



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 12:37 PM UTC

PDB ID : 8SZT / pdb_00008szt
Title : Structure of Kdac1 from Acinetobacter baumannii
Authors : Watson, P.R.; Christianson, D.W.
Deposited on : 2023-05-30
Resolution : 2.50 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

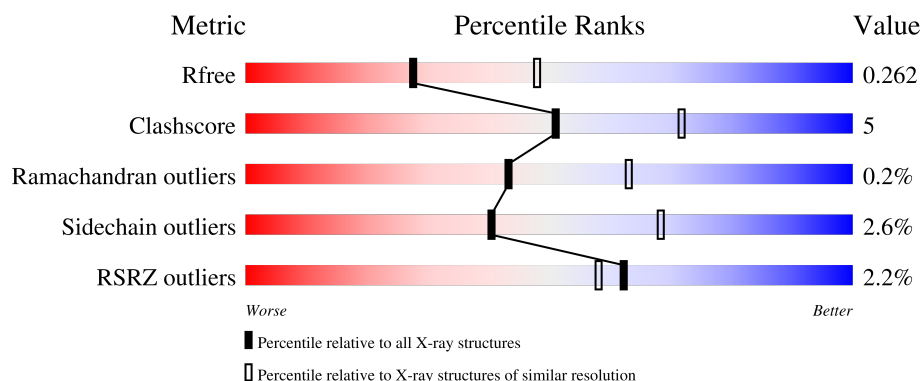
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

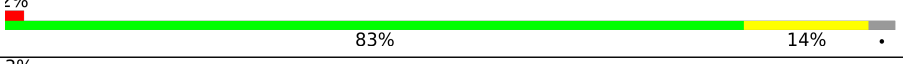
The reported resolution of this entry is 2.50 Å.

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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	372	 3% 87% 11% ..
1	B	372	 2% 83% 14% .
1	C	372	 3% 84% 13% ..
1	D	372	 % 83% 15% .

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10819 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 Histone deacetylase domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	360	Total	C	N	O	S	0	0	0
			2661	1702	448	499	12			
1	A	365	Total	C	N	O	S	0	0	0
			2693	1719	456	506	12			
1	C	365	Total	C	N	O	S	0	0	0
			2705	1727	456	510	12			
1	D	363	Total	C	N	O	S	0	0	0
			2686	1714	452	508	12			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	GLY	-	expression tag	UNP R8YPJ7
B	0	HIS	-	expression tag	UNP R8YPJ7
B	4	ILE	VAL	conflict	UNP R8YPJ7
B	66	TYR	HIS	conflict	UNP R8YPJ7
B	101	VAL	ILE	conflict	UNP R8YPJ7
B	340	ILE	MET	conflict	UNP R8YPJ7
A	-1	GLY	-	expression tag	UNP R8YPJ7
A	0	HIS	-	expression tag	UNP R8YPJ7
A	4	ILE	VAL	conflict	UNP R8YPJ7
A	66	TYR	HIS	conflict	UNP R8YPJ7
A	101	VAL	ILE	conflict	UNP R8YPJ7
A	340	ILE	MET	conflict	UNP R8YPJ7
C	-1	GLY	-	expression tag	UNP R8YPJ7
C	0	HIS	-	expression tag	UNP R8YPJ7
C	4	ILE	VAL	conflict	UNP R8YPJ7
C	66	TYR	HIS	conflict	UNP R8YPJ7
C	101	VAL	ILE	conflict	UNP R8YPJ7
C	340	ILE	MET	conflict	UNP R8YPJ7
D	-1	GLY	-	expression tag	UNP R8YPJ7
D	0	HIS	-	expression tag	UNP R8YPJ7
D	4	ILE	VAL	conflict	UNP R8YPJ7

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Chain	Residue	Modelled	Actual	Comment	Reference
D	66	TYR	HIS	conflict	UNP R8YPJ7
D	101	VAL	ILE	conflict	UNP R8YPJ7
D	340	ILE	MET	conflict	UNP R8YPJ7

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

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

- Molecule 3 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	2	Total K 2 2	0	0
3	A	2	Total K 2 2	0	0
3	C	2	Total K 2 2	0	0
3	D	2	Total K 2 2	0	0

