



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:50 AM GMT

PDB ID : 3CDK
Title : Crystal structure of the co-expressed succinyl-CoA transferase A and B complex from *Bacillus subtilis*
Authors : Kim, Y.; Zhou, M.; Stols, L.; Eschenfeldt, W.; Donnelly, M.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2008-02-27
Resolution : 2.59 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

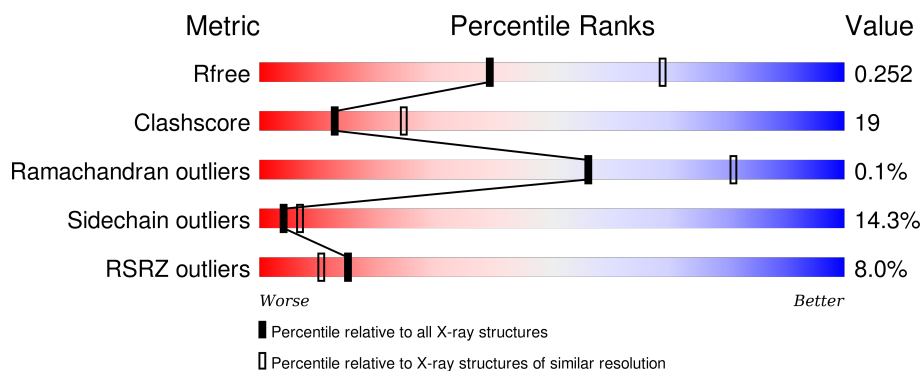
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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	241	
1	C	241	
2	B	219	
2	D	219	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6720 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 Succinyl-CoA:3-ketoacid-coenzyme A transferase subunit A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	229	Total	C	N	O	S	0	3	0
			1752	1104	309	334	5			
1	C	229	Total	C	N	O	S	0	0	0
			1725	1090	302	327	6			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	EXPRESSION TAG	UNP P42315
A	1	ASN	-	EXPRESSION TAG	UNP P42315
A	2	ALA	-	EXPRESSION TAG	UNP P42315
C	0	SER	-	EXPRESSION TAG	UNP P42315
C	1	ASN	-	EXPRESSION TAG	UNP P42315
C	2	ALA	-	EXPRESSION TAG	UNP P42315

- Molecule 2 is a protein called Succinyl-CoA:3-ketoacid-coenzyme A transferase subunit B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	212	Total	C	N	O	S	0	2	0
			1613	1008	276	316	13			
2	D	200	Total	C	N	O	S	0	1	0
			1515	951	256	295	13			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	SER	-	EXPRESSION TAG	UNP P42316
B	-1	ASN	-	EXPRESSION TAG	UNP P42316
B	0	ALA	-	EXPRESSION TAG	UNP P42316
D	-2	SER	-	EXPRESSION TAG	UNP P42316
D	-1	ASN	-	EXPRESSION TAG	UNP P42316
D	0	ALA	-	EXPRESSION TAG	UNP P42316

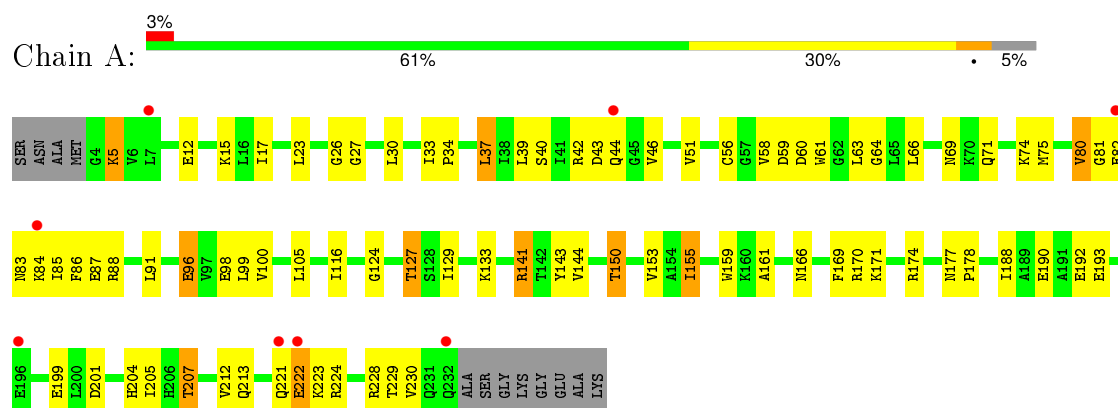
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	48	Total 48	O 48	0	0
3	B	26	Total 26	O 26	0	0
3	C	27	Total 27	O 27	0	0
3	D	14	Total 14	O 14	0	0

