Full wwPDB X-ray Structure Validation Report ⓘ

Aug 6, 2023 – 08:43 PM EDT

PDB ID	:	1KHN
Title	:	E. COLI ALKALINE PHOSPHATASE MUTANT (D153HD330N) ZINC FORM
Authors	:	Le Du, M.H.; Lamoure, C.; Muller, B.H.; Bulgakov, O.V.; Lajeunesse, E.; Menez, A.; Boulain, J.C.
Deposited on	:	2001-11-30
Resolution	:	2.60 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:

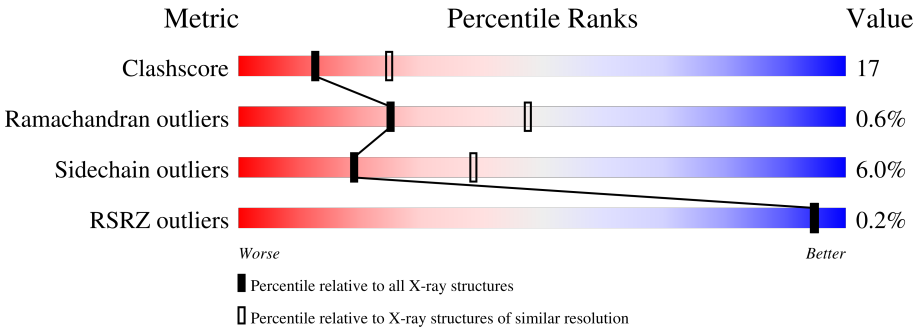
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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 <=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	449	66% 31%
1	B	449	68% 28%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 6837 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 Alkaline phosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	444	Total	C	N	O	S	0	0	0
			3262	2015	579	657	11			
1	B	444	Total	C	N	O	S	0	0	0
			3258	2013	579	655	11			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	15	ASN	ASP	conflict	UNP P00634
A	35	ASN	ASP	conflict	UNP P00634
A	153	HIS	ASP	engineered mutation	UNP P00634
A	176	GLN	GLU	conflict	UNP P00634
A	228	GLU	GLN	conflict	UNP P00634
A	230	GLU	GLN	conflict	UNP P00634
A	330	ASN	ASP	engineered mutation	UNP P00634
B	15	ASN	ASP	conflict	UNP P00634
B	35	ASN	ASP	conflict	UNP P00634
B	153	HIS	ASP	engineered mutation	UNP P00634
B	176	GLN	GLU	conflict	UNP P00634
B	228	GLU	GLN	conflict	UNP P00634
B	230	GLU	GLN	conflict	UNP P00634
B	330	ASN	ASP	engineered mutation	UNP P00634

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 3 is water.

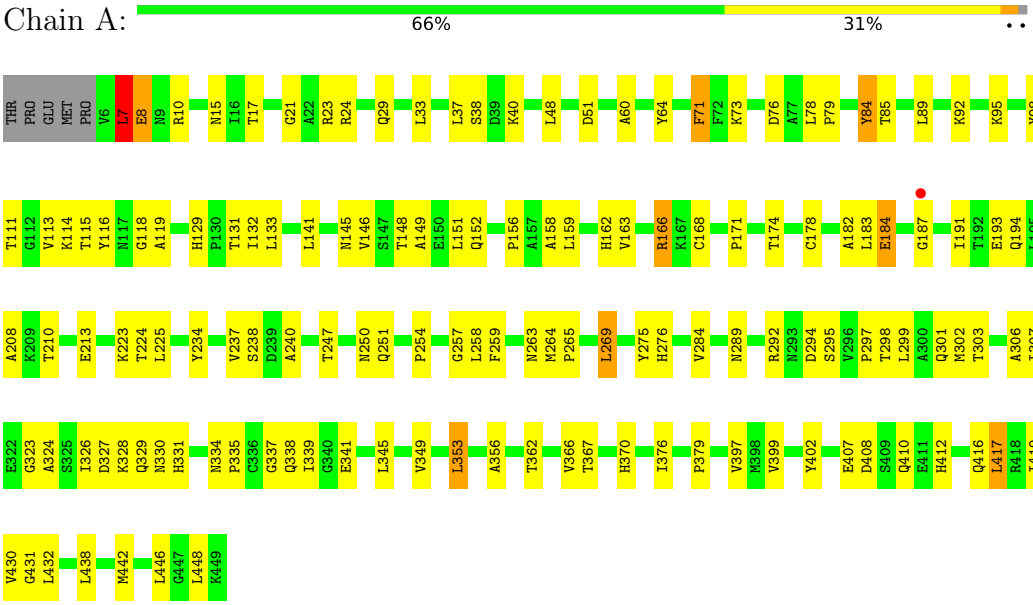
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	181	Total O 181 181	0	0
3	B	130	Total O 130 130	0	0

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.

