



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:40 PM GMT

PDB ID : 4MB0
Title : Crystal structure of TON1374
Authors : Kim, M.-K.; An, Y.J.; Cha, S.-S.
Deposited on : 2013-08-19
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

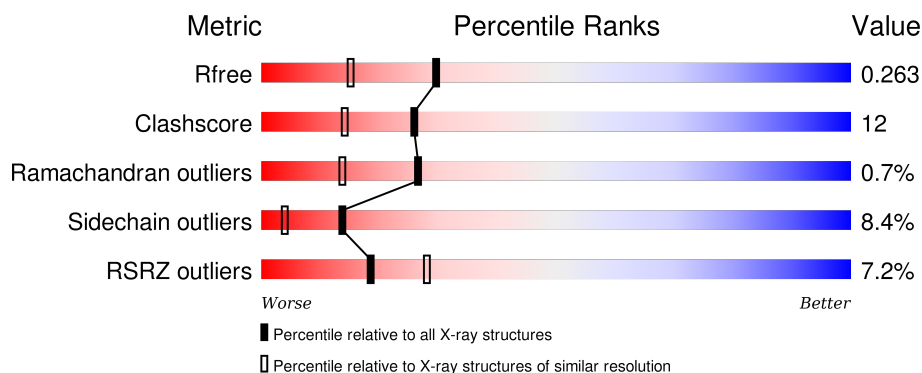
1 Overall quality at a glance

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}	91344	1833 (1.96-1.96)
Clashscore	102246	1953 (1.96-1.96)
Ramachandran outliers	100387	1936 (1.96-1.96)
Sidechain outliers	100360	1936 (1.96-1.96)
RSRZ outliers	91569	1835 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	261	<div> <div>5%</div> <div>75% 17% 5% .</div> </div>
1	B	261	<div> <div>8%</div> <div>69% 23% 5% .</div> </div>
1	C	261	<div> <div>6%</div> <div>74% 20% . .</div> </div>
1	D	261	<div> <div>10%</div> <div>74% 18% . . .</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ACT	D	301	-	-	-	X

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 8358 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 phosphopantothenate synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	253	Total	C	N	O	S	0	1	0
			2027	1281	364	375	7			
1	B	257	Total	C	N	O	S	0	0	0
			2038	1292	365	374	7			
1	C	253	Total	C	N	O	S	0	0	0
			2016	1275	360	374	7			
1	D	252	Total	C	N	O	S	0	0	0
			2004	1267	357	373	7			

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	C	O	0	0
			4	2	2		

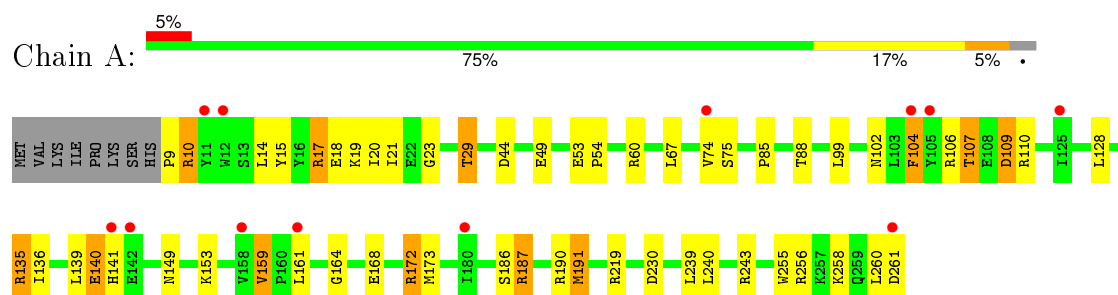
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	52	Total 52	O 52	0	0
3	B	87	Total 87	O 87	0	0
3	C	67	Total 67	O 67	0	0
3	D	63	Total 63	O 63	0	0

