



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2026 – 06:59 PM UTC

PDB ID : 5E2J / pdb_00005e2j
Title : Crystal structure of single mutant thermostable endoglucanase (D468A) from Alicyclobacillus acidocaldarius
Authors : Hsiao, Y.Y.; Wang, H.J.; Tseng, C.P.
Deposited on : 2015-10-01
Resolution : 2.10 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

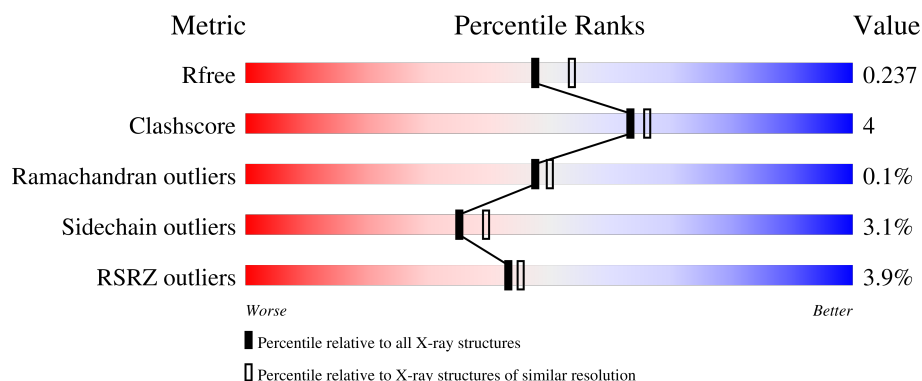
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

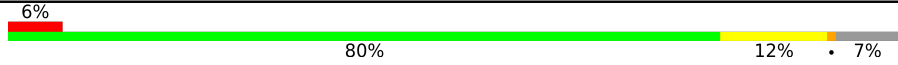
The reported resolution of this entry is 2.10 Å.

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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	570	
1	B	570	

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
4	MPD	A	604	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 8617 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 Endoglucanase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	530	Total	C	N	O	S	0	0	0
			4101	2617	723	738	23			
1	B	529	Total	C	N	O	S	0	0	0
			4090	2611	719	737	23			

There are 68 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MET	-	expression tag	UNP Q9AJS0
A	-21	ALA	-	expression tag	UNP Q9AJS0
A	-20	SER	-	expression tag	UNP Q9AJS0
A	-19	MET	-	expression tag	UNP Q9AJS0
A	-18	THR	-	expression tag	UNP Q9AJS0
A	-17	GLY	-	expression tag	UNP Q9AJS0
A	-16	GLY	-	expression tag	UNP Q9AJS0
A	-15	GLN	-	expression tag	UNP Q9AJS0
A	-14	GLN	-	expression tag	UNP Q9AJS0
A	-13	MET	-	expression tag	UNP Q9AJS0
A	-12	GLY	-	expression tag	UNP Q9AJS0
A	-11	ARG	-	expression tag	UNP Q9AJS0
A	-10	GLY	-	expression tag	UNP Q9AJS0
A	-9	SER	-	expression tag	UNP Q9AJS0
A	-8	GLU	-	expression tag	UNP Q9AJS0
A	-7	PHE	-	expression tag	UNP Q9AJS0
A	-6	GLU	-	expression tag	UNP Q9AJS0
A	-5	LEU	-	expression tag	UNP Q9AJS0
A	-4	ARG	-	expression tag	UNP Q9AJS0
A	-3	ARG	-	expression tag	UNP Q9AJS0
A	-2	ARG	-	expression tag	UNP Q9AJS0
A	-1	PHE	-	expression tag	UNP Q9AJS0
A	0	ASP	-	expression tag	UNP Q9AJS0
A	468	ALA	ASP	engineered mutation	UNP Q9AJS0
A	538	ALA	-	expression tag	UNP Q9AJS0

