



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 01:56 PM UTC

PDB ID : 2GTD / pdb_00002gtd
Title : Crystal Structure of a Type III Pantothenate Kinase: Insight into the Catalysis of an Essential Coenzyme A Biosynthetic Enzyme Universally Distributed in Bacteria
Authors : Yang, K.; Eyobo, Y.; Brand, A.L.; Martynowski, D.; Tomchick, D.
Deposited on : 2006-04-27
Resolution : 2.00 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

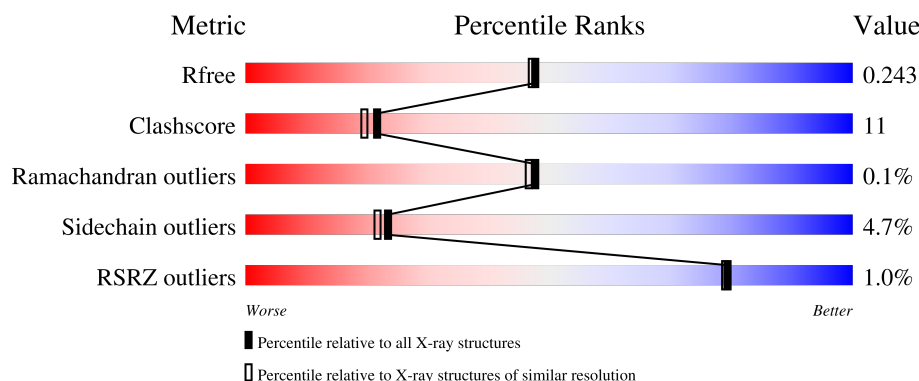
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	251	 80% 17% ..
1	B	251	 2% 83% 14% ..
1	C	251	 80% 18% ..
1	D	251	 82% 14% ..
1	E	251	 2% 85% 12% ..

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Mol	Chain	Length	Quality of chain
1	F	251	2% 80% 15% ••

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12919 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 Type III Pantothenate Kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	248	Total	C	N	O	S	0	0	0
			1929	1244	323	353	9			
1	B	248	Total	C	N	O	S	0	0	0
			1929	1244	323	353	9			
1	C	248	Total	C	N	O	S	0	0	0
			1929	1244	323	353	9			
1	D	248	Total	C	N	O	S	0	0	0
			1929	1244	323	353	9			
1	E	248	Total	C	N	O	S	0	0	0
			1929	1244	323	353	9			
1	F	248	Total	C	N	O	S	0	0	0
			1929	1244	323	353	9			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP Q9WZY5
A	2	ASP	-	cloning artifact	UNP Q9WZY5
A	3	PRO	-	cloning artifact	UNP Q9WZY5
B	1	MET	-	initiating methionine	UNP Q9WZY5
B	2	ASP	-	cloning artifact	UNP Q9WZY5
B	3	PRO	-	cloning artifact	UNP Q9WZY5
C	1	MET	-	initiating methionine	UNP Q9WZY5
C	2	ASP	-	cloning artifact	UNP Q9WZY5
C	3	PRO	-	cloning artifact	UNP Q9WZY5
D	1	MET	-	initiating methionine	UNP Q9WZY5
D	2	ASP	-	cloning artifact	UNP Q9WZY5
D	3	PRO	-	cloning artifact	UNP Q9WZY5
E	1	MET	-	initiating methionine	UNP Q9WZY5
E	2	ASP	-	cloning artifact	UNP Q9WZY5
E	3	PRO	-	cloning artifact	UNP Q9WZY5
F	1	MET	-	initiating methionine	UNP Q9WZY5
F	2	ASP	-	cloning artifact	UNP Q9WZY5

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Chain	Residue	Modelled	Actual	Comment	Reference
F	3	PRO	-	cloning artifact	UNP Q9WZY5

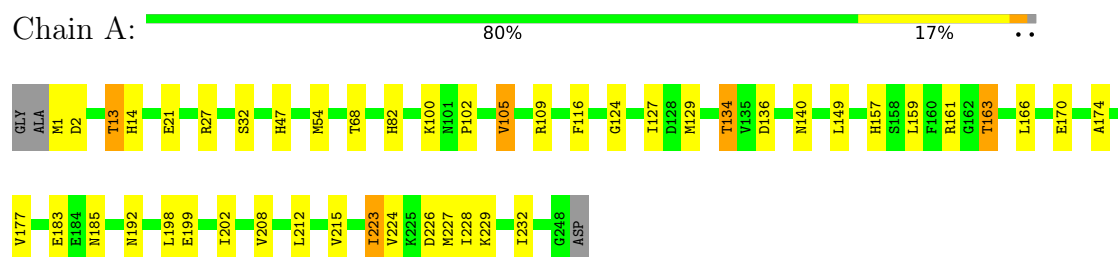
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	227	Total	O	0	0
			227	227		
2	B	223	Total	O	0	0
			223	223		
2	C	245	Total	O	0	0
			245	245		
2	D	219	Total	O	0	0
			219	219		
2	E	200	Total	O	0	0
			200	200		
2	F	231	Total	O	0	0
			231	231		

