Full wwPDB X-ray Structure Validation Report ⓘ

Jun 11, 2024 – 08:42 PM EDT

PDB ID	:	1TDT
Title	:	THREE-DIMENSIONAL STRUCTURE OF TETRAHYDRODIPICOLINATE-N-SUCCINYLTRANSFERASE
Authors	:	Beaman, T.W.; Binder, D.W.; Blanchard, J.S.; Roderick, S.L.
Deposited on	:	1996-11-19
Resolution	:	2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

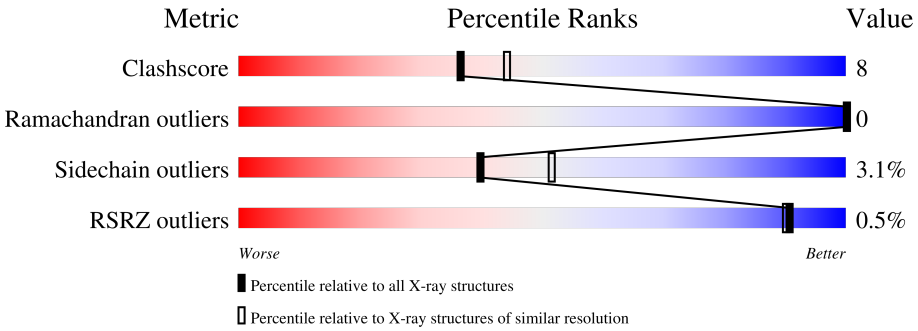
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	259	% 73% 22%
1	B	259	74% 20%
1	C	259	73% 21%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 5995 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 TETRAHYDRODIPICOLINATE-N-SUCCINYLTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	256	Total	C	N	O	S	0	0	0
			1928	1215	331	374	8			
1	B	255	Total	C	N	O	S	0	0	0
			1933	1218	334	373	8			
1	C	253	Total	C	N	O	S	0	0	0
			1919	1209	331	371	8			

- Molecule 2 is water.

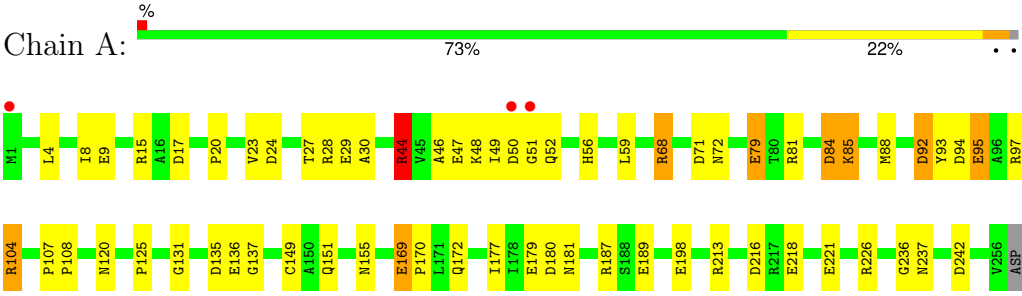
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	80	Total	O	0	0
			80	80		
2	B	58	Total	O	0	0
			58	58		
2	C	77	Total	O	0	0
			77	77		

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.

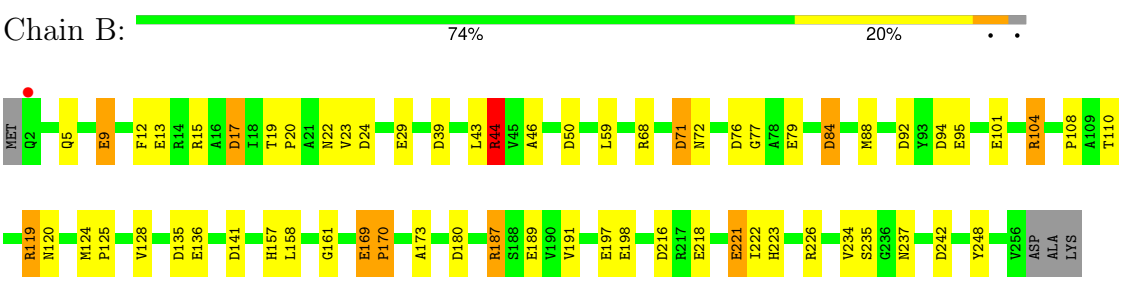
• Molecule 1: TETRAHYDRODIPICOLINATE-N-SUCCINYLTRANSFERASE