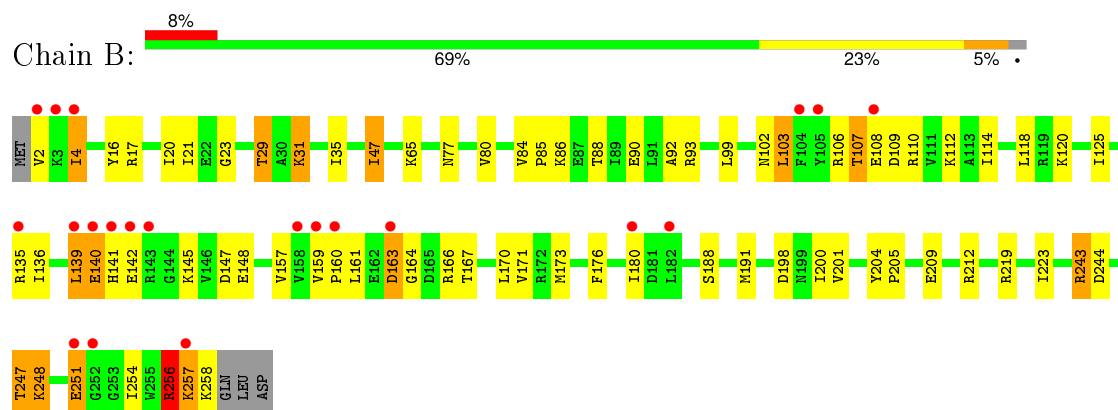
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

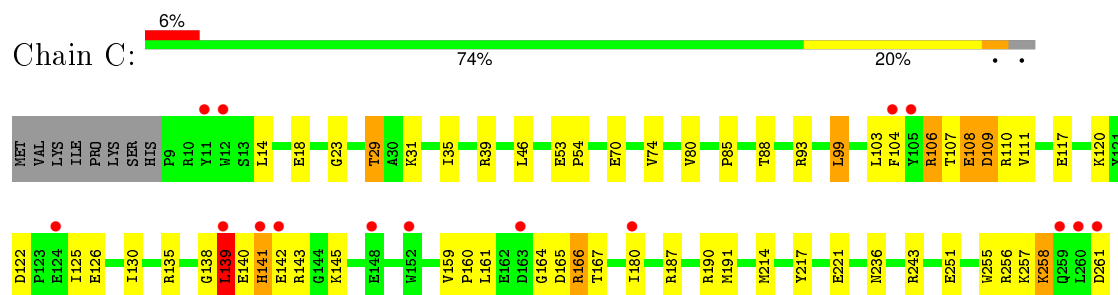
- Molecule 1: phosphopantothenate synthetase



- Molecule 1: phosphopantothenate synthetase

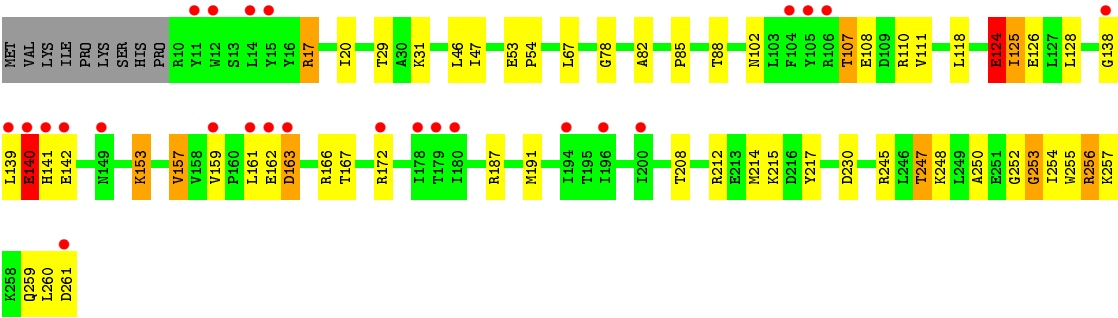


- Molecule 1: phosphopantothenate synthetase



- Molecule 1: phosphopantothenate synthetase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	266.14Å 60.96Å 75.35Å 90.00° 92.38° 90.00°	Depositor
Resolution (Å)	33.26 – 1.96 33.24 – 1.96	Depositor EDS
% Data completeness (in resolution range)	96.5 (33.26-1.96) 96.6 (33.24-1.96)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.45 (at 1.97Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.211 , 0.264 0.214 , 0.263	Depositor DCC
R_{free} test set	4227 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	30.6	Xtriage
Anisotropy	0.333	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.2	EDS
Estimated twinning fraction	0.014 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 84239 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8358	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.06% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: ACT

