



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 08:56 AM UTC

PDB ID : 3CDK / pdb_00003cdk
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

<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

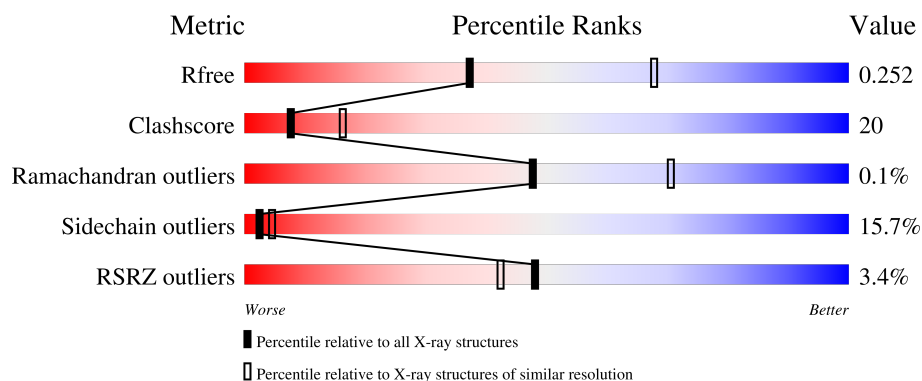
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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}	180053	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (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 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	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

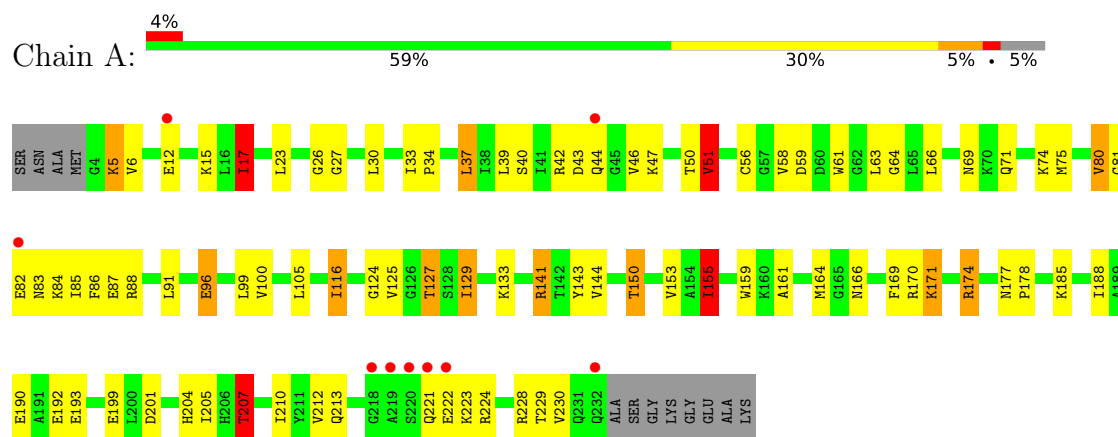
- Molecule 3 is water.

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

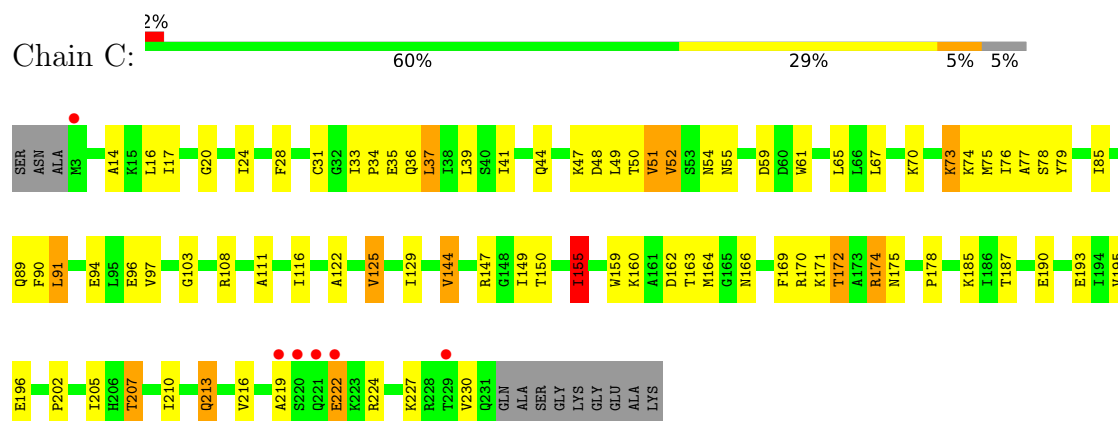
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, 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.