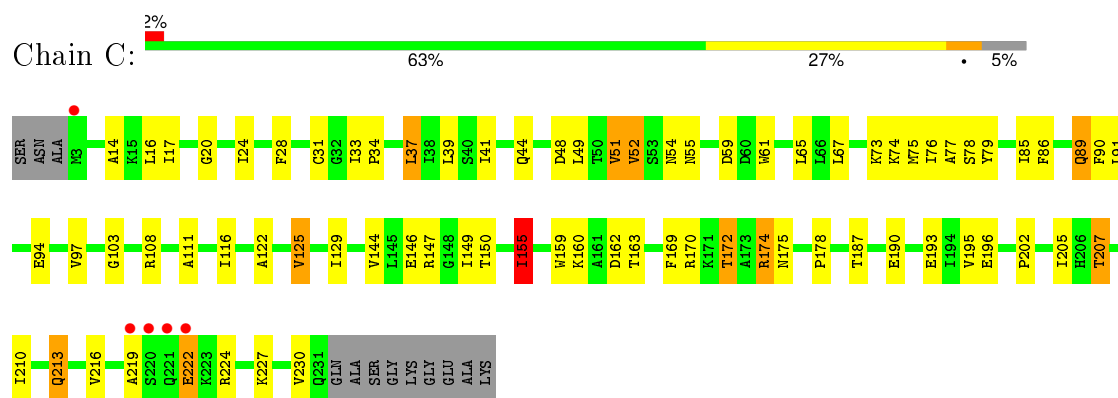
3 Residue-property plots [i](#)

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.

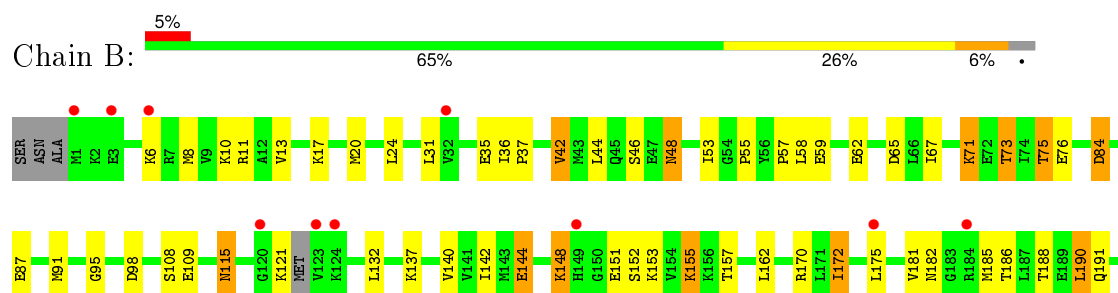
- Molecule 1: Succinyl-CoA:3-ketoacid-coenzyme A transferase subunit A

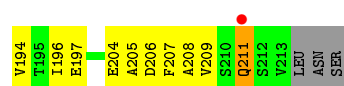


- Molecule 1: Succinyl-CoA:3-ketoacid-coenzyme A transferase subunit A

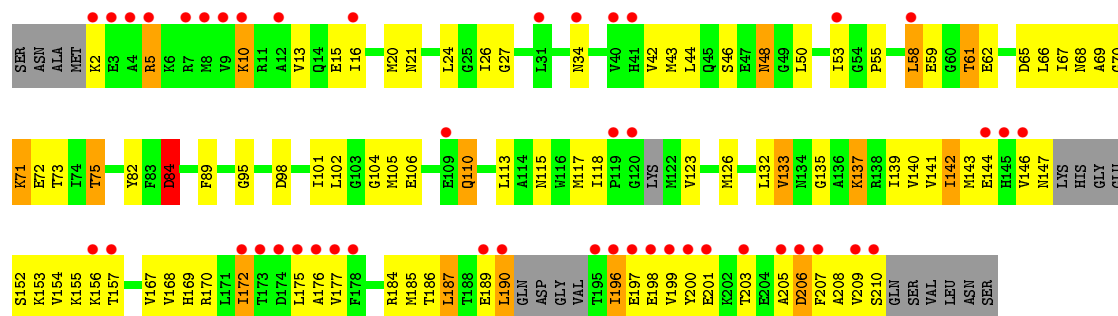


- Molecule 2: Succinyl-CoA:3-ketoacid-coenzyme A transferase subunit B





● Molecule 2: Succinyl-CoA:3-ketoacid-coenzyme A transferase subunit B



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	69.34Å 70.40Å 98.00Å 90.00° 106.31° 90.00°	Depositor
Resolution (Å)	48.30 – 2.59 48.30 – 2.59	Depositor EDS
% Data completeness (in resolution range)	99.2 (48.30-2.59) 99.2 (48.30-2.59)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.52 (at 2.58Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.193 , 0.253 0.193 , 0.252	Depositor DCC
R_{free} test set	1417 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	39.5	Xtriage
Anisotropy	0.636	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 49.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 28070 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6720	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.01% 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

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	0.79	2/1777 (0.1%)	0.81	0/2395
1	C	0.76	3/1750 (0.2%)	0.80	2/2359 (0.1%)
2	B	0.69	0/1632	0.78	0/2197
2	D	0.98	8/1531 (0.5%)	0.74	1/2059 (0.0%)
All	All	0.81	13/6690 (0.2%)	0.79	3/9010 (0.0%)