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Cl 1 1	0	0
4	A	1	Total Cl 1 1	0	0
4	C	1	Total Cl 1 1	0	0
4	D	1	Total Cl 1 1	0	0

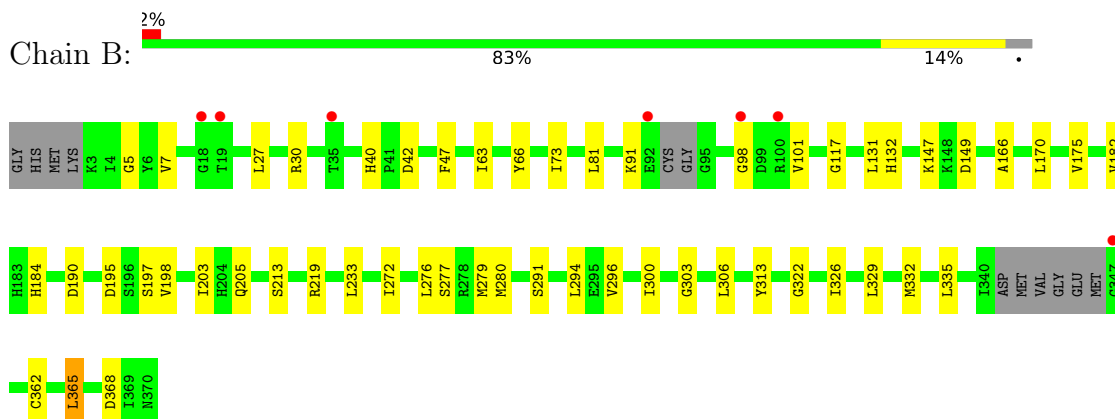
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	16	Total 16	O 16	0	0
5	A	18	Total 18	O 18	0	0
5	C	13	Total 13	O 13	0	0
5	D	11	Total 11	O 11	0	0

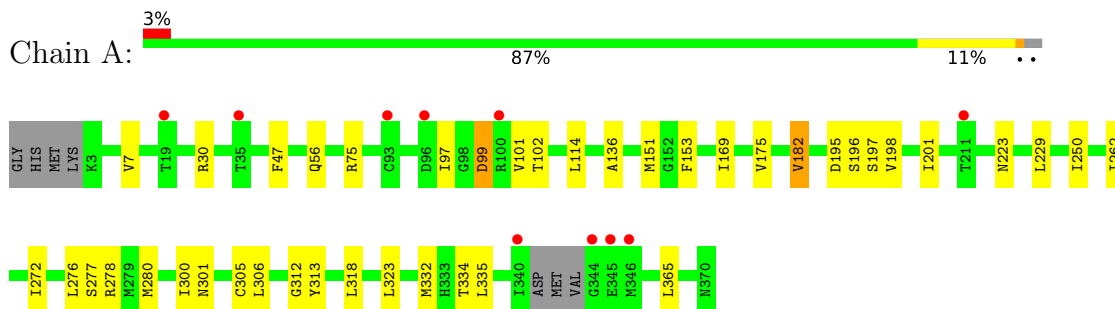
3 Residue-property plots [i](#)

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.

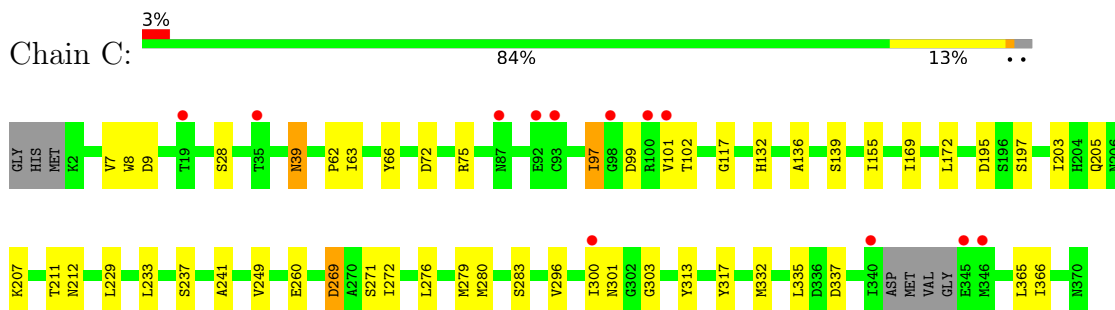
- Molecule 1: Histone deacetylase domain-containing protein



