



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:47 AM UTC

PDB ID : 2WKN / pdb_00002wkn
Title : gamma lactamase from Delftia acidovorans
Authors : Isupov, M.N.; Line, K.; Gonsalvez, I.S.; Gange-Harris, P.; Lanzotti, M.; Saneei, V.; Littlechild, J.A.
Deposited on : 2009-06-16
Resolution : 2.08 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

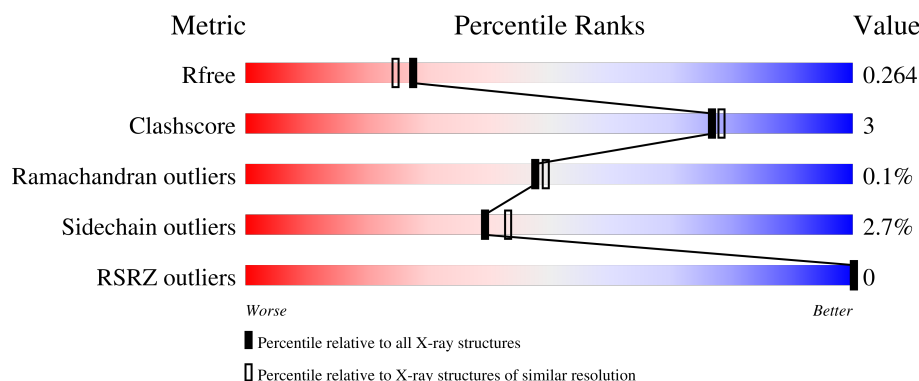
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.08 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	8172 (2.10-2.06)
Clashscore	190562	8714 (2.10-2.06)
Ramachandran outliers	187476	8641 (2.10-2.06)
Sidechain outliers	187428	8642 (2.10-2.06)
RSRZ outliers	180081	8177 (2.10-2.06)

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	409	92% 7% .
1	B	409	89% 9% .
1	C	409	94% 6%
1	D	409	92% 7%
1	E	409	91% 8% .

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Mol	Chain	Length	Quality of chain
1	F	409	 94% 6%
1	G	409	 91% 8% •
1	H	409	 92% 7% •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 26920 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 FORMAMIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	408	Total	C	N	O	S	0	5	0
			3149	2010	531	594	14			
1	B	408	Total	C	N	O	S	0	1	0
			3127	1994	529	590	14			
1	C	408	Total	C	N	O	S	0	1	0
			3130	1996	529	591	14			
1	D	408	Total	C	N	O	S	0	2	0
			3134	2001	530	589	14			
1	E	408	Total	C	N	O	S	0	1	0
			3127	1994	529	590	14			
1	F	408	Total	C	N	O	S	0	1	0
			3129	1995	529	591	14			
1	G	408	Total	C	N	O	S	0	0	0
			3124	1992	529	589	14			
1	H	408	Total	C	N	O	S	0	2	0
			3136	2000	529	593	14			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Zn	0	0
			2	2		
2	B	2	Total	Zn	0	0
			2	2		
2	C	2	Total	Zn	0	0
			2	2		
2	D	2	Total	Zn	0	0
			2	2		
2	E	2	Total	Zn	0	0
			2	2		
2	F	2	Total	Zn	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	2	Total 2	Zn 2	0	0
2	H	2	Total 2	Zn 2	0	0

- Molecule 3 is water.

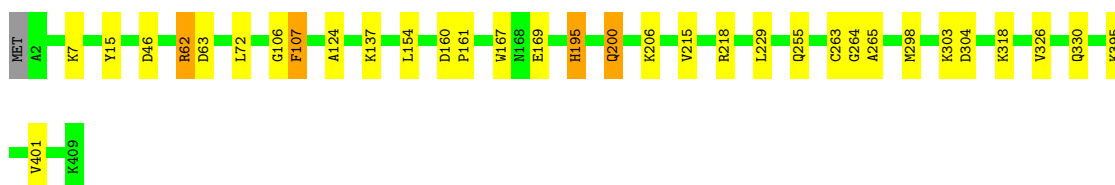
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	332	Total 332	O 332	0	0
3	B	219	Total 219	O 219	0	0
3	C	197	Total 197	O 197	0	0
3	D	269	Total 269	O 269	0	0
3	E	249	Total 249	O 249	0	0
3	F	250	Total 250	O 250	0	0
3	G	166	Total 166	O 166	0	0
3	H	166	Total 166	O 166	0	0

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.

