASN	-	expression tag	UNP B8FWJ2
C	342	ALA	-	expression tag	UNP B8FWJ2
D	340	SER	-	expression tag	UNP B8FWJ2
D	341	ASN	-	expression tag	UNP B8FWJ2
D	342	ALA	-	expression tag	UNP B8FWJ2
E	340	SER	-	expression tag	UNP B8FWJ2
E	341	ASN	-	expression tag	UNP B8FWJ2
E	342	ALA	-	expression tag	UNP B8FWJ2
F	340	SER	-	expression tag	UNP B8FWJ2
F	341	ASN	-	expression tag	UNP B8FWJ2

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Chain	Residue	Modelled	Actual	Comment	Reference
F	342	ALA	-	expression tag	UNP B8FWJ2

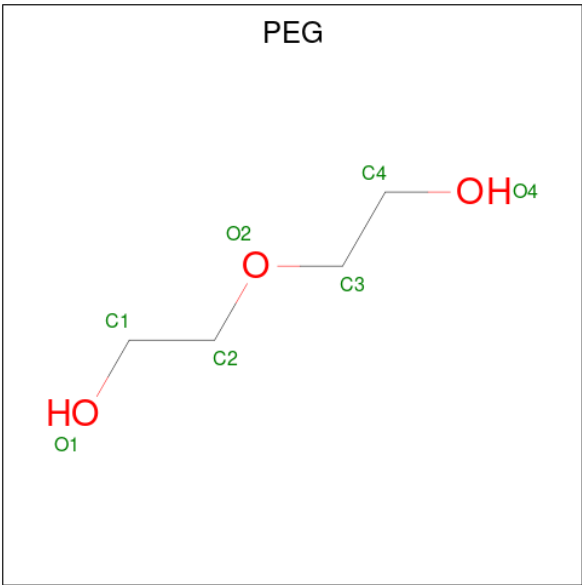
- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	4	Total Ca 4 4	0	0
2	B	1	Total Ca 1 1	0	0
2	C	1	Total Ca 1 1	0	0
2	D	2	Total Ca 2 2	0	0
2	E	1	Total Ca 1 1	0	0
2	F	1	Total Ca 1 1	0	0

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Cl 1 1	0	0
3	F	1	Total Cl 1 1	0	0

- Molecule 4 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



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

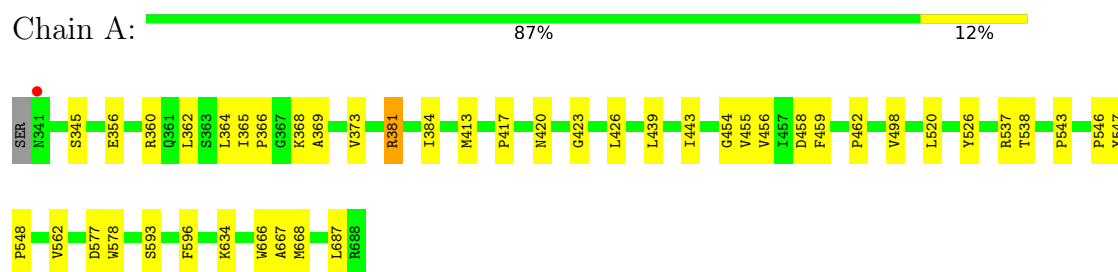
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	54	Total	O	0	0
			54	54		
5	B	59	Total	O	0	0
			59	59		
5	C	23	Total	O	0	0
			23	23		
5	D	32	Total	O	0	0
			32	32		
5	E	27	Total	O	0	0
			27	27		
5	F	100	Total	O	0	0
			100	100		

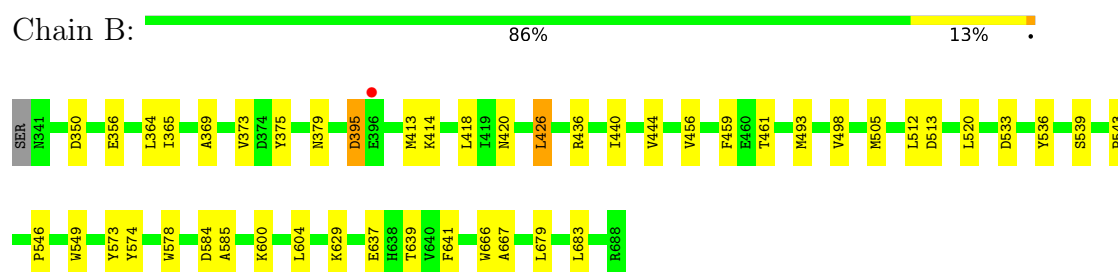
3 Residue-property plots

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.

