



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 11:29 AM UTC

PDB ID : 3ZNB / pdb_00003znb
Title : METALLO-BETA-LACTAMASE (ZN, HG-BOUND FORM)
Authors : Concha, N.O.; Herzberg, O.
Deposited on : 1997-10-15
Resolution : 2.70 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

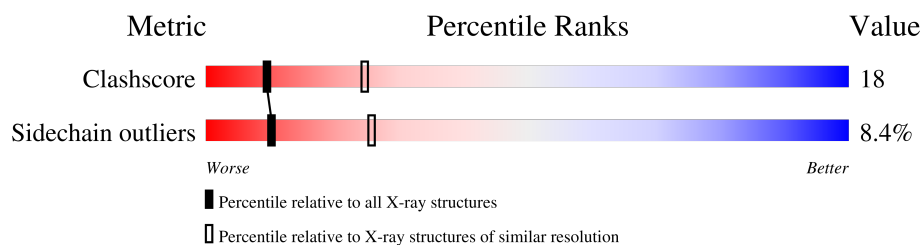
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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))
Clashscore	190562	3843 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	232	
1	B	232	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3480 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 METALLO-BETA-LACTAMASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	226	Total	C	N	O	S	0	0	0
			1697	1076	281	332	8			
1	B	227	Total	C	N	O	S	0	0	0
			1712	1082	282	340	8			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	79	THR	MET	conflict	UNP P25910
A	85	ALA	THR	conflict	UNP P25910
A	113	LYS	ARG	conflict	UNP P25910
B	79	THR	MET	conflict	UNP P25910
B	85	ALA	THR	conflict	UNP P25910
B	113	LYS	ARG	conflict	UNP P25910

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

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

- Molecule 3 is MERCURY (II) ION (CCD ID: HG) (formula: Hg).

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

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Na 1	0	0
4	B	1	Total 1	Na 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	25	Total 25	O 25	0	0
5	B	40	Total 40	O 40	0	0

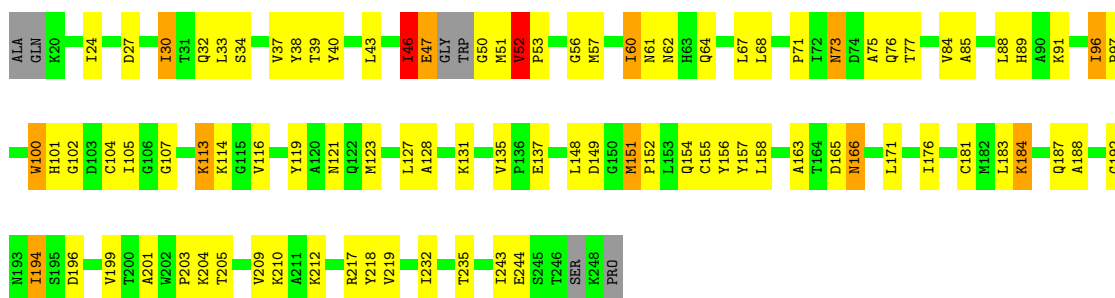
3 Residue-property plots [i](#)

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. 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. 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.

Note EDS was not executed.

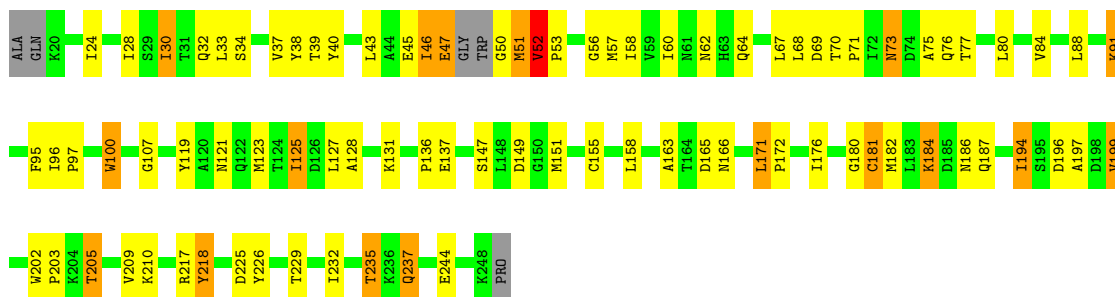
• Molecule 1: METALLO-BETA-LACTAMASE

Chain A: 



