



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 01:57 PM UTC

PDB ID : 2V9O / pdb_00002v9o
Title : L-RHAMNULOSE-1-PHOSPHATE ALDOLASE FROM ESCHERICHIA COLI (MUTANT A87M- T109F-E192A)
Authors : Grueninger, D.; Schulz, G.E.
Deposited on : 2007-08-24
Resolution : 1.95 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

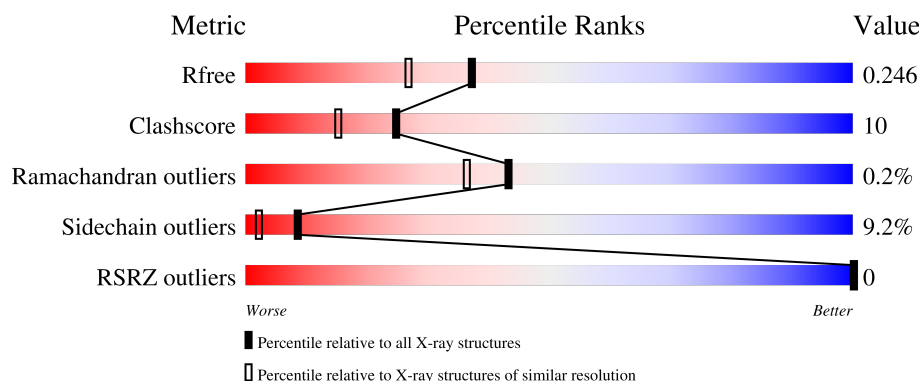
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	274	 69% 25% 5%
1	E	274	 73% 22%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4540 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 RHAMNULOSE-1-PHOSPHATE ALDOLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	274	Total	C	N	O	S	0	5	0
			2154	1381	364	394	15			
1	E	273	Total	C	N	O	S	0	4	0
			2141	1373	364	391	13			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	87	MET	ALA	engineered mutation	UNP P32169
A	109	PHE	THR	engineered mutation	UNP P32169
A	192	ALA	GLU	engineered mutation	UNP P32169
E	87	MET	ALA	engineered mutation	UNP P32169
E	109	PHE	THR	engineered mutation	UNP P32169
E	192	ALA	GLU	engineered mutation	UNP P32169

- 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	E	4	Total	Zn	0	0
			4	4		

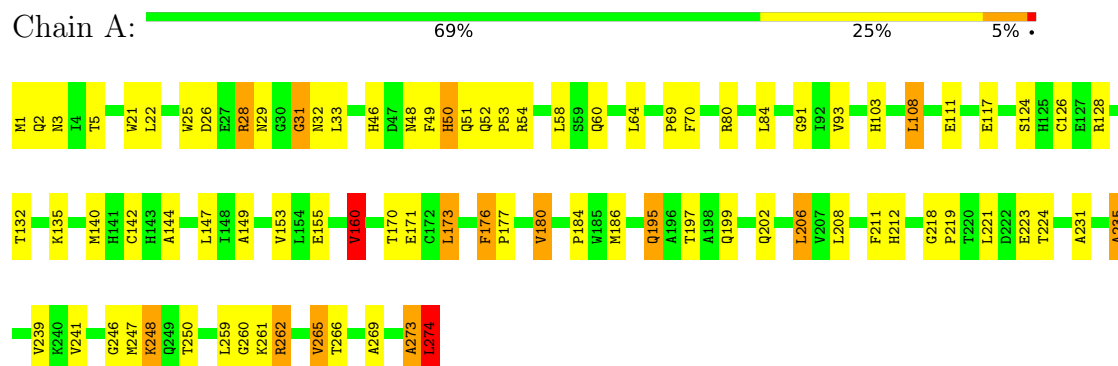
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	124	Total	O	0	0
			124	124		
3	E	115	Total	O	0	0
			115	115		