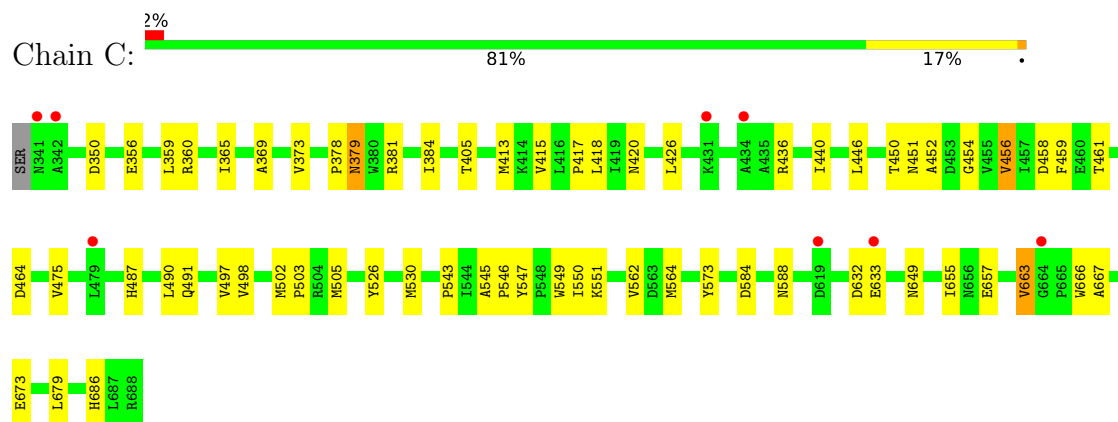
- Molecule 1: Glycoside hydrolase family 18



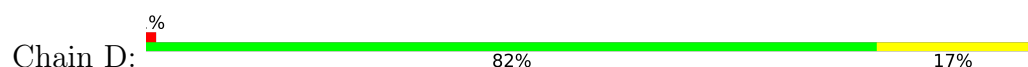
- Molecule 1: Glycoside hydrolase family 18

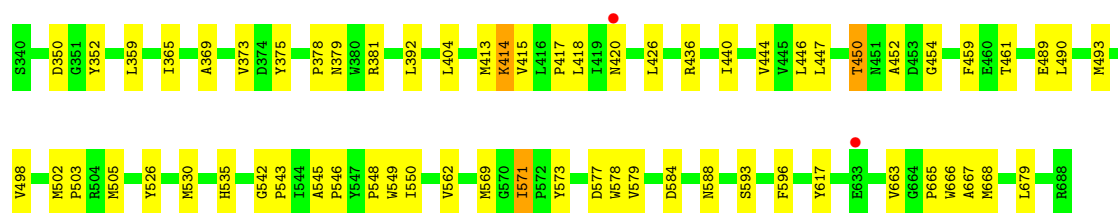


- Molecule 1: Glycoside hydrolase family 18

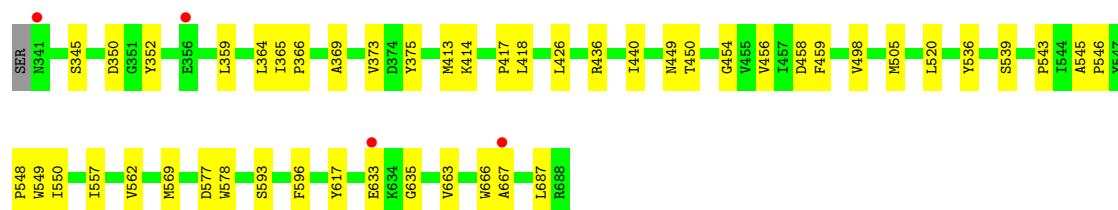
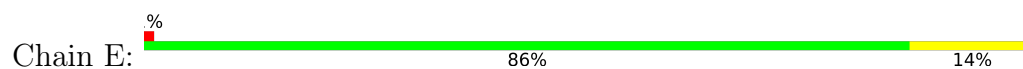