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	0.36	0/2057	0.49	0/2774
1	B	0.40	0/2069	0.54	0/2792
1	C	0.37	0/2046	0.51	0/2760
1	D	0.38	0/2032	0.51	0/2741
All	All	0.38	0/8204	0.51	0/11067

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	3
1	D	0	6
All	All	0	11

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	191	MET	Mainchain
1	A	9	PRO	Peptide
1	B	139	LEU	Peptide
1	B	198	ASP	Mainchain
1	B	251	GLU	Peptide
1	D	124	GLU	Peptide

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Mol	Chain	Res	Type	Group
1	D	140	GLU	Peptide
1	D	252	GLY	Peptide
1	D	253	GLY	Peptide
1	D	255	TRP	Peptide
1	D	260	LEU	Peptide

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	2027	0	2087	46	1
1	B	2038	0	2107	77	0
1	C	2016	0	2075	53	1
1	D	2004	0	2062	50	1
2	D	4	0	3	0	0
3	A	52	0	0	2	0
3	B	87	0	0	2	2
3	C	67	0	0	2	0
3	D	63	0	0	1	2
All	All	8358	0	8334	204	4

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 12.

All (204) 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:141:HIS:CB	1:B:142:GLU:HA	1.49	1.35
1:D:140:GLU:HB2	1:D:141:HIS:CA	1.76	1.15
1:B:243:ARG:HH11	1:B:243:ARG:HG2	0.97	1.10
1:C:256:ARG:NH1	1:D:191:MET:O	1.87	1.07
1:D:140:GLU:HB2	1:D:141:HIS:C	1.78	1.04
1:A:136:ILE:HG23	1:A:173:MET:SD	1.98	1.03
1:B:141:HIS:CB	1:B:142:GLU:CA	2.36	1.02
1:D:107:THR:HG22	1:D:110:ARG:H	1.25	1.01
1:B:243:ARG:NH1	1:B:243:ARG:HG2	1.76	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:243:ARG:HH11	1:B:243:ARG:CG	1.78	0.93
1:A:243:ARG:HD3	1:B:247:THR:HG22	1.49	0.92
1:D:140:GLU:CB	1:D:141:HIS:HA	1.99	0.92
1:D:140:GLU:CB	1:D:141:HIS:CA	2.50	0.90
1:A:260:LEU:O	1:A:261:ASP:HB2	1.72	0.90
1:D:140:GLU:HB2	1:D:141:HIS:HA	1.56	0.86
1:D:140:GLU:CG	1:D:141:HIS:HA	2.10	0.81
1:B:243:ARG:NH1	1:B:244:ASP:OD1	2.13	0.81
1:A:256:ARG:NH1	1:B:191:MET:O	2.14	0.81
1:C:74:VAL:HG23	1:C:99:LEU:CD1	2.11	0.80
1:C:85:PRO:O	1:C:88:THR:HG22	1.82	0.80
1:D:159:VAL:O	1:D:159:VAL:HG23	1.81	0.79
1:C:191:MET:O	1:D:256:ARG:NH1	2.14	0.79
1:B:256:ARG:HA	1:B:257:LYS:HB3	1.63	0.78
1:B:31:LYS:N	1:B:31:LYS:HD3	1.99	0.78
1:D:141:HIS:CB	1:D:142:GLU:HA	2.13	0.76
1:C:138:GLY:HA2	1:C:139:LEU:CB	2.17	0.75
1:C:139:LEU:O	1:C:141:HIS:N	2.19	0.73
1:A:243:ARG:CD	1:B:247:THR:HG22	2.20	0.71
1:B:23:GLY:C	1:B:29:THR:HG23	2.12	0.70
1:B:256:ARG:HA	1:B:257:LYS:CB	2.21	0.70
1:D:107:THR:HG22	1:D:110:ARG:N	2.04	0.69
1:C:74:VAL:CG2	1:C:99:LEU:HD13	2.22	0.69
1:C:74:VAL:CG2	1:C:99:LEU:CD1	2.71	0.68