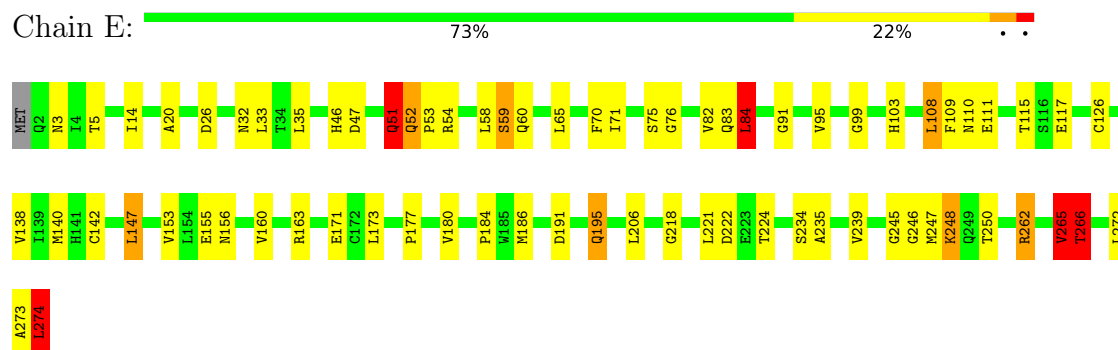
3 Residue-property plots

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: RHAMNULOSE-1-PHOSPHATE ALDOLASE



• Molecule 1: RHAMNULOSE-1-PHOSPHATE ALDOLASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 4	Depositor
Cell constants a, b, c, α , β , γ	86.19Å 86.19Å 89.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.95 – 1.95 44.95 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.7 (44.95-1.95) 99.7 (44.95-1.95)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.04 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.203 , 0.249 0.200 , 0.246	Depositor DCC
R_{free} test set	3347 reflections (7.00%)	wwPDB-VP
Wilson B-factor (Å ²)	30.4	Xtriage
Anisotropy	0.330	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 44.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.045 for -h,-l,-k 0.035 for -h,l,k 0.035 for l,-k,h 0.047 for -l,-k,-h 0.079 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4540	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 38.46 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.6720e-04.*

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	1.42	11/2222 (0.5%)	1.40	22/3023 (0.7%)
1	E	1.34	8/2208 (0.4%)	1.37	25/3005 (0.8%)
All	All	1.38	19/4430 (0.4%)	1.39	47/6028 (0.8%)

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	272	LEU	N-CA	7.49	1.55	1.46
1	A	144	ALA	CA-CB	-7.13	1.44	1.53
1	E	245	GLY	C-O	-6.88	1.15	1.24
1	E	171	GLU	C-O	-6.51	1.15	1.24
1	A	235	ALA	CA-CB	-6.50	1.42	1.53
1	A	22	LEU	C-O	6.04	1.31	1.24
1	A	208	LEU	C-O	-5.76	1.17	1.23
1	A	274	LEU	N-CA	5.70	1.57	1.46
1	E	84	LEU	CG-CD1	-5.69	1.33	1.52
1	A	160	VAL	CA-CB	5.55	1.61	1.54
1	E	234	SER	N-CA	-5.54	1.39	1.46
1	E	138	VAL	N-CA	5.53	1.52	1.46
1	E	266	THR	CA-CB	5.35	1.60	1.53
1	E	75	SER	CA-C	5.16	1.59	1.52
1	A	273	ALA	C-O	-5.08	1.18	1.24
1	A	231	ALA	CA-CB	5.05	1.61	1.53
1	A	149	ALA	CA-C	-5.03	1.46	1.52
1	A	70	PHE	N-CA	-5.02	1.39	1.45
1	A	241	VAL	CA-CB	5.01	1.60	1.54