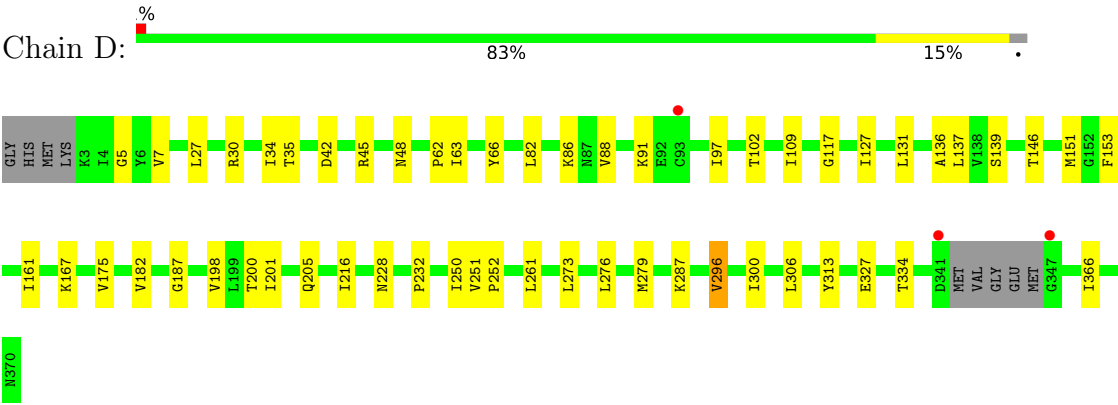
- Molecule 1: Histone deacetylase domain-containing protein



- Molecule 1: Histone deacetylase domain-containing protein



- Molecule 1: Histone deacetylase domain-containing protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	113.95Å 181.69Å 155.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.51 – 2.50 60.51 – 2.50	Depositor EDS
% Data completeness (in resolution range)	97.9 (60.51-2.50) 93.0 (60.51-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.201 , 0.262 0.201 , 0.262	Depositor DCC
R_{free} test set	1981 reflections (3.54%)	wwPDB-VP
Wilson B-factor (Å ²)	17.3	Xtriage
Anisotropy	0.367	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 34.7	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	0.90	EDS
Total number of atoms	10819	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.91% 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: CL, K, 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	0.37	0/2757	0.54	0/3756
1	B	0.37	0/2724	0.56	0/3711
1	C	0.34	0/2769	0.54	0/3772
1	D	0.35	0/2749	0.55	0/3746
All	All	0.36	0/10999	0.55	0/14985

There are no bond length outliers.

There are no bond angle outliers.

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	2693	0	2562	27	0
1	B	2661	0	2526	30	0
1	C	2705	0	2572	28	0
1	D	2686	0	2551	28	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	18	0	0	0	0
5	B	16	0	0	0	0
5	C	13	0	0	1	0
5	D	11	0	0	0	0
All	All	10819	0	10211	108	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 5.