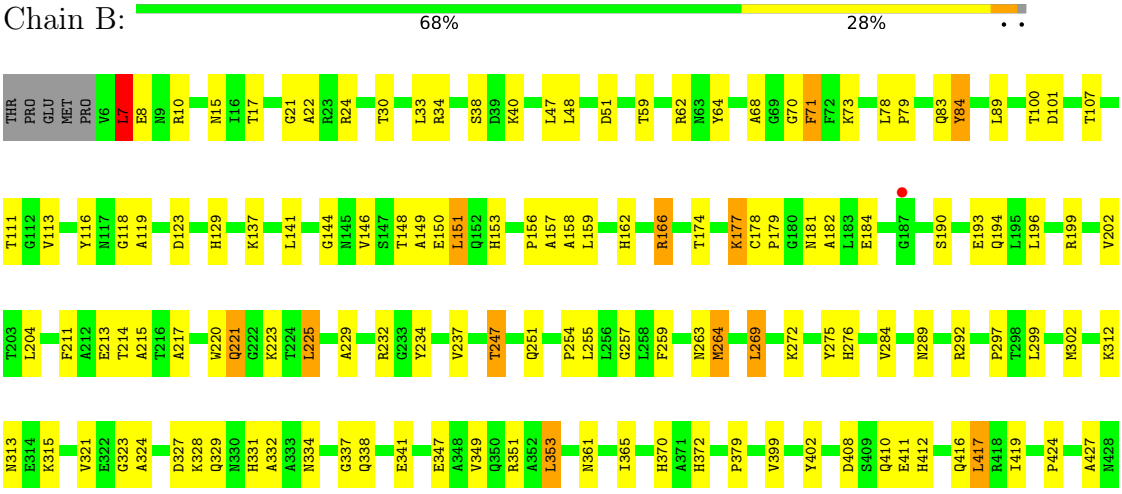
• Molecule 1: Alkaline phosphatase

Chain A:



• Molecule 1: Alkaline phosphatase

Chain B:




V429
L432
T433
D434
D437
Y440
L448
K449

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, α, β, γ	163.51Å 163.51Å 138.03Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 2.60 19.85 – 2.20	Depositor EDS
% Data completeness (in resolution range)	96.3 (10.00-2.60) 54.4 (19.85-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	0.00 (at 2.19Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.207 , 0.248 0.196 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	40.6	Xtriage
Anisotropy	0.714	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 90.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	6837	wwPDB-VP
Average B, all atoms (Å^2)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.50% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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

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	0.47	0/3316	0.76	3/4501 (0.1%)
1	B	0.48	0/3312	0.76	2/4496 (0.0%)
All	All	0.47	0/6628	0.76	5/8997 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	323	GLY	N-CA-C	-6.11	97.82	113.10
1	A	7	LEU	CA-CB-CG	5.86	128.77	115.30
1	A	417	LEU	CA-CB-CG	5.65	128.30	115.30
1	B	417	LEU	CA-CB-CG	5.62	128.23	115.30
1	B	323	GLY	N-CA-C	-5.14	100.24	113.10

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	3262	0	3200	111	0
1	B	3258	0	3196	118	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	181	0	0	5	0
3	B	130	0	0	4	0
All	All	6837	0	6396	216	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 17.