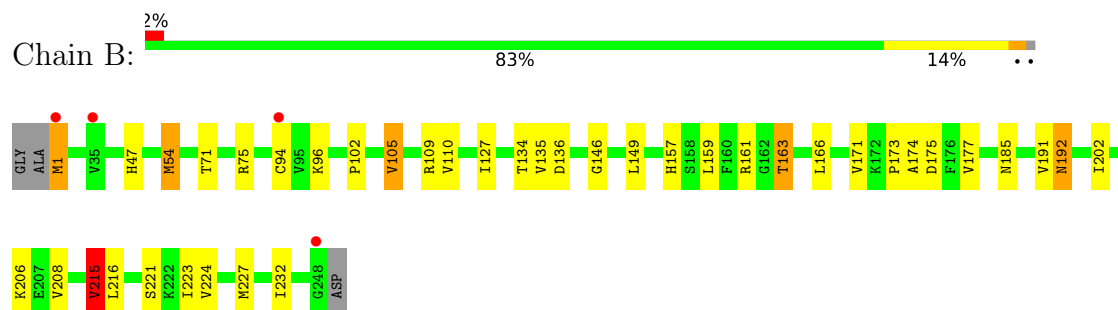
3 Residue-property plots [i](#)

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.

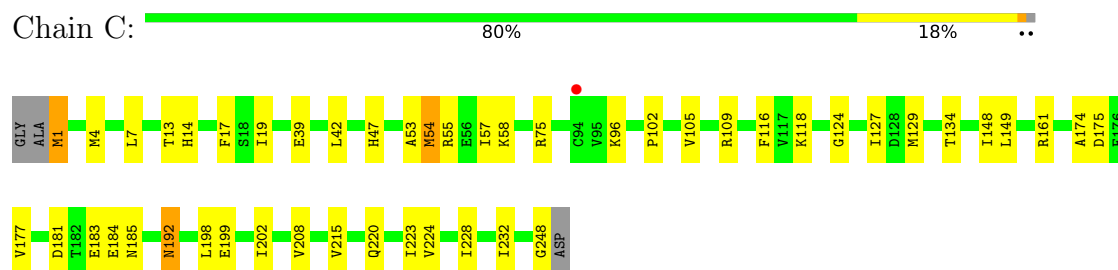
• Molecule 1: Type III Pantothenate Kinase



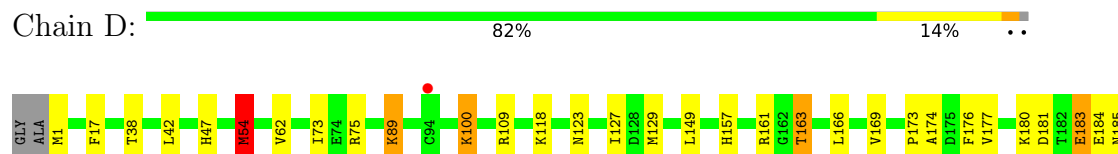
• Molecule 1: Type III Pantothenate Kinase



• Molecule 1: Type III Pantothenate Kinase

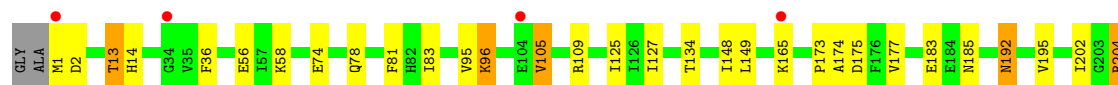
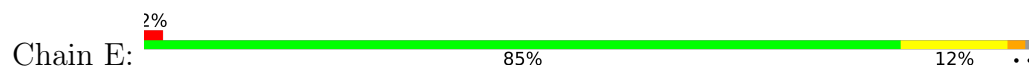