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Chain	Residue	Modelled	Actual	Comment	Reference
A	539	ALA	-	expression tag	UNP Q9AJS0
A	540	LEU	-	expression tag	UNP Q9AJS0
A	541	GLU	-	expression tag	UNP Q9AJS0
A	542	HIS	-	expression tag	UNP Q9AJS0
A	543	HIS	-	expression tag	UNP Q9AJS0
A	544	HIS	-	expression tag	UNP Q9AJS0
A	545	HIS	-	expression tag	UNP Q9AJS0
A	546	HIS	-	expression tag	UNP Q9AJS0
A	547	HIS	-	expression tag	UNP Q9AJS0
B	-22	MET	-	expression tag	UNP Q9AJS0
B	-21	ALA	-	expression tag	UNP Q9AJS0
B	-20	SER	-	expression tag	UNP Q9AJS0
B	-19	MET	-	expression tag	UNP Q9AJS0
B	-18	THR	-	expression tag	UNP Q9AJS0
B	-17	GLY	-	expression tag	UNP Q9AJS0
B	-16	GLY	-	expression tag	UNP Q9AJS0
B	-15	GLN	-	expression tag	UNP Q9AJS0
B	-14	GLN	-	expression tag	UNP Q9AJS0
B	-13	MET	-	expression tag	UNP Q9AJS0
B	-12	GLY	-	expression tag	UNP Q9AJS0
B	-11	ARG	-	expression tag	UNP Q9AJS0
B	-10	GLY	-	expression tag	UNP Q9AJS0
B	-9	SER	-	expression tag	UNP Q9AJS0
B	-8	GLU	-	expression tag	UNP Q9AJS0
B	-7	PHE	-	expression tag	UNP Q9AJS0
B	-6	GLU	-	expression tag	UNP Q9AJS0
B	-5	LEU	-	expression tag	UNP Q9AJS0
B	-4	ARG	-	expression tag	UNP Q9AJS0
B	-3	ARG	-	expression tag	UNP Q9AJS0
B	-2	ARG	-	expression tag	UNP Q9AJS0
B	-1	PHE	-	expression tag	UNP Q9AJS0
B	0	ASP	-	expression tag	UNP Q9AJS0
B	468	ALA	ASP	engineered mutation	UNP Q9AJS0
B	538	ALA	-	expression tag	UNP Q9AJS0
B	539	ALA	-	expression tag	UNP Q9AJS0
B	540	LEU	-	expression tag	UNP Q9AJS0
B	541	GLU	-	expression tag	UNP Q9AJS0
B	542	HIS	-	expression tag	UNP Q9AJS0
B	543	HIS	-	expression tag	UNP Q9AJS0
B	544	HIS	-	expression tag	UNP Q9AJS0
B	545	HIS	-	expression tag	UNP Q9AJS0
B	546	HIS	-	expression tag	UNP Q9AJS0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	547	HIS	-	expression tag	UNP Q9AJS0

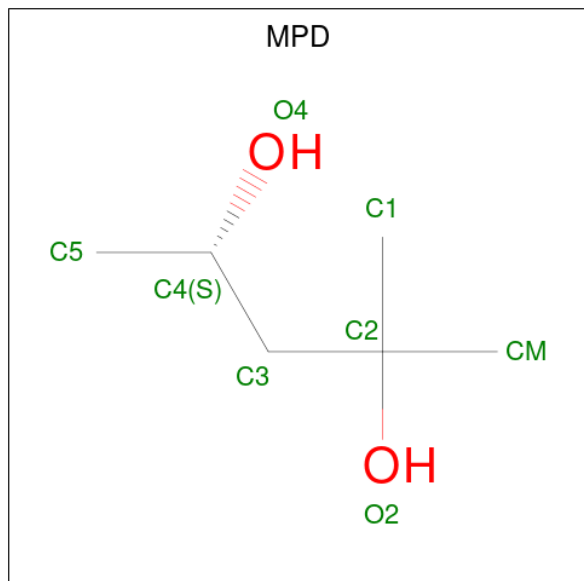
- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Ca 1 1	0	0
2	B	1	Total Ca 1 1	0	0