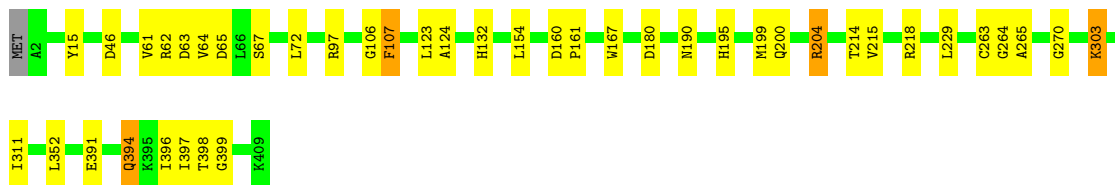
- Molecule 1: FORMAMIDASE

Chain A:  92% 7%

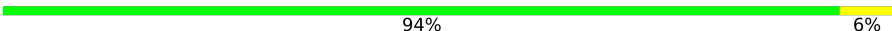


- Molecule 1: FORMAMIDASE

Chain B:  89% 9%



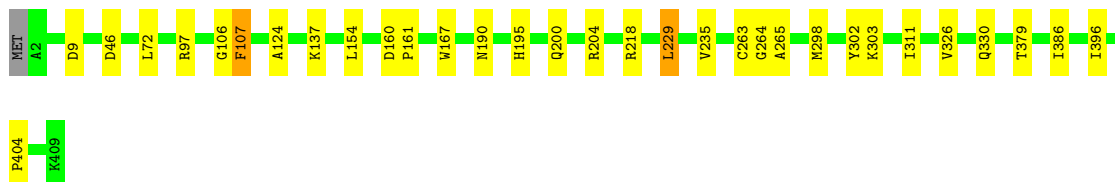
- Molecule 1: FORMAMIDASE

Chain C:  94% 6%



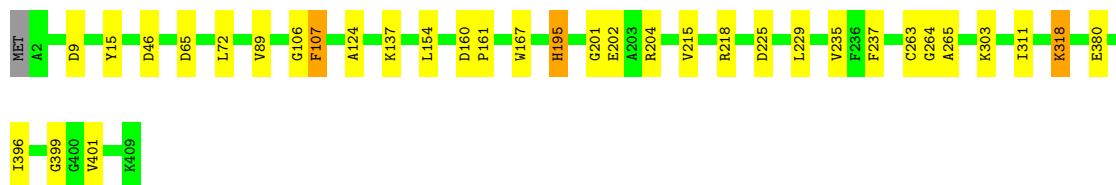
- Molecule 1: FORMAMIDASE

Chain D:  92% 7%

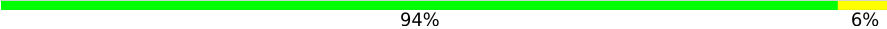


- Molecule 1: FORMAMIDASE

Chain E:  91% 8% .



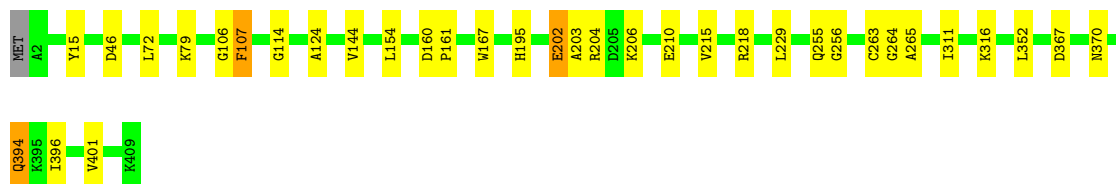
- Molecule 1: FORMAMIDASE

Chain F:  94% 6%



- Molecule 1: FORMAMIDASE

Chain G:  91% 8% .



- Molecule 1: FORMAMIDASE

Chain H:  92% 7% .