• Molecule 1: Type III Pantothenate Kinase

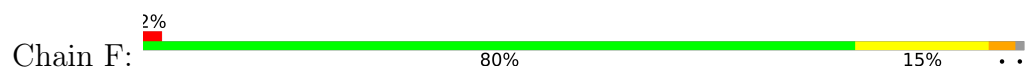




• Molecule 1: Type III Pantothenate Kinase



• Molecule 1: Type III Pantothenate Kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	75.11Å 137.79Å 75.22Å 90.00° 109.23° 90.00°	Depositor
Resolution (Å)	31.56 – 2.00 31.56 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (31.56-2.00) 97.7 (31.56-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.59 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.188 , 0.245 0.187 , 0.243	Depositor DCC
R_{free} test set	4763 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	22.1	Xtriage
Anisotropy	0.045	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 41.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.138 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12919	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.03% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.81	0/1971	0.92	0/2668
1	B	0.79	0/1971	0.91	1/2668 (0.0%)
1	C	0.79	0/1971	0.91	0/2668
1	D	0.84	0/1971	0.92	1/2668 (0.0%)
1	E	0.81	1/1971 (0.1%)	0.92	2/2668 (0.1%)
1	F	0.80	0/1971	0.93	3/2668 (0.1%)
All	All	0.81	1/11826 (0.0%)	0.92	7/16008 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	105	VAL	CA-CB	5.18	1.60	1.54

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	215	VAL	CB-CA-C	-7.84	99.14	110.83
1	F	212	LEU	CA-C-N	-5.71	113.90	119.78
1	F	212	LEU	C-N-CA	-5.71	113.90	119.78
1	F	224	VAL	CB-CA-C	-5.45	104.03	110.84
1	E	105	VAL	CB-CA-C	5.32	118.07	111.15
1	D	54	MET	N-CA-C	5.20	117.34	111.11
1	E	105	VAL	N-CA-CB	5.09	116.08	110.53

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	1929	0	1946	47	0
1	B	1929	0	1946	44	0
1	C	1929	0	1946	49	0
1	D	1929	0	1946	57	0
1	E	1929	0	1946	41	0
1	F	1929	0	1946	38	0
2	A	227	0	0	6	0
2	B	223	0	0	7	0
2	C	245	0	0	11	0
2	D	219	0	0	8	0
2	E	200	0	0	6	0
2	F	231	0	0	9	0
All	All	12919	0	11676	252	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 11.