1:C:138:GLY:HA2	1:C:139:LEU:HG	1.74	0.68
1:A:74:VAL:CG2	1:A:99:LEU:HD12	2.23	0.68
1:A:107:THR:HG22	1:A:110:ARG:H	1.57	0.68
1:D:17:ARG:O	1:D:20:ILE:HG22	1.94	0.68
1:D:141:HIS:CB	1:D:142:GLU:CA	2.73	0.67
1:D:85:PRO:O	1:D:88:THR:HG22	1.96	0.66
1:B:2:VAL:O	1:B:2:VAL:HG13	1.95	0.66
1:A:159:VAL:HG23	1:A:159:VAL:O	1.95	0.65
1:C:138:GLY:HA2	1:C:139:LEU:CG	2.27	0.65
1:A:74:VAL:CG2	1:A:99:LEU:CD1	2.75	0.65
1:A:159:VAL:O	1:A:159:VAL:CG2	2.45	0.65
1:B:219:ARG:HG2	1:B:223:ILE:HD12	1.77	0.64
1:A:74:VAL:HG22	1:A:99:LEU:HD12	1.80	0.64
1:C:159:VAL:HG23	1:C:159:VAL:O	1.97	0.64
1:B:140:GLU:HB2	1:B:141:HIS:O	1.98	0.64
1:C:74:VAL:HG23	1:C:99:LEU:HD12	1.80	0.63
1:B:107:THR:O	1:B:110:ARG:N	2.32	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:135:ARG:NH1	1:B:142:GLU:HG2	2.14	0.63
1:D:153:LYS:HD2	3:D:454:HOH:O	1.99	0.62
1:C:190:ARG:HH21	1:D:259:GLN:NE2	1.98	0.62
1:A:135:ARG:NH1	1:A:140:GLU:O	2.32	0.62
1:A:243:ARG:HD3	1:B:247:THR:CG2	2.25	0.61
1:B:159:VAL:HG23	1:B:159:VAL:O	2.01	0.61
1:C:243:ARG:HD3	1:D:247:THR:HG22	1.81	0.61
1:B:212:ARG:HD2	3:B:448:HOH:O	2.02	0.60
1:C:140:GLU:O	1:C:141:HIS:CB	2.50	0.60
1:A:260:LEU:O	1:A:261:ASP:CB	2.49	0.59
1:A:191:MET:HE1	1:B:257:LYS:HD3	1.85	0.59
1:C:142:GLU:HA	1:C:145:LYS:HE3	1.85	0.58
1:A:23:GLY:C	1:A:29:THR:HG23	2.23	0.58
1:C:23:GLY:C	1:C:29:THR:HG23	2.24	0.58
1:D:250:ALA:O	1:D:253:GLY:HA2	2.03	0.57
1:A:53:GLU:HB3	1:A:54:PRO:HD3	1.86	0.57
1:B:219:ARG:NH1	3:B:424:HOH:O	2.24	0.57
1:D:162:GLU:HG3	1:D:163:ASP:N	2.19	0.57
1:C:85:PRO:HD2	3:C:320:HOH:O	2.04	0.57
1:B:147:ASP:OD1	1:B:148:GLU:N	2.38	0.57
1:C:85:PRO:O	1:C:88:THR:CG2	2.52	0.56
1:D:140:GLU:HG3	1:D:141:HIS:HA	1.85	0.56
1:A:74:VAL:HG22	1:A:99:LEU:CD1	2.35	0.56
1:A:15:TYR:CZ	1:A:19:LYS:HE2	2.40	0.56
1:A:17:ARG:O	1:A:21:ILE:HG13	2.05	0.55
1:D:159:VAL:O	1:D:159:VAL:CG2	2.51	0.55
1:C:93:ARG:HH21	1:C:122:ASP:CG	2.10	0.55
1:B:107:THR:O	1:B:109:ASP:N	2.40	0.55
1:C:164:GLY:HA3	1:C:187:ARG:HB3	1.89	0.55
1:B:243:ARG:NH1	1:B:243:ARG:CG	2.49	0.55
1:C:258:LYS:HG2	1:D:187:ARG:HH22	1.71	0.55
1:A:164:GLY:HA3	1:A:187:ARG:HB3	1.89	0.55
1:C:138:GLY:HA2	1:C:139:LEU:HB2	1.89	0.54
1:C:109:ASP:OD1	1:C:109:ASP:N	2.39	0.53
1:D:139:LEU:O	1:D:140:GLU:C	2.46	0.53
1:A:10:ARG:HB3	1:A:10:ARG:CZ	2.36	0.53
1:B:17:ARG:O	1:B:21:ILE:HG12	2.09	0.53
1:B:180:ILE:HD13	1:B:204:TYR:HE1	1.74	0.53
1:A:149:ASN:HA	1:A:153:LYS:HD3	1.91	0.53
1:C:107:THR:O	1:C:110:ARG:N	2.42	0.52