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	63.87Å 94.62Å 152.30Å 105.71° 90.27° 108.46°	Depositor
Resolution (Å)	24.78 – 2.08 24.78 – 2.08	Depositor EDS
% Data completeness (in resolution range)	90.8 (24.78-2.08) 79.9 (24.78-2.08)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.31 (at 2.08Å)	Xtriage
Refinement program	REFMAC 5.5.0088	Depositor
R, R_{free}	0.220 , 0.269 0.215 , 0.264	Depositor DCC
R_{free} test set	3032 reflections (1.79%)	wwPDB-VP
Wilson B-factor (Å ²)	32.3	Xtriage
Anisotropy	0.683	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 34.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.179 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	26920	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.25% 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: ZN

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	0.73	1/3251 (0.0%)	0.88	3/4422 (0.1%)
1	B	0.77	0/3217	0.91	5/4376 (0.1%)
1	C	0.75	0/3220	0.88	0/4381
1	D	0.74	0/3227	0.86	0/4389
1	E	0.71	1/3217 (0.0%)	0.85	0/4376
1	F	0.70	0/3219	0.88	0/4379
1	G	0.74	1/3211 (0.0%)	0.86	1/4368 (0.0%)
1	H	0.73	1/3229 (0.0%)	0.87	0/4392
All	All	0.74	4/25791 (0.0%)	0.87	9/35083 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	215	VAL	CA-CB	6.45	1.57	1.54
1	H	215	VAL	CA-CB	6.39	1.57	1.54
1	A	215	VAL	CA-CB	5.21	1.57	1.54
1	E	215	VAL	CA-CB	5.05	1.57	1.54

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	62	ARG	NE-CZ-NH2	10.12	128.30	119.20
1	A	62	ARG	NE-CZ-NH1	-9.34	112.16	121.50
1	B	397	ILE	N-CA-C	6.32	116.35	107.37
1	A	62	ARG	CD-NE-CZ	6.18	133.05	124.40
1	B	214	THR	CA-C-N	5.90	123.95	120.24
1	B	214	THR	C-N-CA	5.90	123.95	120.24
1	G	370	ASN	N-CA-C	5.48	117.66	108.73
1	B	215	VAL	CA-C-N	-5.16	115.07	120.38
1	B	215	VAL	C-N-CA	-5.16	115.07	120.38

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	3149	0	3056	18	1
1	B	3127	0	3025	27	0
1	C	3130	0	3026	14	1
1	D	3134	0	3042	27	0
1	E	3127	0	3025	20	0
1	F	3129	0	3024	11	0
1	G	3124	0	3020	21	0
1	H	3136	0	3032	16	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0
2	G	2	0	0	0	0
2	H	2	0	0	0	0
3	A	332	0	0	3	0
3	B	219	0	0	4	0
3	C	197	0	0	3	0
3	D	269	0	0	5	0
3	E	249	0	0	3	0
3	F	250	0	0	2	0
3	G	166	0	0	3	0
3	H	166	0	0	0	0
All	All	26920	0	24250	141	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 3.

