



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:05 PM UTC

PDB ID : 4O26 / pdb_00004o26
Title : Crystal structure of the TRBD domain of TERT and the CR4/5 of TR
Authors : Huang, J.; Wu, J.; Lei, M.
Deposited on : 2013-12-16
Resolution : 3.00 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
Mogul : 2022.3.0, CSD as543be (2022)
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

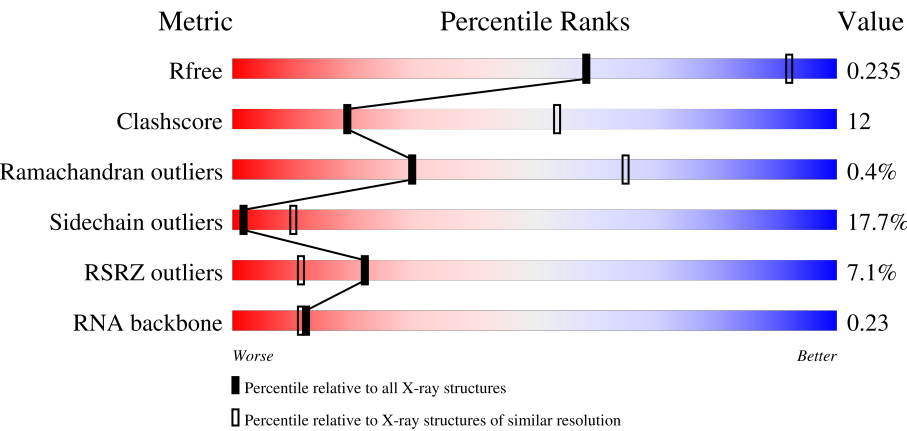
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)
RNA backbone	3983	1109 (3.20-2.80)

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	257	3% 61% 28% 7%
1	B	257	5% 53% 26% 9% 11%
2	E	51	14% 33% 29% 29% 8%
2	F	51	22% 33% 37% 22% 8%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SO4	A	601	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5940 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 Telomerase reverse transcriptase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	240	Total	C	N	O	S	0	0	0
			1965	1261	371	320	13			
1	B	228	Total	C	N	O	S	0	0	0
			1884	1214	354	304	12			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	316	GLY	-	expression tag	UNP Q1PS67
A	317	SER	-	expression tag	UNP Q1PS67
B	316	GLY	-	expression tag	UNP Q1PS67
B	317	SER	-	expression tag	UNP Q1PS67

- Molecule 2 is a RNA chain called Telomerase TR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	47	Total	C	N	O	P	0	0	0
			1000	445	174	334	47			
2	F	47	Total	C	N	O	P	0	0	0
			1000	445	174	334	47			

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

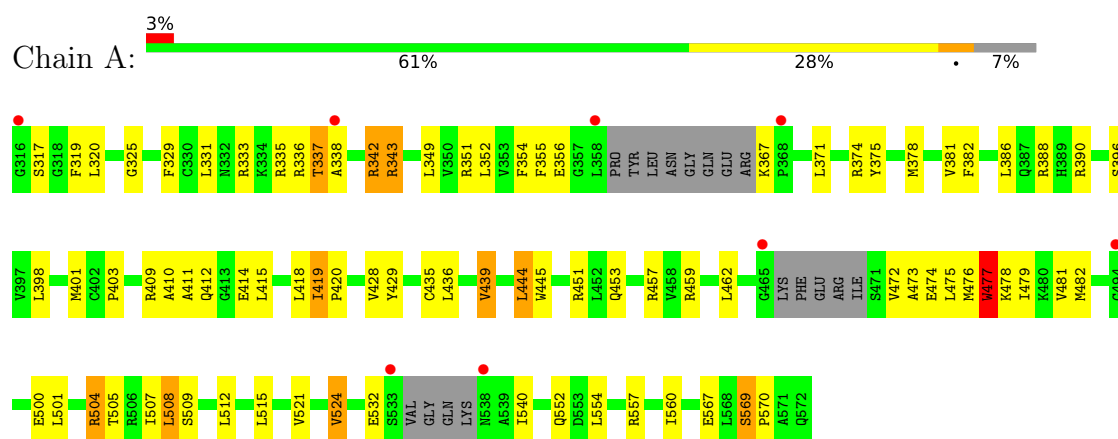
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	21	Total 21	O 21	0	0
4	B	8	Total 8	O 8	0	0
4	E	2	Total 2	O 2	0	0