All (252) 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:D:129:MET:CE	1:D:216:LEU:HD11	1.61	1.29
1:B:173:PRO:CB	1:B:227:MET:HE1	1.69	1.22
1:B:221:SER:HB3	2:B:469:HOH:O	1.45	1.15
1:D:224:VAL:HG12	1:D:227:MET:HE3	1.33	1.09
1:E:127:ILE:HD11	1:E:202:ILE:HD11	1.35	1.07
1:D:129:MET:HE3	1:D:216:LEU:HD11	1.34	1.06
1:A:157:HIS:CE1	1:A:161:ARG:HD3	1.91	1.06
1:B:127:ILE:HD11	1:B:202:ILE:HD11	1.38	1.03
1:B:173:PRO:HB3	1:B:227:MET:CE	1.92	1.00
1:C:127:ILE:HD11	1:C:202:ILE:HD11	1.41	0.99
1:F:127:ILE:HD11	1:F:202:ILE:HD11	1.46	0.98
1:B:173:PRO:HB3	1:B:227:MET:HE1	0.95	0.94
1:B:163:THR:CG2	1:B:166:LEU:H	1.80	0.94
1:E:204:ARG:HG3	1:F:177:VAL:HG23	1.50	0.94
1:B:174:ALA:H	1:B:192:ASN:HD21	1.14	0.93
1:C:102:PRO:O	1:C:105:VAL:HG12	1.70	0.92
1:B:47:HIS:HD2	2:B:465:HOH:O	1.51	0.92
1:E:174:ALA:H	1:E:192:ASN:HD21	1.16	0.90
1:F:174:ALA:H	1:F:192:ASN:HD21	1.16	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:174:ALA:H	1:C:192:ASN:HD21	1.17	0.88
1:A:174:ALA:H	1:A:192:ASN:ND2	1.71	0.87
1:F:248:GLY:HA3	2:F:474:HOH:O	1.73	0.86
1:D:157:HIS:CE1	1:D:161:ARG:HD2	2.10	0.86
1:A:174:ALA:H	1:A:192:ASN:HD21	0.92	0.86
1:E:224:VAL:HA	1:E:227:MET:HE3	1.59	0.84
1:F:78:GLN:HG2	2:F:356:HOH:O	1.78	0.83
1:A:174:ALA:N	1:A:192:ASN:HD21	1.76	0.83
1:D:163:THR:CG2	1:D:166:LEU:H	1.90	0.83
1:E:13:THR:HG22	1:E:14:HIS:ND1	1.94	0.83
1:D:176:PHE:CZ	1:D:184:GLU:HB2	2.14	0.82
1:D:224:VAL:HG12	1:D:227:MET:CE	2.09	0.82
1:F:13:THR:HG23	1:F:14:HIS:ND1	1.94	0.82
1:D:129:MET:CE	1:D:216:LEU:CD1	2.54	0.82
1:A:163:THR:HG21	1:A:166:LEU:HG	1.62	0.81
1:A:127:ILE:HD11	1:A:202:ILE:HD11	1.61	0.81
1:B:173:PRO:CB	1:B:227:MET:CE	2.54	0.81
1:A:68:THR:O	2:A:474:HOH:O	1.98	0.81
1:A:163:THR:CG2	1:A:166:LEU:H	1.94	0.79
1:E:58:LYS:HE2	2:E:366:HOH:O	1.83	0.79
1:A:208:VAL:HG21	1:B:177:VAL:HG21	1.65	0.78
1:E:204:ARG:HG3	1:F:177:VAL:CG2	2.12	0.78
1:D:47:HIS:HA	1:D:54:MET:HE2	1.65	0.77
1:F:118:LYS:HD3	1:F:244:HIS:HB3	1.66	0.77
1:A:134:THR:HG22	2:A:291:HOH:O	1.84	0.76
1:E:185:ASN:HD21	1:F:109:ARG:HH22	1.32	0.76
1:A:13:THR:CG2	1:A:14:HIS:ND1	2.49	0.76
1:D:163:THR:HG21	1:D:166:LEU:HB2	1.67	0.76
1:D:163:THR:HG23	1:D:166:LEU:H	1.48	0.76
1:C:109:ARG:HH22	1:D:185:ASN:ND2	1.86	0.74
1:D:129:MET:HE3	1:D:216:LEU:CD1	2.14	0.74
1:A:170:GLU:HG3	2:A:472:HOH:O	1.89	0.73
1:C:4:MET:HE3	1:C:19:ILE:HG21	1.71	0.73
1:D:174:ALA:H	1:D:192:ASN:HD21	1.35	0.72
1:D:129:MET:HE1	1:D:216:LEU:HD11	1.68	0.72
1:B:127:ILE:HD11	1:B:202:ILE:CD1	2.16	0.72
1:A:82:HIS:ND1	2:A:334:HOH:O	2.23	0.72
1:E:109:ARG:HH22	1:F:185:ASN:ND2	1.88	0.72
1:C:39:GLU:OE1	1:C:75:ARG:NH1	2.23	0.71
1:A:157:HIS:CE1	1:A:161:ARG:CD	2.72	0.71
1:B:163:THR:HG21	1:B:166:LEU:H	1.56	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:183:GLU:HG3	2:E:355:HOH:O	1.92	0.69
1:C:129:MET:HE2	1:C:198:LEU:CD1	2.23	0.69
1:D:163:THR:HG21	1:D:166:LEU:CB	2.21	0.69