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Zn 1 1	0	0
3	B	1	Total Zn 1 1	0	0

- Molecule 4 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 8 6 2	0	0
4	A	1	Total C O 8 6 2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			8	6	2		

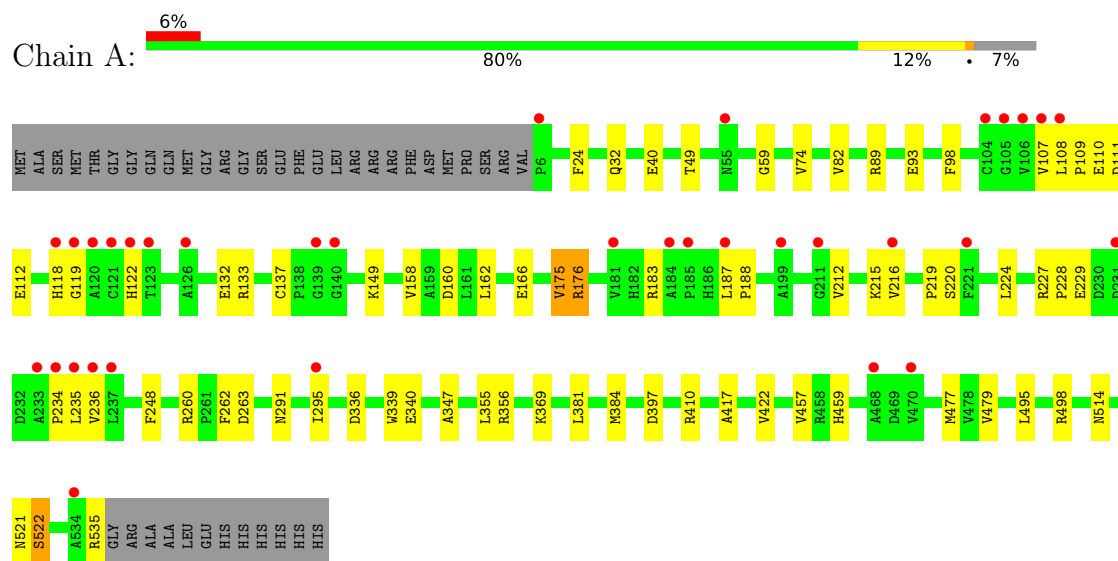
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	206	Total	O	0	0
			206	206		
5	B	192	Total	O	0	0
			192	192		

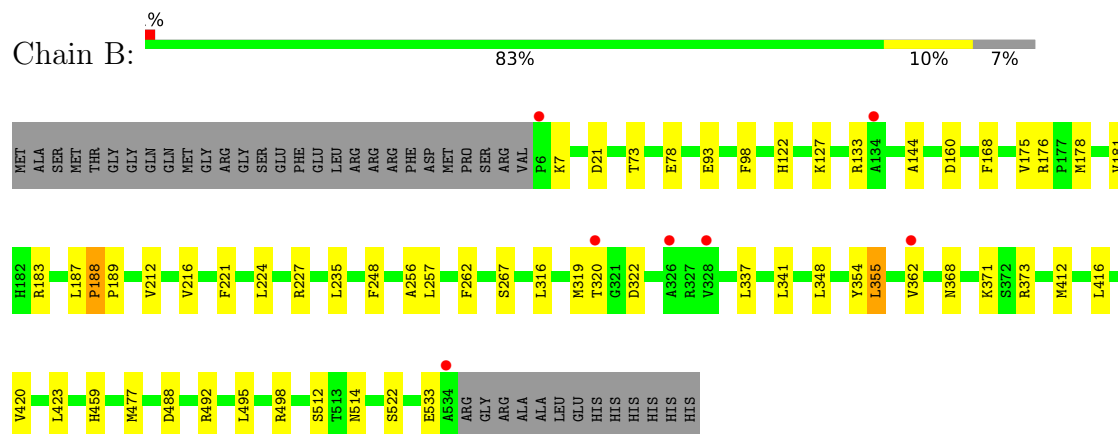
3 Residue-property plots [i](#)