All (141) 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:62:ARG:HD2	1:B:63:ASP:OD1	1.61	1.00
1:A:62:ARG:HD2	1:A:63:ASP:OD1	1.74	0.86
1:F:264:GLY:HA2	1:F:265:ALA:HB3	1.59	0.83
1:G:264:GLY:HA2	1:G:265:ALA:HB3	1.60	0.81
1:H:264:GLY:HA2	1:H:265:ALA:HB3	1.61	0.81
1:E:264:GLY:HA2	1:E:265:ALA:HB3	1.64	0.80
1:C:264:GLY:HA2	1:C:265:ALA:HB3	1.63	0.79
1:D:229:LEU:HD21	1:D:235[B]:VAL:CG2	2.15	0.76
1:D:264:GLY:HA2	1:D:265:ALA:HB3	1.66	0.76
1:B:62:ARG:CD	1:B:63:ASP:OD1	2.34	0.76
1:B:303:LYS:HE3	1:D:303[B]:LYS:HG3	1.67	0.75
1:A:264:GLY:HA2	1:A:265:ALA:HB3	1.71	0.72
1:B:264:GLY:HA2	1:B:265:ALA:HB3	1.71	0.71
1:B:399:GLY:HA3	3:B:2208:HOH:O	1.95	0.67
1:A:137:LYS:HE2	3:A:2092:HOH:O	1.96	0.64
1:D:97:ARG:HD2	3:D:2063:HOH:O	1.97	0.63
1:D:229:LEU:HD21	1:D:235[B]:VAL:HG22	1.79	0.63
1:D:229:LEU:HD21	1:D:235[B]:VAL:HG21	1.80	0.61
1:A:304:ASP:HB3	3:A:2308:HOH:O	2.02	0.60
1:G:79:LYS:NZ	3:G:2043:HOH:O	2.36	0.59
1:D:204:ARG:HD3	1:D:204:ARG:C	2.29	0.58
1:C:311:ILE:HD13	1:D:218:ARG:HA	1.86	0.57
1:G:264:GLY:CA	1:G:265:ALA:HB3	2.34	0.57
1:F:264:GLY:CA	1:F:265:ALA:HB3	2.33	0.57
1:E:72:LEU:HD21	1:E:154:LEU:HD13	1.86	0.56
1:B:107:PHE:O	1:B:124:ALA:HA	2.07	0.55
1:B:160:ASP:HB2	1:B:161:PRO:HD2	1.89	0.54
1:E:218:ARG:HA	1:G:311:ILE:HD13	1.90	0.54
1:F:218:ARG:HA	1:H:311:ILE:HD13	1.91	0.53
1:D:9:ASP:OD2	1:E:9:ASP:OD2	2.27	0.53
1:B:72:LEU:HD21	1:B:154:LEU:HD13	1.90	0.53
1:C:106:GLY:O	1:C:263:CYS:HB3	2.08	0.52
1:A:264:GLY:CA	1:A:265:ALA:HB3	2.40	0.52
1:G:114:GLY:O	3:G:2062:HOH:O	2.19	0.52
1:A:72:LEU:HD21	1:A:154:LEU:HD13	1.92	0.52
1:C:72:LEU:HD21	1:C:154:LEU:HD13	1.91	0.52
1:E:264:GLY:CA	1:E:265:ALA:HB3	2.36	0.52
1:C:41:LYS:NZ	3:C:2016:HOH:O	2.42	0.52
1:D:264:GLY:CA	1:D:265:ALA:HB3	2.38	0.52
1:D:160:ASP:HB2	1:D:161:PRO:HD2	1.92	0.51
1:F:72:LEU:HD21	1:F:154:LEU:HD13	1.91	0.51
1:B:398:THR:HG22	1:B:399:GLY:H	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:160:ASP:HB2	1:E:161:PRO:HD2	1.92	0.51
1:H:326:VAL:O	1:H:330:GLN:HG2	2.11	0.51
1:D:72:LEU:HD21	1:D:154:LEU:HD13	1.92	0.51
1:A:160:ASP:HB2	1:A:161:PRO:HD2	1.92	0.51
1:G:160:ASP:HB2	1:G:161:PRO:HD2	1.92	0.51
1:H:160:ASP:HB2	1:H:161:PRO:HD2	1.93	0.51
1:F:341:LYS:HE2	3:F:2220:HOH:O	2.11	0.50
1:H:264:GLY:CA	1:H:265:ALA:HB3	2.35	0.50
1:B:65:ASP:OD1	1:B:67[A]:SER:OG	2.24	0.50
1:G:72:LEU:HD21	1:G:154:LEU:HD13	1.93	0.50
1:G:106:GLY:O	1:G:263:CYS:HB3	2.12	0.49
1:A:7:LYS:HE3	3:A:2003:HOH:O	2.12	0.49
1:G:316:LYS:NZ	3:G:2134:HOH:O	2.44	0.49
1:B:132:HIS:HD1	1:C:39:GLU:CD	2.20	0.49
1:H:72:LEU:HD21	1:H:154:LEU:HD13	1.93	0.49
1:B:61:VAL:O	1:B:64:VAL:HG12	2.12	0.49
1:G:167:TRP:CD2	1:G:218:ARG:HD3	2.47	0.49
1:C:160:ASP:HB2	1:C:161:PRO:HD2	1.95	0.49
1:F:160:ASP:HB2	1:F:161:PRO:HD2	1.94	0.49
1:G:107:PHE:O	1:G:124:ALA:HA	2.13	0.49
1:H:167:TRP:CD2	1:H:218:ARG:HD3	2.49	0.48
1:C:41:LYS:NZ	3:C:2017:HOH:O	2.44	0.48
1:B:106:GLY:O	1:B:263:CYS:HB3	2.13	0.48
1:B:167:TRP:CD2	1:B:218:ARG:HD3	2.48	0.48
1:C:264:GLY:CA	1:C:265:ALA:HB3	2.38	0.48
1:D:167:TRP:CD2	1:D:218:ARG:HD3	2.48	0.48
1:E:107:PHE:O	1:E:124:ALA:HA	2.14	0.47
1:E:311:ILE:HD13	1:G:218:ARG:HA	1.97	0.47
1:H:107:PHE:O	1:H:124:ALA:HA	2.15	0.47