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	273	ALA	N-CA-C	9.04	122.43	111.40
1	A	49	PHE	N-CA-C	8.05	121.42	110.55
1	E	14	ILE	N-CA-C	-7.65	102.99	110.72
1	A	260	GLY	N-CA-C	-7.54	102.55	113.86
1	E	273	ALA	CA-C-N	7.22	134.69	121.70
1	E	273	ALA	C-N-CA	7.22	134.69	121.70
1	A	274	LEU	N-CA-C	7.20	131.16	111.00
1	E	272	LEU	N-CA-C	7.13	125.99	110.80
1	E	155	GLU	N-CA-C	-7.05	99.39	109.96
1	A	31	GLY	N-CA-C	-6.87	104.60	112.29
1	E	273	ALA	N-CA-C	-6.67	96.60	110.80
1	E	274	LEU	CA-CB-CG	6.59	139.37	116.30
1	E	51	GLN	N-CA-C	6.25	118.06	109.18
1	E	265	VAL	CB-CA-C	6.22	120.40	110.82
1	A	170	THR	N-CA-C	6.02	118.61	111.33
1	A	147	LEU	CD1-CG-CD2	-5.93	97.76	110.80
1	A	273	ALA	CA-C-O	-5.88	114.26	120.55
1	E	46	HIS	N-CA-C	5.62	118.13	111.33
1	A	273	ALA	O-C-N	-5.61	115.38	122.17
1	E	47	ASP	CA-C-N	-5.60	113.72	122.73
1	E	47	ASP	C-N-CA	-5.60	113.72	122.73
1	E	70	PHE	CA-C-N	-5.57	115.91	123.10
1	E	70	PHE	C-N-CA	-5.57	115.91	123.10
1	E	273	ALA	O-C-N	5.44	129.83	122.59
1	A	176	PHE	CA-C-N	5.33	124.84	119.19
1	A	176	PHE	C-N-CA	5.33	124.84	119.19
1	A	21	TRP	N-CA-C	-5.32	105.38	111.07
1	A	246	GLY	N-CA-C	5.29	117.44	112.04
1	A	265	VAL	CB-CA-C	5.26	118.80	110.81
1	E	52[A]	GLN	CA-C-N	-5.23	114.41	119.90
1	E	52[A]	GLN	C-N-CA	-5.23	114.41	119.90
1	E	52[B]	GLN	CA-C-N	-5.23	114.41	119.90
1	E	52[B]	GLN	C-N-CA	-5.23	114.41	119.90
1	E	156	ASN	CA-C-N	5.19	130.32	123.00
1	E	156	ASN	C-N-CA	5.19	130.32	123.00
1	E	76	GLY	N-CA-C	-5.18	108.48	115.36
1	A	70	PHE	CA-C-N	-5.16	116.44	123.10
1	A	70	PHE	C-N-CA	-5.16	116.44	123.10
1	E	246	GLY	N-CA-C	5.09	118.49	112.33
1	E	82	VAL	N-CA-C	5.07	115.68	110.36
1	A	49	PHE	CA-C-N	5.06	128.04	120.90
1	A	49	PHE	C-N-CA	5.06	128.04	120.90
1	A	28	ARG	CG-CD-NE	-5.03	100.92	112.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	222	ASP	N-CA-C	5.03	119.67	111.37
1	A	197	THR	N-CA-C	-5.02	105.72	111.14
1	A	273	ALA	CA-C-N	-5.00	112.69	121.70
1	A	273	ALA	C-N-CA	-5.00	112.69	121.70

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	2154	0	2133	46	0
1	E	2141	0	2125	39	0
2	A	2	0	0	0	0
2	E	4	0	0	0	0
3	A	124	0	0	10	1
3	E	115	0	0	5	1
All	All	4540	0	4258	83	2

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 10.