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: Endoglucanase



• Molecule 1: Endoglucanase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	55.42Å 144.37Å 158.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.96 – 2.10 19.96 – 2.10	Depositor EDS
% Data completeness (in resolution range)	97.2 (19.96-2.10) 96.8 (19.96-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 2.09Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.202 , 0.237 0.204 , 0.237	Depositor DCC
R_{free} test set	5675 reflections (7.54%)	wwPDB-VP
Wilson B-factor (Å ²)	41.3	Xtriage
Anisotropy	0.132	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 51.5	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	0.96	EDS
Total number of atoms	8617	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.53% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, MPD, 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	0.29	0/4227	0.73	4/5781 (0.1%)
1	B	0.29	0/4216	0.71	2/5767 (0.0%)
All	All	0.29	0/8443	0.72	6/11548 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	188	PRO	CA-C-N	5.43	126.63	119.84
1	A	188	PRO	C-N-CA	5.43	126.63	119.84
1	B	188	PRO	CA-C-N	5.29	126.46	119.84
1	B	188	PRO	C-N-CA	5.29	126.46	119.84
1	A	522	SER	CA-C-N	-5.03	113.86	119.19
1	A	522	SER	C-N-CA	-5.03	113.86	119.19

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	4101	0	3962	40	0
1	B	4090	0	3949	26	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	16	0	28	6	0
4	B	8	0	14	0	0
5	A	206	0	0	1	0
5	B	192	0	0	2	0
All	All	8617	0	7953	65	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 4.

