



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:48 AM UTC

PDB ID : 1PTM / pdb_00001ptm
Title : Crystal structure of E.coli PdxA
Authors : Sivaraman, J.; Li, Y.; Banks, J.; Cane, D.E.; Matte, A.; Cygler, M.; Montreal-Kingston Bacterial Structural Genomics Initiative (BSGI)
Deposited on : 2003-06-23
Resolution : 1.96 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

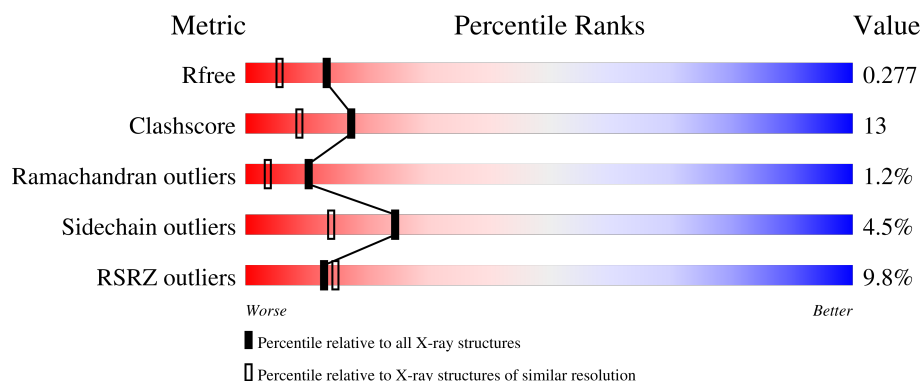
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.96 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	329	 6% 74% 21% 5%
1	B	329	 13% 77% 19% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5129 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 4-hydroxythreonine-4-phosphate dehydrogenase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	329	Total	C	N	O	S	Se	0	0	0
			2444	1554	427	451	4	8			
1	B	328	Total	C	N	O	S	Se	0	0	0
			2415	1536	423	445	4	7			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	UNP P19624
A	49	MSE	MET	modified residue	UNP P19624
A	151	MSE	MET	modified residue	UNP P19624
A	152	MSE	MET	modified residue	UNP P19624
A	218	MSE	MET	modified residue	UNP P19624
A	238	MSE	MET	modified residue	UNP P19624
A	264	MSE	MET	modified residue	UNP P19624
A	324	MSE	MET	modified residue	UNP P19624
B	1	MSE	MET	modified residue	UNP P19624
B	49	MSE	MET	modified residue	UNP P19624
B	151	MSE	MET	modified residue	UNP P19624
B	152	MSE	MET	modified residue	UNP P19624
B	218	MSE	MET	modified residue	UNP P19624
B	238	MSE	MET	modified residue	UNP P19624
B	264	MSE	MET	modified residue	UNP P19624
B	324	MSE	MET	modified residue	UNP P19624

- 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 PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		

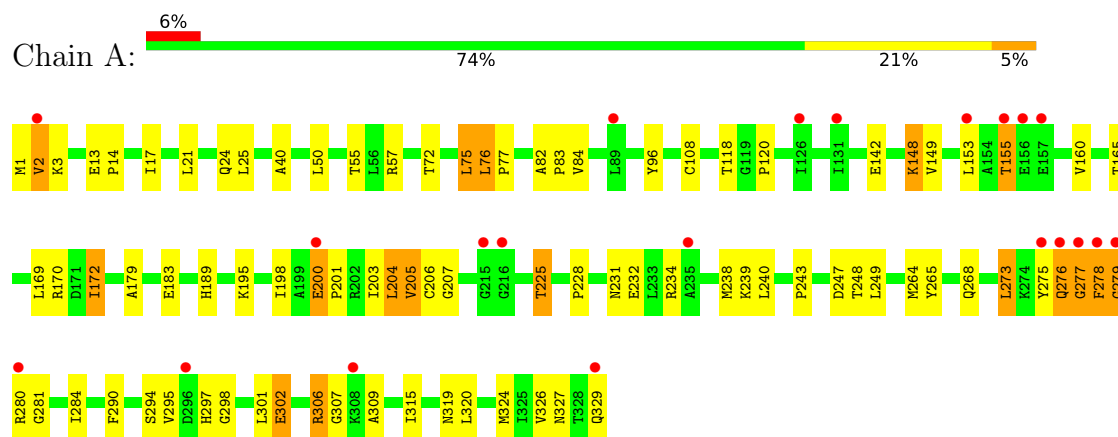
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	132	Total	O	0	0
			132	132		
4	B	131	Total	O	0	0
			131	131		