3:B:2159:HOH:O	1:D:303[A]:LYS:HE2	2.14	0.46
1:C:304:ASP:HB3	3:C:2178:HOH:O	2.15	0.46
1:B:264:GLY:HA3	1:B:265:ALA:O	2.15	0.46
1:F:167:TRP:CD2	1:F:218:ARG:HD3	2.50	0.46
1:A:167:TRP:CD2	1:A:218:ARG:HD3	2.50	0.46
1:B:394:GLN:O	1:B:396:ILE:HG12	2.16	0.46
1:B:97:ARG:HG3	1:B:270:GLY:HA2	1.97	0.46
1:E:137:LYS:HE2	3:E:2060:HOH:O	2.16	0.46
1:E:225:ASP:OD1	1:G:367:ASP:OD2	2.34	0.46
1:G:144:VAL:HG11	1:G:352:LEU:HD11	1.98	0.46
1:B:123:LEU:HD21	3:B:2217:HOH:O	2.15	0.45
1:A:326:VAL:O	1:A:330[B]:GLN:HG2	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:107:PHE:O	1:C:124:ALA:HA	2.16	0.45
1:B:62:ARG:HH21	1:B:63:ASP:CG	2.23	0.45
1:D:107:PHE:O	1:D:124:ALA:HA	2.16	0.45
1:H:106:GLY:O	1:H:263:CYS:HB3	2.16	0.45
1:D:106:GLY:O	1:D:263:CYS:HB3	2.17	0.45
1:E:89:VAL:HB	1:E:235:VAL:CG1	2.47	0.45
1:E:318:LYS:HB3	3:E:2212:HOH:O	2.16	0.45
1:H:204:ARG:HD3	1:H:204:ARG:C	2.42	0.45
1:A:218:ARG:HA	1:B:311:ILE:HD13	1.97	0.45
1:B:15:TYR:HA	1:B:195:HIS:CD2	2.52	0.44
1:D:326:VAL:O	1:D:330:GLN:HG2	2.16	0.44
1:G:15:TYR:HA	1:G:195:HIS:CD2	2.52	0.44
1:E:235:VAL:CG1	1:E:237:PHE:CE1	3.01	0.44
1:E:235:VAL:HG11	1:E:237:PHE:CE1	2.52	0.44
1:A:107:PHE:CD2	1:A:107:PHE:N	2.85	0.43
1:A:107:PHE:O	1:A:124:ALA:HA	2.18	0.43
1:A:200[B]:GLN:OE1	1:A:200[B]:GLN:HA	2.17	0.43
1:B:190:ASN:OD1	1:B:190:ASN:C	2.61	0.43
1:D:107:PHE:CD2	1:D:107:PHE:N	2.87	0.43
1:E:89:VAL:HB	1:E:235:VAL:HG12	1.99	0.43
1:F:27:ASP:HA	3:F:2043:HOH:O	2.19	0.43
1:G:352:LEU:HD12	1:G:352:LEU:HA	1.89	0.43
1:H:255:GLN:HG2	1:H:256:GLY:N	2.33	0.43
1:E:65:ASP:OD2	3:E:2074:HOH:O	2.21	0.43
1:A:106:GLY:O	1:A:263:CYS:HB3	2.19	0.42
1:A:255:GLN:HG3	1:A:265:ALA:HB2	2.01	0.42
1:B:303:LYS:HE3	1:D:303[B]:LYS:CG	2.43	0.42
1:D:302:TYR:O	1:D:303[A]:LYS:HD2	2.18	0.42
1:D:379:THR:HG22	1:D:386:ILE:HD11	2.00	0.42
1:D:404:PRO:HD3	3:D:2263:HOH:O	2.19	0.42
1:F:106:GLY:O	1:F:263:CYS:HB3	2.20	0.42
1:F:107:PHE:CD2	1:F:107:PHE:N	2.88	0.42
1:D:137:LYS:HE2	3:D:2065:HOH:O	2.19	0.42
1:E:167:TRP:CD2	1:E:218:ARG:HD3	2.55	0.42
1:B:199:MET:HA	3:B:2121:HOH:O	2.19	0.41
1:C:167:TRP:CD2	1:C:218:ARG:HD3	2.55	0.41
1:G:202:GLU:HG3	1:G:203:ALA:N	2.35	0.41
1:B:199:MET:HB3	1:B:204:ARG:HB2	2.02	0.41
1:C:218:ARG:HA	1:D:311:ILE:HD13	2.02	0.41
1:D:330:GLN:HB3	3:D:2227:HOH:O	2.20	0.41
1:H:167:TRP:CD2	1:H:218:ARG:CD	3.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:15:TYR:HA	1:A:195:HIS:CD2	2.55	0.41
1:E:106:GLY:O	1:E:263:CYS:HB3	2.20	0.41
1:E:380:GLU:OE2	1:H:303:LYS:NZ	2.49	0.41
1:G:255:GLN:HG2	1:G:256:GLY:N	2.35	0.41
1:H:15:TYR:HA	1:H:195:HIS:CD2	2.55	0.41
1:A:298:MET:HE3	1:D:298:MET:HE3	2.02	0.41
1:F:107:PHE:O	1:F:124:ALA:HA	2.20	0.41
1:G:107:PHE:CD2	1:G:107:PHE:N	2.86	0.41
1:H:107:PHE:CD2	1:H:107:PHE:N	2.86	0.41
1:H:395:LYS:HB3	1:H:395:LYS:HE2	1.67	0.41
1:B:264:GLY:CA	1:B:265:ALA:HB3	2.47	0.41
1:E:15:TYR:HA	1:E:195:HIS:CD2	2.56	0.41
1:G:394:GLN:HG2	1:G:396:ILE:HD12	2.03	0.41
1:G:206:LYS:O	1:G:210:GLU:HG2	2.20	0.41
1:D:190:ASN:OD1	1:D:190:ASN:C	2.64	0.40
1:C:289:LYS:NZ	3:D:2227:HOH:O	2.54	0.40
1:B:352:LEU:HA	1:B:352:LEU:HD12	1.82	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:GLU:OE1	1:C:178:ASP:OD1[1_545]	1.66	0.54