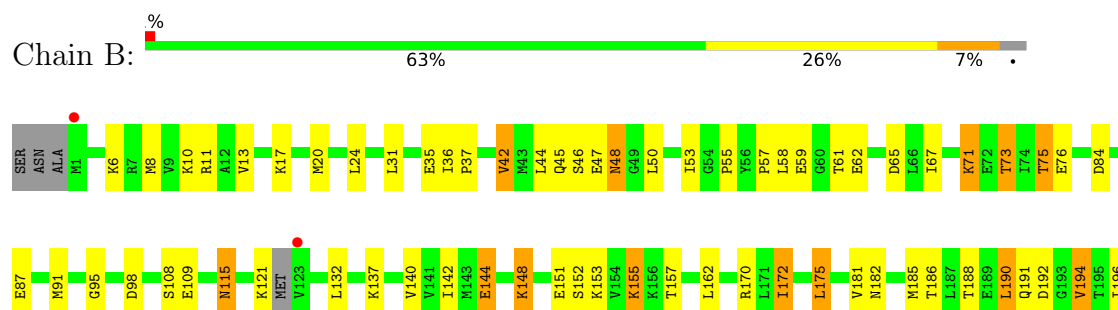
• Molecule 1: Succinyl-CoA:3-ketoacid-coenzyme A transferase subunit A



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

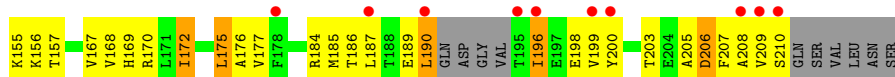


• Molecule 2: Succinyl-CoA:3-ketoacid-coenzyme A transferase subunit B





● 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) 99.1 (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 (5.00%)	wwPDB-VP
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 , 47.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	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, 0.333 respectively for untwinned datasets, and 0.375, 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	0.95	4/1777 (0.2%)	1.10	6/2395 (0.3%)
1	C	0.92	0/1750	1.09	7/2359 (0.3%)
2	B	0.84	0/1632	1.06	5/2197 (0.2%)
2	D	0.94	4/1531 (0.3%)	1.05	3/2059 (0.1%)
All	All	0.91	8/6690 (0.1%)	1.07	21/9010 (0.2%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	198	GLU	CD-OE2	13.82	1.51	1.25
2	D	175	LEU	C-N	9.65	1.45	1.33
2	D	175	LEU	C-O	7.72	1.31	1.23
2	D	190	LEU	C-O	7.38	1.38	1.23
1	A	125	VAL	CA-CB	5.78	1.60	1.54
1	A	17	ILE	CA-CB	5.54	1.60	1.54
1	A	207	THR	CA-CB	5.41	1.60	1.53
1	A	51	VAL	CA-CB	5.09	1.60	1.54

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	192	ASP	N-CA-C	6.73	118.61	111.28
1	C	174	ARG	CB-CG-CD	-6.33	96.75	111.30
1	C	155	ILE	N-CA-C	-6.30	96.65	107.24
2	D	16	ILE	N-CA-C	6.27	116.86	107.51
1	A	155	ILE	CB-CA-C	6.18	119.54	110.77
2	D	117	MET	N-CA-C	6.01	118.26	108.76
1	C	171	LYS	CB-CA-C	-5.95	109.70	116.54
1	C	50	THR	CA-C-N	-5.83	114.58	122.91
1	C	50	THR	C-N-CA	-5.83	114.58	122.91
2	D	16	ILE	CB-CA-C	-5.63	103.43	111.31
1	A	171	LYS	CB-CA-C	-5.58	110.12	116.54
2	B	209	VAL	CB-CA-C	5.53	118.36	110.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	47	LYS	N-CA-C	5.51	117.01	109.18
1	A	6	VAL	N-CA-C	5.41	116.03	108.89
2	B	194	VAL	N-CA-C	5.40	117.33	108.97
1	A	174	ARG	CB-CG-CD	-5.23	99.27	111.30
1	C	144	VAL	CB-CA-C	-5.17	102.96	110.81
1	A	50	THR	N-CA-C	-5.14	100.35	108.73
1	C	47	LYS	N-CA-C	5.14	117.66	109.65
2	B	208	ALA	CA-C-N	-5.11	116.50	122.93
2	B	208	ALA	C-N-CA	-5.11	116.50	122.93

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	77	0
1	C	1725	0	1761	57	0
2	B	1613	0	1627	56	0
2	D	1515	0	1531	98	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	268	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 20.