Chain A:



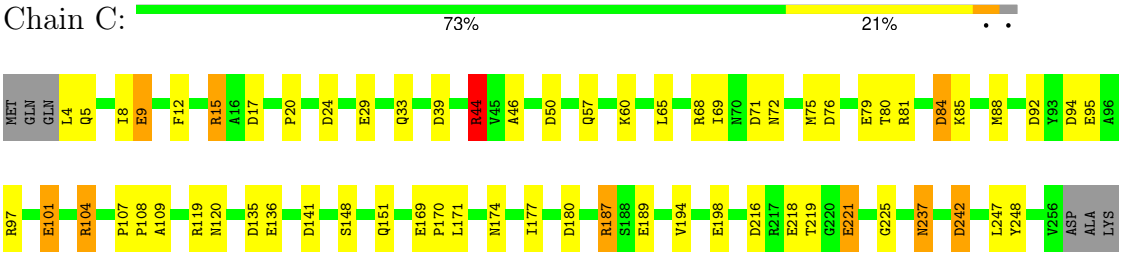
• Molecule 1: TETRAHYDRODIPICOLINATE-N-SUCCINYLTRANSFERASE

Chain B:



• Molecule 1: TETRAHYDRODIPICOLINATE-N-SUCCINYLTRANSFERASE

Chain C:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	60.60Å 107.90Å 70.70Å 90.00° 113.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.20 23.02 – 2.20	Depositor EDS
% Data completeness (in resolution range)	90.0 ((Not available)-2.20) 90.6 (23.02-2.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.47 (at 2.19Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.170 , 0.230 0.160 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	19.1	Xtriage
Anisotropy	0.308	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 81.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.019 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	5995	wwPDB-VP
Average B, all atoms (Å^2)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.28% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (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	0.85	11/1962 (0.6%)	1.25	30/2671 (1.1%)
1	B	0.86	12/1967 (0.6%)	1.26	28/2675 (1.0%)
1	C	0.84	10/1953 (0.5%)	1.23	29/2657 (1.1%)
All	All	0.85	33/5882 (0.6%)	1.24	87/8003 (1.1%)

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	29	GLU	CD-OE2	6.85	1.33	1.25
1	C	79	GLU	CD-OE2	6.71	1.33	1.25
1	B	13	GLU	CD-OE2	6.62	1.32	1.25
1	A	79	GLU	CD-OE2	6.38	1.32	1.25
1	B	79	GLU	CD-OE2	6.32	1.32	1.25
1	B	221	GLU	CD-OE2	6.29	1.32	1.25
1	C	189	GLU	CD-OE2	6.25	1.32	1.25
1	A	189	GLU	CD-OE2	6.21	1.32	1.25
1	B	101	GLU	CD-OE2	6.04	1.32	1.25
1	B	197	GLU	CD-OE2	6.02	1.32	1.25
1	A	9	GLU	CD-OE2	5.98	1.32	1.25
1	C	198	GLU	CD-OE2	5.81	1.32	1.25
1	A	218	GLU	CD-OE2	5.72	1.31	1.25
1	A	169	GLU	CD-OE2	5.71	1.31	1.25
1	B	169	GLU	CD-OE2	5.66	1.31	1.25
1	C	218	GLU	CD-OE2	5.61	1.31	1.25
1	C	136	GLU	CD-OE2	5.49	1.31	1.25
1	B	29	GLU	CD-OE2	5.47	1.31	1.25
1	B	136	GLU	CD-OE2	5.44	1.31	1.25
1	A	198	GLU	CD-OE1	5.43	1.31	1.25
1	A	29	GLU	CD-OE2	5.43	1.31	1.25
1	B	9	GLU	CD-OE2	5.25	1.31	1.25
1	C	221	GLU	CD-OE2	5.24	1.31	1.25
1	A	101	GLU	CD-OE2	5.22	1.31	1.25
1	C	101	GLU	CD-OE2	5.22	1.31	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	136	GLU	CD-OE2	5.19	1.31	1.25
1	B	95	GLU	CD-OE2	5.15	1.31	1.25
1	C	95	GLU	CD-OE1	5.14	1.31	1.25
1	A	95	GLU	CD-OE2	5.12	1.31	1.25
1	B	218	GLU	CD-OE1	5.09	1.31	1.25
1	B	198	GLU	CD-OE2	5.09	1.31	1.25
1	A	179	GLU	CD-OE2	5.06	1.31	1.25
1	C	9	GLU	CD-OE2	5.03	1.31	1.25