All (83) 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:E:140:MET:HE2	1:E:142[B]:CYS:SG	1.69	1.30
1:A:140:MET:HE2	1:A:142[B]:CYS:SG	1.82	1.19
1:A:140:MET:CE	1:A:142[B]:CYS:SG	2.57	0.93
1:E:91:GLY:HA3	1:E:108:LEU:HD22	1.51	0.90
1:A:91:GLY:HA3	1:A:108:LEU:HD22	1.56	0.86
1:E:140:MET:CE	1:E:142[B]:CYS:SG	2.61	0.85
1:A:64:LEU:HD23	1:A:126[B]:CYS:SG	2.22	0.79
1:A:273:ALA:O	1:A:274:LEU:HD22	1.83	0.78
1:E:262[A]:ARG:CG	1:E:262[A]:ARG:HH11	2.00	0.75
1:A:219:PRO:HG2	1:A:223:GLU:OE2	1.88	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:ARG:NH1	3:A:2047:HOH:O	2.22	0.71
1:E:274:LEU:HD13	3:E:2076:HOH:O	1.90	0.71
1:E:117:GLU:OE1	3:E:2070:HOH:O	2.10	0.69
1:A:262:ARG:NH1	3:A:2120:HOH:O	2.21	0.68
1:A:1:MET:HG3	1:A:2:GLN:O	1.94	0.68
1:E:140:MET:HE2	1:E:142[B]:CYS:HG	1.55	0.66
1:E:163:ARG:HG2	1:E:274:LEU:HA	1.77	0.65
1:A:84:LEU:HD11	3:A:2047:HOH:O	1.98	0.63
1:A:54:ARG:NH2	1:A:111:GLU:HA	2.14	0.62
1:A:247[B]:MET:HE1	3:A:2076:HOH:O	2.01	0.61
1:A:273:ALA:HB2	1:E:53:PRO:HD2	1.83	0.61
1:E:247[B]:MET:HE1	3:E:2078:HOH:O	2.04	0.56
1:E:262[A]:ARG:CG	1:E:262[A]:ARG:NH1	2.68	0.56
1:E:218:GLY:HA3	1:E:224:THR:OG1	2.06	0.55
1:A:50:HIS:HD2	3:A:2058:HOH:O	1.90	0.55
1:A:266:THR:HG23	1:A:266:THR:O	2.07	0.54
1:E:20:ALA:HB1	1:E:26:ASP:OD1	2.08	0.54
1:E:60:GLN:HE22	1:E:195:GLN:HE22	1.56	0.53
1:E:262[A]:ARG:HH11	1:E:262[A]:ARG:HG3	1.73	0.53
1:A:235:ALA:O	1:A:239:VAL:HG23	2.09	0.53
1:E:177:PRO:HB3	1:E:265:VAL:HG13	1.92	0.52
1:A:132:THR:O	1:A:135:LYS:HD3	2.10	0.50
1:E:247[B]:MET:HE2	3:E:2107:HOH:O	2.11	0.50
1:A:3:ASN:OD1	1:A:5:THR:HG23	2.10	0.50
1:A:117:GLU:OE1	3:A:2062:HOH:O	2.18	0.50
1:A:212:HIS:CD2	1:A:212:HIS:C	2.89	0.50
1:E:59:SER:HB2	1:E:191:ASP:OD2	2.12	0.50
1:E:60:GLN:NE2	1:E:195:GLN:HE22	2.09	0.50
1:E:65:LEU:HG	1:E:126:CYS:SG	2.54	0.48
1:A:60:GLN:NE2	1:A:195:GLN:HE22	2.11	0.48
1:E:54:ARG:NH2	1:E:111:GLU:HA	2.28	0.48
1:A:50:HIS:CD2	3:A:2058:HOH:O	2.65	0.47
1:E:262[A]:ARG:HH11	1:E:262[A]:ARG:HG2	1.75	0.47
1:A:184:PRO:HD2	1:A:186:MET:HE3	1.97	0.46
1:E:184:PRO:HG2	1:E:186:MET:HE3	1.96	0.46
1:A:28:ARG:HG2	1:A:29:ASN:H	1.81	0.46
1:E:147:LEU:HD13	1:E:235:ALA:HB2	1.97	0.46
1:E:262[A]:ARG:NH1	1:E:262[A]:ARG:HG3	2.30	0.46
1:A:124:SER:O	1:A:128:ARG:HG2	2.15	0.46
1:A:262:ARG:HG3	1:A:262:ARG:HH11	1.79	0.46
1:A:60:GLN:HE22	1:A:195:GLN:HE22	1.64	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:247[A]:MET:SD	1:A:250:THR:HB	2.57	0.44
1:A:84:LEU:HD21	3:A:2047:HOH:O	2.17	0.44
1:E:247[B]:MET:CE	3:E:2107:HOH:O	2.65	0.44
1:E:266:THR:HG23	1:E:266:THR:O	2.18	0.44
1:A:266:THR:O	1:A:266:THR:CG2	2.65	0.44
1:A:155:GLU:HB3	1:A:160:VAL:HG21	1.99	0.44
1:E:35:LEU:HD23	1:E:71:ILE:CG2	2.48	0.44
1:E:83:GLN:HG3	1:E:84:LEU:HD13	2.00	0.43
1:A:69:PRO:HA	1:A:93:VAL:O	2.18	0.43
1:A:180:VAL:HB	1:A:206:LEU:HB3	1.99	0.43
1:E:3:ASN:OD1	1:E:5:THR:HG23	2.18	0.43
1:A:248:LYS:HE3	1:A:248:LYS:HB3	1.73	0.43
1:A:269:ALA:HB1	1:E:51:GLN:HA	2.00	0.43
1:E:195:GLN:HE21	1:E:195:GLN:HB2	1.55	0.43
1:A:25:TRP:HB3	1:A:142[A]:CYS:HB2	2.00	0.43
1:A:184:PRO:O	1:A:186:MET:HG3	2.19	0.43
1:E:54:ARG:O	1:E:103:HIS:HA	2.18	0.42
1:E:109:PHE:O	1:E:110:ASN:HB2	2.19	0.42
1:A:218:GLY:HA3	1:A:224:THR:OG1	2.18	0.42
1:A:247[B]:MET:HE2	3:A:2111:HOH:O	2.19	0.42
1:E:52[A]:GLN:H	1:E:52[A]:GLN:CD	2.27	0.42
1:A:199:GLN:O	1:A:202:GLN:HB2	2.20	0.42
1:E:235:ALA:O	1:E:239:VAL:HG23	2.20	0.41
1:E:95:VAL:CG1	1:E:99:GLY:HA2	2.50	0.41
1:A:173:LEU:HG	1:A:259:LEU:HD23	2.02	0.41
1:E:248:LYS:HB3	1:E:248:LYS:HE3	1.74	0.41
1:A:53:PRO:HG3	3:A:2035:HOH:O	2.21	0.41
1:A:54:ARG:O	1:A:103:HIS:HA	2.20	0.41
1:A:50:HIS:O	1:A:51:GLN:C	2.64	0.41
1:A:26:ASP:OD2	1:A:31:GLY:HA3	2.21	0.40
1:A:176:PHE:HA	1:A:177:PRO:HD2	1.95	0.40
1:E:247[A]:MET:SD	1:E:250:THR:HB	2.62	0.40