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	198	GLU	CD-OE2	23.59	1.51	1.25
2	D	201	GLU	CG-CD	8.08	1.64	1.51
2	D	190	LEU	C-O	7.88	1.38	1.23
2	D	201	GLU	CD-OE2	7.28	1.33	1.25
2	D	201	GLU	CD-OE1	7.08	1.33	1.25
1	C	222	GLU	CD-OE2	6.13	1.32	1.25
2	D	197	GLU	CD-OE1	5.88	1.32	1.25
2	D	152	SER	CB-OG	5.41	1.49	1.42
1	A	98	GLU	CG-CD	5.35	1.59	1.51
2	D	198	GLU	CD-OE1	-5.33	1.19	1.25
1	C	222	GLU	CG-CD	5.24	1.59	1.51
1	C	146	GLU	CG-CD	5.20	1.59	1.51
1	A	222	GLU	CB-CG	5.12	1.61	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	174	ARG	CB-CG-CD	-5.71	96.75	111.60
1	C	155	ILE	N-CA-C	-5.31	96.65	111.00
2	D	84	ASP	CB-CG-OD1	5.03	122.83	118.30

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	1752	0	1781	75	0
1	C	1725	0	1761	54	0
2	B	1613	0	1627	56	0
2	D	1515	0	1531	89	0
3	A	48	0	0	9	0
3	B	26	0	0	0	0
3	C	27	0	0	4	0
3	D	14	0	0	0	0
All	All	6720	0	6700	256	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 19.