- Molecule 1: Glycoside hydrolase family 18

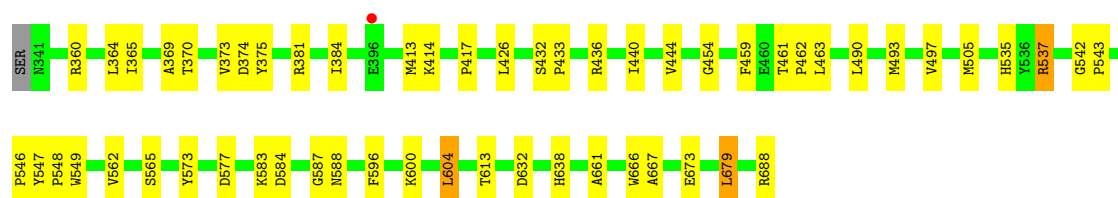
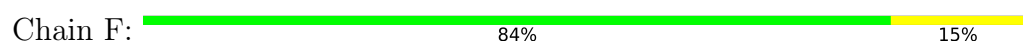




● Molecule 1: Glycoside hydrolase family 18



● Molecule 1: Glycoside hydrolase family 18



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	79.93Å 346.62Å 193.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.14 – 2.64 39.14 – 2.64	Depositor EDS
% Data completeness (in resolution range)	92.0 (39.14-2.64) 91.9 (39.14-2.64)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.19 (at 2.65Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.3_1479)	Depositor
R, R_{free}	0.184 , 0.236 0.192 , 0.238	Depositor DCC
R_{free} test set	3641 reflections (4.54%)	wwPDB-VP
Wilson B-factor (Å ²)	39.3	Xtriage
Anisotropy	0.242	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 34.1	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	16700	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.15% 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, CL, CA

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	0/2797	0.69	0/3797
1	B	0.29	0/2799	0.68	0/3800
1	C	0.27	0/2785	0.70	0/3783
1	D	0.28	0/2781	0.71	1/3781 (0.0%)
1	E	0.27	0/2804	0.70	1/3809 (0.0%)
1	F	0.29	0/2810	0.68	0/3814
All	All	0.28	0/16776	0.69	2/22784 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	352	TYR	CB-CA-C	-5.09	110.31	117.23
1	D	352	TYR	CB-CA-C	-5.06	110.34	117.23