1:C:46:LEU:HD11	1:D:46:LEU:HD13	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:255:TRP:CZ3	1:C:108:GLU:HG3	2.44	0.52
1:B:147:ASP:OD1	1:B:147:ASP:C	2.48	0.52
1:D:53:GLU:HB3	1:D:54:PRO:HD3	1.91	0.52
1:D:29:THR:HG22	1:D:78:GLY:O	2.10	0.51
1:A:191:MET:HE2	1:B:257:LYS:HB2	1.91	0.51
1:D:162:GLU:CG	1:D:163:ASP:N	2.72	0.51
1:B:125:ILE:O	1:B:125:ILE:HG23	2.10	0.51
1:B:102:ASN:OD1	1:B:166:ARG:NH1	2.43	0.51
1:B:106:ARG:O	1:B:107:THR:OG1	2.29	0.51
1:B:4:ILE:HD13	1:B:4:ILE:H	1.76	0.50
1:B:103:LEU:HD21	1:B:114:ILE:CD1	2.42	0.50
1:A:190:ARG:HD3	1:B:257:LYS:O	2.12	0.49
1:C:106:ARG:HB2	1:C:142:GLU:CD	2.32	0.49
1:D:124:GLU:O	1:D:125:ILE:O	2.29	0.49
1:B:2:VAL:CG1	1:B:2:VAL:O	2.60	0.49
1:B:107:THR:O	1:B:108:GLU:C	2.51	0.48
1:B:107:THR:HG22	1:B:109:ASP:H	1.77	0.48
1:C:243:ARG:HD3	1:D:247:THR:CG2	2.43	0.48
1:B:77:ASN:HA	1:B:103:LEU:HD23	1.96	0.48
1:C:190:ARG:HD3	1:D:257:LYS:O	2.13	0.48
1:B:85:PRO:O	1:B:88:THR:HG22	2.13	0.48
1:B:159:VAL:CG2	1:B:159:VAL:O	2.62	0.48
1:C:93:ARG:NH2	1:C:122:ASP:OD2	2.39	0.47
1:B:140:GLU:N	1:B:141:HIS:HA	2.29	0.47
1:C:142:GLU:HA	1:C:145:LYS:CE	2.44	0.47
1:A:240:LEU:HD23	3:A:318:HOH:O	2.14	0.47
1:B:136:ILE:HB	1:B:139:LEU:HD12	1.96	0.47
1:A:136:ILE:HB	1:A:139:LEU:HD12	1.97	0.47
1:C:236:ASN:OD1	1:D:254:ILE:HD12	2.15	0.47
1:B:140:GLU:HG2	1:B:141:HIS:HA	1.96	0.47
1:C:93:ARG:HD2	3:C:353:HOH:O	2.15	0.47
1:B:17:ARG:O	1:B:20:ILE:HG22	2.15	0.47
1:A:240:LEU:CD2	3:A:318:HOH:O	2.62	0.47
1:B:140:GLU:CG	1:B:141:HIS:HA	2.46	0.46
1:C:141:HIS:O	1:C:143:ARG:N	2.48	0.46
1:D:139:LEU:HD23	1:D:139:LEU:HA	1.83	0.46
1:B:31:LYS:H	1:B:31:LYS:HD3	1.78	0.46
1:D:141:HIS:CB	1:D:142:GLU:HB3	2.46	0.46
1:D:128:LEU:HD12	1:D:128:LEU:N	2.30	0.46
1:D:107:THR:O	1:D:111:VAL:HG23	2.14	0.46
1:C:165:ASP:OD1	1:C:165:ASP:N	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:135:ARG:HH11	1:B:142:GLU:HG2	1.80	0.46
1:A:14:LEU:O	1:A:18:GLU:HG2	2.16	0.46
1:B:170:LEU:O	1:B:173:MET:HB2	2.16	0.46
1:B:65:LYS:HG2	1:B:176:PHE:CE2	2.51	0.46
1:C:108:GLU:H	1:C:108:GLU:CD	2.18	0.46
1:A:239:LEU:HD12	1:B:254:ILE:HD11	1.97	0.45
1:C:255:TRP:O	1:C:256:ARG:HB2	2.17	0.45
1:A:135:ARG:HD2	1:A:141:HIS:CD2	2.51	0.45
1:C:166:ARG:O	1:C:167:THR:C	2.54	0.45
1:B:107:THR:HG22	1:B:109:ASP:HB2	1.99	0.45
1:C:190:ARG:NH2	1:D:259:GLN:NE2	2.62	0.45
1:B:80:VAL:HG22	1:B:200:ILE:CD1	2.46	0.45
1:B:80:VAL:HG21	1:B:160:PRO:HG3	1.99	0.45
1:A:44:ASP:OD1	1:A:49:GLU:HA	2.17	0.45
1:D:162:GLU:OE2	1:D:163:ASP:N	2.48	0.45