1:D:157:HIS:HE1	1:D:161:ARG:HD2	1.54	0.69
1:E:185:ASN:ND2	1:F:109:ARG:HH22	1.91	0.68
1:D:181:ASP:OD1	1:D:184:GLU:HG2	1.94	0.67
1:A:157:HIS:HE1	1:A:161:ARG:HD3	1.55	0.67
1:C:223:ILE:HG13	1:C:224:VAL:HG13	1.75	0.67
1:E:127:ILE:CD1	1:E:202:ILE:HD11	2.20	0.67
1:C:1:MET:SD	2:C:280:HOH:O	2.51	0.67
1:D:75:ARG:NH2	2:D:440:HOH:O	2.26	0.67
1:A:13:THR:HG23	1:A:14:HIS:ND1	2.09	0.67
1:A:223:ILE:HD12	1:A:224:VAL:HG13	1.76	0.67
1:A:226:ASP:OD1	1:A:227:MET:HG3	1.94	0.67
1:B:47:HIS:CD2	2:B:465:HOH:O	2.36	0.67
1:A:185:ASN:ND2	1:B:109:ARG:HH22	1.93	0.66
1:B:127:ILE:CD1	1:B:202:ILE:HD11	2.22	0.66
1:D:183:GLU:HB3	2:D:346:HOH:O	1.96	0.66
1:A:177:VAL:HG21	1:B:208:VAL:HG21	1.77	0.66
1:B:71:THR:HG22	1:B:75:ARG:HE	1.61	0.66
1:D:118:LYS:HE3	2:D:309:HOH:O	1.96	0.66
1:E:13:THR:HG22	1:E:14:HIS:CE1	2.31	0.65
1:C:199:GLU:OE1	2:C:487:HOH:O	2.14	0.65
1:A:163:THR:HG21	1:A:166:LEU:CG	2.26	0.65
1:B:159:LEU:O	1:B:163:THR:HB	1.96	0.65
1:C:109:ARG:HH22	1:D:185:ASN:HD21	1.43	0.65
1:C:118:LYS:HG2	2:C:333:HOH:O	1.96	0.65
1:F:94:CYS:HB2	2:F:416:HOH:O	1.95	0.64
1:F:13:THR:CG2	1:F:14:HIS:ND1	2.61	0.64
1:F:192:ASN:H	1:F:192:ASN:HD22	1.46	0.63
1:A:215:VAL:HG22	1:A:232:ILE:HB	1.80	0.63
1:D:129:MET:HE2	1:D:216:LEU:HD11	1.71	0.63
1:C:134:THR:HG22	2:C:269:HOH:O	1.99	0.63
1:C:127:ILE:HD11	1:C:202:ILE:CD1	2.23	0.62
1:D:229:LYS:HB3	2:D:456:HOH:O	1.98	0.62
1:E:109:ARG:HH22	1:F:185:ASN:HD21	1.48	0.62
1:E:2:ASP:HB3	2:E:275:HOH:O	2.00	0.61
1:E:174:ALA:H	1:E:192:ASN:ND2	1.95	0.61
1:C:4:MET:HE2	1:C:57:ILE:HG13	1.82	0.61
1:E:165:LYS:HE2	2:E:328:HOH:O	2.00	0.61
1:C:183:GLU:HG3	2:C:315:HOH:O	1.99	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:173:PRO:HG3	1:D:227:MET:CE	2.31	0.60
1:A:47:HIS:HA	1:A:54:MET:HE2	1.82	0.60
1:E:74:GLU:O	1:E:78:GLN:HG2	2.00	0.60
1:F:3:PRO:HG2	2:F:447:HOH:O	2.01	0.60
1:C:4:MET:HE2	1:C:57:ILE:CG1	2.31	0.60
1:B:163:THR:HG21	1:B:166:LEU:CB	2.32	0.60
1:D:207:GLU:HB2	2:D:465:HOH:O	2.01	0.59
1:B:71:THR:CG2	1:B:75:ARG:HE	2.15	0.59
1:E:173:PRO:HG3	1:E:227:MET:CE	2.33	0.58
1:A:47:HIS:HA	1:A:54:MET:CE	2.33	0.58
1:D:1:MET:N	2:D:251:HOH:O	2.29	0.58
1:E:173:PRO:HG3	1:E:227:MET:HE1	1.85	0.58
1:E:109:ARG:NH1	1:E:148:ILE:HD11	2.20	0.57
1:A:109:ARG:HH22	1:B:185:ASN:ND2	2.03	0.57
1:A:129:MET:HE2	1:A:198:LEU:CD1	2.35	0.57
1:F:172:LYS:HD2	2:F:431:HOH:O	2.04	0.57
1:B:163:THR:HG21	1:B:166:LEU:HB2	1.86	0.57
1:B:192:ASN:HD22	1:B:192:ASN:H	1.52	0.56
1:F:47:HIS:HA	1:F:54:MET:CE	2.35	0.56
1:B:71:THR:HG21	1:B:75:ARG:HH21	1.69	0.56
1:F:13:THR:HG23	1:F:14:HIS:CE1	2.41	0.56
1:E:223:ILE:HG13	1:E:224:VAL:HG23	1.86	0.56
1:C:185:ASN:ND2	1:D:109:ARG:HH22	2.03	0.56
1:F:92:ASN:C	1:F:92:ASN:HD22	2.14	0.56
1:D:127:ILE:HD11	1:D:202:ILE:HD11	1.87	0.55
1:A:159:LEU:O	1:A:163:THR:HB	2.06	0.55
1:C:183:GLU:CG	2:C:315:HOH:O	2.55	0.55
1:A:140:ASN:HB2	2:A:288:HOH:O	2.07	0.55
1:C:13:THR:HG23	1:E:36:PHE:CD1	2.42	0.55