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	2732	0	2618	27	0
1	B	2734	0	2622	26	0
1	C	2720	0	2601	43	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2716	0	2582	40	0
1	E	2739	0	2608	26	0
1	F	2745	0	2634	31	0
2	A	4	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	2	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	1	0	0	0	0
3	F	1	0	0	1	0
4	D	7	0	10	0	0
5	A	54	0	0	0	0
5	B	59	0	0	1	0
5	C	23	0	0	0	0
5	D	32	0	0	0	0
5	E	27	0	0	0	0
5	F	100	0	0	0	0
All	All	16700	0	15675	189	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 (189) 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:369:ALA:HB1	1:B:413:MSE:HE3	1.41	1.02
1:A:369:ALA:HB1	1:A:413:MSE:HE3	1.50	0.91
1:C:369:ALA:HB1	1:C:413:MSE:HE3	1.51	0.91
1:F:369:ALA:HB1	1:F:413:MSE:HE3	1.60	0.81
1:E:369:ALA:HB1	1:E:413:MSE:HE3	1.61	0.81
1:C:365:ILE:HD13	1:C:413:MSE:HE1	1.63	0.81
1:A:373:VAL:HB	1:A:413:MSE:HE2	1.68	0.75
1:B:373:VAL:HB	1:B:413:MSE:HE2	1.67	0.74
1:B:365:ILE:HD13	1:B:413:MSE:HE1	1.69	0.73
1:F:373:VAL:HB	1:F:413:MSE:HE2	1.69	0.73
1:D:373:VAL:HB	1:D:413:MSE:HE2	1.72	0.71
1:D:545:ALA:HB1	1:D:550:ILE:HD11	1.73	0.71
1:C:543:PRO:HB2	1:C:546:PRO:HG3	1.74	0.70
1:B:456:VAL:HG12	1:B:498:VAL:HB	1.72	0.69
1:B:356:GLU:OE1	1:F:360[A]:ARG:NH2	2.25	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:543:PRO:HB2	1:D:546:PRO:HG3	1.74	0.69
1:B:543:PRO:HB2	1:B:546:PRO:HG3	1.76	0.68
1:C:545:ALA:HB1	1:C:550:ILE:HD11	1.76	0.67
1:D:350:ASP:HA	1:D:379:ASN:HD21	1.57	0.67
1:D:369:ALA:HB1	1:D:413:MSE:HE3	1.77	0.66
1:E:545:ALA:HB1	1:E:550:ILE:HD11	1.77	0.66
1:E:373:VAL:HB	1:E:413:MSE:HE2	1.76	0.66
1:E:543:PRO:HB2	1:E:546:PRO:HG3	1.76	0.66
1:F:490:LEU:HA	1:F:493:MSE:HE3	1.78	0.66
1:C:450:THR:HG23	1:C:452:ALA:H	1.62	0.65
1:C:373:VAL:HB	1:C:413:MSE:HE2	1.78	0.65
1:C:498:VAL:HG22	1:C:526:TYR:HB2	1.81	0.63
1:D:450:THR:HG23	1:D:452:ALA:H	1.63	0.62
1:D:444:VAL:HG22	1:D:493:MSE:HE1	1.83	0.61
1:C:418:LEU:HA	1:C:456:VAL:HG23	1.82	0.61
1:F:370:THR:O	1:F:688:ARG:NH2	2.34	0.60
1:F:374:ASP:OD2	1:F:688:ARG:NH1	2.33	0.60
1:A:543:PRO:HB2	1:A:546:PRO:HG3	1.84	0.60
1:C:381:ARG:HD3	1:C:384:ILE:HG13	1.84	0.60
1:D:350:ASP:HA	1:D:379:ASN:ND2	2.17	0.59
1:A:381:ARG:HD3	1:A:384:ILE:HG13	1.86	0.58
1:A:365:ILE:HD13	1:A:413:MSE:HE1	1.86	0.58
1:D:505:MSE:HE2	1:D:549:TRP:HE1	1.68	0.57
1:E:456:VAL:HG22	1:E:498:VAL:HB	1.87	0.57
1:F:543:PRO:HB2	1:F:546:PRO:HG3	1.87	0.57
1:A:356:GLU:OE2	1:C:360:ARG:NH2	2.37	0.57