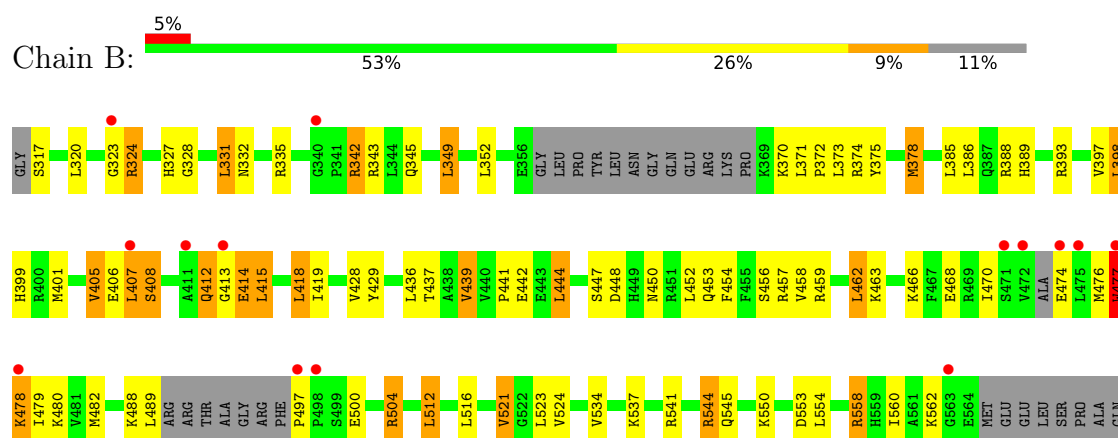
3 Residue-property plots [i](#)

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.

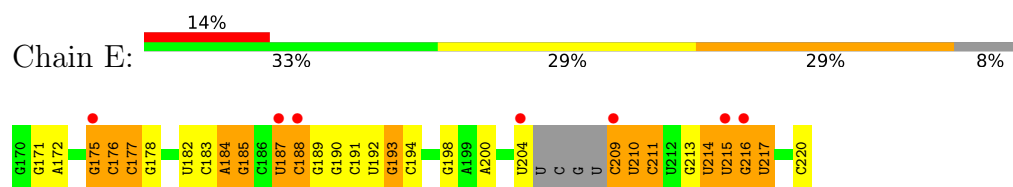
• Molecule 1: Telomerase reverse transcriptase



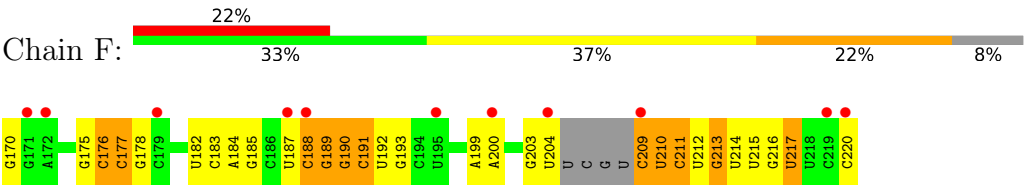
• Molecule 1: Telomerase reverse transcriptase



• Molecule 2: Telomerase TR



• Molecule 2: Telomerase TR



4 Data and refinement statistics

Property	Value	Source
Space group	P 62 2 2	Depositor
Cell constants a, b, c, α , β , γ	118.09Å 118.09Å 358.02Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	41.97 – 3.00 41.97 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.9 (41.97-3.00) 99.8 (41.97-3.00)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.85 (at 3.01Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309, REFMAC	Depositor
R, R_{free}	0.221 , 0.261 0.228 , 0.235	Depositor DCC
R_{free} test set	1543 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	74.8	Xtriage
Anisotropy	0.173	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 96.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5940	wwPDB-VP
Average B, all atoms (Å ²)	101.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.39% 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

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

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	0.49	0/2012	0.93	8/2702 (0.3%)
1	B	0.40	0/1929	0.96	8/2588 (0.3%)
2	E	0.21	0/1114	0.41	0/1732
2	F	0.15	0/1114	0.41	0/1732
All	All	0.37	0/6169	0.78	16/8754 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	470	ILE	CB-CA-C	-8.73	102.92	111.30
1	B	478	LYS	N-CA-C	-8.26	100.30	110.88
1	A	569	SER	CA-C-N	7.46	127.90	119.92
1	A	569	SER	C-N-CA	7.46	127.90	119.92
1	B	470	ILE	N-CA-C	-7.37	105.93	111.90
1	B	448	ASP	N-CA-C	-7.23	104.71	113.97
1	A	477	TRP	N-CA-C	7.12	119.99	107.61
1	B	477	TRP	N-CA-C	7.04	119.85	107.61
1	A	325	GLY	N-CA-C	6.74	118.91	112.04
1	B	497	PRO	CA-C-N	6.25	127.66	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	497	PRO	C-N-CA	6.25	127.66	119.84
1	A	419	ILE	CA-C-N	5.34	125.33	120.21
1	A	419	ILE	C-N-CA	5.34	125.33	120.21
1	A	478	LYS	N-CA-C	-5.20	99.72	110.80
1	B	412	GLN	N-CA-C	5.17	116.80	108.32
1	A	411	ALA	N-CA-C	-5.06	107.28	113.50

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	477	TRP	Peptide
1	B	477	TRP	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1965	0	1996	49	0
1	B	1884	0	1920	48	0
2	E	1000	0	508	20	1
2	F	1000	0	508	12	1
3	A	40	0	0	4	0
3	B	20	0	0	1	0
4	A	21	0	0	0	0
4	B	8	0	0	0	0
4	E	2	0	0	0	0
All	All	5940	0	4932	123	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 12.