All (2) 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)
3:E:2054:HOH:O	3:E:2054:HOH:O[3_555]	1.66	0.54
3:A:2051:HOH:O	3:A:2051:HOH:O[3_655]	1.71	0.49

5.3 Torsion angles

5.3.1 Protein backbone

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	277/274 (101%)	265 (96%)	12 (4%)	0	100	100
1	E	275/274 (100%)	264 (96%)	10 (4%)	1 (0%)	30	21
All	All	552/548 (101%)	529 (96%)	22 (4%)	1 (0%)	43	36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	115	THR

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	233/228 (102%)	209 (90%)	24 (10%)	7	1
1	E	231/228 (101%)	210 (91%)	21 (9%)	9	2
All	All	464/456 (102%)	419 (90%)	45 (10%)	8	2

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	ASN
1	A	33	LEU
1	A	46[A]	HIS
1	A	46[B]	HIS
1	A	48	ASN

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Mol	Chain	Res	Type
1	A	50	HIS
1	A	52	GLN
1	A	58	LEU
1	A	108	LEU
1	A	153	VAL
1	A	160	VAL
1	A	171[A]	GLU
1	A	171[B]	GLU
1	A	173	LEU
1	A	180	VAL
1	A	195	GLN
1	A	206	LEU
1	A	211	PHE
1	A	221	LEU
1	A	248	LYS
1	A	261	LYS
1	A	262	ARG
1	A	265	VAL
1	A	274	LEU
1	E	32	ASN
1	E	33	LEU
1	E	51	GLN
1	E	58	LEU
1	E	59	SER
1	E	84	LEU
1	E	108	LEU
1	E	147	LEU
1	E	153	VAL
1	E	160	VAL
1	E	173	LEU
1	E	180	VAL
1	E	195	GLN
1	E	206	LEU
1	E	221	LEU
1	E	248	LYS
1	E	262[A]	ARG
1	E	262[B]	ARG
1	E	265	VAL
1	E	266	THR
1	E	274	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	GLN
1	A	156	ASN
1	A	195	GLN
1	E	50	HIS
1	E	51	GLN
1	E	156	ASN
1	E	195	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	274/274 (100%)	-1.06	0 100 100	14, 29, 41, 52	5 (1%)
1	E	273/274 (99%)	-1.05	0 100 100	18, 29, 40, 44	4 (1%)
All	All	547/548 (99%)	-1.05	0 100 100	14, 29, 40, 52	9 (1%)

There are no RSRZ outliers to report.

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	E	1276	1/1	0.99	0.03	61,61,61,61	0
2	ZN	A	1276	1/1	1.00	0.03	55,55,55,55	0
2	ZN	E	1275	1/1	1.00	0.02	35,35,35,35	0
2	ZN	A	1275	1/1	1.00	0.02	29,29,29,29	0
2	ZN	E	1277	1/1	1.00	0.04	39,39,39,39	0
2	ZN	E	1278	1/1	1.00	0.03	50,50,50,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.