1:F:444:VAL:HG22	1:F:493:MSE:HE1	1.87	0.57
1:D:378:PRO:HD2	1:D:417:PRO:HA	1.86	0.56
1:D:415:VAL:HG12	1:D:417:PRO:HD3	1.89	0.55
1:B:629:LYS:HG2	1:B:639:THR:HG22	1.89	0.55
1:F:381:ARG:HD3	1:F:384:ILE:HG12	1.88	0.55
1:F:505:MSE:HE2	1:F:549:TRP:HE1	1.72	0.54
1:C:420:ASN:HB3	1:C:458:ASP:HB3	1.90	0.54
1:B:513:ASP:OD2	5:B:853:HOH:O	2.18	0.54
1:B:418:LEU:HA	1:B:456:VAL:HG23	1.89	0.54
1:E:426:LEU:HD21	1:E:459:PHE:CG	2.44	0.53
1:E:418:LEU:HD11	1:E:458:ASP:HB2	1.90	0.53
1:C:632:ASP:OD1	1:C:633:GLU:N	2.41	0.53
1:C:379:ASN:HB3	1:C:418:LEU:HD23	1.91	0.52
1:E:345:SER:HB2	1:E:687:LEU:HD13	1.92	0.52
1:F:365:ILE:HD13	1:F:413:MSE:HE1	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:503:PRO:HG3	1:D:530:MSE:O	2.10	0.51
1:A:423:GLY:HA2	1:A:462:PRO:HG2	1.91	0.51
1:C:551:LYS:NZ	1:C:657:GLU:OE1	2.42	0.51
1:F:577:ASP:HB2	1:F:596:PHE:HE2	1.75	0.51
1:D:461:THR:HG23	1:D:502:MSE:HG3	1.93	0.51
1:D:571:ILE:HG22	1:D:665:PRO:HB3	1.93	0.50
1:B:584:ASP:OD1	1:B:585:ALA:N	2.43	0.50
1:D:577:ASP:HB2	1:D:596:PHE:HE2	1.76	0.50
1:C:490:LEU:HD22	1:C:497:VAL:HG22	1.93	0.50
1:F:375:TYR:HD1	1:F:414:LYS:HB2	1.76	0.50
1:D:379:ASN:HB3	1:D:418:LEU:HD23	1.93	0.49
1:F:436:ARG:O	1:F:440:ILE:HG12	2.11	0.49
1:C:405:THR:HG22	1:C:415:VAL:HG11	1.92	0.49
1:E:505:MSE:HE2	1:E:549:TRP:HE1	1.76	0.49
1:F:583:LYS:HE3	1:F:587:GLY:HA2	1.95	0.49
1:F:426:LEU:HD21	1:F:459:PHE:CG	2.48	0.49
1:A:577:ASP:HB2	1:A:596:PHE:CE2	2.47	0.49
1:B:600:LYS:HE3	1:B:604:LEU:HD13	1.94	0.49
1:C:505:MSE:HE2	1:C:549:TRP:HE1	1.76	0.49
1:C:649:ASN:OD1	1:C:686:HIS:NE2	2.30	0.48
1:B:395:ASP:N	1:B:395:ASP:OD1	2.44	0.48
1:E:569:MSE:HG3	1:E:663:VAL:HG23	1.95	0.48
1:C:655:ILE:HD11	1:C:663:VAL:HG22	1.95	0.48
1:B:505:MSE:HE2	1:B:549:TRP:HE1	1.78	0.48
1:C:405:THR:HA	1:C:415:VAL:HG21	1.95	0.48
1:D:436:ARG:O	1:D:440:ILE:HG12	2.14	0.47
1:A:426:LEU:HD21	1:A:459:PHE:CG	2.49	0.47
1:B:375:TYR:HE1	1:B:414:LYS:HE2	1.78	0.47
1:C:446:LEU:O	1:C:450:THR:HG22	2.14	0.47
1:A:498:VAL:HG22	1:A:526:TYR:HB2	1.95	0.47
1:C:418:LEU:HD11	1:C:458:ASP:HB2	1.97	0.47
1:D:381:ARG:HH11	1:D:420:ASN:HB2	1.79	0.47
1:C:379:ASN:HA	1:C:418:LEU:HB3	1.97	0.47
1:C:456:VAL:HG12	1:C:498:VAL:HB	1.96	0.47
1:E:436:ARG:O	1:E:440:ILE:HG12	2.15	0.47
1:D:381:ARG:HE	1:D:420:ASN:HB2	1.80	0.47
1:D:584:ASP:HB2	1:D:588:ASN:H	1.80	0.47
1:C:666:TRP:CD2	1:C:667:ALA:HB2	2.51	0.46
1:C:546:PRO:O	1:C:550:ILE:HG12	2.15	0.46
1:C:461:THR:HG23	1:C:502:MSE:HG3	1.97	0.46