All (87) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	104	ARG	NE-CZ-NH1	10.41	125.51	120.30
1	A	104	ARG	NE-CZ-NH1	8.60	124.60	120.30
1	A	180	ASP	CB-CG-OD2	-8.18	110.94	118.30
1	C	44	ARG	NE-CZ-NH2	-7.95	116.33	120.30
1	B	135	ASP	CB-CG-OD1	7.76	125.28	118.30
1	C	104	ARG	NE-CZ-NH2	-7.67	116.47	120.30
1	B	104	ARG	NE-CZ-NH2	-7.51	116.55	120.30
1	C	216	ASP	CB-CG-OD2	-7.49	111.56	118.30
1	A	24	ASP	CB-CG-OD2	-7.48	111.57	118.30
1	A	71	ASP	CB-CG-OD2	-7.45	111.59	118.30
1	B	242	ASP	CB-CG-OD2	-7.45	111.59	118.30
1	C	92	ASP	CB-CG-OD2	-7.45	111.60	118.30
1	C	71	ASP	CB-CG-OD2	-7.40	111.64	118.30
1	B	135	ASP	CB-CG-OD2	-7.39	111.65	118.30
1	B	180	ASP	CB-CG-OD2	-7.28	111.75	118.30
1	C	104	ARG	NE-CZ-NH1	7.27	123.94	120.30
1	A	242	ASP	CB-CG-OD2	-7.25	111.78	118.30
1	C	216	ASP	CB-CG-OD1	7.04	124.64	118.30
1	A	180	ASP	CB-CG-OD1	7.02	124.62	118.30
1	A	84	ASP	CB-CG-OD2	-6.97	112.03	118.30
1	B	180	ASP	CB-CG-OD1	6.92	124.53	118.30
1	A	221	GLU	CB-CA-C	-6.87	96.67	110.40
1	B	24	ASP	CB-CG-OD2	-6.84	112.14	118.30
1	B	71	ASP	CB-CG-OD2	-6.82	112.16	118.30
1	B	84	ASP	CB-CG-OD2	-6.79	112.19	118.30
1	A	135	ASP	CB-CG-OD1	6.73	124.36	118.30
1	C	94	ASP	CB-CG-OD2	-6.62	112.34	118.30
1	B	44	ARG	NE-CZ-NH1	6.58	123.59	120.30
1	A	44	ARG	NE-CZ-NH1	6.54	123.57	120.30
1	B	17	ASP	CB-CG-OD2	-6.50	112.45	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	94	ASP	CB-CG-OD2	-6.47	112.47	118.30
1	B	216	ASP	CB-CG-OD2	-6.46	112.49	118.30
1	B	92	ASP	CB-CG-OD2	-6.44	112.51	118.30
1	A	71	ASP	CB-CG-OD1	6.42	124.08	118.30
1	C	92	ASP	CB-CG-OD1	6.37	124.04	118.30
1	B	71	ASP	CB-CG-OD1	6.37	124.03	118.30
1	A	92	ASP	CB-CG-OD2	-6.36	112.57	118.30
1	A	242	ASP	CB-CG-OD1	6.36	124.03	118.30
1	A	50	ASP	CB-CG-OD2	-6.30	112.63	118.30
1	A	135	ASP	CB-CG-OD2	-6.30	112.63	118.30
1	C	135	ASP	CB-CG-OD2	-6.29	112.64	118.30
1	A	94	ASP	CB-CG-OD1	6.29	123.96	118.30
1	C	50	ASP	CB-CG-OD2	-6.24	112.68	118.30
1	A	84	ASP	CB-CG-OD1	6.21	123.89	118.30
1	C	17	ASP	CB-CG-OD2	-6.19	112.73	118.30
1	C	71	ASP	CB-CG-OD1	6.15	123.84	118.30
1	C	39	ASP	CB-CG-OD1	6.11	123.80	118.30
1	A	24	ASP	CB-CG-OD1	6.08	123.77	118.30
1	A	94	ASP	CB-CG-OD2	-6.06	112.84	118.30
1	C	141	ASP	CB-CG-OD2	-6.05	112.86	118.30
1	C	24	ASP	CB-CG-OD2	-6.00	112.90	118.30
1	C	135	ASP	CB-CG-OD1	6.00	123.69	118.30
1	B	39	ASP	CB-CG-OD2	-5.92	112.97	118.30
1	C	17	ASP	CB-CG-OD1	5.91	123.62	118.30
1	B	84	ASP	CB-CG-OD1	5.89	123.60	118.30
1	A	216	ASP	CB-CG-OD1	5.83	123.55	118.30
1	B	242	ASP	CB-CG-OD1	5.81	123.53	118.30
1	A	104	ARG	NE-CZ-NH2	-5.77	117.41	120.30
1	B	141	ASP	CB-CG-OD2	-5.75	113.12	118.30
1	C	44	ARG	NE-CZ-NH1	5.71	123.15	120.30
1	C	84	ASP	CB-CG-OD2	-5.67	113.20	118.30
1	B	50	ASP	CB-CG-OD2	-5.58	113.28	118.30
1	A	81	ARG	NE-CZ-NH1	5.57	123.09	120.30
1	C	242	ASP	CB-CG-OD2	-5.57	113.29	118.30
1	B	94	ASP	CB-CG-OD1	5.54	123.29	118.30
1	A	17	ASP	CB-CG-OD2	-5.52	113.33	118.30
1	C	39	ASP	CB-CG-OD2	-5.51	113.34	118.30
1	C	15	ARG	NE-CZ-NH1	5.50	123.05	120.30
1	A	216	ASP	CB-CG-OD2	-5.41	113.43	118.30
1	A	15	ARG	NE-CZ-NH1	5.41	123.00	120.30
1	B	50	ASP	CB-CG-OD1	5.37	123.14	118.30
1	C	187	ARG	NE-CZ-NH2	-5.37	117.62	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	76	ASP	CB-CG-OD2	-5.34	113.49	118.30
1	C	141	ASP	CB-CG-OD1	5.34	123.10	118.30
1	C	180	ASP	CB-CG-OD2	-5.31	113.52	118.30
1	B	92	ASP	CB-CG-OD1	5.30	123.07	118.30
1	B	216	ASP	CB-CG-OD1	5.29	123.06	118.30
1	A	50	ASP	CB-CG-OD1	5.28	123.05	118.30
1	A	15	ARG	NE-CZ-NH2	-5.24	117.68	120.30
1	C	24	ASP	CB-CG-OD1	5.24	123.01	118.30
1	A	213	ARG	NE-CZ-NH1	5.23	122.92	120.30
1	A	17	ASP	CB-CG-OD1	5.21	122.99	118.30
1	B	39	ASP	CB-CG-OD1	5.21	122.99	118.30
1	A	81	ARG	NE-CZ-NH2	-5.18	117.71	120.30
1	C	180	ASP	CB-CG-OD1	5.03	122.83	118.30
1	B	77	GLY	N-CA-C	-5.02	100.56	113.10
1	C	94	ASP	CB-CG-OD1	5.01	122.81	118.30