All (65) 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:381:LEU:HD11	4:A:604:MPD:HM1	1.72	0.72
1:A:176:ARG:NH2	1:B:221:PHE:O	2.23	0.71
1:A:340:GLU:HA	4:A:604:MPD:H12	1.72	0.71
1:A:410:ARG:HG2	4:A:604:MPD:HM3	1.73	0.69
1:B:168:PHE:HB3	1:B:533:GLU:HG3	1.78	0.65
1:B:73:THR:HG22	1:B:78:GLU:HG2	1.79	0.64
1:B:320:THR:HG23	1:B:322:ASP:H	1.62	0.64
1:A:122:HIS:HB3	1:A:235:LEU:HD11	1.78	0.64
1:A:187:LEU:HD22	1:A:262:PHE:HB3	1.81	0.62
1:A:108:LEU:N	1:A:118:HIS:O	2.35	0.59
1:A:339:TRP:O	4:A:604:MPD:H31	2.01	0.59
1:A:118:HIS:ND1	1:A:119:GLY:O	2.34	0.59
1:A:89:ARG:HH22	1:A:175:VAL:HA	1.67	0.59
1:B:183:ARG:NH1	5:B:703:HOH:O	2.38	0.56
1:A:535:ARG:NH2	5:A:709:HOH:O	2.39	0.55
1:B:495:LEU:HA	1:B:498:ARG:HD2	1.88	0.54
1:A:339:TRP:HA	1:A:347:ALA:HB3	1.89	0.53
1:A:160:ASP:HB2	1:A:522:SER:HB2	1.91	0.52
1:A:89:ARG:NH2	1:A:175:VAL:HA	2.24	0.52
1:B:160:ASP:HB2	1:B:522:SER:HB2	1.91	0.52
1:A:495:LEU:HA	1:A:498:ARG:HD2	1.92	0.51
1:B:187:LEU:HD22	1:B:262:PHE:HB3	1.92	0.51
1:A:118:HIS:HB2	1:A:119:GLY:HA3	1.91	0.50
1:A:459:HIS:O	1:A:514:ASN:ND2	2.44	0.50
1:B:122:HIS:HB3	1:B:235:LEU:HD11	1.92	0.50
1:A:183:ARG:NH2	1:A:263:ASP:OD2	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:PHE:CE2	1:A:59:GLY:HA3	2.49	0.48
1:B:257:LEU:HD21	1:B:319:MET:HA	1.95	0.48
1:A:397:ASP:OD1	1:A:397:ASP:N	2.47	0.47
1:A:212:VAL:HG21	1:A:248:PHE:CG	2.50	0.47
1:B:98:PHE:CD1	1:B:477:MET:HG2	2.50	0.47
1:B:492:ARG:NH1	5:B:714:HOH:O	2.48	0.47
1:B:256:ALA:O	1:B:267:SER:OG	2.31	0.46
1:B:488:ASP:OD2	1:B:512:SER:OG	2.31	0.46
1:B:337:LEU:O	1:B:373:ARG:NH2	2.44	0.46
1:A:236:VAL:HG11	1:A:295:ILE:HD11	1.98	0.46
1:A:227:ARG:HD3	1:A:228:PRO:HD2	1.98	0.45
1:B:355:LEU:HB3	1:B:420:VAL:HG11	1.99	0.45
1:B:316:LEU:O	1:B:320:THR:HG22	2.17	0.45
1:A:93:GLU:OE2	1:A:176:ARG:NH1	2.49	0.45
1:B:178:MET:O	1:B:181:VAL:HG22	2.16	0.45
1:B:341:LEU:HG	1:B:348:LEU:HD23	1.99	0.45
1:B:459:HIS:O	1:B:514:ASN:ND2	2.51	0.44
1:B:412:MET:HE2	1:B:416:LEU:HD11	2.00	0.44
1:B:368:ASN:HA	1:B:371:LYS:HD3	2.00	0.44
1:A:109:PRO:HG2	1:A:112:GLU:HB2	1.99	0.43
1:A:166:GLU:OE2	1:A:356:ARG:NH2	2.37	0.43
1:A:107:VAL:HG22	1:A:119:GLY:H	1.84	0.43
1:A:137:CYS:O	1:A:149:LYS:NZ	2.51	0.43
1:B:93:GLU:OE1	1:B:176:ARG:HD3	2.19	0.42
1:B:188:PRO:HA	1:B:189:PRO:HD3	1.89	0.42
1:A:410:ARG:HA	4:A:604:MPD:HM3	2.02	0.42
1:A:98:PHE:CD1	1:A:477:MET:HG2	2.54	0.42
1:A:158:VAL:O	1:A:162:LEU:HG	2.19	0.42
1:B:212:VAL:HG21	1:B:248:PHE:CG	2.55	0.41
1:A:109:PRO:HB2	1:A:111:ASP:OD1	2.19	0.41
1:B:354:TYR:CE1	1:B:362:VAL:HG21	2.55	0.41
1:A:32:GLN:H	1:A:49:THR:HG1	1.67	0.41
1:A:410:ARG:HA	4:A:604:MPD:CM	2.51	0.41
1:A:457:VAL:HG11	1:A:479:VAL:HG11	2.02	0.41
1:A:219:PRO:HD3	1:A:234:PRO:O	2.21	0.41
1:A:417:ALA:HB1	1:A:422:VAL:HB	2.03	0.41
1:A:384:MET:HE1	1:A:410:ARG:NH1	2.36	0.40
1:A:369:LYS:HE2	1:A:369:LYS:HB3	1.86	0.40
1:A:215:LYS:NZ	1:A:291:ASN:OD1	2.50	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	528/570 (93%)	510 (97%)	18 (3%)	0	100	100
1	B	527/570 (92%)	512 (97%)	14 (3%)	1 (0%)	43	44
All	All	1055/1140 (92%)	1022 (97%)	32 (3%)	1 (0%)	48	50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	144	ALA

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	414/446 (93%)	398 (96%)	16 (4%)	28	31
1	B	413/446 (93%)	403 (98%)	10 (2%)	43	49
All	All	827/892 (93%)	801 (97%)	26 (3%)	35	39

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	GLU
1	A	74	VAL
1	A	82	VAL
1	A	110	GLU
1	A	132	GLU