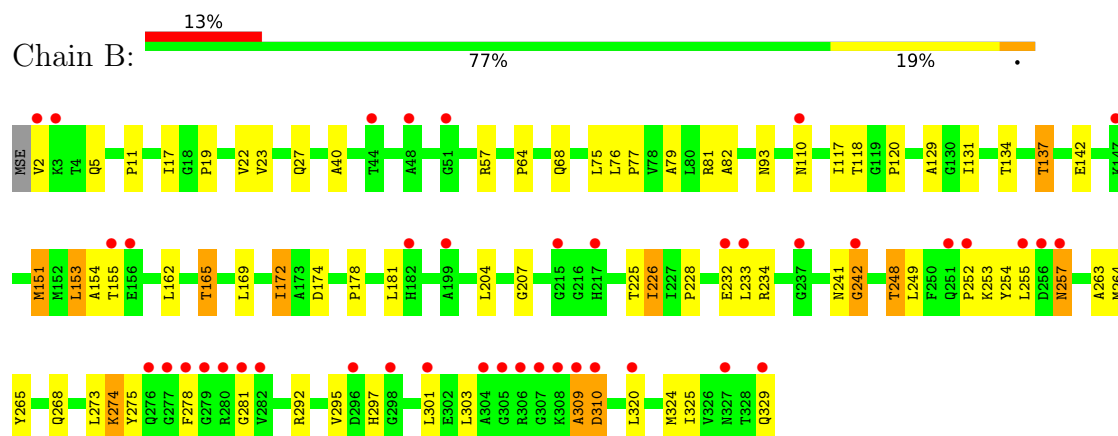
3 Residue-property plots [i](#)

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: 4-hydroxythreonine-4-phosphate dehydrogenase



- Molecule 1: 4-hydroxythreonine-4-phosphate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	75.48Å 79.23Å 114.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.00 – 1.96 45.00 – 1.96	Depositor EDS
% Data completeness (in resolution range)	(Not available) (45.00-1.96) 98.2 (45.00-1.96)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.63 (at 1.95Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.217 , 0.267 0.226 , 0.277	Depositor DCC
R_{free} test set	2668 reflections (2.91%)	wwPDB-VP
Wilson B-factor (Å ²)	18.9	Xtriage
Anisotropy	0.298	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 38.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.012 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5129	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.27% 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: PO4, 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	0.52	1/2484 (0.0%)	1.03	12/3380 (0.4%)
1	B	0.53	0/2456	1.04	13/3344 (0.4%)
All	All	0.53	1/4940 (0.0%)	1.03	25/6724 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1	MSE	SE-CE	-5.26	1.79	1.95

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	172	ILE	N-CA-C	9.02	119.08	110.42
1	B	77	PRO	N-CA-C	7.09	121.95	111.03
1	A	172	ILE	N-CA-C	7.04	117.17	110.42
1	A	279	GLY	N-CA-C	-6.75	97.19	113.18
1	B	131	ILE	N-CA-C	6.43	114.98	107.77
1	A	205	VAL	N-CA-C	6.37	117.09	108.17
1	B	207	GLY	N-CA-C	-6.29	104.13	112.82
1	B	226	ILE	N-CA-C	6.01	116.19	110.42
1	A	77	PRO	N-CA-C	5.95	120.42	111.14
1	A	207	GLY	N-CA-C	-5.87	103.16	112.66
1	A	225	THR	N-CA-C	5.67	119.72	112.12
1	B	248	THR	N-CA-C	-5.62	107.07	114.04
1	B	281	GLY	N-CA-C	5.60	117.91	111.36
1	A	165	THR	N-CA-C	5.53	116.93	108.30
1	B	129	ALA	N-CA-C	-5.52	106.52	113.20
1	B	155	THR	N-CA-C	-5.48	99.13	110.80
1	B	79	ALA	N-CA-C	5.45	119.11	109.96
1	A	155	THR	N-CA-C	-5.44	99.21	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(^o)	Ideal(^o)
1	B	118	THR	N-CA-C	5.39	118.60	109.76
1	A	84	VAL	N-CA-C	5.39	116.34	108.53
1	A	76	LEU	N-CA-C	-5.21	98.23	109.01
1	A	232	GLU	N-CA-C	-5.12	105.70	111.28
1	B	76	LEU	N-CA-C	-5.03	98.59	109.01
1	B	165	THR	N-CA-C	5.03	116.15	108.30
1	A	75	LEU	N-CA-C	5.01	117.14	109.07