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	1928	0	1883	36	0
1	B	1933	0	1900	34	0
1	C	1919	0	1885	33	0
2	A	80	0	0	5	0
2	B	58	0	0	4	0
2	C	77	0	0	5	0
All	All	5995	0	5668	90	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 8.

All (90) 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:15:ARG:HG2	1:B:15:ARG:HH11	1.47	0.80
1:A:44:ARG:HD2	1:A:47:GLU:HB3	1.62	0.80
1:B:88:MET:HE2	1:C:80:THR:HB	1.66	0.77
1:A:172:GLN:HB2	2:A:323:HOH:O	1.85	0.75
1:A:226:ARG:HD2	2:A:316:HOH:O	1.90	0.72
1:C:5:GLN:O	1:C:9:GLU:HG3	1.90	0.70
1:A:108:PRO:HD2	1:A:125:PRO:HB3	1.73	0.70
1:C:174:ASN:HB3	2:C:272:HOH:O	1.96	0.65
1:B:12:PHE:O	1:B:15:ARG:HG2	1.98	0.64
1:C:85:LYS:NZ	2:C:284:HOH:O	2.31	0.64
1:A:72:ASN:OD1	1:C:104:ARG:HD3	1.98	0.64
1:B:88:MET:CE	1:B:110:THR:HG23	2.29	0.62
1:A:68:ARG:NH1	2:A:301:HOH:O	2.31	0.62
1:B:88:MET:HE1	2:C:291:HOH:O	1.99	0.62
1:A:51:GLY:O	1:A:52:GLN:NE2	2.32	0.61
1:C:12:PHE:O	1:C:15:ARG:NH1	2.34	0.61
1:C:75:MET:HE1	1:C:84:ASP:HB3	1.85	0.59
1:B:12:PHE:O	1:B:15:ARG:NH1	2.35	0.59
1:B:17:ASP:O	1:B:22:ASN:ND2	2.35	0.59
1:B:226:ARG:NH2	2:B:279:HOH:O	2.35	0.59
1:B:15:ARG:HG2	1:B:15:ARG:NH1	2.18	0.58
1:A:84:ASP:OD1	1:C:104:ARG:NH1	2.33	0.58
1:B:128:VAL:HG23	2:B:317:HOH:O	2.03	0.57
1:C:4:LEU:O	1:C:8:ILE:HG12	2.05	0.56
1:B:104:ARG:HD3	1:C:72:ASN:OD1	2.05	0.56
1:A:79:GLU:CG	1:C:88:MET:HE2	2.35	0.56
1:A:104:ARG:NH1	1:B:84:ASP:OD1	2.34	0.55
1:C:219:THR:OG1	1:C:221:GLU:HB2	2.07	0.55
1:A:20:PRO:HD3	1:A:68:ARG:O	2.06	0.55
1:B:44:ARG:HD2	1:B:46:ALA:O	2.07	0.54
1:B:19:THR:HB	1:B:20:PRO:HD2	1.90	0.54
1:A:44:ARG:HD3	1:A:46:ALA:O	2.08	0.54
1:A:85:LYS:NZ	2:A:291:HOH:O	2.39	0.54
1:B:169:GLU:HA	1:B:170:PRO:C	2.29	0.53
1:A:104:ARG:HD3	1:B:72:ASN:OD1	2.08	0.53
1:B:88:MET:HE1	1:B:110:THR:HG23	1.89	0.52
1:C:4:LEU:HD21	1:C:33:GLN:HG2	1.91	0.52
1:C:169:GLU:HA	1:C:170:PRO:C	2.30	0.52
1:C:44:ARG:HD2	1:C:46:ALA:O	2.09	0.52
1:B:170:PRO:HG2	1:B:173:ALA:HB2	1.92	0.51
1:B:237:ASN:ND2	1:B:248:TYR:HA	2.25	0.51
1:C:97:ARG:O	1:C:101:GLU:HG3	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:GLY:C	1:A:237:ASN:HD22	2.14	0.50
1:A:79:GLU:HG2	1:C:88:MET:HE2	1.94	0.49
1:C:20:PRO:HD3	1:C:68:ARG:O	2.12	0.49
1:A:92:ASP:HB2	2:A:310:HOH:O	2.13	0.49
1:A:48:LYS:HA	1:A:52:GLN:O	2.13	0.49