• Molecule 1: METALLO-BETA-LACTAMASE

Chain B: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	78.20 Å 78.20 Å 140.60 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.70	Depositor
% Data completeness (in resolution range)	90.0 (8.00-2.70)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.200 , 0.268	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3480	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, HG, NA

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	1.26	5/1733 (0.3%)	1.36	17/2358 (0.7%)
1	B	1.22	3/1747 (0.2%)	1.38	16/2384 (0.7%)
All	All	1.24	8/3480 (0.2%)	1.37	33/4742 (0.7%)

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	1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	96	ILE	CA-C	8.63	1.59	1.52
1	B	181	CYS	CB-SG	6.66	2.03	1.81
1	B	194	ILE	CA-CB	6.22	1.61	1.54
1	A	46	ILE	CA-CB	5.89	1.62	1.54
1	A	188	ALA	CA-CB	-5.74	1.44	1.53
1	B	205	THR	CA-CB	-5.72	1.44	1.53
1	A	201	ALA	CA-CB	-5.40	1.44	1.53
1	A	96	ILE	CA-CB	-5.22	1.50	1.54

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	52	VAL	N-CA-C	9.46	119.36	108.96
1	A	52	VAL	N-CA-C	8.24	118.03	108.96
1	A	217	ARG	N-CA-C	-8.13	102.01	111.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	100	TRP	N-CA-C	8.00	122.61	113.01
1	B	137	GLU	N-CA-C	7.99	123.14	113.23
1	B	171	LEU	CA-C-N	7.98	127.54	119.24
1	B	171	LEU	C-N-CA	7.98	127.54	119.24
1	A	137	GLU	N-CA-C	7.26	122.23	113.23
1	B	37	VAL	N-CA-C	7.19	118.83	108.48
1	A	135	VAL	N-CA-C	7.12	115.84	107.73
1	B	217	ARG	N-CA-C	-7.11	103.13	111.03
1	B	128	ALA	N-CA-C	-7.11	103.53	111.28
1	A	100	TRP	N-CA-C	6.70	121.04	113.01
1	A	171	LEU	CA-C-N	6.64	128.15	119.84
1	A	171	LEU	C-N-CA	6.64	128.15	119.84
1	A	37	VAL	N-CA-C	6.62	118.01	108.48
1	B	51	MET	N-CA-C	-6.55	96.85	110.80
1	A	51	MET	N-CA-C	-6.13	97.75	110.80
1	B	235	THR	N-CA-C	-5.85	104.91	111.28
1	A	116	VAL	N-CA-C	5.68	116.39	108.89
1	A	33	LEU	N-CA-C	-5.60	107.00	113.88
1	A	60	ILE	N-CA-C	5.46	116.44	108.53
1	A	166	ASN	N-CA-C	5.45	118.24	110.24
1	A	128	ALA	N-CA-C	-5.44	105.35	111.28
1	B	73	ASN	N-CA-C	5.40	115.67	108.38
1	B	58	ILE	N-CA-C	-5.38	98.16	109.34
1	B	199	VAL	N-CA-C	5.37	115.58	110.42
1	A	73	ASN	N-CA-C	5.24	115.45	108.38
1	B	197	ALA	N-CA-C	5.19	118.57	110.32
1	B	136	PRO	N-CA-C	-5.15	103.02	111.21
1	A	105	ILE	CB-CA-C	-5.03	103.04	111.29
1	B	237	GLN	N-CA-C	-5.03	105.71	111.14
1	A	154	GLN	N-CA-C	5.00	116.88	108.73

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	218	TYR	Sidechain

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	1697	0	1628	58	0
1	B	1712	0	1645	62	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	25	0	0	3	0
5	B	40	0	0	12	0
All	All	3480	0	3273	120	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 18.