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	2444	0	2498	69	1
1	B	2415	0	2448	57	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	5	0	0	0	0
4	A	132	0	0	2	0
4	B	131	0	0	6	0
All	All	5129	0	4946	123	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 13.

All (123) 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:306:ARG:HH11	1:A:306:ARG:HG3	0.88	1.05
1:A:306:ARG:HH11	1:A:306:ARG:CG	1.78	0.96
1:A:306:ARG:HG3	1:A:306:ARG:NH1	1.69	0.96
1:B:137:THR:HG22	1:B:292:ARG:HH12	1.33	0.93
1:A:160:VAL:HG23	1:A:273:LEU:HD23	1.56	0.85
1:A:160:VAL:CG2	1:A:273:LEU:HD23	2.16	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:5:GLN:CD	1:B:329:GLN:HE22	1.94	0.75
1:A:55:THR:CG2	1:A:72:THR:HG22	2.18	0.73
1:A:179:ALA:O	1:A:183:GLU:HG3	1.90	0.72
1:B:169:LEU:HD22	1:B:172:ILE:HD12	1.71	0.71
1:A:55:THR:HG22	1:A:72:THR:HG22	1.72	0.69
1:B:234:ARG:NH2	1:B:242:GLY:HA2	2.12	0.65
1:B:325:ILE:O	1:B:329:GLN:HG3	1.94	0.65
1:B:82:ALA:H	1:B:93:ASN:HD21	1.46	0.64
1:B:204:LEU:HD23	1:B:241:ASN:HB3	1.80	0.64
1:B:153:LEU:HD12	1:B:273:LEU:HD11	1.81	0.63
1:A:275:TYR:HD1	1:A:275:TYR:O	1.81	0.63
1:B:234:ARG:HH22	1:B:242:GLY:HA2	1.64	0.62
1:B:264:MSE:H	1:B:268:GLN:NE2	1.98	0.61
1:A:40:ALA:HB2	1:A:75:LEU:CD2	2.31	0.61
1:A:40:ALA:HB2	1:A:75:LEU:HD21	1.83	0.60
1:B:153:LEU:HD13	1:B:154:ALA:N	2.17	0.60
1:A:14:PRO:HG3	1:A:96:TYR:CE2	2.38	0.58
1:B:273:LEU:C	1:B:273:LEU:HD13	2.28	0.58
1:B:178:PRO:HG3	1:B:225:THR:HG23	1.86	0.57
1:B:153:LEU:HD12	1:B:273:LEU:CD1	2.34	0.57
1:B:241:ASN:O	1:B:242:GLY:O	2.22	0.57
1:B:320:LEU:HD23	1:B:324:MSE:SE	2.55	0.57
1:A:264:MSE:H	1:A:268:GLN:NE2	2.03	0.56
1:A:275:TYR:O	1:A:275:TYR:CD1	2.59	0.56
1:A:276:GLN:NE2	1:B:275:TYR:OH	2.39	0.56
1:B:17:ILE:HD11	1:B:297:HIS:CD2	2.41	0.56
1:A:155:THR:HG23	1:A:273:LEU:CD1	2.36	0.55
1:B:309:ALA:O	1:B:310:ASP:HB2	2.06	0.55
1:B:27:GLN:NE2	4:B:638:HOH:O	2.39	0.55
1:A:279:GLY:O	1:A:281:GLY:N	2.40	0.55