All (256) 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:42[B]:ARG:HH21	1:A:42[B]:ARG:HG2	1.04	1.15
1:A:105:LEU:HB2	3:A:253:HOH:O	1.47	1.14
2:D:13:VAL:O	2:D:16:ILE:HG22	1.51	1.09
2:D:139:ILE:H	2:D:169:HIS:CD2	1.73	1.06
1:A:83:ASN:HB3	1:A:86:PHE:HB2	1.39	1.02
2:B:48:ASN:HD22	2:B:48:ASN:H	1.01	0.99
2:D:48:ASN:HD22	2:D:48:ASN:H	1.09	0.96
1:A:5:LYS:HE3	1:A:212:VAL:O	1.66	0.96
2:D:55:PRO:O	2:D:75:THR:HB	1.68	0.93
2:D:16:ILE:HG23	2:D:16:ILE:O	1.70	0.89
1:A:42[B]:ARG:HG2	1:A:42[B]:ARG:NH2	1.83	0.88
2:D:139:ILE:H	2:D:169:HIS:HD2	0.91	0.87
1:C:90:PHE:CE2	2:D:123:VAL:HG11	2.10	0.87
2:B:142:ILE:HG22	2:B:172:ILE:HG12	1.56	0.87
2:B:148:LYS:HD3	2:B:148:LYS:H	1.39	0.86
1:A:124:GLY:O	1:A:127:THR:HB	1.77	0.83
1:A:207:THR:HB	3:A:271:HOH:O	1.80	0.81
2:B:148:LYS:HD3	2:B:148:LYS:N	1.98	0.78
2:D:200:TYR:CD1	2:D:207:PHE:CE1	2.72	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:185:MET:HE3	2:B:207:PHE:HB3	1.66	0.77
1:A:80:VAL:HG21	1:A:86:PHE:CE1	2.20	0.77
1:A:83:ASN:HB3	1:A:86:PHE:CB	2.15	0.76
2:B:48:ASN:ND2	2:B:48:ASN:H	1.78	0.75
2:D:139:ILE:N	2:D:169:HIS:HD2	1.76	0.75
2:D:200:TYR:HD1	2:D:207:PHE:CZ	2.05	0.74
1:A:37:LEU:HD22	1:A:190:GLU:HB2	1.69	0.74
1:A:40:SER:O	1:A:44:GLN:HG2	1.87	0.74
2:B:190:LEU:HG	2:B:194:VAL:HG23	1.70	0.74
2:B:115:ASN:C	2:B:115:ASN:HD22	1.90	0.73
2:D:185:MET:HE2	2:D:207:PHE:HB3	1.69	0.73
1:A:17:ILE:HD12	1:A:153:VAL:HG11	1.72	0.72
1:C:172:THR:HG23	3:C:246:HOH:O	1.90	0.71
2:B:148:LYS:H	2:B:148:LYS:CD	2.00	0.71
1:A:155:ILE:HD11	1:A:188:ILE:HD12	1.73	0.71
2:D:48:ASN:HD22	2:D:48:ASN:N	1.85	0.70
1:A:166:ASN:HB3	1:A:205:ILE:HD13	1.72	0.70
1:C:147:ARG:HD3	3:C:255:HOH:O	1.91	0.70
2:D:200:TYR:HD1	2:D:207:PHE:CE1	2.07	0.70
1:A:105:LEU:CB	3:A:253:HOH:O	2.21	0.70
1:A:201:ASP:HB3	1:A:204:HIS:CD2	2.27	0.69
2:B:75:THR:CG2	2:B:76:GLU:N	2.55	0.69
1:C:187:THR:H	1:C:213:GLN:HG2	1.57	0.69
1:C:202:PRO:HA	1:C:205:ILE:HD12	1.76	0.68
2:B:142:ILE:HG22	2:B:172:ILE:CG1	2.24	0.68
1:A:127:THR:CG2	1:A:129:ILE:HB	2.24	0.67
2:B:11:ARG:HB3	2:B:172:ILE:HD13	1.76	0.67
1:C:111:ALA:HB1	1:C:116:ILE:HB	1.76	0.67
2:D:15:GLU:OE2	2:D:170:ARG:NH1	2.19	0.67
1:A:17:ILE:CD1	1:A:153:VAL:HG11	2.26	0.66
1:C:169:PHE:HE2	1:C:207:THR:CG2	2.08	0.65
1:C:90:PHE:HE2	2:D:123:VAL:HG11	1.56	0.65
1:C:73:LYS:NZ	3:C:257:HOH:O	2.24	0.65
1:C:172:THR:HG21	2:D:67:ILE:H	1.62	0.65
2:D:20:MET:O	2:D:42:VAL:HG13	1.97	0.65
2:D:58:LEU:O	2:D:61:THR:HG22	1.97	0.64
2:D:139:ILE:N	2:D:169:HIS:CD2	2.58	0.64
2:D:200:TYR:CE1	2:D:207:PHE:CE1	2.86	0.64
1:C:89:GLN:HG2	1:C:94:GLU:HB3	1.80	0.64
2:D:48:ASN:ND2	2:D:48:ASN:H	1.89	0.64
2:B:11:ARG:HB3	2:B:172:ILE:CD1	2.29	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:48:ASN:HD22	2:B:48:ASN:N	1.83	0.63
2:D:203:THR:HB	2:D:207:PHE:HE2	1.64	0.63
1:C:187:THR:H	1:C:213:GLN:CG	2.12	0.62
2:B:115:ASN:ND2	2:B:115:ASN:C	2.52	0.62
1:A:80:VAL:HG12	1:A:81:GLY:N	2.14	0.61
1:C:122:ALA:O	1:C:125:VAL:HG13	2.00	0.61
1:A:80:VAL:CG2	1:A:86:PHE:CZ	2.84	0.61
2:B:211:GLN:NE2	2:B:211:GLN:HA	2.15	0.61
1:A:42[B]:ARG:HH21	1:A:42[B]:ARG:CG	1.96	0.60
2:D:156:LYS:HG3	2:D:205:ALA:HA	1.83	0.60
2:B:67:ILE:HD12	2:B:71:LYS:HA	1.84	0.60
1:A:30:LEU:HD13	1:A:224:ARG:HD3	1.84	0.60
2:B:75:THR:HG23	2:B:76:GLU:N	2.17	0.60
1:A:169:PHE:HE1	1:A:207:THR:HG23	1.67	0.60
1:A:51:VAL:HG22	1:A:75:MET:SD	2.42	0.59
1:C:54:ASN:ND2	1:C:78:SER:OG	2.35	0.59
2:D:16:ILE:CG2	2:D:16:ILE:O	2.44	0.59
1:C:90:PHE:HD1	1:C:97:VAL:HG21	1.68	0.59
2:D:200:TYR:HA	2:D:207:PHE:HZ	1.68	0.59
2:D:13:VAL:O	2:D:16:ILE:CG2	2.41	0.59
2:B:55:PRO:O	2:B:75:THR:HB	2.01	0.59
2:B:211:GLN:HE21	2:B:211:GLN:HA	1.68	0.59
2:D:70:GLY:O	2:D:71:LYS:HB2	2.01	0.59
1:C:169:PHE:CE2	1:C:207:THR:CG2	2.85	0.59
1:A:127:THR:HG22	1:A:129:ILE:HB	1.84	0.59