1:A:230:ASP:C	1:A:230:ASP:OD1	2.55	0.44
1:D:126:GLU:HG2	1:D:128:LEU:HD12	2.00	0.44
1:C:214:MET:O	1:C:217:TYR:HB2	2.17	0.44
1:D:162:GLU:HG2	1:D:167:THR:OG1	2.17	0.44
1:B:167:THR:O	1:B:171:VAL:HG23	2.18	0.44
1:B:107:THR:C	1:B:109:ASP:N	2.71	0.44
1:B:84:VAL:N	1:B:85:PRO:CD	2.81	0.44
1:B:140:GLU:H	1:B:140:GLU:HG2	1.63	0.44
1:C:103:LEU:HD13	1:C:110:ARG:HG2	2.00	0.44
1:C:180:ILE:HD12	1:C:180:ILE:N	2.33	0.44
1:D:126:GLU:CD	1:D:128:LEU:HD11	2.38	0.43
1:D:138:GLY:HA3	1:D:172:ARG:HH22	1.82	0.43
1:A:75:SER:OG	1:A:102:ASN:HB3	2.18	0.43
1:C:14:LEU:O	1:C:18:GLU:HG2	2.18	0.43
1:B:219:ARG:HG2	1:B:223:ILE:CD1	2.46	0.43
1:A:168:GLU:OE2	1:A:172:ARG:NH1	2.47	0.43
1:C:111:VAL:CG1	1:C:130:ILE:HG13	2.49	0.43
1:B:248:LYS:CD	1:B:248:LYS:C	2.86	0.43
1:B:164:GLY:HA2	1:B:188:SER:HB2	2.01	0.43
1:D:157:VAL:HG23	1:D:159:VAL:HG13	2.00	0.43
1:B:108:GLU:O	1:B:112:LYS:HB2	2.19	0.43
1:C:117:GLU:O	1:C:120:LYS:HB3	2.19	0.43
1:B:205:PRO:O	1:B:209:GLU:HG3	2.18	0.43
1:B:257:LYS:O	1:B:258:LYS:C	2.57	0.43
1:D:85:PRO:O	1:D:88:THR:CG2	2.67	0.42
1:A:240:LEU:C	1:A:240:LEU:HD23	2.40	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:157:VAL:O	1:B:157:VAL:HG23	2.19	0.42
1:A:109:ASP:OD1	1:A:109:ASP:N	2.51	0.42
1:A:53:GLU:N	1:A:54:PRO:CD	2.83	0.42
1:A:17:ARG:HG2	1:B:35:ILE:HD13	2.01	0.42
1:C:53:GLU:HB3	1:C:54:PRO:HD3	2.01	0.42
1:A:186:SER:O	1:A:190:ARG:HG3	2.20	0.42
1:B:142:GLU:HB3	1:B:145:LYS:HG3	2.02	0.42
1:B:16:TYR:CE2	1:B:201:VAL:HG11	2.55	0.42
1:C:139:LEU:HA	1:C:139:LEU:HD23	1.82	0.41
1:C:138:GLY:CA	1:C:139:LEU:CB	2.95	0.41
1:D:250:ALA:O	1:D:253:GLY:CA	2.67	0.41
1:C:35:ILE:HG22	1:C:39:ARG:NH1	2.35	0.41
1:B:163:ASP:OD2	1:B:163:ASP:C	2.57	0.41
1:C:139:LEU:HB3	1:C:140:GLU:H	1.56	0.41
1:C:80:VAL:HG21	1:C:160:PRO:HG2	2.02	0.41
1:A:243:ARG:NE	1:B:247:THR:HG22	2.36	0.41
1:A:191:MET:CE	1:B:257:LYS:HB2	2.51	0.41
1:D:29:THR:CG2	1:D:82:ALA:HB2	2.51	0.41
1:B:23:GLY:HA3	1:B:29:THR:CG2	2.51	0.41
1:B:92:ALA:HB2	1:B:99:LEU:HD13	2.02	0.41
1:D:208:THR:O	1:D:212:ARG:HG3	2.21	0.41
1:A:67:LEU:HA	1:A:67:LEU:HD23	1.92	0.41
1:D:230:ASP:OD1	1:D:230:ASP:C	2.59	0.41
1:C:122:ASP:OD1	1:C:125:ILE:N	2.54	0.41
1:D:214:MET:O	1:D:217:TYR:HB2	2.21	0.41
1:D:67:LEU:HA	1:D:67:LEU:HD23	1.92	0.41
1:B:159:VAL:HA	1:B:160:PRO:HD2	1.93	0.40
1:A:104:PHE:N	1:A:104:PHE:CD1	2.89	0.40
1:B:86:LYS:HE2	1:B:90:GLU:OE2	2.22	0.40
1:B:47:ILE:HD13	1:B:47:ILE:HA	1.72	0.40
1:A:85:PRO:O	1:A:88:THR:HG22	2.21	0.40