All (216) 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)
3:A:1267:HOH:O	1:B:7:LEU:HD11	1.78	0.82
1:B:111:THR:HG22	1:B:113:VAL:HG12	1.62	0.82
1:A:48:LEU:HD13	1:A:321:VAL:HB	1.60	0.82
1:A:149:ALA:HA	1:A:263:ASN:HD22	1.45	0.81
1:B:111:THR:CG2	1:B:113:VAL:HG12	2.11	0.80
1:B:38:SER:OG	1:B:40:LYS:HG2	1.83	0.78
1:A:111:THR:CG2	1:A:113:VAL:HG12	2.12	0.78
1:B:7:LEU:HD13	1:B:10:ARG:HG3	1.66	0.76
1:B:48:LEU:HD13	1:B:321:VAL:HB	1.67	0.76
1:A:10:ARG:HB2	1:A:71:PHE:CD1	2.22	0.74
1:A:148:THR:HA	1:A:264:MET:HE2	1.71	0.73
1:B:10:ARG:O	1:B:24:ARG:HD3	1.88	0.72
1:B:419:ILE:HG13	1:B:429:VAL:HG13	1.70	0.72
1:A:334:ASN:HD22	1:A:337:GLY:H	1.38	0.72
1:A:111:THR:HG21	1:A:113:VAL:HG12	1.71	0.71
1:A:213:GLU:O	1:A:225:LEU:HD12	1.89	0.71
1:A:7:LEU:HB3	1:A:10:ARG:NE	2.06	0.71
1:B:149:ALA:HA	1:B:263:ASN:HD22	1.56	0.70
1:A:269:LEU:HD23	1:A:289:ASN:ND2	2.08	0.69
1:B:334:ASN:HD22	1:B:337:GLY:H	1.40	0.68
1:A:337:GLY:O	1:A:341:GLU:HG2	1.93	0.67
1:A:240:ALA:HA	1:A:297:PRO:HG3	1.77	0.67
1:B:15:ASN:O	1:B:21:GLY:HA3	1.96	0.66
1:B:118:GLY:HA2	1:B:166:ARG:HG2	1.78	0.65
1:B:10:ARG:HB2	1:B:71:PHE:CD1	2.33	0.64
1:A:10:ARG:HB2	1:A:71:PHE:HD1	1.63	0.64
1:A:111:THR:HG22	1:A:113:VAL:HG12	1.81	0.63
1:B:7:LEU:HD13	1:B:10:ARG:CG	2.28	0.62
1:A:328:LYS:HE2	3:A:1235:HOH:O	1.99	0.62
1:B:137:LYS:HG2	3:B:1144:HOH:O	2.01	0.61
1:A:264:MET:HE1	1:A:302:MET:SD	2.41	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98:TYR:HE1	1:B:68:ALA:HB2	1.65	0.60
1:B:7:LEU:CD1	1:B:10:ARG:HG3	2.32	0.60
1:B:193:GLU:HG2	1:B:225:LEU:HD21	1.82	0.60
1:A:10:ARG:HD2	1:B:432:LEU:O	2.02	0.60
1:A:141:LEU:HD21	1:A:446:LEU:HD22	1.82	0.60
1:B:419:ILE:CG1	1:B:429:VAL:HG13	2.32	0.60
1:A:116:TYR:CZ	1:A:119:ALA:HB2	2.38	0.58
1:A:98:TYR:CE1	1:B:68:ALA:HB2	2.39	0.58
1:A:107:THR:O	1:A:111:THR:HB	2.04	0.58
1:B:199:ARG:HD3	1:B:251:GLN:NE2	2.19	0.58
1:B:292:ARG:NH2	1:B:297:PRO:O	2.38	0.57
3:A:1006:HOH:O	1:B:123:ASP:HB2	2.05	0.57
1:A:419:ILE:HG13	1:A:429:VAL:HG13	1.87	0.57
1:A:334:ASN:ND2	1:A:337:GLY:H	2.02	0.57
1:B:89:LEU:O	1:B:116:TYR:HA	2.05	0.56
1:A:38:SER:OG	1:A:40:LYS:HG2	2.06	0.56
1:B:190:SER:OG	1:B:193:GLU:HG3	2.05	0.56
1:B:334:ASN:ND2	1:B:337:GLY:H	2.03	0.56
1:B:129:HIS:O	1:B:162:HIS:CE1	2.59	0.56
1:A:7:LEU:HD21	1:A:79:PRO:HA	1.88	0.56
1:A:146:VAL:HG23	1:A:302:MET:HE3	1.87	0.56
1:B:51:ASP:O	1:B:327:ASP:HB2	2.06	0.56
1:B:84:TYR:HD1	1:B:84:TYR:O	1.87	0.56