1:A:169:PHE:HE1	1:A:207:THR:CG2	2.16	0.58
2:D:186:THR:CG2	2:D:208:ALA:O	2.51	0.58
2:D:190:LEU:HD21	2:D:196:ILE:HG23	1.84	0.58
2:D:186:THR:HG23	2:D:208:ALA:O	2.04	0.58
1:C:37:LEU:HD22	1:C:190:GLU:HB2	1.86	0.58
2:D:200:TYR:CD1	2:D:207:PHE:CZ	2.90	0.57
1:A:229:THR:O	2:B:73:THR:HG22	2.05	0.57
1:C:51:VAL:HG22	1:C:75:MET:SD	2.44	0.57
1:A:83:ASN:OD1	1:A:85:ILE:HB	2.04	0.57
1:C:172:THR:HG22	2:D:65:ASP:O	2.05	0.57
2:D:186:THR:HA	2:D:208:ALA:O	2.05	0.57
2:B:108:SER:HA	2:B:155:LYS:O	2.05	0.57
1:C:67:LEU:CD2	1:C:85:ILE:HG22	2.35	0.56
2:D:133:VAL:HG13	2:D:167:VAL:HG11	1.86	0.56
2:B:57:PRO:HG3	2:B:73:THR:HG23	1.87	0.56
2:B:98:ASP:OD1	2:B:137:LYS:HE3	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:ARG:HD3	3:A:248:HOH:O	2.05	0.56
2:D:203:THR:HG22	2:D:205:ALA:H	1.71	0.56
2:B:191:GLN:O	2:B:194:VAL:HG22	2.06	0.56
1:C:219:ALA:HB1	1:C:222:GLU:OE1	2.06	0.56
2:D:102:LEU:O	2:D:141:VAL:HA	2.06	0.55
1:C:103:GLY:HA2	2:D:89:PHE:CG	2.42	0.55
2:B:46:SER:OG	2:B:48:ASN:ND2	2.39	0.55
2:D:156:LYS:HE3	2:D:206:ASP:HB2	1.89	0.55
1:A:169:PHE:CE1	1:A:207:THR:HG23	2.42	0.55
1:C:227:LYS:HB2	2:D:72:GLU:HA	1.88	0.55
1:A:27:GLY:HA3	1:A:33:ILE:HG22	1.89	0.55
1:C:20:GLY:HA2	1:C:48:ASP:O	2.07	0.54
2:B:144:GLU:O	2:B:153:LYS:HD3	2.08	0.54
2:B:155:LYS:HA	2:B:204:GLU:HB2	1.89	0.54
2:D:117:MET:HG2	2:D:118:ILE:N	2.23	0.53
2:D:42:VAL:CG1	2:D:43:MET:N	2.71	0.53
2:D:177:VAL:HB	2:D:189:GLU:HB3	1.90	0.53
2:D:200:TYR:CD1	2:D:207:PHE:HE1	2.25	0.53
1:A:26:GLY:HA3	1:A:177:ASN:OD1	2.08	0.53
1:A:100:VAL:HG12	3:A:253:HOH:O	2.08	0.53
1:C:34:PRO:CG	1:C:155:ILE:HG23	2.38	0.53
1:A:12:GLU:HA	1:A:15:LYS:HE3	1.91	0.53
2:D:142:ILE:HG22	2:D:172:ILE:HG22	1.90	0.53
1:A:221:GLN:HG3	3:A:287:HOH:O	2.07	0.53
1:A:83:ASN:CB	1:A:86:PHE:HB2	2.24	0.53
2:B:109:GLU:OE2	2:B:206:ASP:HB2	2.08	0.53
1:C:162:ASP:HA	1:C:195:VAL:O	2.09	0.53
2:B:175:LEU:O	2:B:194:VAL:HG21	2.07	0.53
2:D:68:ASN:ND2	2:D:72:GLU:OE1	2.35	0.52
2:D:154:VAL:HB	2:D:203:THR:HG23	1.91	0.52
1:A:133:LYS:HB2	1:A:144:VAL:CG2	2.40	0.52
2:B:142:ILE:CG2	2:B:172:ILE:HG12	2.36	0.52
1:C:24:ILE:HG12	1:C:52:VAL:HG13	1.92	0.52
1:A:222:GLU:HA	3:A:288:HOH:O	2.08	0.51
2:D:21:ASN:N	2:D:98:ASP:OD2	2.36	0.51
2:B:20:MET:O	2:B:42:VAL:HG22	2.10	0.51
1:A:69:ASN:ND2	1:A:71:GLN:OE1	2.38	0.51
1:A:42[B]:ARG:NH2	1:A:43:ASP:OD1	2.44	0.51
1:A:58:VAL:O	1:A:64:GLY:HA3	2.11	0.51
2:B:185:MET:HE1	2:B:205:ALA:HB3	1.93	0.51
1:A:80:VAL:HG21	1:A:86:PHE:CZ	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:ASN:CB	1:A:205:ILE:HD13	2.41	0.51
1:A:34:PRO:HG2	1:A:155:ILE:HG23	1.92	0.51
2:D:2:LYS:HE3	2:D:5:ARG:HD2	1.92	0.50
2:D:98:ASP:O	2:D:137:LYS:HG3	2.11	0.50
1:A:37:LEU:CD2	1:A:190:GLU:HB2	2.40	0.50
2:B:75:THR:HG23	2:B:76:GLU:H	1.76	0.50
2:B:8:MET:O	2:B:172:ILE:HD11	2.12	0.50
2:B:95:GLY:O	2:B:137:LYS:HE2	2.12	0.50
1:C:55:ASN:ND2	1:C:79:TYR:HB3	2.27	0.50
2:D:61:THR:O	2:D:61:THR:HG23	2.12	0.49
2:B:62:GLU:HB2	2:B:73:THR:HG21	1.94	0.49
1:A:141:ARG:HD3	1:A:143:TYR:OH	2.13	0.49
1:A:192:GLU:OE2	1:A:223:LYS:NZ	2.41	0.49
1:A:127:THR:HG22	1:A:129:ILE:N	2.20	0.48
1:C:52:VAL:HB	1:C:76:ILE:HB	1.94	0.48
1:C:163:THR:HG23	1:C:196:GLU:HA	1.95	0.48
1:A:34:PRO:CG	1:A:155:ILE:HG23	2.43	0.48
1:A:87:GLU:O	1:A:91:LEU:HD23	2.13	0.48
2:B:91:MET:HE3	1:C:210:ILE:HD13	1.94	0.48
2:D:104:GLY:HA2	2:D:115:ASN:HB3	1.95	0.48
1:A:230:VAL:HG21	2:B:59:GLU:HG3	1.94	0.48
1:A:201:ASP:HB3	1:A:204:HIS:CG	2.48	0.48
1:A:80:VAL:CG1	1:A:81:GLY:N	2.76	0.47
2:D:50:LEU:HD13	2:D:82:TYR:CE1	2.48	0.47
2:B:8:MET:HA	2:B:172:ILE:HD11	1.97	0.47
1:C:34:PRO:HG2	1:C:155:ILE:HG23	1.96	0.47
2:B:13:VAL:HG11	2:B:35:GLU:O	2.14	0.47
2:B:46:SER:HB2	2:B:53:ILE:HD11	1.96	0.47
1:A:116:ILE:HD13	1:A:116:ILE:N	2.30	0.47
2:D:146:VAL:HG12	2:D:147:ASN:N	2.30	0.47