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Mol	Chain	Res	Type
1	A	133	ARG
1	A	175	VAL
1	A	176	ARG
1	A	216	VAL
1	A	220	SER
1	A	224	LEU
1	A	229	GLU
1	A	260	ARG
1	A	336	ASP
1	A	355	LEU
1	A	521	ASN
1	B	7	LYS
1	B	21	ASP
1	B	127	LYS
1	B	133	ARG
1	B	175	VAL
1	B	216	VAL
1	B	224	LEU
1	B	227	ARG
1	B	355	LEU
1	B	423	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	GLN
1	A	32	GLN
1	A	432	GLN
1	B	32	GLN
1	B	459	HIS
1	B	485	HIS
1	B	494	GLN

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

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 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	MPD	A	603	-	7,7,7	0.67	0	9,10,10	0.58	0
4	MPD	B	603	-	7,7,7	0.65	0	9,10,10	0.60	0
4	MPD	A	604	-	7,7,7	0.77	0	9,10,10	0.63	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	MPD	A	603	-	-	2/5/5/5	-
4	MPD	B	603	-	-	2/5/5/5	-
4	MPD	A	604	-	-	2/5/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	603	MPD	C2-C3-C4-C5
4	B	603	MPD	C2-C3-C4-C5
4	A	604	MPD	C2-C3-C4-C5
4	A	603	MPD	O2-C2-C3-C4

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Mol	Chain	Res	Type	Atoms
4	B	603	MPD	O2-C2-C3-C4
4	A	604	MPD	C2-C3-C4-O4

There are no ring outliers.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	604	MPD	6	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	530/570 (92%)	0.33	34 (6%) 25 27	26, 48, 93, 117	0
1	B	529/570 (92%)	0.22	7 (1%) 75 77	32, 55, 90, 128	0
All	All	1059/1140 (92%)	0.27	41 (3%) 43 45	26, 51, 91, 128	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	6	PRO	4.8
1	A	139	GLY	4.5
1	B	534	ALA	4.3
1	A	105	GLY	3.6
1	A	233	ALA	3.4
1	A	470	VAL	3.3
1	A	237	LEU	3.2
1	A	236	VAL	3.2
1	A	231	ASP	3.1
1	A	122	HIS	3.0
1	A	468	ALA	2.9
1	A	234	PRO	2.9
1	A	120	ALA	2.8
1	A	534	ALA	2.8
1	A	118	HIS	2.7
1	A	126	ALA	2.7
1	A	221	PHE	2.7
1	A	184	ALA	2.7
1	B	320	THR	2.6
1	A	104	CYS	2.6
1	A	187	LEU	2.6
1	A	107	VAL	2.5
1	A	119	GLY	2.5
1	B	328	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	362	VAL	2.4
1	A	123	THR	2.4
1	A	181	VAL	2.3
1	A	216	VAL	2.3
1	A	199	ALA	2.3
1	A	121	CYS	2.2
1	B	134	ALA	2.2
1	A	185	PRO	2.2
1	A	235	LEU	2.2
1	A	6	PRO	2.1
1	A	140	GLY	2.1
1	A	108	LEU	2.1
1	A	106	VAL	2.1
1	B	326	ALA	2.1
1	A	295	ILE	2.1
1	A	55	ASN	2.0
1	A	211	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](#)

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	MPD	B	603	8/8	0.69	0.22	88,93,94,96	0
4	MPD	A	603	8/8	0.79	0.19	63,67,71,75	0
4	MPD	A	604	8/8	0.81	0.18	69,71,73,77	0
3	ZN	A	602	1/1	0.93	0.07	77,77,77,77	0
2	CA	A	601	1/1	0.97	0.03	41,41,41,41	0
3	ZN	B	602	1/1	0.99	0.03	45,45,45,45	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CA	B	601	1/1	0.99	0.04	55,55,55,55	0

6.5 Other polymers [i](#)

There are no such residues in this entry.