1:F:109:ARG:NH2	1:F:136:ASP:OD1	2.36	0.55
1:E:96:LYS:HE3	2:E:433:HOH:O	2.06	0.54
1:C:13:THR:HG23	1:E:36:PHE:CG	2.43	0.54
1:A:229:LYS:HG2	2:A:392:HOH:O	2.06	0.54
1:D:47:HIS:HA	1:D:54:MET:CE	2.37	0.54
1:A:163:THR:HG21	1:A:166:LEU:CB	2.37	0.54
1:E:81:PHE:HB3	1:E:83:ILE:HD12	1.90	0.54
1:A:102:PRO:O	1:A:105:VAL:HG13	2.08	0.54
1:C:208:VAL:HG21	1:D:177:VAL:HG21	1.89	0.54
1:C:4:MET:HE3	1:C:19:ILE:CG2	2.35	0.53
1:D:174:ALA:H	1:D:192:ASN:ND2	2.05	0.53
1:E:192:ASN:HD22	1:E:192:ASN:H	1.57	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:192:ASN:HD22	1:F:192:ASN:N	2.06	0.53
1:C:1:MET:N	2:C:250:HOH:O	2.28	0.52
1:C:58:LYS:HE3	2:C:387:HOH:O	2.09	0.52
1:D:224:VAL:CG1	1:D:227:MET:HE3	2.23	0.52
1:D:62:VAL:HG21	1:D:73:ILE:HG21	1.92	0.50
1:B:109:ARG:NH2	1:B:136:ASP:OD1	2.34	0.50
1:A:13:THR:HG22	1:A:14:HIS:ND1	2.24	0.50
1:B:134:THR:HG22	2:B:283:HOH:O	2.11	0.50
1:C:174:ALA:H	1:C:192:ASN:ND2	1.97	0.50
1:E:127:ILE:HD11	1:E:202:ILE:CD1	2.25	0.50
1:C:127:ILE:CD1	1:C:202:ILE:HD11	2.29	0.50
1:A:13:THR:HG23	1:A:14:HIS:CE1	2.47	0.49
1:F:224:VAL:HG13	1:F:227:MET:HE3	1.94	0.49
1:B:192:ASN:HD22	1:B:192:ASN:N	2.10	0.49
1:D:173:PRO:HG3	1:D:227:MET:HE1	1.94	0.49
1:F:127:ILE:CD1	1:F:202:ILE:HD11	2.30	0.49
1:C:129:MET:CE	1:C:198:LEU:CD1	2.89	0.49
1:A:185:ASN:HD21	1:B:109:ARG:HH22	1.60	0.49
1:D:176:PHE:CE1	1:D:184:GLU:HB2	2.48	0.49
1:C:181:ASP:OD1	1:C:184:GLU:HG3	2.13	0.49
1:F:74:GLU:HG2	2:F:425:HOH:O	2.13	0.49
1:C:13:THR:HG22	1:C:14:HIS:CE1	2.48	0.48
1:A:32:SER:HB3	1:F:36:PHE:O	2.13	0.48
1:E:13:THR:CG2	1:E:14:HIS:CE1	2.96	0.48
1:A:199:GLU:HG3	1:A:228:ILE:HD13	1.95	0.48
1:D:163:THR:HG21	1:D:166:LEU:H	1.71	0.48
1:E:134:THR:HG22	1:E:148:ILE:HG12	1.95	0.48
1:F:123:ASN:HB3	1:F:212:LEU:HD22	1.96	0.48
1:C:185:ASN:HD21	1:D:109:ARG:HH22	1.63	0.47
1:C:4:MET:HE1	1:C:53:ALA:HB1	1.95	0.47
1:B:71:THR:CG2	1:B:75:ARG:HH21	2.28	0.47
1:B:1:MET:N	2:B:250:HOH:O	2.31	0.47
1:B:163:THR:HG21	1:B:166:LEU:HG	1.95	0.47
1:C:192:ASN:H	1:C:192:ASN:HD22	1.63	0.47
1:E:174:ALA:N	1:E:192:ASN:HD21	1.99	0.47
1:E:195:VAL:HG22	1:E:224:VAL:HG11	1.97	0.47
1:F:47:HIS:HA	1:F:54:MET:HE1	1.96	0.47
1:E:224:VAL:HG22	1:E:227:MET:CE	2.45	0.47
1:B:102:PRO:O	1:B:105:VAL:HG13	2.14	0.46
1:B:135:VAL:O	1:B:146:GLY:HA3	2.15	0.46
1:D:163:THR:HG21	1:D:166:LEU:CG	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:173:PRO:HG3	1:D:227:MET:HE2	1.96	0.46
1:A:163:THR:HG22	1:A:166:LEU:H	1.79	0.46
1:A:223:ILE:CD1	1:A:224:VAL:HG13	2.45	0.46
2:C:286:HOH:O	1:D:204:ARG:HG2	2.14	0.46
1:D:176:PHE:CZ	1:D:184:GLU:CB	2.93	0.46
1:B:105:VAL:HG22	1:B:110:VAL:HG23	1.97	0.46
1:D:174:ALA:N	1:D:192:ASN:HD21	2.10	0.46
1:D:174:ALA:HB1	1:D:176:PHE:CE1	2.51	0.46
1:B:174:ALA:H	1:B:192:ASN:ND2	1.97	0.45
1:A:116:PHE:CZ	1:A:124:GLY:HA3	2.51	0.45
1:D:89:LYS:HE3	1:D:89:LYS:HB2	1.56	0.45
1:D:173:PRO:CG	1:D:227:MET:HE2	2.46	0.45