All (268) 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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:H	2:D:48:ASN:HD22	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
1:C:90:PHE:CE2	2:D:123:VAL:HG11	2.10	0.87
2:D:139:ILE:H	2:D:169:HIS:HD2	0.91	0.87
2:B:142:ILE:HG22	2:B:172:ILE:HG12	1.56	0.87
2:B:148:LYS:H	2:B:148:LYS:HD3	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:B:115:ASN:C	2:B:115:ASN:HD22	1.90	0.77
2:D:200:TYR:CD1	2:D:207:PHE:CE1	2.72	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:H	2:B:48:ASN:ND2	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
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
1:A:155:ILE:HD11	1:A:188:ILE:HD12	1.73	0.71
2:B:148:LYS:H	2:B:148:LYS:CD	2.00	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
2:B:185:MET:HE1	2:B:205:ALA:HB3	1.75	0.68
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
2:B:91:MET:HE3	1:C:210:ILE:HD13	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
2:D:185:MET:HE3	2:D:207:PHE:HB3	1.77	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:B:11:ARG:HB3	2:B:172:ILE:CD1	2.29	0.63
2:B:48:ASN:HD22	2:B:48:ASN:N	1.83	0.63
2:B:115:ASN:C	2:B:115:ASN:ND2	2.52	0.63
2:D:203:THR:HB	2:D:207:PHE:HE2	1.64	0.63
2:B:185:MET:HE1	2:B:205:ALA:CB	2.28	0.62
1:C:187:THR:H	1:C:213:GLN:CG	2.12	0.62
2:D:48:ASN:H	2:D:48:ASN:ND2	1.89	0.61
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:HA	2:B:211:GLN:NE2	2.15	0.61
1:A:42[B]:ARG:HH21	1:A:42[B]:ARG:CG	1.96	0.60
1:A:30:LEU:HD13	1:A:224:ARG:HD3	1.84	0.60
2:B:67:ILE:HD12	2:B:71:LYS:HA	1.84	0.60
2:D:156:LYS:HG3	2:D:205:ALA:HA	1.83	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:O	2:D:16:ILE:CG2	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:B:55:PRO:O	2:B:75:THR:HB	2.01	0.59
2:B:211:GLN:HA	2:B:211:GLN:HE21	1.68	0.59
2:D:13:VAL:O	2:D:16:ILE:CG2	2.41	0.59
2:D:70:GLY:O	2:D:71:LYS:HB2	2.01	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:C:37:LEU:HD22	1:C:190:GLU:HB2	1.86	0.58
2:D:186:THR:HG23	2:D:208:ALA:O	2.04	0.58
2:D:17:LYS:H	2:D:20:MET:HE3	1.69	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
2:D:200:TYR:CD1	2:D:207:PHE:CZ	2.90	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:B:108:SER:HA	2:B:155:LYS:O	2.05	0.57
2:D:186:THR:HA	2:D:208:ALA:O	2.05	0.57
1:C:67:LEU:CD2	1:C:85:ILE:HG22	2.35	0.56
2:B:57:PRO:HG3	2:B:73:THR:HG23	1.87	0.56
2:D:133:VAL:HG13	2:D:167:VAL:HG11	1.86	0.56
1:A:174:ARG:HD3	3:A:248:HOH:O	2.05	0.56
2:B:98:ASP:OD1	2:B:137:LYS:HE3	2.06	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
2:B:185:MET:HE3	2:B:207:PHE:HB3	1.87	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:D:17:LYS:HG3	2:D:20:MET:HE3	1.87	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:56:CYS:HB2	1:A:75:MET:HE2	1.88	0.55
1:A:27:GLY:HA3	1:A:33:ILE:HG22	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
2:B:144:GLU:O	2:B:153:LYS:HD3	2.08	0.54
1:C:20:GLY:HA2	1:C:48:ASP:O	2.07	0.54
2:B:155:LYS:HA	2:B:204:GLU:HB2	1.89	0.54
2:D:17:LYS:HE2	2:D:20:MET:HE2	1.88	0.54
1:A:210:ILE:HD13	2:D:91:MET:HE3	1.91	0.53
2:D:42:VAL:CG1	2:D:43:MET:N	2.71	0.53
2:D:117:MET:HG2	2:D:118:ILE:N	2.23	0.53
2:D:177:VAL:HB	2:D:189:GLU:HB3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:GLY:HA3	1:A:177:ASN:OD1	2.08	0.53
2:D:200:TYR:CD1	2:D:207:PHE:HE1	2.25	0.53
1:A:12:GLU:HA	1:A:15:LYS:HE3	1.91	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