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	411/409 (100%)	392 (95%)	19 (5%)	0	100	100
1	B	407/409 (100%)	384 (94%)	23 (6%)	0	100	100
1	C	407/409 (100%)	385 (95%)	22 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	408/409 (100%)	390 (96%)	18 (4%)	0	100	100
1	E	407/409 (100%)	385 (95%)	20 (5%)	2 (0%)	24	21
1	F	407/409 (100%)	383 (94%)	24 (6%)	0	100	100
1	G	406/409 (99%)	384 (95%)	22 (5%)	0	100	100
1	H	408/409 (100%)	386 (95%)	21 (5%)	1 (0%)	43	44
All	All	3261/3272 (100%)	3089 (95%)	169 (5%)	3 (0%)	48	49

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	398	THR
1	E	201	GLY
1	E	399	GLY

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	334/330 (101%)	323 (97%)	11 (3%)	33	35
1	B	330/330 (100%)	321 (97%)	9 (3%)	39	43
1	C	330/330 (100%)	323 (98%)	7 (2%)	47	52
1	D	331/330 (100%)	325 (98%)	6 (2%)	51	58
1	E	330/330 (100%)	320 (97%)	10 (3%)	36	39
1	F	330/330 (100%)	318 (96%)	12 (4%)	31	32
1	G	329/330 (100%)	322 (98%)	7 (2%)	47	52
1	H	331/330 (100%)	321 (97%)	10 (3%)	36	39
All	All	2645/2640 (100%)	2573 (97%)	72 (3%)	39	43