1:A:301:LEU:H	1:A:301:LEU:HD22	1.72	0.54
1:B:248:THR:HG23	4:B:540:HOH:O	2.07	0.54
1:A:279:GLY:C	1:A:281:GLY:H	2.16	0.53
1:B:234:ARG:NH2	1:B:242:GLY:CA	2.72	0.53
1:A:264:MSE:H	1:A:268:GLN:HE22	1.57	0.53
1:B:81:ARG:H	1:B:93:ASN:ND2	2.06	0.53
1:B:181:LEU:HD11	1:B:226:ILE:HG12	1.90	0.53
1:B:110:ASN:ND2	4:B:435:HOH:O	2.41	0.53
1:B:151:MSE:HG2	1:B:165:THR:CG2	2.38	0.53
1:A:284:ILE:HD11	1:A:320:LEU:HD21	1.90	0.52
1:A:169:LEU:HD12	1:A:172:ILE:HD12	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:225:THR:C	1:A:228:PRO:HD2	2.34	0.52
1:B:117:ILE:HD12	1:B:117:ILE:N	2.25	0.52
1:B:301:LEU:HD22	4:B:491:HOH:O	2.10	0.51
1:A:160:VAL:HG23	1:A:273:LEU:CD2	2.36	0.51
1:B:234:ARG:HH21	1:B:242:GLY:N	2.08	0.51
1:A:2:VAL:HG23	1:A:3:LYS:N	2.24	0.51
1:B:249:LEU:C	1:B:249:LEU:HD23	2.36	0.51
1:B:309:ALA:O	1:B:310:ASP:CB	2.59	0.51
1:B:120:PRO:HD3	1:B:295:VAL:HG21	1.92	0.51
1:A:118:THR:OG1	1:A:294:SER:HB2	2.11	0.50
1:B:252:PRO:C	1:B:254:TYR:N	2.69	0.50
1:A:189:HIS:CD2	1:A:238:MSE:HG2	2.46	0.50
1:A:21:LEU:HD21	1:A:309:ALA:HB1	1.93	0.50
1:A:189:HIS:CD2	1:A:201:PRO:HG2	2.47	0.50
1:A:203:ILE:HB	1:A:240:LEU:HD23	1.94	0.49
1:A:204:LEU:HD22	1:A:249:LEU:HG	1.93	0.49
1:A:297:HIS:HD2	1:A:298:GLY:O	1.96	0.49
1:A:249:LEU:C	1:A:249:LEU:HD23	2.38	0.49
1:A:25:LEU:CD2	1:A:315:ILE:HD13	2.43	0.49
1:A:148:LYS:HG3	1:A:149:VAL:N	2.28	0.49
1:A:24:GLN:HE22	1:A:309:ALA:HB3	1.78	0.48
1:A:326:VAL:O	1:A:329:GLN:O	2.31	0.48
1:B:11:PRO:HG3	1:B:22:VAL:HG21	1.95	0.48
1:A:326:VAL:HG13	1:A:327:ASN:N	2.28	0.48
1:A:200:GLU:OE1	1:A:239:LYS:HG3	2.15	0.47
1:A:273:LEU:O	1:A:276:GLN:HB3	2.15	0.47
1:A:82:ALA:HB1	1:A:83:PRO:HD2	1.98	0.46
1:A:195:LYS:HB3	1:A:324:MSE:CE	2.45	0.46
1:A:326:VAL:HB	4:A:537:HOH:O	2.15	0.46
1:B:225:THR:C	1:B:228:PRO:HD2	2.41	0.46
1:B:228:PRO:O	1:B:232:GLU:HG3	2.16	0.46
1:A:155:THR:HG23	1:A:273:LEU:HD11	1.96	0.46
1:B:301:LEU:HD22	1:B:301:LEU:H	1.81	0.45
1:A:301:LEU:HD22	1:A:301:LEU:N	2.32	0.45