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:B:20:MET:O	2:B:42:VAL:HG22	2.10	0.51
2:D:21:ASN:N	2:D:98:ASP:OD2	2.36	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
1:A:80:VAL:HG21	1:A:86:PHE:CZ	2.46	0.51
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
1:A:192:GLU:OE2	1:A:223:LYS:NZ	2.41	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
2:B:45:GLN:HG3	2:B:91:MET:HE1	1.94	0.50
1:C:55:ASN:ND2	1:C:79:TYR:HB3	2.27	0.50
2:D:186:THR:HG23	2:D:208:ALA:C	2.36	0.50
2:D:61:THR:O	2:D:61:THR:HG23	2.12	0.49
1:A:141:ARG:HD3	1:A:143:TYR:OH	2.13	0.49
2:B:62:GLU:HB2	2:B:73:THR:HG21	1.94	0.49
1:A:164:MET:HE1	2:D:43:MET:HE2	1.95	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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
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:14:ALA:C	1:C:16:LEU:H	2.21	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:133:LYS:HB2	1:A:144:VAL:HG22	1.97	0.46
1:A:169:PHE:CE1	1:A:207:THR:CG2	2.96	0.46
1:C:34:PRO:HG3	1:C:155:ILE:HG23	1.98	0.46
2:D:155:LYS:O	2:D:205:ALA:HB2	2.15	0.46
1:C:28:PHE:CE2	2:D:69:ALA:HA	2.51	0.46
1:C:67:LEU:HD23	1:C:85:ILE:HG22	1.97	0.46
2:B:186:THR:HA	2:B:208:ALA:O	2.16	0.46
2:B:140:VAL:HG22	2:B:170:ARG:HB3	1.97	0.45
1:C:17:ILE:HD12	1:C:41:ILE:HD11	1.99	0.45
2:D:71:LYS:HD3	2:D:71:LYS:N	2.32	0.45
1:A:178:PRO:HG2	2:B:84:ASP:HB3	1.98	0.45
2:B:57:PRO:CG	2:B:73:THR:HG23	2.46	0.45
2:D:203:THR:HG22	2:D:205:ALA:N	2.30	0.45
1:C:174:ARG:HD2	3:C:267:HOH:O	2.15	0.45
1:A:80:VAL:CG1	1:A:81:GLY:N	2.76	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:117:MET:HE2	2:D:119:PRO:HA	1.99	0.44
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:B:196:ILE:HD12	2:B:196:ILE:H	1.82	0.44
2:D:186:THR:HG22	2:D:187:LEU:H	1.81	0.44
2:B:162:LEU:HD23	2:B:162:LEU:HA	1.82	0.44
2:D:143:MET:HG2	2:D:144:GLU:O	2.18	0.44
1:C:14:ALA:C	1:C:16:LEU:N	2.73	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:C:230:VAL:HG21	2:D:59:GLU:HG3	1.99	0.43
2:D:10:LYS:HE2	2:D:10:LYS:HA	1.99	0.43
2:D:42:VAL:HG12	2:D:43:MET:N	2.33	0.43
2:D:203:THR:HB	2:D:207:PHE:CE2	2.48	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
1:A:199:GLU:HA	1:A:199:GLU:OE1	2.18	0.43
2:D:196:ILE:O	2:D:199:VAL:HG23	2.19	0.43
1:C:162:ASP:CG	1:C:166:ASN:HB2	2.44	0.43
2:D:46:SER:HB2	2:D:53:ILE:HD11	2.00	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
1:A:171:LYS:HB2	2:B:65:ASP:O	2.19	0.43
2:D:5:ARG:HA	2:D:8:MET:HE3	2.00	0.43
2:D:177:VAL:HB	2:D:189:GLU:CB	2.48	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
1:C:34:PRO:HB2	1:C:37:LEU:HB2	2.01	0.42
1:C:111:ALA:CB	1:C:116:ILE:HB	2.48	0.42
2:D:117:MET:HG3	2:D:123:VAL:HG13	2.02	0.42
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
2:D:110:GLN:H	2:D:110:GLN:HG3	1.74	0.41
2:D:186:THR:CG2	2:D:209:VAL:HA	2.50	0.41
1:A:58:VAL:HG23	1:A:61:TRP:H	1.86	0.41
2:B:172:ILE:O	2:B:172:ILE:HG13	2.16	0.41
1:C:67:LEU:HD21	1:C:85:ILE:HG22	2.02	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:A:228:ARG:HG3	2:B:62:GLU:HG3	2.03	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
2:D:91:MET:HB3	2:D:91:MET:HE2	1.66	0.41
2:D:126:MET:HE3	2:D:126:MET:HB3	1.90	0.41
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
1:C:31:CYS:HB2	2:D:67:ILE:CD1	2.52	0.40
1:A:159:TRP:HB2	1:A:170:ARG:HB2	2.03	0.40
2:D:101:ILE:HA	2:D:140:VAL:O	2.21	0.40
1:C:162:ASP:C	1:C:164:MET:H	2.29	0.40
1:C:36:GLN:O	1:C:37:LEU:C	2.65	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