1:B:111:THR:HG21	1:B:113:VAL:HG12	1.88	0.56
1:A:335:PRO:O	1:A:339:ILE:HG13	2.05	0.55
1:A:208:ALA:HB2	1:A:258:LEU:HB3	1.89	0.55
1:A:7:LEU:HB3	1:A:10:ARG:HE	1.71	0.55
1:A:370:HIS:CE1	1:A:412:HIS:CE1	2.95	0.55
1:A:275:TYR:O	1:A:276:HIS:HB2	2.07	0.55
1:A:33:LEU:HD12	1:B:427:ALA:HB1	1.89	0.54
1:A:116:TYR:CE1	1:A:119:ALA:HB2	2.43	0.54
1:B:419:ILE:HG13	1:B:429:VAL:CG1	2.38	0.54
1:A:292:ARG:NH2	1:A:297:PRO:O	2.42	0.53
1:A:298:THR:OG1	1:A:301:GLN:HG3	2.07	0.53
1:B:264:MET:HE1	1:B:302:MET:SD	2.47	0.53
1:A:129:HIS:O	1:A:162:HIS:CE1	2.62	0.53
1:B:247:THR:O	1:B:312:LYS:NZ	2.41	0.53
1:A:345:LEU:O	1:A:349:VAL:HG23	2.09	0.53
1:A:199:ARG:HD3	1:A:251:GLN:NE2	2.24	0.52
1:A:51:ASP:O	1:A:327:ASP:HB2	2.10	0.52
1:A:432:LEU:O	1:B:10:ARG:HD2	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:129:HIS:O	1:B:162:HIS:HE1	1.92	0.52
1:B:370:HIS:CE1	1:B:412:HIS:CE1	2.98	0.52
1:B:313:ASN:OD1	1:B:315:LYS:HE2	2.10	0.52
1:A:338:GLN:NE2	1:A:402:TYR:OH	2.43	0.52
1:B:370:HIS:CE1	1:B:412:HIS:ND1	2.79	0.51
1:A:10:ARG:O	1:A:24:ARG:HD3	2.11	0.51
1:A:48:LEU:HB2	1:A:366:VAL:HG22	1.91	0.51
1:A:427:ALA:HB1	1:B:33:LEU:HD12	1.93	0.51
1:B:40:LYS:NZ	1:B:361:ASN:HD22	2.09	0.51
1:B:269:LEU:HD23	1:B:289:ASN:ND2	2.26	0.50
1:A:7:LEU:HD22	1:A:10:ARG:CZ	2.42	0.50
1:A:15:ASN:O	1:A:21:GLY:HA3	2.10	0.50
1:A:299:LEU:HD11	1:A:321:VAL:CG1	2.42	0.50
1:A:367:THR:HB	1:A:419:ILE:HD12	1.92	0.50
1:B:47:LEU:HD12	1:B:365:ILE:HB	1.93	0.50
1:B:229:ALA:O	1:B:234:TYR:HB2	2.11	0.50
1:B:182:ALA:HB1	1:B:184:GLU:OE1	2.12	0.50
1:B:100:THR:OG1	1:B:101:ASP:N	2.46	0.49
1:B:379:PRO:HA	1:B:399:VAL:HG21	1.94	0.49
1:B:149:ALA:HB2	1:B:324:ALA:CB	2.43	0.49
1:B:148:THR:HA	1:B:264:MET:HE3	1.95	0.49
1:B:100:THR:HG22	3:B:1010:HOH:O	2.11	0.49
1:B:177:LYS:C	1:B:179:PRO:HD3	2.33	0.48
1:A:299:LEU:HD11	1:A:321:VAL:HG13	1.95	0.48
1:B:40:LYS:O	1:B:424:PRO:HB3	2.13	0.48
1:A:174:THR:HG23	1:A:178:CYS:HB2	1.95	0.48
1:B:62:ARG:HD2	3:B:1053:HOH:O	2.13	0.48
1:B:337:GLY:O	1:B:341:GLU:HG2	2.12	0.48
1:A:132:ILE:HD13	1:A:442:MET:HB2	1.96	0.48
1:A:199:ARG:HA	1:A:234:TYR:OH	2.14	0.48
1:A:328:LYS:HG3	3:A:1235:HOH:O	2.14	0.48
1:B:178:CYS:N	1:B:179:PRO:HD3	2.29	0.48
1:B:263:ASN:HD21	1:B:328:LYS:NZ	2.11	0.48
1:A:265:PRO:O	1:A:292:ARG:NH1	2.47	0.48
1:A:295:SER:O	1:A:297:PRO:HD3	2.13	0.48
1:A:145:ASN:HD21	1:A:156:PRO:HB2	1.78	0.47
1:B:347:GLU:O	1:B:351:ARG:HG2	2.14	0.47
1:B:213:GLU:O	1:B:225:LEU:HD12	2.15	0.47