1:A:169:PHE:CE1	1:A:207:THR:CG2	2.96	0.46
1:A:133:LYS:HB2	1:A:144:VAL:HG22	1.97	0.46
2:D:155:LYS:O	2:D:205:ALA:HB2	2.15	0.46
1:C:34:PRO:HG3	1:C:155:ILE:HG23	1.98	0.46
1:C:67:LEU:HD23	1:C:85:ILE:HG22	1.97	0.46
1:C:28:PHE:CE2	2:D:69:ALA:HA	2.51	0.46
2:B:186:THR:HA	2:B:208:ALA:O	2.16	0.46
2:D:71:LYS:HD3	2:D:71:LYS:N	2.32	0.45
1:C:17:ILE:HD12	1:C:41:ILE:HD11	1.99	0.45
2:B:140:VAL:HG22	2:B:170:ARG:HB3	1.97	0.45
2:B:185:MET:HE2	2:B:205:ALA:HB1	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:PRO:HG2	2:B:84:ASP:HB3	1.98	0.45
2:D:203:THR:HG22	2:D:205:ALA:N	2.30	0.45
1:A:56:CYS:HB2	1:A:75:MET:HE2	1.98	0.45
2:B:57:PRO:CG	2:B:73:THR:HG23	2.46	0.45
1:C:174:ARG:HD2	3:C:267:HOH:O	2.15	0.45
2:D:133:VAL:HA	2:D:139:ILE:CD1	2.46	0.45
2:B:75:THR:HG22	2:B:76:GLU:N	2.32	0.45
2:D:186:THR:HG23	2:D:208:ALA:C	2.36	0.45
2:D:209:VAL:HG22	2:D:210:SER:N	2.32	0.44
2:D:26:ILE:HD12	2:D:27:GLY:H	1.83	0.44
2:D:48:ASN:N	2:D:48:ASN:ND2	2.58	0.44
2:D:144:GLU:O	2:D:153:LYS:HD3	2.17	0.44
2:D:186:THR:HG22	2:D:187:LEU:H	1.81	0.44
2:B:196:ILE:H	2:B:196:ILE:HD12	1.82	0.44
2:D:143:MET:HG2	2:D:144:GLU:O	2.18	0.44
2:B:162:LEU:HA	2:B:162:LEU:HD23	1.82	0.44
1:C:54:ASN:O	1:C:77:ALA:HB1	2.18	0.43
1:C:159:TRP:CE3	1:C:170:ARG:HG3	2.53	0.43
2:D:203:THR:HB	2:D:207:PHE:CE2	2.48	0.43
2:D:42:VAL:HG12	2:D:43:MET:N	2.33	0.43
2:D:10:LYS:HE2	2:D:10:LYS:HA	1.99	0.43
1:C:230:VAL:HG21	2:D:59:GLU:HG3	1.99	0.43
1:C:169:PHE:HE2	1:C:207:THR:HG22	1.83	0.43
1:C:91:LEU:HD11	2:D:123:VAL:HG23	1.99	0.43
2:D:176:ALA:HB1	2:D:189:GLU:O	2.17	0.43
1:A:161:ALA:HA	1:A:166:ASN:O	2.19	0.43
2:D:34:ASN:N	2:D:34:ASN:HD22	2.17	0.43
2:D:95:GLY:HA2	2:D:135:GLY:O	2.18	0.43
2:D:196:ILE:O	2:D:199:VAL:HG23	2.19	0.43
1:A:199:GLU:OE1	1:A:199:GLU:HA	2.18	0.43
1:C:14:ALA:C	1:C:16:LEU:H	2.21	0.43
2:D:106:GLU:OE2	2:D:155:LYS:HE3	2.19	0.43
2:D:175:LEU:HD21	2:D:199:VAL:HG22	2.01	0.43
2:D:46:SER:HB2	2:D:53:ILE:HD11	2.00	0.43
2:D:177:VAL:HB	2:D:189:GLU:CB	2.48	0.43
1:A:171:LYS:HB2	2:B:65:ASP:O	2.19	0.43
2:B:20:MET:O	2:B:42:VAL:CG2	2.67	0.42
1:C:61:TRP:CD1	1:C:224:ARG:HD3	2.54	0.42
2:D:117:MET:HG3	2:D:123:VAL:HG13	2.02	0.42
1:C:111:ALA:CB	1:C:116:ILE:HB	2.48	0.42
1:C:34:PRO:HB2	1:C:37:LEU:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:227:LYS:HD2	2:D:72:GLU:HG2	2.01	0.42
1:A:39:LEU:HD11	1:A:61:TRP:HH2	1.84	0.42
1:A:74:LYS:HG3	1:A:96[B]:GLU:HB3	2.02	0.42
1:C:178:PRO:HG2	2:D:84:ASP:HB2	2.01	0.42
1:A:150:THR:HG22	3:A:270:HOH:O	2.18	0.42
1:C:219:ALA:CB	1:C:222:GLU:OE1	2.68	0.41
2:D:113:LEU:HD13	2:D:141:VAL:HG21	2.02	0.41
1:C:14:ALA:C	1:C:16:LEU:N	2.73	0.41
2:D:110:GLN:HG3	2:D:110:GLN:H	1.74	0.41
2:D:186:THR:CG2	2:D:209:VAL:HA	2.50	0.41
2:B:172:ILE:HG13	2:B:172:ILE:O	2.16	0.41
1:C:67:LEU:HD21	1:C:85:ILE:HG22	2.02	0.41
1:A:58:VAL:HG23	1:A:61:TRP:H	1.86	0.41
1:C:108:ARG:HB3	1:C:149:ILE:HD12	2.03	0.41
1:A:39:LEU:HD11	1:A:61:TRP:CH2	2.55	0.41
1:A:59:ASP:HB2	1:A:85:ILE:CD1	2.50	0.41
2:D:200:TYR:HE1	2:D:207:PHE:CE1	2.37	0.41
1:A:127:THR:HG21	1:A:129:ILE:HB	2.00	0.41
2:D:15:GLU:OE2	2:D:170:ARG:HD2	2.21	0.41
2:D:26:ILE:HD13	2:D:70:GLY:HA3	2.02	0.41
1:A:228:ARG:HG3	2:B:62:GLU:HG3	2.03	0.41
1:C:31:CYS:HB2	2:D:67:ILE:CD1	2.52	0.40
1:A:174:ARG:HB2	3:A:248:HOH:O	2.21	0.40
2:B:36:ILE:HA	2:B:37:PRO:HD3	1.87	0.40
2:B:91:MET:HB3	2:B:91:MET:HE3	1.86	0.40
2:D:101:ILE:HA	2:D:140:VAL:O	2.21	0.40
1:A:159:TRP:HB2	1:A:170:ARG:HB2	2.03	0.40
1:C:160:LYS:HG2	1:C:193:GLU:HB2	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	230/241 (95%)	216 (94%)	14 (6%)	0	100	100
1	C	227/241 (94%)	216 (95%)	10 (4%)	1 (0%)	39	65
2	B	210/219 (96%)	198 (94%)	12 (6%)	0	100	100
2	D	193/219 (88%)	182 (94%)	11 (6%)	0	100	100
All	All	860/920 (94%)	812 (94%)	47 (6%)	1 (0%)	56	81