1:D:365:ILE:HD11	1:D:668:MSE:HE1	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:345:SER:HB2	1:A:687:LEU:HD13	1.97	0.46
1:D:426:LEU:HD21	1:D:459:PHE:CG	2.50	0.46
1:E:577:ASP:HB2	1:E:596:PHE:HE2	1.81	0.46
1:C:350:ASP:HA	1:C:379:ASN:OD1	2.15	0.46
1:D:578:TRP:CD1	1:D:593:SER:HB3	2.51	0.46
1:E:548:PRO:HG2	1:E:617:TYR:HB3	1.98	0.46
1:C:378:PRO:HD2	1:C:417:PRO:HA	1.98	0.45
1:C:436:ARG:O	1:C:440:ILE:HG12	2.17	0.45
1:F:632:ASP:OD2	1:F:638:HIS:NE2	2.42	0.45
1:D:573:TYR:CE1	1:D:679:LEU:HD11	2.51	0.45
1:A:417:PRO:HD2	1:A:454:GLY:O	2.17	0.45
1:A:364:LEU:HD12	1:A:364:LEU:HA	1.82	0.45
1:D:666:TRP:HA	1:D:667:ALA:HA	1.79	0.45
1:F:600:LYS:HE3	1:F:604:LEU:HD13	1.99	0.45
1:F:537:ARG:HG3	3:F:702:CL:CL	2.54	0.45
1:B:436:ARG:O	1:B:440:ILE:HG12	2.17	0.45
1:D:359:LEU:HB3	1:E:359:LEU:HD13	1.97	0.45
1:E:375:TYR:HE1	1:E:414:LYS:HE3	1.83	0.45
1:F:547:TYR:HB3	1:F:548:PRO:HD3	1.99	0.45
1:D:666:TRP:CD2	1:D:667:ALA:HB2	2.52	0.44
1:E:633:GLU:C	1:E:635:GLY:H	2.24	0.44
1:A:443:ILE:HG23	1:A:455:VAL:HG21	1.99	0.44
1:C:503:PRO:HG3	1:C:530:MSE:O	2.16	0.44
1:D:440:ILE:O	1:D:444:VAL:HG23	2.17	0.44
1:E:578:TRP:CD1	1:E:593:SER:HB3	2.52	0.44
1:B:444:VAL:HG22	1:B:493:MSE:HE1	1.99	0.44
1:A:547:TYR:HB3	1:A:548:PRO:HD3	1.99	0.44
1:E:417:PRO:HD2	1:E:454:GLY:O	2.17	0.44
1:E:557:ILE:HG22	1:E:562:VAL:HG21	1.99	0.44
1:A:537:ARG:HD2	1:A:538:THR:HG23	2.00	0.44
1:B:461:THR:HG21	1:B:512:LEU:HD23	2.00	0.44
1:B:426:LEU:HD21	1:B:459:PHE:CD1	2.52	0.43
1:C:464:ASP:HB3	1:C:475:VAL:HG13	1.99	0.43
1:D:666:TRP:CG	1:D:667:ALA:HB2	2.53	0.43
1:D:392:LEU:HD11	1:D:446:LEU:HD13	1.99	0.43
1:D:548:PRO:HG2	1:D:617:TYR:HB3	2.00	0.43
1:E:666:TRP:CD2	1:E:667:ALA:HB2	2.52	0.43
1:F:432:SER:HA	1:F:433:PRO:HD2	1.91	0.43
1:F:535:HIS:CE1	1:F:542:GLY:HA3	2.54	0.43
1:A:666:TRP:HA	1:A:667:ALA:HA	1.85	0.43
1:D:490:LEU:HA	1:D:493:MSE:HE3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:584:ASP:OD1	1:F:588:ASN:N	2.51	0.43
1:A:360:ARG:HH21	1:C:356:GLU:HG2	1.83	0.43
1:A:634:LYS:HB3	1:A:634:LYS:HE2	1.78	0.43
1:E:546:PRO:O	1:E:550:ILE:HG12	2.18	0.43
1:A:365:ILE:HD11	1:A:668:MSE:SE	2.68	0.43
1:C:547:TYR:CZ	1:C:551:LYS:HD2	2.54	0.43
1:A:578:TRP:CD1	1:A:593:SER:HB3	2.54	0.42
1:A:577:ASP:HB2	1:A:596:PHE:HE2	1.84	0.42
1:F:577:ASP:HB2	1:F:596:PHE:CE2	2.54	0.42
1:B:666:TRP:CD2	1:B:667:ALA:HB2	2.54	0.42
1:C:417:PRO:HD2	1:C:454:GLY:O	2.19	0.42
1:C:487:HIS:O	1:C:491:GLN:HB2	2.19	0.42