1:B:313:ASN:OD1	1:B:315:LYS:HB2	2.13	0.47
1:A:264:MET:CE	1:A:302:MET:SD	3.02	0.47
1:A:168:CYS:O	1:A:191:ILE:HG13	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:ARG:HD2	1:B:440:TYR:CD2	2.50	0.47
1:A:23:ARG:HD2	1:B:440:TYR:CE2	2.50	0.47
1:A:118:GLY:HA2	1:A:166:ARG:HG2	1.96	0.47
1:A:149:ALA:HB2	1:A:324:ALA:CB	2.44	0.47
1:A:303:THR:O	1:A:307:ILE:HG13	2.15	0.47
1:B:144:GLY:HA2	1:B:202:VAL:O	2.14	0.47
1:B:214:THR:HA	1:B:223:LYS:O	2.14	0.47
1:B:338:GLN:NE2	1:B:402:TYR:OH	2.48	0.47
1:A:183:LEU:HD12	1:A:187:GLY:O	2.15	0.47
1:A:370:HIS:CE1	1:A:412:HIS:ND1	2.83	0.47
1:A:379:PRO:HA	1:A:399:VAL:HG21	1.97	0.47
1:A:199:ARG:HD3	1:A:251:GLN:HE22	1.80	0.46
1:A:110:SER:O	1:A:131:THR:HA	2.15	0.46
1:A:10:ARG:HH22	1:A:29:GLN:CD	2.18	0.46
1:A:37:LEU:HD22	1:A:427:ALA:HB2	1.97	0.46
1:A:416:GLN:HG2	1:B:59:THR:OG1	2.16	0.46
1:B:7:LEU:HD22	1:B:10:ARG:CZ	2.45	0.46
1:A:306:ALA:O	1:A:310:LEU:HG	2.15	0.46
1:B:8:GLU:O	1:B:71:PHE:CZ	2.69	0.46
1:B:48:LEU:HG	1:B:349:VAL:HG22	1.97	0.46
1:A:237:VAL:O	1:A:257:GLY:HA2	2.16	0.46
1:A:329:GLN:HE21	1:A:334:ASN:HB3	1.81	0.46
1:B:24:ARG:NH1	1:B:70:GLY:O	2.48	0.46
1:B:83:GLN:HG2	1:B:416:GLN:HE21	1.81	0.46
1:B:217:ALA:O	1:B:221:GLN:HB3	2.16	0.46
1:A:107:THR:HG21	1:A:115:THR:HG21	1.97	0.46
1:B:275:TYR:O	1:B:276:HIS:HB2	2.16	0.46
1:B:116:TYR:CZ	1:B:119:ALA:HB2	2.52	0.45
1:B:149:ALA:HA	1:B:263:ASN:ND2	2.29	0.45
1:B:149:ALA:CA	1:B:263:ASN:HD22	2.25	0.45
1:B:148:THR:HG23	1:B:299:LEU:HD13	1.99	0.45
1:A:133:LEU:O	1:A:133:LEU:HD23	2.17	0.45
1:B:84:TYR:CD1	1:B:84:TYR:C	2.90	0.45
1:A:89:LEU:HD13	1:B:17:THR:HG22	1.99	0.45
1:A:95:LYS:HD3	1:A:95:LYS:HA	1.81	0.45
1:B:153:HIS:O	1:B:156:PRO:HD2	2.17	0.45
1:B:190:SER:O	1:B:194:GLN:HG3	2.16	0.44
1:B:199:ARG:HD3	1:B:251:GLN:HE22	1.80	0.44
1:B:237:VAL:O	1:B:257:GLY:HA2	2.17	0.44
1:B:107:THR:O	1:B:111:THR:HB	2.17	0.44
1:A:149:ALA:CA	1:A:263:ASN:HD22	2.23	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:326:ILE:O	1:A:330:ASN:HB2	2.17	0.44
1:B:146:VAL:HG22	1:B:321:VAL:HG13	2.00	0.44
1:A:84:TYR:HD1	1:A:84:TYR:O	2.00	0.44
1:A:250:ASN:O	1:A:254:PRO:HG3	2.18	0.44
1:B:272:LYS:HE2	1:B:332:ALA:O	2.18	0.44
1:A:110:SER:HB3	1:A:159:LEU:HD23	2.00	0.44
1:A:193:GLU:HG3	1:A:225:LEU:HD21	2.00	0.44
1:B:30:THR:O	1:B:34:ARG:HB2	2.18	0.44
1:A:234:TYR:CD2	1:A:254:PRO:HB2	2.53	0.43
1:B:7:LEU:HD22	1:B:10:ARG:NH2	2.33	0.43
1:B:331:HIS:ND1	1:B:410:GLN:O	2.49	0.43