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	175	ASN

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	182/186 (98%)	161 (88%)	21 (12%)	7	12
1	C	179/186 (96%)	158 (88%)	21 (12%)	7	12
2	B	175/179 (98%)	144 (82%)	31 (18%)	2	3
2	D	164/179 (92%)	137 (84%)	27 (16%)	3	4
All	All	700/730 (96%)	600 (86%)	100 (14%)	4	7

All (100) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LYS
1	A	23	LEU
1	A	37	LEU
1	A	46	VAL
1	A	60	ASP
1	A	63	LEU
1	A	66	LEU
1	A	80	VAL
1	A	82	GLU
1	A	84	LYS

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Mol	Chain	Res	Type
1	A	88	ARG
1	A	96[A]	GLU
1	A	96[B]	GLU
1	A	99	LEU
1	A	127	THR
1	A	141	ARG
1	A	150	THR
1	A	155	ILE
1	A	193	GLU
1	A	207	THR
1	A	213	GLN
2	B	6	LYS
2	B	10	LYS
2	B	17	LYS
2	B	24	LEU
2	B	31	LEU
2	B	42	VAL
2	B	44	LEU
2	B	48	ASN
2	B	58	LEU
2	B	71	LYS
2	B	73	THR
2	B	75	THR
2	B	84	ASP
2	B	87	GLU
2	B	115	ASN
2	B	121	LYS
2	B	132	LEU
2	B	144	GLU
2	B	148	LYS
2	B	151	GLU
2	B	152	SER
2	B	155	LYS
2	B	157	THR
2	B	172	ILE
2	B	181	VAL
2	B	182	ASN
2	B	188	THR
2	B	190	LEU
2	B	197	GLU
2	B	209	VAL
2	B	211	GLN