All (4) 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:139:LEU:O	1:A:139:LEU:O[2_554]	1.40	0.80
3:B:468:HOH:O	3:D:451:HOH:O[1_554]	1.79	0.41
3:B:431:HOH:O	3:D:430:HOH:O[1_554]	2.02	0.18
1:C:217:TYR:OH	1:D:261:ASP:OD1[4_555]	2.12	0.08

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	252/261 (97%)	244 (97%)	8 (3%)	0	100	100
1	B	255/261 (98%)	242 (95%)	10 (4%)	3 (1%)	16	5
1	C	251/261 (96%)	239 (95%)	10 (4%)	2 (1%)	24	11
1	D	250/261 (96%)	236 (94%)	12 (5%)	2 (1%)	24	11
All	All	1008/1044 (97%)	961 (95%)	40 (4%)	7 (1%)	26	14

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	139	LEU
1	D	125	ILE
1	D	140	GLU
1	B	256	ARG
1	B	257	LYS
1	C	141	HIS
1	B	107	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	212/220 (96%)	194 (92%)	18 (8%)	13	4
1	B	213/220 (97%)	197 (92%)	16 (8%)	17	5
1	C	211/220 (96%)	193 (92%)	18 (8%)	13	4

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	209/220 (95%)	190 (91%)	19 (9%)	12	3
All	All	845/880 (96%)	774 (92%)	71 (8%)	14	4

All (71) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ARG
1	A	17	ARG
1	A	20	ILE
1	A	29	THR
1	A	60	ARG
1	A	104	PHE
1	A	106	ARG
1	A	107	THR
1	A	109	ASP
1	A	128	LEU
1	A	135	ARG
1	A	140	GLU
1	A	159	VAL
1	A	161	LEU
1	A	172	ARG
1	A	187	ARG
1	A	219	ARG
1	A	258	LYS
1	B	4	ILE
1	B	29	THR
1	B	31	LYS
1	B	47	ILE
1	B	93	ARG
1	B	103	LEU
1	B	118	LEU
1	B	120	LYS
1	B	140	GLU
1	B	161	LEU
1	B	163	ASP
1	B	243	ARG
1	B	247	THR
1	B	248	LYS
1	B	251	GLU
1	B	256	ARG
1	C	29	THR
1	C	31	LYS