All (108) 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:30:ARG:HB2	1:A:335:LEU:HD21	1.66	0.77
1:B:40:HIS:HD2	1:B:42:ASP:H	1.30	0.77
1:D:42:ASP:OD1	1:D:45:ARG:NH2	2.21	0.74
1:C:39:ASN:H	1:C:39:ASN:HD22	1.42	0.68
1:A:75:ARG:CZ	1:A:169:ILE:HD11	2.25	0.67
1:B:40:HIS:CD2	1:B:42:ASP:H	2.13	0.65
1:D:88:VAL:HA	1:D:91:LYS:HD2	1.80	0.62
1:C:72:ASP:HA	1:C:75:ARG:NH1	2.17	0.60
1:A:75:ARG:NH2	1:A:169:ILE:HD11	2.17	0.59
1:C:97:ILE:HD13	1:C:155:ILE:HD12	1.84	0.58
1:C:335:LEU:HD21	1:D:30:ARG:HB2	1.86	0.58
1:A:276:LEU:HD12	1:A:313:TYR:CD2	2.39	0.57
1:A:56:GLN:HG3	1:A:323:LEU:HD11	1.87	0.56
1:B:272:ILE:HD13	1:B:280:MET:HA	1.87	0.56
1:D:102:THR:HG23	1:D:153:PHE:HA	1.88	0.55
1:C:62:PRO:O	1:C:63:ILE:HD13	2.07	0.55
1:A:300:ILE:HG13	1:A:301:ASN:N	2.21	0.55
1:C:276:LEU:HD12	1:C:313:TYR:CD2	2.42	0.54
1:B:276:LEU:HD12	1:B:313:TYR:CD2	2.42	0.54
1:C:195:ASP:OD1	1:C:197:SER:OG	2.22	0.54
1:D:62:PRO:O	1:D:63:ILE:HD13	2.08	0.54
1:D:287:LYS:NZ	1:D:327:GLU:HG2	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:66:TYR:O	1:C:117:GLY:HA3	2.08	0.53
1:B:66:TYR:O	1:B:117:GLY:HA3	2.09	0.53
1:C:296:VAL:O	1:C:300:ILE:HG12	2.09	0.52
1:D:175:VAL:O	1:D:198:VAL:HA	2.09	0.52
1:B:205:GLN:HA	1:B:279:MET:HE2	1.92	0.51
1:A:196:SER:HB3	1:A:223:ASN:O	2.09	0.51
1:C:7:VAL:HB	1:C:136:ALA:HA	1.93	0.51
1:B:175:VAL:O	1:B:198:VAL:HA	2.10	0.51
1:B:195:ASP:OD1	1:B:197:SER:OG	2.28	0.51
1:B:205:GLN:HB2	1:B:279:MET:HE2	1.92	0.51
1:A:182:VAL:HG11	1:A:277:SER:HB3	1.93	0.50
1:C:205:GLN:HA	1:C:279:MET:HE2	1.94	0.49
1:C:283:SER:OG	1:C:337:ASP:OD2	2.21	0.49
1:C:97:ILE:HG22	1:C:102:THR:HB	1.94	0.49
1:B:335:LEU:HD21	1:A:30:ARG:HB2	1.94	0.49
1:C:269:ASP:OD1	1:C:269:ASP:N	2.46	0.49
1:A:229:LEU:HD13	1:A:365:LEU:HD12	1.94	0.48
1:A:102:THR:HG23	1:A:153:PHE:HA	1.96	0.48
1:C:249:VAL:HA	1:C:366:ILE:HD11	1.94	0.48
1:B:362:CYS:HA	1:B:365:LEU:HD23	1.96	0.47
1:D:5:GLY:HA3	1:D:131:LEU:HD13	1.96	0.47
1:D:276:LEU:HD12	1:D:313:TYR:CD2	2.50	0.47
1:C:75:ARG:NH2	1:C:169:ILE:HD11	2.30	0.47
1:B:132:HIS:NE2	1:B:303:GLY:HA3	2.30	0.47
1:A:175:VAL:O	1:A:198:VAL:HA	2.15	0.47
1:D:97:ILE:CG2	1:D:102:THR:HB	2.43	0.47
1:D:200:THR:OG1	1:D:228:ASN:OD1	2.23	0.47
1:A:151:MET:HE2	1:A:151:MET:HB3	1.82	0.46
1:A:272:ILE:HD13	1:A:280:MET:HB3	1.97	0.46
1:A:201:ILE:HG21	1:A:250:ILE:HG12	1.98	0.46
1:D:66:TYR:O	1:D:117:GLY:HA3	2.15	0.46
1:A:195:ASP:OD1	1:A:197:SER:OG	2.32	0.46
1:B:219:ARG:NH1	1:B:365:LEU:HD12	2.31	0.46
1:B:47:PHE:CZ	1:B:322:GLY:HA3	2.52	0.45
1:C:300:ILE:HG13	1:C:301:ASN:H	1.80	0.45