1:B:157:HIS:CE1	1:B:161:ARG:HD2	2.51	0.45
1:C:129:MET:HE2	1:C:198:LEU:HD11	1.96	0.45
1:D:17:PHE:CZ	1:D:42:LEU:HD11	2.51	0.45
1:C:181:ASP:HB3	1:D:100:LYS:HD2	1.99	0.45
1:A:109:ARG:NH2	1:A:136:ASP:OD1	2.49	0.45
1:A:163:THR:HG21	1:A:166:LEU:H	1.78	0.45
1:E:192:ASN:HD22	1:E:192:ASN:N	2.14	0.45
1:A:2:ASP:HB3	1:A:21:GLU:O	2.17	0.44
1:C:4:MET:CE	1:C:19:ILE:HG21	2.45	0.44
1:E:224:VAL:HG12	1:E:224:VAL:O	2.17	0.44
1:B:163:THR:HG21	1:B:166:LEU:N	2.30	0.44
1:C:177:VAL:HG21	1:D:208:VAL:HG21	2.00	0.44
1:F:17:PHE:CZ	1:F:42:LEU:HD11	2.52	0.44
1:C:13:THR:CG2	1:E:36:PHE:CG	3.01	0.43
1:C:215:VAL:HG22	1:C:232:ILE:HD12	2.01	0.43
1:C:248:GLY:HA3	2:C:483:HOH:O	2.18	0.43
1:C:47:HIS:HA	1:C:54:MET:CE	2.48	0.43
1:C:96:LYS:HD3	2:C:491:HOH:O	2.19	0.43
1:A:163:THR:HG21	1:A:166:LEU:HB2	1.99	0.43
1:F:78:GLN:CG	2:F:356:HOH:O	2.53	0.43
1:F:86:ILE:HD12	1:F:247:PHE:CE1	2.53	0.43
1:A:127:ILE:CD1	1:A:202:ILE:HD11	2.42	0.43
1:B:171:VAL:HG13	1:B:191:VAL:HG22	2.01	0.43
1:C:109:ARG:NH1	1:C:148:ILE:HD11	2.34	0.43
1:B:47:HIS:HE1	2:B:463:HOH:O	2.01	0.42
1:C:17:PHE:CZ	1:C:42:LEU:HD11	2.54	0.42
1:F:13:THR:HG22	2:F:259:HOH:O	2.18	0.42
1:D:38:THR:HB	2:D:265:HOH:O	2.18	0.42
1:B:96:LYS:HA	1:B:96:LYS:HD2	1.91	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:220:GLN:O	1:C:223:ILE:HG12	2.19	0.42
1:A:212:LEU:H	1:A:212:LEU:HG	1.57	0.42
1:A:129:MET:HE2	1:A:198:LEU:HD11	2.00	0.42
1:C:116:PHE:CZ	1:C:124:GLY:HA3	2.55	0.42
1:B:47:HIS:HA	1:B:54:MET:CE	2.49	0.42
1:C:175:ASP:OD1	1:C:175:ASP:N	2.53	0.41
1:D:123:ASN:HB3	1:D:212:LEU:HD22	2.01	0.41
1:E:177:VAL:HG21	1:F:208:VAL:HG21	2.01	0.41
1:E:224:VAL:HG22	1:E:227:MET:HE1	2.03	0.41
1:F:62:VAL:HG21	1:F:73:ILE:HG21	2.02	0.41
1:C:148:ILE:HG13	1:D:185:ASN:HB3	2.03	0.41
1:D:180:LYS:H	1:D:184:GLU:HG3	1.85	0.41
1:E:74:GLU:OE1	2:E:439:HOH:O	2.22	0.41
1:F:105:VAL:HG22	1:F:110:VAL:HG23	2.02	0.41
1:F:195:VAL:HG13	1:F:228:ILE:HG12	2.01	0.41
1:B:216:LEU:HD13	2:B:416:HOH:O	2.21	0.41
1:E:195:VAL:CG2	1:E:224:VAL:HG11	2.50	0.41
1:F:39:GLU:HB2	1:F:76:PHE:HA	2.03	0.41
1:A:223:ILE:H	1:A:223:ILE:HG13	1.71	0.41
1:D:163:THR:HG21	1:D:166:LEU:HG	2.02	0.41
1:D:222:LYS:HB2	1:D:222:LYS:HE2	1.69	0.41
1:F:69:GLN:NE2	2:F:299:HOH:O	2.54	0.41
1:B:215:VAL:HG13	1:B:232:ILE:HB	2.03	0.41
1:B:163:THR:HG21	1:B:166:LEU:CG	2.50	0.40
1:C:192:ASN:HD22	1:C:192:ASN:N	2.18	0.40
1:D:204:ARG:HD2	2:D:465:HOH:O	2.20	0.40
1:E:125:ILE:HB	1:E:214:VAL:HG22	2.03	0.40
1:D:169:VAL:HG22	1:D:183:GLU:HG3	2.03	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	246/251 (98%)	243 (99%)	3 (1%)	0	100	100
1	B	246/251 (98%)	241 (98%)	5 (2%)	0	100	100
1	C	246/251 (98%)	242 (98%)	4 (2%)	0	100	100
1	D	246/251 (98%)	243 (99%)	3 (1%)	0	100	100
1	E	246/251 (98%)	242 (98%)	3 (1%)	1 (0%)	30	27
1	F	246/251 (98%)	243 (99%)	3 (1%)	0	100	100
All	All	1476/1506 (98%)	1454 (98%)	21 (1%)	1 (0%)	48	46