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	230/241 (95%)	216 (94%)	14 (6%)	0	100	100
1	C	227/241 (94%)	216 (95%)	10 (4%)	1 (0%)	30	51
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%)	48	70

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	175	ASN

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	182/186 (98%)	157 (86%)	25 (14%)	3	7
1	C	179/186 (96%)	153 (86%)	26 (14%)	3	6
2	B	175/179 (98%)	141 (81%)	34 (19%)	1	2
2	D	164/179 (92%)	139 (85%)	25 (15%)	3	5
All	All	700/730 (96%)	590 (84%)	110 (16%)	2	4

All (110) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LYS
1	A	17	ILE
1	A	23	LEU
1	A	37	LEU
1	A	46	VAL
1	A	51	VAL
1	A	63	LEU
1	A	66	LEU
1	A	80	VAL
1	A	82	GLU
1	A	84	LYS
1	A	88	ARG
1	A	96[A]	GLU
1	A	96[B]	GLU
1	A	99	LEU
1	A	116	ILE
1	A	127	THR
1	A	129	ILE
1	A	141	ARG
1	A	150	THR
1	A	155	ILE

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Mol	Chain	Res	Type
1	A	185	LYS
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	47	GLU
2	B	48	ASN
2	B	50	LEU
2	B	58	LEU
2	B	61	THR
2	B	71	LYS
2	B	73	THR
2	B	75	THR
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	175	LEU
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
1	C	33	ILE
1	C	35	GLU
1	C	37	LEU
1	C	39	LEU

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Mol	Chain	Res	Type
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	70	LYS
1	C	73	LYS
1	C	74	LYS
1	C	91	LEU
1	C	96	GLU
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	185	LYS
1	C	207	THR
1	C	213	GLN
1	C	216	VAL
1	C	222	GLU
2	D	3	GLU
2	D	5	ARG
2	D	6	LYS
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	73	THR
2	D	75	THR
2	D	84	ASP
2	D	105	MET
2	D	110	GLN
2	D	132	LEU
2	D	133	VAL
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	196	ILE
2	D	206	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	204	HIS
2	B	21	ASN
2	B	48	ASN
2	B	110	GLN
2	B	115	ASN
2	B	134	ASN
2	B	211	GLN
1	C	18	HIS
1	C	44	GLN
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 oligosaccharides 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%)	-0.12	9 (3%)	43	38	15, 37, 71, 117	3 (1%)
1	C	229/241 (95%)	-0.12	6 (2%)	57	51	23, 44, 73, 131	0
2	B	212/219 (96%)	-0.17	3 (1%)	73	69	21, 45, 74, 97	2 (0%)
2	D	200/219 (91%)	0.32	12 (6%)	27	22	17, 74, 140, 181	1 (0%)
All	All	870/920 (94%)	-0.03	30 (3%)	48	42	15, 46, 106, 181	6 (0%)

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	219	ALA	6.3
1	C	220	SER	5.2
1	C	221	GLN	4.5
2	D	190	LEU	3.5
2	D	210	SER	3.4
2	D	196	ILE	3.2
1	C	222	GLU	3.2
1	A	219	ALA	3.1
1	A	222	GLU	3.0
2	D	209	VAL	3.0
1	A	220	SER	2.8
2	D	199	VAL	2.8
1	C	3	MET	2.8
1	A	221	GLN	2.7
2	B	123	VAL	2.7
1	A	232	GLN	2.6
2	D	208	ALA	2.5
2	B	213	VAL	2.4
2	D	195	THR	2.2
1	A	218	GLY	2.2
2	D	58	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
2	D	187	LEU	2.1
1	A	12	GLU	2.1
1	A	82	GLU	2.1
2	D	16	ILE	2.1
2	B	1	MET	2.1
1	A	44	GLN	2.1
2	D	178	PHE	2.0
1	C	229	THR	2.0
2	D	200	TYR	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.