1:A:79:GLU:HG3	1:C:88:MET:HE2	1.94	0.49
1:C:194:VAL:HG22	1:C:225:GLY:HA2	1.95	0.48
1:A:56:HIS:HB2	1:A:59:LEU:HD12	1.95	0.48
1:B:124:MET:HB3	1:B:125:PRO:HD2	1.96	0.48
1:B:19:THR:N	1:B:22:ASN:OD1	2.36	0.48
1:A:95:GLU:O	1:A:99:GLN:HG3	2.14	0.47
1:C:75:MET:HE2	2:C:311:HOH:O	2.14	0.47
1:A:68:ARG:HE	1:A:68:ARG:HB3	1.37	0.46
1:C:242:ASP:OD1	1:C:242:ASP:N	2.48	0.46
1:B:189:GLU:HG3	1:B:191:VAL:HG13	1.97	0.46
1:B:88:MET:CE	1:C:80:THR:HB	2.42	0.45
1:B:15:ARG:NH1	1:B:15:ARG:CG	2.79	0.45
1:A:169:GLU:HA	1:A:170:PRO:C	2.36	0.45
1:A:120:ASN:O	1:A:137:GLY:HA2	2.17	0.45
1:A:131:GLY:O	1:A:149:CYS:HA	2.17	0.45
1:A:23:VAL:CG1	1:A:28:ARG:NH2	2.80	0.44
1:A:155:ASN:O	1:A:181:ASN:HA	2.17	0.44
1:B:108:PRO:HD2	1:B:125:PRO:HB3	1.98	0.44
1:C:57:GLN:HG2	1:C:171:LEU:O	2.17	0.44
1:A:236:GLY:O	1:A:237:ASN:ND2	2.50	0.43
1:B:157:HIS:HD2	1:B:158:LEU:N	2.17	0.43
1:A:88:MET:HE3	1:C:107:PRO:HG2	2.00	0.43
1:B:43:LEU:HD13	1:B:59:LEU:HD21	2.00	0.43
1:B:119:ARG:O	1:B:120:ASN:HB2	2.18	0.43
1:B:187:ARG:NH2	2:B:281:HOH:O	2.42	0.43
1:A:107:PRO:HA	1:A:108:PRO:HA	1.70	0.42
1:C:60:LYS:NZ	1:C:170:PRO:O	2.52	0.42
1:A:44:ARG:NH1	1:A:151:GLN:OE1	2.52	0.42
1:A:93:TYR:CE2	1:A:97:ARG:HG2	2.54	0.42
1:A:4:LEU:O	1:A:8:ILE:HG12	2.20	0.42
1:C:187:ARG:NH2	2:C:294:HOH:O	2.53	0.42
1:C:237:ASN:ND2	1:C:248:TYR:HA	2.35	0.42
1:C:108:PRO:O	1:C:109:ALA:C	2.58	0.41
1:C:65:LEU:O	1:C:69:ILE:HG23	2.21	0.41
1:C:151:GLN:HB2	1:C:177:ILE:HD13	2.03	0.41
1:A:49:ILE:N	1:A:52:GLN:O	2.40	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:119:ARG:O	1:C:120:ASN:HB2	2.21	0.41
1:A:187:ARG:NH1	1:B:161:GLY:O	2.54	0.41
1:B:223:HIS:HB3	2:B:278:HOH:O	2.21	0.41
1:B:88:MET:HE3	1:B:110:THR:HG23	2.00	0.40
1:A:27:THR:O	1:A:30:ALA:HB3	2.22	0.40
1:B:5:GLN:O	1:B:9:GLU:HG3	2.20	0.40
1:B:234:VAL:CG2	1:C:247:LEU:HD22	2.52	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	254/259 (98%)	244 (96%)	10 (4%)	0	100	100
1	B	253/259 (98%)	245 (97%)	8 (3%)	0	100	100
1	C	251/259 (97%)	241 (96%)	10 (4%)	0	100	100
All	All	758/777 (98%)	730 (96%)	28 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	203/213 (95%)	199 (98%)	4 (2%)	55 69
1	B	205/213 (96%)	195 (95%)	10 (5%)	25 31
1	C	204/213 (96%)	199 (98%)	5 (2%)	47 60
All	All	612/639 (96%)	593 (97%)	19 (3%)	40 51