All (120) 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:181:CYS:CB	1:B:181:CYS:SG	2.03	1.46
1:B:24:ILE:HD13	1:B:30:ILE:HD11	1.61	0.82
1:A:24:ILE:HD13	1:A:30:ILE:HD11	1.66	0.78
1:B:46:ILE:HD12	1:B:47:GLU:HB2	1.66	0.78
1:B:46:ILE:HD12	1:B:47:GLU:H	1.52	0.74
1:A:184:LYS:HD2	1:A:235:THR:HG23	1.70	0.73
1:B:57:MET:HG3	1:B:68:LEU:HD12	1.71	0.73
1:B:184:LYS:HD2	1:B:235:THR:HG23	1.71	0.71
1:B:237:GLN:NE2	1:B:237:GLN:OE1	2.25	0.69
1:A:46:ILE:CG1	1:A:47:GLU:H	2.06	0.69
1:B:46:ILE:CG1	1:B:47:GLU:H	2.03	0.69
1:B:123:MET:HE3	5:B:266:HOH:O	1.92	0.68
1:A:27:ASP:HB2	5:A:272:HOH:O	1.94	0.68
1:B:181:CYS:SG	5:B:259:HOH:O	2.50	0.68
1:A:181:CYS:SG	5:A:262:HOH:O	2.52	0.67
1:A:100:TRP:CH2	1:A:131:LYS:HE2	2.29	0.67
1:B:186:ASN:HB2	5:B:289:HOH:O	1.94	0.67
1:B:100:TRP:CH2	1:B:131:LYS:HE2	2.30	0.67
1:B:46:ILE:CD1	1:B:47:GLU:H	2.09	0.66
1:B:196:ASP:HB3	5:B:274:HOH:O	1.96	0.65
1:A:71:PRO:HB2	1:A:76:GLN:HB2	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:62:ASN:HD22	1:B:149:ASP:CG	2.06	0.63
1:B:47:GLU:O	1:B:50:GLY:N	2.32	0.63
1:A:47:GLU:O	1:A:50:GLY:N	2.33	0.61
1:B:45:GLU:HA	1:B:51:MET:HA	1.82	0.60
1:A:24:ILE:HG21	1:A:30:ILE:HG13	1.83	0.59
1:B:91:LYS:HA	5:B:276:HOH:O	2.03	0.59
1:B:71:PRO:HB2	1:B:76:GLN:HB2	1.85	0.59
1:B:46:ILE:CD1	1:B:47:GLU:HB2	2.32	0.58
1:A:100:TRP:CD1	1:A:127:LEU:HD13	2.39	0.58
1:A:57:MET:HG3	1:A:68:LEU:HD12	1.87	0.57
1:B:24:ILE:HG21	1:B:30:ILE:HG13	1.86	0.56
1:A:46:ILE:HD12	1:A:47:GLU:HB2	1.87	0.56
1:B:24:ILE:HG21	1:B:30:ILE:CG1	2.35	0.56
1:A:46:ILE:HD12	1:A:47:GLU:H	1.71	0.55
1:A:163:ALA:HA	1:A:196:ASP:O	2.07	0.55
1:B:199:VAL:O	1:B:203:PRO:HD3	2.08	0.54
1:A:96:ILE:HD13	1:A:155:CYS:SG	2.48	0.54
1:A:24:ILE:HG21	1:A:30:ILE:CG1	2.37	0.54
1:B:225:ASP:HA	5:B:286:HOH:O	2.08	0.54
1:B:210:LYS:HB2	1:B:232:ILE:HG21	1.90	0.53
1:B:226:TYR:N	5:B:286:HOH:O	2.41	0.53
1:B:46:ILE:CG1	1:B:47:GLU:N	2.72	0.53
1:B:229:THR:HG21	5:B:278:HOH:O	2.08	0.52
1:A:176:ILE:HG12	1:A:218:TYR:HB2	1.92	0.52
1:A:84:VAL:O	1:A:88:LEU:HB2	2.10	0.52
1:B:176:ILE:HG12	1:B:218:TYR:HB2	1.92	0.52
1:A:123:MET:HE1	1:A:157:TYR:CE2	2.45	0.51
1:B:73:ASN:HD21	1:B:75:ALA:HB3	1.75	0.50
1:A:46:ILE:CD1	1:A:47:GLU:H	2.24	0.50
1:B:123:MET:HB3	1:B:165:ASP:HB3	1.92	0.50
1:B:43:LEU:HD13	5:B:272:HOH:O	2.10	0.50
1:A:148:LEU:N	1:A:151:MET:O	2.40	0.50
1:B:64:GLN:NE2	1:B:149:ASP:HB2	2.27	0.50
1:A:205:THR:O	1:A:209:VAL:HG23	2.12	0.50
1:B:24:ILE:CG2	1:B:30:ILE:HG12	2.43	0.49
1:A:210:LYS:HB2	1:A:232:ILE:HG21	1.95	0.49
1:B:46:ILE:HD12	1:B:47:GLU:CB	2.40	0.49
1:A:199:VAL:O	1:A:203:PRO:HD3	2.13	0.48
1:B:100:TRP:CD1	1:B:127:LEU:HD13	2.48	0.48