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Mol	Chain	Res	Type
1	C	70	GLU
1	C	99	LEU
1	C	104	PHE
1	C	106	ARG
1	C	108	GLU
1	C	109	ASP
1	C	126	GLU
1	C	135	ARG
1	C	139	LEU
1	C	161	LEU
1	C	166	ARG
1	C	221	GLU
1	C	251	GLU
1	C	257	LYS
1	C	258	LYS
1	C	261	ASP
1	D	17	ARG
1	D	31	LYS
1	D	47	ILE
1	D	102	ASN
1	D	107	THR
1	D	108	GLU
1	D	118	LEU
1	D	124	GLU
1	D	140	GLU
1	D	153	LYS
1	D	157	VAL
1	D	161	LEU
1	D	163	ASP
1	D	166	ARG
1	D	215	LYS
1	D	245	ARG
1	D	247	THR
1	D	248	LYS
1	D	256	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	141	HIS
1	D	259	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

1 ligand is modelled in this entry.

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	$\# Z > 2$
2	ACT	D	301	-	1,3,3	1.39	0	0,3,3	0.00	-

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ACT	D	301	-	-	0/0/0/0	0/0/0/0

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

There are no such residues in this entry.

5.8 Polymer linkage issues

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	253/261 (96%)	0.27	12 (4%) 35 46	25, 40, 73, 110	0
1	B	257/261 (98%)	0.40	21 (8%) 14 23	19, 38, 87, 123	0
1	C	253/261 (96%)	0.29	15 (5%) 26 36	24, 42, 84, 140	0
1	D	252/261 (96%)	0.48	25 (9%) 9 15	25, 42, 88, 146	0
All	All	1015/1044 (97%)	0.36	73 (7%) 18 28	19, 41, 84, 146	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	139	LEU	6.9
1	A	105	TYR	6.5
1	C	11	TYR	5.1
1	B	105	TYR	4.9
1	D	141	HIS	4.8
1	D	142	GLU	4.5
1	D	105	TYR	4.5
1	D	140	GLU	4.4
1	C	261	ASP	4.3
1	B	142	GLU	4.2
1	C	260	LEU	4.2
1	A	141	HIS	4.0
1	D	261	ASP	4.0
1	B	140	GLU	3.9
1	A	104	PHE	3.8
1	B	104	PHE	3.7
1	C	104	PHE	3.7
1	B	251	GLU	3.7
1	B	2	VAL	3.5
1	C	142	GLU	3.5
1	B	257	LYS	3.4

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Mol	Chain	Res	Type	RSRZ
1	C	105	TYR	3.4
1	D	104	PHE	3.4
1	B	163	ASP	3.4
1	D	11	TYR	3.3
1	B	141	HIS	3.3
1	D	12	TRP	3.3
1	B	3	LYS	3.3
1	D	172	ARG	3.2
1	D	15	TYR	3.2
1	B	159	VAL	3.1
1	C	259	GLN	3.1
1	A	11	TYR	3.1
1	D	180	ILE	3.1
1	A	12	TRP	3.0
1	D	14	LEU	3.0
1	D	159	VAL	2.9
1	D	196	ILE	2.8
1	B	108	GLU	2.8
1	B	135	ARG	2.8
1	C	141	HIS	2.7
1	D	163	ASP	2.7
1	D	179	THR	2.7
1	B	4	ILE	2.6
1	C	12	TRP	2.6
1	B	252	GLY	2.6
1	C	180	ILE	2.6
1	A	125	ILE	2.6
1	A	180	ILE	2.6
1	A	261	ASP	2.5
1	D	162	GLU	2.5
1	C	124	GLU	2.4
1	B	139	LEU	2.4
1	A	158	VAL	2.4
1	B	160	PRO	2.4
1	B	180	ILE	2.3
1	D	200	ILE	2.3
1	C	152	TRP	2.3
1	B	182	LEU	2.3
1	D	161	LEU	2.3
1	D	138	GLY	2.2
1	A	161	LEU	2.2
1	C	148	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	143	ARG	2.2
1	D	106	ARG	2.2
1	A	74	VAL	2.2
1	A	142	GLU	2.1
1	C	163	ASP	2.1
1	D	194	ILE	2.1
1	D	139	LEU	2.1
1	D	149	ASN	2.0
1	B	158	VAL	2.0
1	D	178	ILE	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	ACT	D	301	4/4	0.74	0.24	2.04	49,49,61,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.