1:F:565:SER:O	1:F:661:ALA:HB2	2.20	0.42
1:A:439:LEU:O	1:A:443:ILE:HG13	2.19	0.42
1:C:456:VAL:HA	1:C:498:VAL:O	2.19	0.42
1:F:490:LEU:HD22	1:F:497:VAL:HG22	2.01	0.42
1:C:505:MSE:HE2	1:C:549:TRP:NE1	2.34	0.42
1:F:573:TYR:CE1	1:F:679:LEU:HD11	2.54	0.42
1:D:447:LEU:HD22	1:D:452:ALA:HB3	2.01	0.42
1:F:673:GLU:H	1:F:673:GLU:CD	2.27	0.42
1:D:498:VAL:HG22	1:D:526:TYR:HB2	2.02	0.42
1:A:365:ILE:HA	1:A:366:PRO:HD3	1.79	0.41
1:B:536:TYR:CZ	1:B:539:SER:HB3	2.55	0.41
1:D:417:PRO:HD2	1:D:454:GLY:O	2.19	0.41
1:C:666:TRP:HA	1:C:667:ALA:HA	1.77	0.41
1:D:502:MSE:HE2	1:D:502:MSE:HA	2.02	0.41
1:D:535:HIS:CE1	1:D:542:GLY:HA3	2.55	0.41
1:F:666:TRP:CD2	1:F:667:ALA:HB2	2.55	0.41
1:B:578:TRP:HZ3	1:B:641:PHE:HB2	1.85	0.41
1:A:456:VAL:HG22	1:A:498:VAL:HB	2.01	0.41
1:C:573:TYR:CE1	1:C:679:LEU:HD11	2.56	0.41
1:E:365:ILE:HA	1:E:366:PRO:HD3	1.83	0.41
1:A:362:LEU:HD23	1:A:362:LEU:HA	1.94	0.41
1:B:350:ASP:OD1	1:B:350:ASP:N	2.53	0.41
1:C:426:LEU:HD21	1:C:459:PHE:CG	2.55	0.41
1:F:461:THR:OG1	1:F:462:PRO:HA	2.21	0.41
1:B:573:TYR:CE1	1:B:679:LEU:HD11	2.56	0.41
1:C:584:ASP:N	1:C:588:ASN:O	2.36	0.41
1:C:673:GLU:H	1:C:673:GLU:CD	2.28	0.41
1:D:375:TYR:CD1	1:D:414:LYS:HB2	2.56	0.41
1:E:350:ASP:OD1	1:E:350:ASP:N	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:449:ASN:O	1:E:449:ASN:ND2	2.54	0.41
1:E:577:ASP:HB2	1:E:596:PHE:CE2	2.56	0.41
1:F:417:PRO:HD2	1:F:454:GLY:O	2.20	0.41
1:E:536:TYR:CZ	1:E:539:SER:HB3	2.56	0.41
1:D:489:GLU:HG3	1:D:493:MSE:HE2	2.03	0.40
1:D:569:MSE:HG3	1:D:663:VAL:HG23	2.03	0.40
1:B:379:ASN:HA	1:B:418:LEU:HB3	2.02	0.40
1:B:533:ASP:OD1	1:B:574:TYR:OH	2.33	0.40
1:A:420:ASN:HB3	1:A:458:ASP:HB3	2.02	0.40
1:B:420:ASN:OD1	1:B:420:ASN:N	2.53	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	346/349 (99%)	332 (96%)	14 (4%)	0	100	100
1	B	346/349 (99%)	336 (97%)	10 (3%)	0	100	100
1	C	346/349 (99%)	330 (95%)	16 (5%)	0	100	100
1	D	347/349 (99%)	333 (96%)	14 (4%)	0	100	100
1	E	348/349 (100%)	330 (95%)	18 (5%)	0	100	100
1	F	347/349 (99%)	335 (96%)	12 (4%)	0	100	100
All	All	2080/2094 (99%)	1996 (96%)	84 (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	291/283 (103%)	287 (99%)	4 (1%)	59	75
1	B	292/283 (103%)	286 (98%)	6 (2%)	47	67
1	C	288/283 (102%)	281 (98%)	7 (2%)	43	64
1	D	287/283 (101%)	281 (98%)	6 (2%)	47	67
1	E	291/283 (103%)	288 (99%)	3 (1%)	68	80
1	F	293/283 (104%)	286 (98%)	7 (2%)	43	64
All	All	1742/1698 (103%)	1709 (98%)	33 (2%)	50	70