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	95	VAL

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	211/212 (100%)	201 (95%)	10 (5%)	23	22
1	B	211/212 (100%)	199 (94%)	12 (6%)	18	15
1	C	211/212 (100%)	203 (96%)	8 (4%)	29	29
1	D	211/212 (100%)	203 (96%)	8 (4%)	29	29
1	E	211/212 (100%)	202 (96%)	9 (4%)	26	25
1	F	211/212 (100%)	198 (94%)	13 (6%)	16	13
All	All	1266/1272 (100%)	1206 (95%)	60 (5%)	23	22

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	13	THR
1	A	27	ARG
1	A	100	LYS

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Mol	Chain	Res	Type
1	A	105	VAL
1	A	134	THR
1	A	149	LEU
1	A	163	THR
1	A	183	GLU
1	A	223	ILE
1	B	1	MET
1	B	54	MET
1	B	94	CYS
1	B	105	VAL
1	B	149	LEU
1	B	163	THR
1	B	175	ASP
1	B	192	ASN
1	B	206	LYS
1	B	215	VAL
1	B	223	ILE
1	B	224	VAL
1	C	1	MET
1	C	7	LEU
1	C	54	MET
1	C	55	ARG
1	C	149	LEU
1	C	161	ARG
1	C	192	ASN
1	C	228	ILE
1	D	54	MET
1	D	89	LYS
1	D	100	LYS
1	D	149	LEU
1	D	163	THR
1	D	183	GLU
1	D	216	LEU
1	D	222	LYS
1	E	1	MET
1	E	13	THR
1	E	56	GLU
1	E	96	LYS
1	E	105	VAL
1	E	149	LEU
1	E	175	ASP
1	E	192	ASN

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Mol	Chain	Res	Type
1	E	204	ARG
1	F	2	ASP
1	F	13	THR
1	F	58	LYS
1	F	92	ASN
1	F	105	VAL
1	F	118	LYS
1	F	149	LEU
1	F	177	VAL
1	F	183	GLU
1	F	192	ASN
1	F	207	GLU
1	F	224	VAL
1	F	225	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	157	HIS
1	A	185	ASN
1	A	192	ASN
1	B	14	HIS
1	B	37	GLN
1	B	157	HIS
1	B	185	ASN
1	B	192	ASN
1	C	37	GLN
1	C	69	GLN
1	C	92	ASN
1	C	157	HIS
1	C	185	ASN
1	C	192	ASN
1	D	92	ASN
1	D	101	ASN
1	D	157	HIS
1	D	185	ASN
1	D	192	ASN
1	E	37	GLN
1	E	92	ASN
1	E	98	ASN
1	E	185	ASN
1	E	192	ASN

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Mol	Chain	Res	Type
1	F	45	HIS
1	F	92	ASN
1	F	98	ASN
1	F	185	ASN
1	F	192	ASN

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](#)

There are no ligands in this entry.

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	248/251 (98%)	-0.03	0	100	100	10, 22, 36, 41	0
1	B	248/251 (98%)	0.02	4 (1%)	70	70	11, 22, 39, 45	0
1	C	248/251 (98%)	-0.18	1 (0%)	88	88	12, 19, 33, 44	0
1	D	248/251 (98%)	-0.09	1 (0%)	88	88	11, 20, 35, 45	0
1	E	248/251 (98%)	-0.02	4 (1%)	70	70	12, 21, 35, 48	0
1	F	248/251 (98%)	-0.08	5 (2%)	65	64	11, 20, 36, 65	0
All	All	1488/1506 (98%)	-0.06	15 (1%)	79	79	10, 21, 36, 65	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	248	GLY	5.7
1	F	1	MET	4.0
1	C	94	CYS	3.1
1	F	224	VAL	2.8
1	B	248	GLY	2.7
1	F	2	ASP	2.4
1	B	94	CYS	2.4
1	B	35	VAL	2.3
1	D	94	CYS	2.2
1	E	34	GLY	2.2
1	E	1	MET	2.2
1	F	3	PRO	2.1
1	B	1	MET	2.0
1	E	104	GLU	2.0
1	E	165	LYS	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.