1:A:64:GLN:NE2	1:A:149:ASP:HB2	2.29	0.48
1:B:68:LEU:O	1:B:69:ASP:HB2	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:84:VAL:O	1:B:88:LEU:HB2	2.14	0.47
1:B:96:ILE:HD13	1:B:155:CYS:SG	2.54	0.47
1:B:147:SER:HB2	5:B:283:HOH:O	2.14	0.47
1:B:163:ALA:HA	1:B:196:ASP:O	2.14	0.47
1:A:46:ILE:CG1	1:A:47:GLU:N	2.77	0.47
1:A:123:MET:HB3	1:A:165:ASP:HB3	1.97	0.46
1:B:226:TYR:HD2	5:B:286:HOH:O	1.97	0.46
1:B:32:GLN:NE2	1:B:34:SER:O	2.48	0.46
1:A:24:ILE:CG2	1:A:30:ILE:HG12	2.46	0.45
1:B:97:PRO:HD2	1:B:119:TYR:O	2.15	0.45
1:B:73:ASN:OD1	1:B:76:GLN:HG3	2.16	0.45
1:B:28:ILE:HG21	1:B:80:LEU:HD13	1.99	0.45
1:A:24:ILE:HG23	1:A:30:ILE:HG12	1.99	0.45
1:A:73:ASN:OD1	1:A:76:GLN:HG3	2.17	0.45
1:A:156:TYR:CD2	1:A:212:LYS:HE2	2.52	0.44
1:A:97:PRO:HD2	1:A:119:TYR:O	2.18	0.44
1:A:100:TRP:NE1	1:A:127:LEU:HD13	2.33	0.44
1:A:52:VAL:HG22	1:A:53:PRO:HD2	2.00	0.44
1:A:151:MET:HA	1:A:152:PRO:HD2	1.91	0.44
1:A:62:ASN:HD22	1:A:149:ASP:CG	2.26	0.44
1:A:32:GLN:NE2	1:A:38:TYR:CZ	2.86	0.44
1:A:61:ASN:O	1:A:64:GLN:HB2	2.17	0.44
1:A:192:GLY:O	1:A:194:ILE:HG23	2.18	0.44
1:B:121:ASN:O	1:B:125:ILE:HG13	2.18	0.43
1:A:43:LEU:HD12	1:A:43:LEU:N	2.33	0.43
1:B:24:ILE:HG23	1:B:30:ILE:HG12	2.00	0.43
1:A:77:THR:OG1	1:A:107:GLY:HA3	2.18	0.43
1:B:181:CYS:HB3	5:B:258:HOH:O	2.19	0.43
1:A:52:VAL:HG22	1:A:53:PRO:CD	2.49	0.43
1:B:57:MET:CG	1:B:68:LEU:HD12	2.44	0.43
1:B:32:GLN:NE2	1:B:38:TYR:CZ	2.86	0.42
1:B:202:TRP:HB3	1:B:203:PRO:HD3	1.99	0.42
1:A:40:TYR:CE2	1:A:56:GLY:HA3	2.55	0.42
1:A:52:VAL:HA	1:A:53:PRO:HD3	1.83	0.42
1:B:77:THR:OG1	1:B:107:GLY:HA3	2.19	0.42
1:A:67:LEU:HA	1:A:67:LEU:HD12	1.74	0.42
1:A:88:LEU:HD23	1:A:88:LEU:HA	1.80	0.42
1:A:24:ILE:CD1	1:A:30:ILE:HD11	2.45	0.42
1:B:180:GLY:C	1:B:182:MET:H	2.27	0.42
1:A:85:ALA:O	1:A:89:HIS:HA	2.19	0.42
1:A:46:ILE:CD1	1:A:47:GLU:HB2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:PRO:HB2	1:A:76:GLN:CB	2.49	0.41
1:A:199:VAL:HG12	1:A:243:ILE:HD11	2.01	0.41
1:B:33:LEU:HG	1:B:39:THR:HG22	2.03	0.41
1:B:205:THR:O	1:B:209:VAL:HG23	2.21	0.41
1:B:40:TYR:CE1	1:B:56:GLY:HA3	2.56	0.41
1:B:121:ASN:ND2	1:B:123:MET:HB2	2.35	0.41
1:B:171:LEU:HA	1:B:172:PRO:HD2	1.85	0.41
1:A:113:LYS:HE2	5:A:264:HOH:O	2.20	0.41
1:B:67:LEU:O	1:B:95:PHE:HA	2.21	0.41
1:A:73:ASN:HD21	1:A:75:ALA:HB3	1.86	0.41
1:A:121:ASN:ND2	1:A:123:MET:HB2	2.36	0.41
1:B:70:THR:HG23	1:B:95:PHE:CE1	2.56	0.41
1:A:101:HIS:O	1:A:104:CYS:HB2	2.20	0.40
1:A:183:LEU:HD13	1:A:219:VAL:CG1	2.51	0.40
1:A:32:GLN:NE2	1:A:34:SER:O	2.54	0.40
1:A:102:GLY:C	1:A:104:CYS:H	2.29	0.40
1:B:52:VAL:HG22	1:B:53:PRO:CD	2.52	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