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	ARG
1	A	68	ARG
1	A	85	LYS
1	A	177	ILE
1	B	23	VAL
1	B	44	ARG
1	B	68	ARG
1	B	71	ASP
1	B	119	ARG
1	B	170	PRO
1	B	187	ARG
1	B	221	GLU
1	B	222	ILE
1	B	235	SER
1	C	44	ARG
1	C	76	ASP
1	C	81	ARG
1	C	148	SER
1	C	237	ASN

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

Mol	Chain	Res	Type
1	A	32	ASN
1	A	52	GLN
1	A	57	GLN
1	A	120	ASN
1	A	237	ASN
1	B	32	ASN
1	B	57	GLN
1	B	120	ASN
1	C	32	ASN
1	C	57	GLN

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Mol	Chain	Res	Type
1	C	120	ASN
1	C	157	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 monosaccharides 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, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
1	A	256/259 (98%)	-0.63	3 (1%) 79 77	8, 17, 46, 57	0
1	B	255/259 (98%)	-0.78	1 (0%) 92 91	6, 17, 43, 55	0
1	C	253/259 (97%)	-0.68	0 100 100	6, 17, 42, 61	0
All	All	764/777 (98%)	-0.70	4 (0%) 91 90	6, 17, 44, 61	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	4.4
1	A	51	GLY	3.0
1	A	50	ASP	2.6
1	B	2	GLN	2.2

6.2 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates

There are no monosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.