Full wwPDB X-ray Structure Validation Report ⓘ

Dec 2, 2023 – 01:11 pm GMT

PDB ID	:	2V9O
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:

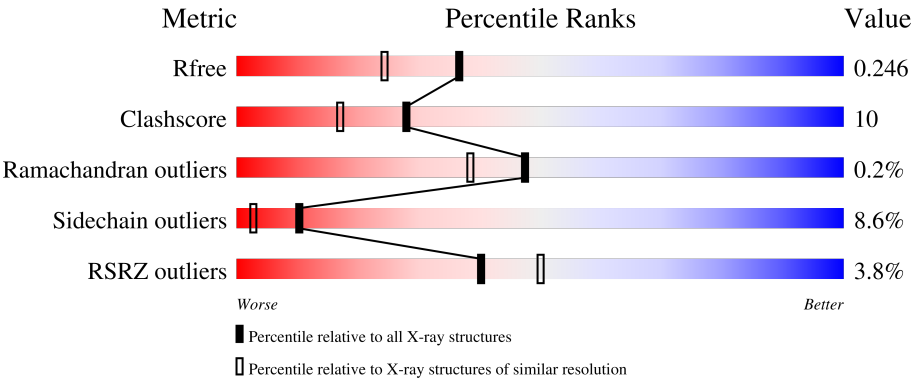
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 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}	130704	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (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 <=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	4% 73% 20% 5%
1	E	274	4% 78% 17%

2 Entry composition (i)

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 (three-letter code: 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.

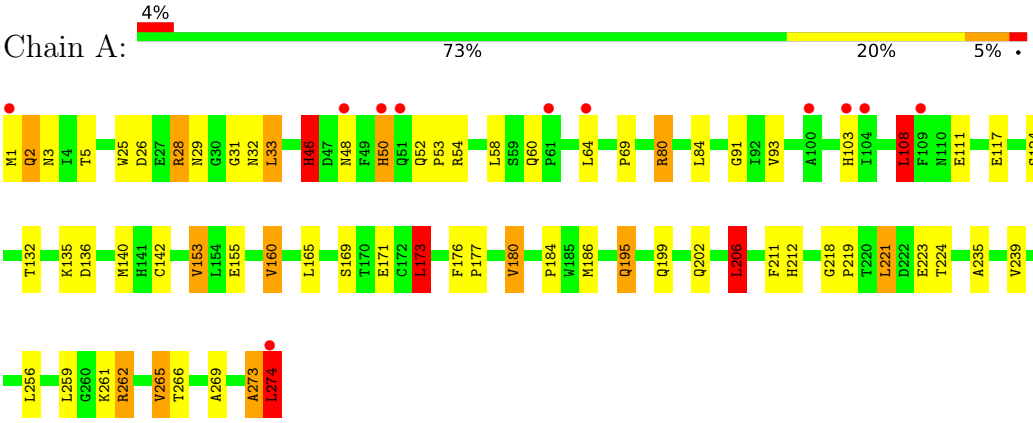
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	124	Total	O	0	0
			124	124		
3	E	115	Total	O	0	0
			115	115		

3 Residue-property plots (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.

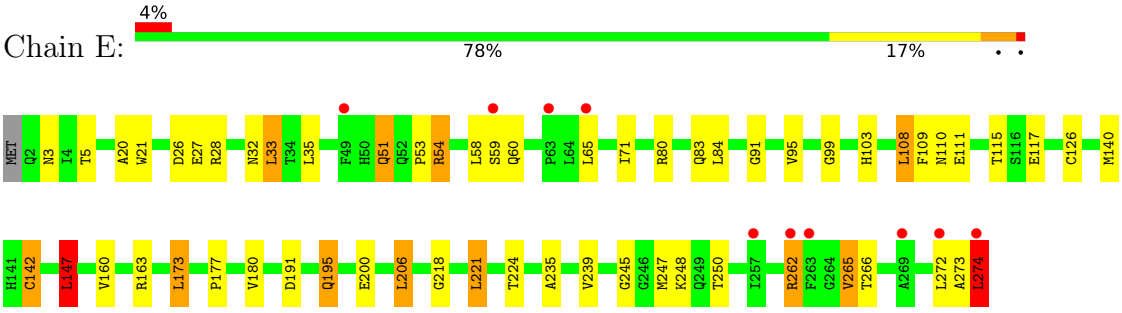
• Molecule 1: RHAMNULOSE-1-PHOSPHATE ALDOLASE

Chain A:



• Molecule 1: RHAMNULOSE-1-PHOSPHATE ALDOLASE

Chain E:



4 Data and refinement statistics (i)

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 44.94 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.7 (44.95-1.95) 99.7 (44.94-1.95)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [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 , 48.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.43$, $< L^2 > = 0.25$	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.*

[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	1.22	8/2222 (0.4%)	1.12	16/3023 (0.5%)
1	E	1.09	5/2208 (0.2%)	1.09	16/3005 (0.5%)
All	All	1.16	13/4430 (0.3%)	1.10	32/6028 (0.5%)

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	46[A]	HIS	CE1-NE2	10.60	1.57	1.32
1	A	46[B]	HIS	CE1-NE2	10.60	1.57	1.32
1	A	46[A]	HIS	CG-CD2	9.05	1.51	1.35
1	A	46[B]	HIS	CG-CD2	9.05	1.51	1.35
1	E	27	GLU	CD-OE1	7.63	1.34	1.25
1	E	142[A]	CYS	CB-SG	-7.53	1.69	1.82
1	E	142[B]	CYS	CB-SG	-7.53	1.69	1.82
1	A	46[A]	HIS	CG-ND1	6.83	1.53	1.38
1	A	46[B]	HIS	CG-ND1	6.83	1.53	1.38
1	A	169	SER	CB-OG	5.55	1.49	1.42
1	A	274	LEU	N-CA	5.36	1.57	1.46
1	E	21	TRP	CE3-CZ3	5.35	1.47	1.38
1	E	245	GLY	C-O	-5.28	1.15	1.23

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	E	274	LEU	CA-CB-CG	10.47	139.37	115.30
1	E	173	LEU	CA-CB-CG	8.02	133.75	115.30
1	A	206	LEU	CA-CB-CG	7.61	132.81	115.30
1	E	147	LEU	CB-CG-CD1	7.57	123.86	111.00
1	A	274	LEU	N-CA-C	7.47	131.16	111.00
1	E	28	ARG	NE-CZ-NH2	-7.01	116.79	120.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	256	LEU	CB-CG-CD1	-6.82	99.41	111.00
1	E	54	ARG	NE-CZ-NH1	6.76	123.68	120.30
1	A	28	ARG	NE-CZ-NH2	-6.35	117.13	120.30
1	A	221	LEU	CA-CB-CG	6.20	129.57	115.30
1	E	54	ARG	NE-CZ-NH2	-6.09	117.26	120.30
1	A	265	VAL	CG1-CB-CG2	6.07	120.60	110.90
1	E	221	LEU	CB-CG-CD1	6.04	121.27	111.00
1	A	153	VAL	CG1-CB-CG2	5.98	120.47	110.90
1	A	273	ALA	CA-C-N	5.98	130.35	117.20
1	A	108	LEU	CB-CG-CD1	5.94	121.10	111.00
1	A	33	LEU	CB-CG-CD2	5.90	121.03	111.00
1	E	33	LEU	CB-CG-CD2	5.72	120.73	111.00
1	E	206	LEU	CA-CB-CG	5.65	128.29	115.30
1	E	272	LEU	N-CA-C	5.55	125.99	111.00
1	A	136	ASP	CB-CG-OD1	5.46	123.21	118.30
1	E	33	LEU	CA-CB-CG	5.41	127.75	115.30
1	E	147	LEU	CA-CB-CG	5.34	127.57	115.30
1	E	273	ALA	N-CA-C	-5.33	96.60	111.00
1	A	165	LEU	CB-CG-CD2	5.27	119.95	111.00
1	E	206	LEU	CB-CG-CD1	5.25	119.93	111.00
1	A	173	LEU	CA-CB-CG	5.23	127.32	115.30
1	E	273	ALA	C-N-CA	5.20	134.69	121.70
1	A	28	ARG	CG-CD-NE	-5.18	100.92	111.80
1	A	221	LEU	CB-CG-CD1	5.12	119.70	111.00
1	E	200	GLU	OE1-CD-OE2	-5.08	117.20	123.30
1	A	206	LEU	CB-CG-CD1	5.03	119.55	111.00