1:B:68:GLN:HA	4:B:526:HOH:O	2.16	0.45
1:A:284:ILE:HD11	1:A:320:LEU:CD2	2.46	0.45
1:A:277:GLY:O	1:A:278:PHE:C	2.60	0.45
1:B:234:ARG:HH21	1:B:242:GLY:H	1.65	0.44
1:A:205:VAL:O	1:A:243:PRO:HA	2.17	0.44
1:B:40:ALA:HB2	1:B:75:LEU:CD2	2.47	0.44
1:A:231:ASN:ND2	1:A:234:ARG:HH11	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:252:PRO:C	1:B:254:TYR:H	2.26	0.44
1:A:301:LEU:H	1:A:301:LEU:CD2	2.31	0.43
1:A:231:ASN:HD22	1:A:231:ASN:HA	1.70	0.43
1:A:169:LEU:HD12	1:A:172:ILE:CD1	2.47	0.43
1:B:151:MSE:HG2	1:B:165:THR:HG22	1.99	0.43
1:B:255:LEU:C	1:B:257:ASN:N	2.75	0.43
1:A:17:ILE:O	1:A:21:LEU:HG	2.18	0.43
1:A:50:LEU:HD22	1:A:307:GLY:HA2	2.01	0.43
1:B:225:THR:O	1:B:228:PRO:HD2	2.19	0.43
1:B:253:LYS:HD2	1:B:254:TYR:CE2	2.53	0.43
1:A:265:TYR:CZ	1:A:268:GLN:HB2	2.54	0.42
1:A:276:GLN:OE1	1:A:276:GLN:O	2.38	0.42
1:A:247:ASP:HB2	1:B:274:LYS:HE2	2.01	0.42
1:B:181:LEU:HD23	1:B:181:LEU:HA	1.90	0.42
1:A:13:GLU:OE1	1:A:14:PRO:HD2	2.20	0.42
1:A:195:LYS:HB3	1:A:324:MSE:HE2	2.00	0.42
1:B:64:PRO:HD2	4:B:651:HOH:O	2.20	0.42
1:A:326:VAL:CG1	1:A:327:ASN:N	2.83	0.41
1:B:153:LEU:C	1:B:153:LEU:CD1	2.92	0.41
1:B:273:LEU:HD13	1:B:273:LEU:O	2.19	0.41
1:B:19:PRO:O	1:B:23:VAL:HG23	2.19	0.41
1:B:151:MSE:HG2	1:B:165:THR:HG21	2.02	0.41
1:A:279:GLY:C	1:A:281:GLY:N	2.79	0.41
1:A:306:ARG:CG	1:A:306:ARG:NH1	2.50	0.41
1:A:206:CYS:SG	1:A:249:LEU:HD13	2.59	0.41
1:A:297:HIS:HE1	4:A:464:HOH:O	2.04	0.41
1:A:108:CYS:HB3	1:A:290:PHE:CG	2.56	0.41
1:A:170:ARG:NH2	1:B:174:ASP:OD1	2.41	0.41
1:B:162:LEU:HD23	1:B:263:ALA:HB3	2.02	0.41
1:B:265:TYR:CZ	1:B:268:GLN:HB2	2.54	0.41
1:A:120:PRO:HD3	1:A:295:VAL:HG21	2.03	0.40
1:B:93:ASN:HD22	1:B:93:ASN:HA	1.77	0.40
1:B:153:LEU:HD13	1:B:153:LEU:C	2.46	0.40
1:A:198:ILE:O	1:A:201:PRO:HD3	2.21	0.40
1:A:302:GLU:H	1:A:302:GLU:HG3	1.53	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:319:ASN:OD1	1:B:110:ASN:OD1[4_555]	2.08	0.12