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Mol	Chain	Res	Type
1	C	33	ILE
1	C	37	LEU
1	C	39	LEU
1	C	44	GLN
1	C	49	LEU
1	C	51	VAL
1	C	52	VAL
1	C	59	ASP
1	C	65	LEU
1	C	74	LYS
1	C	86	PHE
1	C	89	GLN
1	C	125	VAL
1	C	129	ILE
1	C	144	VAL
1	C	150	THR
1	C	155	ILE
1	C	172	THR
1	C	207	THR
1	C	213	GLN
1	C	216	VAL
2	D	5	ARG
2	D	10	LYS
2	D	24	LEU
2	D	44	LEU
2	D	48	ASN
2	D	58	LEU
2	D	61	THR
2	D	62	GLU
2	D	66	LEU
2	D	71	LYS
2	D	73	THR
2	D	75	THR
2	D	84	ASP
2	D	105	MET
2	D	110	GLN
2	D	126	MET
2	D	132	LEU
2	D	133	VAL
2	D	137	LYS
2	D	142	ILE
2	D	157	THR

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Mol	Chain	Res	Type
2	D	168	VAL
2	D	172	ILE
2	D	184	ARG
2	D	187	LEU
2	D	196	ILE
2	D	206	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	204	HIS
2	B	48	ASN
2	B	115	ASN
2	B	211	GLN
1	C	18	HIS
1	C	54	ASN
1	C	69	ASN
1	C	102	GLN
1	C	213	GLN
2	D	34	ASN
2	D	48	ASN
2	D	134	ASN
2	D	145	HIS
2	D	169	HIS

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

There are no ligands in this entry.

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	229/241 (95%)	0.03	8 (3%)	48 40	22, 38, 71, 117	0
1	C	229/241 (95%)	0.03	5 (2%)	65 59	23, 44, 73, 131	0
2	B	212/219 (96%)	0.12	11 (5%)	31 24	24, 45, 74, 97	0
2	D	200/219 (91%)	0.99	46 (23%)	1 0	30, 74, 140, 181	0
All	All	870/920 (94%)	0.27	70 (8%)	15 10	22, 47, 106, 181	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	199	VAL	9.9
2	D	201	GLU	9.8
2	D	4	ALA	7.2
2	D	200	TYR	6.9
2	D	197	GLU	6.9
2	D	178	PHE	6.4
2	D	175	LEU	5.7
2	D	195	THR	5.7
1	C	219	ALA	5.5
2	D	10	LYS	5.4
2	D	196	ILE	5.4
2	D	5	ARG	5.1
2	B	1	MET	4.5
1	C	221	GLN	4.2
2	D	172	ILE	4.1
1	C	220	SER	4.1
2	D	207	PHE	3.9
2	B	123	VAL	3.7
2	D	58	LEU	3.6
2	D	190	LEU	3.6
2	D	40	VAL	3.5

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Mol	Chain	Res	Type	RSRZ
2	D	3	GLU	3.5
2	D	173	THR	3.4
2	D	145	HIS	3.4
2	D	16	ILE	3.3
1	A	232	GLN	3.3
2	D	8	MET	3.2
2	D	156	LYS	3.2
2	B	124	LYS	3.1
2	D	198	GLU	3.0
2	D	174	ASP	3.0
1	A	84	LYS	2.9
1	A	222	GLU	2.8
2	D	206	ASP	2.8
2	D	144	GLU	2.8
2	D	146	VAL	2.8
2	D	41	HIS	2.7
2	D	9	VAL	2.7
2	D	177	VAL	2.6
2	D	205	ALA	2.6
2	D	2	LYS	2.6
1	A	7	LEU	2.5
1	A	196	GLU	2.5
2	B	6	LYS	2.5
2	D	210	SER	2.5
1	A	44	GLN	2.5
1	A	221	GLN	2.4
2	D	203	THR	2.4
1	C	222	GLU	2.4
2	B	3	GLU	2.4
1	A	82	GLU	2.4
2	D	189	GLU	2.4
2	B	211	GLN	2.3
2	B	149	HIS	2.3
2	D	176	ALA	2.3
1	C	3	MET	2.3
2	B	120	GLY	2.3
2	D	109	GLU	2.2
2	B	184	ARG	2.2
2	D	209	VAL	2.2
2	D	119	PRO	2.2
2	D	34	ASN	2.2
2	D	120	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
2	D	7	ARG	2.2
2	D	53	ILE	2.1
2	D	157	THR	2.1
2	D	12	ALA	2.1
2	B	32	VAL	2.1
2	B	175	LEU	2.0
2	D	31	LEU	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.