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	183/196 (93%)	166 (91%)	17 (9%)	8	21
1	B	187/196 (95%)	173 (92%)	14 (8%)	12	31
All	All	370/392 (94%)	339 (92%)	31 (8%)	10	26

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	ILE
1	A	39	THR
1	A	46	ILE
1	A	47	GLU
1	A	52	VAL
1	A	60	ILE
1	A	91	LYS
1	A	113	LYS
1	A	114	LYS
1	A	151	MET
1	A	158	LEU
1	A	166	ASN
1	A	184	LYS
1	A	187	GLN
1	A	194	ILE
1	A	204	LYS
1	A	244	GLU
1	B	30	ILE
1	B	46	ILE
1	B	47	GLU
1	B	52	VAL
1	B	60	ILE
1	B	91	LYS
1	B	125	ILE
1	B	151	MET
1	B	158	LEU
1	B	166	ASN
1	B	184	LYS
1	B	187	GLN
1	B	194	ILE
1	B	244	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	GLN
1	A	64	GLN
1	A	240	ASN
1	B	32	GLN
1	B	62	ASN
1	B	64	GLN

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Mol	Chain	Res	Type
1	B	186	ASN
1	B	187	GLN
1	B	240	ASN
1	B	241	GLN

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 6 ligands modelled in this entry, 6 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 [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section is therefore empty.

6.4 Ligands ⓘ

EDS was not executed - this section is therefore empty.

6.5 Other polymers ⓘ

EDS was not executed - this section is therefore empty.