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	327/329 (99%)	312 (95%)	11 (3%)	4 (1%)	10	4
1	B	326/329 (99%)	305 (94%)	17 (5%)	4 (1%)	10	4
All	All	653/658 (99%)	617 (94%)	28 (4%)	8 (1%)	10	4

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	276	GLN
1	B	242	GLY
1	B	278	PHE
1	B	310	ASP
1	A	280	ARG
1	B	309	ALA
1	A	278	PHE
1	A	277	GLY

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	257/254 (101%)	245 (95%)	12 (5%)	23	12
1	B	252/254 (99%)	241 (96%)	11 (4%)	25	15
All	All	509/508 (100%)	486 (96%)	23 (4%)	24	14

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	VAL
1	A	57	ARG
1	A	76	LEU
1	A	142	GLU
1	A	148	LYS
1	A	153	LEU
1	A	200	GLU
1	A	204	LEU
1	A	248	THR
1	A	273	LEU
1	A	302	GLU
1	A	306	ARG
1	B	2	VAL
1	B	57	ARG
1	B	134	THR
1	B	137	THR
1	B	142	GLU
1	B	151	MSE
1	B	153	LEU
1	B	233	LEU
1	B	257	ASN
1	B	274	LYS
1	B	303	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	27	GLN
1	A	45	ASN
1	A	190	HIS
1	A	231	ASN
1	A	268	GLN
1	A	276	GLN
1	A	297	HIS
1	B	27	GLN
1	B	93	ASN
1	B	110	ASN
1	B	127	ASN
1	B	182	HIS
1	B	231	ASN
1	B	236	GLN
1	B	268	GLN

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Mol	Chain	Res	Type
1	B	283	ASN
1	B	329	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 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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	$\# Z > 2$
3	PO4	A	332	-	4,4,4	1.42	0	6,6,6	0.47	0

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	321/329 (97%)	0.47	21 (6%) 25 28	10, 19, 31, 42	0
1	B	321/329 (97%)	0.69	42 (13%) 7 8	8, 21, 36, 41	0
All	All	642/658 (97%)	0.58	63 (9%) 13 15	8, 20, 34, 42	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	278	PHE	9.1
1	A	278	PHE	6.8
1	B	242	GLY	6.4
1	A	329	GLN	6.3
1	B	309	ALA	5.0
1	A	279	GLY	4.6
1	A	275	TYR	4.5
1	B	281	GLY	4.4
1	B	329	GLN	4.3
1	B	279	GLY	4.3
1	B	296	ASP	4.3
1	A	277	GLY	4.0
1	B	147	LYS	3.8
1	A	216	GLY	3.7
1	B	277	GLY	3.5
1	A	2	VAL	3.4
1	B	2	VAL	3.3
1	B	232	GLU	3.2
1	B	237	GLY	3.1
1	B	305	GLY	3.1
1	A	276	GLN	3.0
1	A	280	ARG	3.0
1	B	182	HIS	2.9
1	B	215	GLY	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	304	ALA	2.7
1	B	255	LEU	2.7
1	B	252	PRO	2.7
1	B	256	ASP	2.6
1	B	257	ASN	2.6
1	A	296	ASP	2.6
1	A	308	LYS	2.6
1	B	3	LYS	2.5
1	B	306	ARG	2.4
1	A	153	LEU	2.4
1	A	200	GLU	2.4
1	A	215	GLY	2.4
1	B	320	LEU	2.4
1	B	155	THR	2.4
1	B	251	GLN	2.4
1	B	307	GLY	2.4
1	B	233	LEU	2.3
1	B	301	LEU	2.3
1	B	110	ASN	2.3
1	A	157	GLU	2.3
1	B	308	LYS	2.3
1	B	48	ALA	2.3
1	A	156	GLU	2.3
1	B	44	THR	2.2
1	A	89	LEU	2.2
1	B	298	GLY	2.2
1	A	235	ALA	2.2
1	B	276	GLN	2.2
1	B	156	GLU	2.2
1	A	155	THR	2.2
1	B	280	ARG	2.2
1	B	310	ASP	2.1
1	A	131	ILE	2.1
1	B	327	ASN	2.1
1	B	199	ALA	2.1
1	A	126	ILE	2.0
1	B	51	GLY	2.0
1	B	282	VAL	2.0
1	B	217	HIS	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
3	PO4	A	332	5/5	0.96	0.10	23,23,23,23	0
2	ZN	A	330	1/1	0.99	0.02	25,25,25,25	0
2	ZN	B	331	1/1	1.00	0.01	20,20,20,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.