1:D:7:VAL:HB	1:D:136:ALA:HA	1.97	0.45
1:D:97:ILE:HG22	1:D:102:THR:HB	1.97	0.45
1:B:203:ILE:HD11	1:B:233:LEU:HD11	1.98	0.45
1:B:5:GLY:HA3	1:B:131:LEU:HD13	1.98	0.45
1:A:99:ASP:O	1:A:101:VAL:HG23	2.17	0.44
1:A:332:MET:HG3	1:A:334:THR:HG23	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:182:VAL:HG11	1:B:277:SER:HB3	1.98	0.44
1:B:272:ILE:HD12	1:A:278:ARG:HD2	1.99	0.44
1:C:172:LEU:HD23	1:C:260:GLU:HB2	2.00	0.44
1:B:7:VAL:HG22	1:B:63:ILE:HB	2.00	0.43
1:A:97:ILE:HG22	1:A:102:THR:HB	2.00	0.43
1:B:147:LYS:HG3	1:B:190:ASP:OD2	2.17	0.43
1:B:166:ALA:HA	1:B:170:LEU:HD12	2.00	0.43
1:D:205:GLN:HA	1:D:279:MET:HE2	2.01	0.43
1:A:7:VAL:HB	1:A:136:ALA:HA	2.01	0.43
1:C:203:ILE:HD11	1:C:233:LEU:HD11	2.01	0.43
1:D:296:VAL:O	1:D:300:ILE:HG12	2.18	0.43
1:C:132:HIS:NE2	1:C:303:GLY:HA3	2.34	0.43
1:C:332:MET:HE3	1:C:332:MET:HB3	1.72	0.43
1:C:272:ILE:HD13	1:C:280:MET:HB3	2.01	0.43
1:B:322:GLY:O	1:B:326:ILE:HG12	2.19	0.43
1:A:114:LEU:HD12	1:A:114:LEU:HA	1.89	0.43
1:C:272:ILE:HG22	1:C:317:TYR:CE2	2.54	0.43
1:B:205:GLN:CB	1:B:279:MET:HE2	2.48	0.42
1:C:8:TRP:CG	1:C:9:ASP:N	2.87	0.42
1:C:237:SER:HB3	1:C:241:ALA:HB3	2.00	0.42
1:C:212:ASN:OD1	1:C:212:ASN:N	2.52	0.42
5:C:511:HOH:O	1:D:34:ILE:HD12	2.18	0.42
1:B:332:MET:HE1	1:D:334:THR:HA	2.01	0.42
1:D:167:LYS:HB3	1:D:167:LYS:HE2	1.82	0.42
1:D:273:LEU:HD23	1:D:273:LEU:HA	1.87	0.42
1:A:56:GLN:NE2	1:A:332:MET:O	2.53	0.42
1:D:48:ASN:HB2	1:D:137:LEU:HD23	2.02	0.42
1:B:81:LEU:HD12	1:B:149:ASP:N	2.35	0.42
1:C:99:ASP:CG	1:C:101:VAL:HG12	2.44	0.41
1:B:184:HIS:ND1	1:B:213:SER:OG	2.53	0.41
1:A:312:GLY:HA3	1:A:318:LEU:CD1	2.51	0.41
1:D:34:ILE:HG13	1:D:35:THR:N	2.35	0.41
1:C:97:ILE:CG2	1:C:102:THR:HB	2.51	0.41
1:B:294:LEU:HD11	1:B:329:LEU:HD23	2.02	0.41
1:A:97:ILE:CG2	1:A:102:THR:HB	2.51	0.41
1:A:262:ILE:O	1:A:305:CYS:HA	2.20	0.41
1:D:151:MET:HE2	1:D:151:MET:HB3	1.80	0.41
1:D:252:PRO:HB2	1:D:366:ILE:HD13	2.02	0.41
1:B:296:VAL:O	1:B:300:ILE:HG12	2.20	0.40
1:B:368:ASP:OD1	1:B:368:ASP:N	2.54	0.40
1:D:82:LEU:O	1:D:86:LYS:HG3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:201:ILE:HG21	1:D:250:ILE:HG12	2.03	0.40
1:A:47:PHE:C	1:A:47:PHE:CD1	3.00	0.40
1:D:146:THR:HG22	1:D:187:GLY:HA2	2.02	0.40
1:D:127:ILE:HD11	1:D:261:LEU:HD13	2.03	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	361/372 (97%)	338 (94%)	23 (6%)	0	100	100
1	B	354/372 (95%)	339 (96%)	13 (4%)	2 (1%)	21	38
1	C	361/372 (97%)	340 (94%)	21 (6%)	0	100	100
1	D	359/372 (96%)	345 (96%)	13 (4%)	1 (0%)	36	55
All	All	1435/1488 (96%)	1362 (95%)	70 (5%)	3 (0%)	43	63