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	ASP
1	A	107	PHE
1	A	195	HIS
1	A	200[A]	GLN
1	A	200[B]	GLN
1	A	206	LYS
1	A	229	LEU
1	A	303	LYS
1	A	318	LYS
1	A	395	LYS
1	A	401	VAL
1	B	46	ASP
1	B	107	PHE
1	B	180	ASP
1	B	200	GLN
1	B	204	ARG
1	B	229	LEU
1	B	303	LYS
1	B	391	GLU
1	B	394	GLN
1	C	46	ASP
1	C	107	PHE
1	C	195	HIS
1	C	229	LEU
1	C	303	LYS
1	C	318	LYS
1	C	401	VAL
1	D	46	ASP
1	D	107	PHE
1	D	195	HIS
1	D	200	GLN
1	D	229	LEU
1	D	396	ILE
1	E	46	ASP
1	E	107	PHE
1	E	195	HIS
1	E	202	GLU
1	E	204	ARG
1	E	229	LEU
1	E	303	LYS
1	E	318	LYS
1	E	396	ILE
1	E	401	VAL

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Mol	Chain	Res	Type
1	F	46	ASP
1	F	107	PHE
1	F	195	HIS
1	F	229	LEU
1	F	303	LYS
1	F	330	GLN
1	F	377	LEU
1	F	394	GLN
1	F	395	LYS
1	F	396	ILE
1	F	398	THR
1	F	401	VAL
1	G	46	ASP
1	G	107	PHE
1	G	202	GLU
1	G	204	ARG
1	G	229	LEU
1	G	394	GLN
1	G	401	VAL
1	H	35	GLU
1	H	39[A]	GLU
1	H	39[B]	GLU
1	H	46	ASP
1	H	107	PHE
1	H	195	HIS
1	H	229	LEU
1	H	391	GLU
1	H	395	LYS
1	H	409	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	111	GLN
1	A	145	ASN
1	B	145	ASN
1	C	111	GLN
1	C	125	GLN
1	C	145	ASN
1	D	111	GLN
1	D	120	HIS
1	D	125	GLN

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Mol	Chain	Res	Type
1	E	111	GLN
1	E	394	GLN
1	F	111	GLN
1	F	145	ASN
1	G	111	GLN
1	G	125	GLN
1	G	145	ASN
1	H	111	GLN
1	H	125	GLN
1	H	145	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 16 ligands modelled in this entry, 16 are monoatomic - leaving 0 for Mogul analysis.

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 ⓘ

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	408/409 (99%)	-1.59	0 100 100	18, 36, 56, 72	5 (1%)
1	B	408/409 (99%)	-1.53	0 100 100	20, 36, 61, 97	1 (0%)
1	C	408/409 (99%)	-1.56	0 100 100	20, 36, 61, 83	1 (0%)
1	D	408/409 (99%)	-1.58	0 100 100	15, 36, 58, 71	2 (0%)
1	E	408/409 (99%)	-1.61	0 100 100	20, 36, 61, 80	1 (0%)
1	F	408/409 (99%)	-1.60	0 100 100	19, 36, 59, 77	1 (0%)
1	G	408/409 (99%)	-1.55	0 100 100	20, 36, 62, 84	0
1	H	408/409 (99%)	-1.59	0 100 100	18, 36, 61, 86	2 (0%)
All	All	3264/3272 (99%)	-1.58	0 100 100	15, 36, 61, 97	13 (0%)

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates

There are no oligosaccharides in this entry.

6.4 Ligands

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
2	ZN	A	411	1/1	0.99	0.03	55,55,55,55	0
2	ZN	B	411	1/1	0.99	0.02	48,48,48,48	0
2	ZN	C	411	1/1	0.99	0.02	56,56,56,56	0
2	ZN	D	411	1/1	0.99	0.03	55,55,55,55	0
2	ZN	G	411	1/1	0.99	0.02	67,67,67,67	0
2	ZN	H	411	1/1	0.99	0.02	59,59,59,59	0
2	ZN	B	412	1/1	1.00	0.03	45,45,45,45	0
2	ZN	D	412	1/1	1.00	0.02	38,38,38,38	0
2	ZN	E	411	1/1	1.00	0.02	61,61,61,61	0
2	ZN	E	412	1/1	1.00	0.01	46,46,46,46	0
2	ZN	F	411	1/1	1.00	0.03	68,68,68,68	0
2	ZN	F	412	1/1	1.00	0.01	46,46,46,46	0
2	ZN	A	412	1/1	1.00	0.02	42,42,42,42	0
2	ZN	G	412	1/1	1.00	0.02	51,51,51,51	0
2	ZN	C	412	1/1	1.00	0.03	47,47,47,47	0
2	ZN	H	412	1/1	1.00	0.03	49,49,49,49	0

6.5 Other polymers [i](#)

There are no such residues in this entry.