1:A:84:TYR:CD1	1:A:84:TYR:C	2.92	0.43
1:A:17:THR:HG22	1:B:89:LEU:HD13	1.99	0.43
1:A:162:HIS:H	1:A:194:GLN:HE22	1.66	0.43
1:B:234:TYR:CD2	1:B:254:PRO:HB2	2.53	0.43
1:B:202:VAL:HG13	1:B:255:LEU:HD23	2.01	0.43
1:B:204:LEU:HB3	1:B:259:PHE:HE2	1.84	0.43
1:B:211:PHE:HA	1:B:225:LEU:HB2	2.01	0.43
1:B:157:ALA:O	1:B:159:LEU:N	2.52	0.42
1:B:263:ASN:ND2	1:B:328:LYS:NZ	2.67	0.42
1:B:379:PRO:HA	1:B:399:VAL:CG2	2.48	0.42
1:B:220:TRP:CZ2	1:B:232:ARG:HD3	2.54	0.42
1:A:60:ALA:HB2	1:A:376:ILE:HD11	2.01	0.42
1:B:199:ARG:HA	1:B:234:TYR:OH	2.19	0.42
1:B:372:HIS:HA	1:B:412:HIS:HA	2.01	0.42
1:B:411:GLU:HG2	3:B:1152:HOH:O	2.19	0.42
1:A:129:HIS:O	1:A:162:HIS:HE1	1.99	0.42
1:B:193:GLU:HG2	1:B:225:LEU:CD2	2.47	0.42
1:B:193:GLU:CG	1:B:225:LEU:HD21	2.49	0.42
1:A:379:PRO:HA	1:A:399:VAL:CG2	2.49	0.42
1:A:24:ARG:HG3	1:B:434:ASP:HB2	2.02	0.42
1:A:240:ALA:CA	1:A:297:PRO:HG3	2.45	0.42
1:B:329:GLN:HE21	1:B:334:ASN:HB3	1.84	0.42
1:B:141:LEU:HD12	1:B:315:LYS:HB3	2.01	0.42
1:A:331:HIS:ND1	1:A:410:GLN:O	2.53	0.42
1:B:150:GLU:H	1:B:150:GLU:HG2	1.68	0.41
1:A:89:LEU:O	1:A:116:TYR:HA	2.20	0.41
1:A:7:LEU:HD12	1:A:76:ASP:O	2.20	0.41
1:A:131:THR:HG22	3:A:1226:HOH:O	2.20	0.41
1:A:182:ALA:HB1	1:A:184:GLU:OE1	2.20	0.41
1:A:238:SER:C	1:A:259:PHE:HD1	2.24	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:430:VAL:HG12	1:A:431:GLY:N	2.34	0.41
1:B:174:THR:HG22	1:B:178:CYS:O	2.20	0.41
1:B:434:ASP:O	1:B:437:ASP:HB2	2.21	0.41
1:A:48:LEU:HG	1:A:349:VAL:HG22	2.02	0.41
1:A:64:TYR:CD1	1:A:64:TYR:C	2.93	0.41
1:A:213:GLU:O	1:A:224:THR:HA	2.20	0.41
1:A:438:LEU:O	1:A:442:MET:HG3	2.20	0.41
1:B:151:LEU:HD13	1:B:151:LEU:HA	1.90	0.41
1:B:269:LEU:HD23	1:B:289:ASN:HD22	1.86	0.41
1:A:152:GLN:NE2	1:A:210:THR:HB	2.36	0.41
1:B:64:TYR:CD1	1:B:64:TYR:C	2.94	0.41
1:A:171:PRO:HD2	1:A:213:GLU:OE1	2.21	0.40
1:A:356:ALA:HA	1:A:362:THR:HB	2.02	0.40
1:B:78:LEU:HA	1:B:79:PRO:HD2	1.77	0.40
1:B:178:CYS:HB3	1:B:181:ASN:OD1	2.21	0.40
1:A:78:LEU:HA	1:A:79:PRO:HD2	1.82	0.40
1:B:215:ALA:HB3	1:B:221:GLN:HA	2.04	0.40
1:B:7:LEU:HB3	1:B:8:GLU:H	1.56	0.40
1:A:114:LYS:HE2	1:B:22:ALA:HB1	2.03	0.40
1:A:353:LEU:HA	1:A:353:LEU:HD12	1.84	0.40
1:B:353:LEU:HD12	1:B:353:LEU:HA	1.88	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	442/449 (98%)	417 (94%)	22 (5%)	3 (1%)	22	43
1	B	442/449 (98%)	415 (94%)	25 (6%)	2 (0%)	29	52
All	All	884/898 (98%)	832 (94%)	47 (5%)	5 (1%)	25	47