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	98	GLY
1	B	91	LYS
1	D	232	PRO

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	272/306 (89%)	269 (99%)	3 (1%)	65	84
1	B	268/306 (88%)	262 (98%)	6 (2%)	45	73
1	C	274/306 (90%)	264 (96%)	10 (4%)	31	58
1	D	272/306 (89%)	263 (97%)	9 (3%)	33	61
All	All	1086/1224 (89%)	1058 (97%)	28 (3%)	40	68

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	27	LEU
1	B	73	ILE
1	B	101	VAL
1	B	291	SER
1	B	306	LEU
1	B	365	LEU
1	A	99	ASP
1	A	182	VAL
1	A	306	LEU
1	C	28	SER
1	C	39	ASN
1	C	97	ILE
1	C	139	SER
1	C	207	LYS
1	C	211	THR
1	C	229	LEU
1	C	269	ASP
1	C	271	SER
1	C	365	LEU
1	D	27	LEU
1	D	109	ILE
1	D	139	SER
1	D	161	ILE
1	D	182	VAL
1	D	216	ILE
1	D	251	VAL
1	D	296	VAL
1	D	306	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	40	HIS
1	B	217	ASN
1	B	316	HIS
1	B	349	ASN
1	A	103	HIS
1	A	158	ASN
1	A	206	ASN
1	A	212	ASN
1	C	39	ASN
1	C	84	ASN
1	C	87	ASN
1	C	103	HIS
1	C	106	ASN
1	D	56	GLN
1	D	206	ASN
1	D	349	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 ⓘ

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 [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	365/372 (98%)	0.00	10 (2%) 56 51	9, 20, 37, 54	0
1	B	360/372 (96%)	0.04	7 (1%) 66 62	9, 20, 39, 60	0
1	C	365/372 (98%)	0.14	12 (3%) 49 45	10, 22, 40, 52	0
1	D	363/372 (97%)	-0.03	3 (0%) 82 80	11, 20, 35, 66	0
All	All	1453/1488 (97%)	0.04	32 (2%) 62 58	9, 21, 38, 66	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	340	ILE	4.7
1	C	93	CYS	4.3
1	A	93	CYS	4.2
1	B	19	THR	4.0
1	D	341	ASP	3.9
1	B	92	GLU	3.8
1	D	93	CYS	3.8
1	A	346	MET	3.4
1	A	344	GLY	3.4
1	B	100	ARG	3.4
1	B	98	GLY	3.4
1	A	19	THR	3.4
1	C	345	GLU	3.1
1	C	100	ARG	3.1
1	C	346	MET	2.9
1	A	345	GLU	2.7
1	D	347	GLY	2.6
1	C	35	THR	2.6
1	B	347	GLY	2.6
1	C	19	THR	2.6
1	C	101	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	340	ILE	2.5
1	B	35	THR	2.5
1	B	18	GLY	2.4
1	A	35	THR	2.4
1	A	96	ASP	2.4
1	C	87	ASN	2.3
1	C	300	ILE	2.3
1	C	92	GLU	2.2
1	A	100	ARG	2.2
1	C	98	GLY	2.1
1	A	211	THR	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	ZN	B	401	1/1	0.82	0.13	70,70,70,70	0
2	ZN	D	401	1/1	0.82	0.14	80,80,80,80	0
2	ZN	A	401	1/1	0.84	0.14	74,74,74,74	0
2	ZN	C	401	1/1	0.89	0.10	65,65,65,65	0
4	CL	A	404	1/1	0.97	0.19	6,6,6,6	0
3	K	A	403	1/1	0.98	0.03	22,22,22,22	0
3	K	C	402	1/1	0.98	0.09	25,25,25,25	0
3	K	C	403	1/1	0.98	0.03	23,23,23,23	0
3	K	D	403	1/1	0.98	0.03	21,21,21,21	0
3	K	B	403	1/1	0.98	0.05	23,23,23,23	0
3	K	D	402	1/1	0.99	0.13	29,29,29,29	0
3	K	A	402	1/1	0.99	0.07	23,23,23,23	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	CL	B	404	1/1	0.99	0.12	0,0,0,0	0
3	K	B	402	1/1	0.99	0.11	30,30,30,30	0
4	CL	C	404	1/1	0.99	0.15	0,0,0,0	0
4	CL	D	404	1/1	0.99	0.18	0,0,0,0	0

6.5 Other polymers [i](#)

There are no such residues in this entry.