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	368	LYS
1	A	381	ARG
1	A	520	LEU
1	A	562	VAL
1	B	364	LEU
1	B	395	ASP
1	B	426	LEU
1	B	520	LEU
1	B	637	GLU
1	B	683	LEU
1	C	359	LEU
1	C	379	ASN
1	C	451	ASN
1	C	456	VAL
1	C	562	VAL
1	C	564	MSE
1	C	663	VAL
1	D	404	LEU
1	D	414	LYS
1	D	450	THR
1	D	562	VAL
1	D	571	ILE

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Mol	Chain	Res	Type
1	D	579	VAL
1	E	364	LEU
1	E	450	THR
1	E	520	LEU
1	F	364	LEU
1	F	463	LEU
1	F	537	ARG
1	F	562	VAL
1	F	604	LEU
1	F	613	THR
1	F	679	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	403	GLN
1	C	535	HIS
1	C	560	GLN
1	D	379	ASN
1	D	684	ASN
1	F	406	ASN

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](#)

Of 13 ligands modelled in this entry, 12 are monoatomic - leaving 1 for Mogul analysis.

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	$\# Z > 2$
4	PEG	D	703	-	6,6,6	0.46	0	5,5,5	0.30	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
4	PEG	D	703	-	-	0/4/4/4	-

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	337/349 (96%)	-0.32	1 (0%) 90 89	23, 39, 62, 91	0
1	B	337/349 (96%)	-0.41	1 (0%) 90 89	22, 35, 61, 111	0
1	C	337/349 (96%)	0.25	8 (2%) 59 55	43, 61, 81, 103	0
1	D	338/349 (96%)	0.12	2 (0%) 85 83	31, 54, 78, 120	0
1	E	337/349 (96%)	0.24	4 (1%) 76 73	33, 59, 83, 103	2 (0%)
1	F	337/349 (96%)	-0.63	1 (0%) 90 89	14, 27, 46, 69	1 (0%)
All	All	2023/2094 (96%)	-0.12	17 (0%) 82 80	14, 48, 76, 120	3 (0%)

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	341	ASN	4.2
1	A	341	ASN	3.6
1	C	341	ASN	3.3
1	C	619	ASP	3.0
1	E	356[A]	GLU	2.5
1	D	633	GLU	2.4
1	C	342	ALA	2.4
1	C	633	GLU	2.4
1	E	667	ALA	2.4
1	D	420	ASN	2.3
1	E	633	GLU	2.2
1	C	434	ALA	2.2
1	B	396	GLU	2.2
1	C	431	LYS	2.2
1	F	396	GLU	2.1
1	C	479	LEU	2.1
1	C	664	GLY	2.1

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($\text{\AA}^2$)	Q<0.9
4	PEG	D	703	7/7	0.85	0.14	37,45,62,62	0
2	CA	A	704	1/1	0.91	0.16	66,66,66,66	0
2	CA	D	702	1/1	0.94	0.05	59,59,59,59	1
2	CA	E	701	1/1	0.94	0.06	57,57,57,57	0
2	CA	D	701	1/1	0.94	0.23	116,116,116,116	0
2	CA	A	701	1/1	0.95	0.19	59,59,59,59	0
2	CA	C	701	1/1	0.95	0.06	47,47,47,47	0
2	CA	A	703	1/1	0.96	0.14	57,57,57,57	0
2	CA	B	701	1/1	0.98	0.06	58,58,58,58	0
2	CA	F	701	1/1	0.99	0.03	40,40,40,40	0
3	CL	A	705	1/1	0.99	0.03	37,37,37,37	0
3	CL	F	702	1/1	0.99	0.07	39,39,39,39	0
2	CA	A	702	1/1	0.99	0.04	30,30,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.