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	7	LEU
1	A	8	GLU
1	B	158	ALA
1	A	92	LYS
1	A	158	ALA

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	333/340 (98%)	311 (93%)	22 (7%)	16	33
1	B	332/340 (98%)	314 (95%)	18 (5%)	22	44
All	All	665/680 (98%)	625 (94%)	40 (6%)	19	39

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LEU
1	A	8	GLU
1	A	71	PHE
1	A	73	LYS
1	A	84	TYR
1	A	85	THR
1	A	151	LEU
1	A	163	VAL
1	A	166	ARG
1	A	184	GLU
1	A	196	LEU
1	A	223	LYS
1	A	247	THR
1	A	269	LEU
1	A	284	VAL
1	A	294	ASP
1	A	353	LEU
1	A	397	VAL

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Mol	Chain	Res	Type
1	A	407	GLU
1	A	408	ASP
1	A	417	LEU
1	A	448	LEU
1	B	7	LEU
1	B	71	PHE
1	B	73	LYS
1	B	84	TYR
1	B	151	LEU
1	B	166	ARG
1	B	177	LYS
1	B	196	LEU
1	B	221	GLN
1	B	225	LEU
1	B	247	THR
1	B	264	MET
1	B	269	LEU
1	B	284	VAL
1	B	353	LEU
1	B	408	ASP
1	B	417	LEU
1	B	448	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	ASN
1	A	83	GLN
1	A	145	ASN
1	A	194	GLN
1	A	235	GLN
1	A	251	GLN
1	A	263	ASN
1	A	289	ASN
1	A	329	GLN
1	A	334	ASN
1	A	338	GLN
1	A	361	ASN
1	B	15	ASN
1	B	83	GLN
1	B	194	GLN
1	B	221	GLN

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Mol	Chain	Res	Type
1	B	235	GLN
1	B	251	GLN
1	B	252	GLN
1	B	263	ASN
1	B	289	ASN
1	B	329	GLN
1	B	334	ASN
1	B	338	GLN
1	B	361	ASN
1	B	375	GLN
1	B	410	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 monosaccharides 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

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, 95^{th} 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	444/449 (98%)	-0.58	1 (0%)	95	95	8, 30, 65, 98	0
1	B	444/449 (98%)	-0.55	1 (0%)	95	95	10, 32, 65, 87	0
All	All	888/898 (98%)	-0.56	2 (0%)	95	95	8, 31, 65, 98	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	187	GLY	2.8
1	A	187	GLY	2.3

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 monosaccharides 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, 95^{th} 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	450	1/1	0.96	0.05	31,31,31,31	0
2	ZN	B	451	1/1	0.96	0.04	23,23,23,23	0
2	ZN	B	452	1/1	0.96	0.06	32,32,32,32	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	ZN	A	452	1/1	0.97	0.04	27,27,27,27	0
2	ZN	A	451	1/1	0.97	0.06	24,24,24,24	0
2	ZN	A	450	1/1	0.99	0.04	27,27,27,27	0

6.5 Other polymers (i)

There are no such residues in this entry.