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	2154	0	2133	49	0
1	E	2141	0	2125	37	0
2	A	2	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	4	0	0	0	0
3	A	124	0	0	14	1
3	E	115	0	0	5	1
All	All	4540	0	4258	84	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 (84) 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:46[A]:HIS:NE2	3:A:2031:HOH:O	1.60	1.31
1:A:46[A]:HIS:HB2	3:A:2032:HOH:O	1.14	1.30
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:E:140:MET:CE	1:E:142[B]:CYS:SG	2.61	0.89
1:A:140:MET:HE3	1:A:142[B]:CYS:SG	2.13	0.87
1:A:91:GLY:HA3	1:A:108:LEU:HD22	1.56	0.86
1:E:140:MET:HE1	1:E:142[B]:CYS:SG	2.18	0.83
1:A:140:MET:HE2	1:A:142[B]:CYS:SG	2.19	0.82
1:E:140:MET:HE2	1:E:142[B]:CYS:SG	2.22	0.80
1:A:64:LEU:HD23	1:A:126[B]:CYS:SG	2.22	0.79
1:A:46[A]:HIS:CB	3:A:2032:HOH:O	1.88	0.79
1:A:273:ALA:O	1:A:274:LEU:HD22	1.83	0.78
1:A:46[A]:HIS:CE1	3:A:2031:HOH:O	2.26	0.76
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
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:A:46[A]:HIS:CD2	3:A:2031:HOH:O	2.23	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: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:247[B]:MET:HE1	3:A:2076:HOH:O	1.99	0.63
1:A:54:ARG:NH2	1:A:111:GLU:HA	2.14	0.62
1:E:247[B]:MET:HE1	3:E:2078:HOH:O	1.99	0.62
1:A:273:ALA:HB2	1:E:53:PRO:HD2	1.83	0.61
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

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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: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: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
1:A:212:HIS:CD2	1:A:212:HIS:C	2.89	0.45
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:247[B]:MET:HE2	3:E:2107:HOH:O	2.18	0.43
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

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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: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 (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	277/274 (101%)	265 (96%)	12 (4%)	0	100 100
1	E	275/274 (100%)	264 (96%)	10 (4%)	1 (0%)	34 22
All	All	552/548 (101%)	529 (96%)	22 (4%)	1 (0%)	47 38

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	115	THR

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	233/228 (102%)	208 (89%)	25 (11%)	6 1
1	E	231/228 (101%)	214 (93%)	17 (7%)	13 4
All	All	464/456 (102%)	422 (91%)	42 (9%)	10 2

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	GLN
1	A	32	ASN
1	A	33	LEU
1	A	46[A]	HIS
1	A	46[B]	HIS
1	A	48	ASN
1	A	50	HIS
1	A	52	GLN
1	A	58	LEU
1	A	80	ARG
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	261	LYS
1	A	262	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
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	80	ARG
1	E	108	LEU
1	E	147	LEU
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	262[A]	ARG
1	E	262[B]	ARG
1	E	265	VAL
1	E	274	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	GLN
1	A	60	GLN
1	A	156	ASN
1	A	164	GLN
1	A	195	GLN
1	E	2	GLN
1	E	50	HIS
1	E	51	GLN
1	E	60	GLN
1	E	156	ASN
1	E	164	GLN
1	E	195	GLN

5.3.3 RNA

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	274/274 (100%)	0.24	11 (4%)	38	48	24, 29, 41, 52	0
1	E	273/274 (99%)	0.23	10 (3%)	41	51	24, 29, 40, 44	0
All	All	547/548 (99%)	0.23	21 (3%)	40	50	24, 29, 40, 52	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	5.5
1	E	274	LEU	4.1
1	A	51	GLN	3.9
1	E	59	SER	3.8
1	A	100	ALA	3.0
1	A	274	LEU	3.0
1	E	262[A]	ARG	3.0
1	E	257	ILE	2.9
1	E	65	LEU	2.8
1	E	269	ALA	2.8
1	A	103	HIS	2.6
1	E	63	PRO	2.5
1	A	109	PHE	2.5
1	E	49	PHE	2.4
1	A	61	PRO	2.4
1	A	64	LEU	2.3
1	E	272	LEU	2.2
1	A	50	HIS	2.2
1	A	104	ILE	2.1
1	E	263	PHE	2.1
1	A	48	ASN	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 monosaccharides 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, 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	E	1278	1/1	0.96	0.05	50,50,50,50	0
2	ZN	E	1276	1/1	0.98	0.05	61,61,61,61	0
2	ZN	A	1275	1/1	0.99	0.09	29,29,29,29	0
2	ZN	E	1277	1/1	0.99	0.04	39,39,39,39	0
2	ZN	E	1275	1/1	0.99	0.09	35,35,35,35	0
2	ZN	A	1276	1/1	1.00	0.12	55,55,55,55	0

6.5 Other polymers (i)

There are no such residues in this entry.