All (123) 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:481:VAL:HG21	1:A:505:THR:HG22	1.53	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:209:C:HO2'	2:F:210:U:H6	1.28	0.80
1:B:332:ASN:OD1	1:B:389:HIS:NE2	2.16	0.77
1:B:323:GLY:O	1:B:545:GLN:NE2	2.19	0.75
1:A:532:GLU:HG2	1:A:540:ILE:HG13	1.70	0.74
1:B:415:LEU:HD12	1:B:554:LEU:HD13	1.70	0.73
1:B:447:SER:HG	1:B:450:ASN:H	1.39	0.71
1:A:482:MET:HE1	1:A:504:ARG:HH21	1.54	0.71
1:A:342:ARG:HD3	1:A:343:ARG:H	1.58	0.69
1:B:477:TRP:HE3	1:B:479:ILE:HB	1.60	0.65
1:B:374:ARG:NH2	1:B:500:GLU:OE1	2.27	0.63
2:F:170:G:H1	2:F:220:C:H42	1.48	0.61
1:B:477:TRP:CE3	1:B:479:ILE:HB	2.35	0.60
1:B:371:LEU:O	2:F:189:G:N1	2.35	0.60
1:A:343:ARG:NH2	3:A:601:SO4:O4	2.34	0.60
2:F:209:C:O2'	2:F:210:U:O5'	2.19	0.60
1:A:390:ARG:NH2	3:A:601:SO4:O2	2.34	0.60
2:E:209:C:O2'	2:E:210:U:OP1	2.18	0.60
2:E:191:C:N4	2:E:192:U:O4	2.36	0.59
1:B:408:SER:HB3	1:B:412:GLN:HA	1.86	0.58
1:A:351:ARG:HD3	1:A:367:LYS:HB3	1.86	0.58
1:B:405:VAL:HG12	1:B:406:GLU:H	1.68	0.58
2:F:176:C:N4	2:F:217:U:O2	2.37	0.57
1:A:342:ARG:HD3	1:A:343:ARG:N	2.19	0.56
1:A:375:TYR:HA	1:A:378:MET:HE2	1.85	0.56
1:B:558:ARG:HH11	1:B:558:ARG:HB3	1.70	0.56
1:A:374:ARG:NH2	1:A:500:GLU:OE1	2.38	0.56
1:B:480:LYS:NZ	2:F:212:U:OP1	2.38	0.56
1:A:451:ARG:NH2	1:B:553:ASP:OD2	2.40	0.55
1:B:349:LEU:HD11	1:B:385:LEU:HD22	1.87	0.55
1:A:388:ARG:HB3	1:A:439:VAL:HA	1.89	0.54
1:B:558:ARG:HH12	1:B:560:ILE:HD13	1.72	0.54
1:A:319:PHE:HB2	1:A:524:VAL:HG22	1.89	0.54
1:A:554:LEU:HD23	1:A:557:ARG:NH1	2.23	0.53
1:B:408:SER:HA	1:B:413:GLY:H	1.73	0.53
1:B:331:LEU:HB3	1:B:389:HIS:CD2	2.43	0.53
1:B:398:LEU:HD13	1:B:544:ARG:CZ	2.39	0.53
2:F:177:C:N3	2:F:217:U:H1'	2.23	0.53
1:B:349:LEU:HD12	1:B:386:LEU:HG	1.91	0.53
2:F:210:U:H2'	2:F:211:C:C6	2.44	0.52
1:A:473:ALA:O	1:A:477:TRP:NE1	2.37	0.52
1:B:441:PRO:HD2	1:B:444:LEU:HD23	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:188:C:O2'	2:F:190:G:N1	2.39	0.52
1:B:442:GLU:OE1	1:B:442:GLU:N	2.40	0.52
1:A:501:LEU:O	1:A:505:THR:HG23	2.09	0.51
1:A:390:ARG:NE	3:A:601:SO4:O2	2.43	0.51
1:A:476:MET:C	1:A:477:TRP:CD1	2.88	0.51
1:A:335:ARG:CZ	1:A:342:ARG:HB3	2.42	0.50
1:A:351:ARG:HG3	1:A:371:LEU:HD21	1.94	0.50
1:B:418:LEU:HD21	1:B:550:LYS:HB3	1.93	0.50
2:E:187:U:H4'	2:E:188:C:OP2	2.12	0.50
1:B:474:GLU:C	1:B:476:MET:H	2.20	0.50
2:E:177:C:N3	2:E:217:U:H1'	2.26	0.50
2:F:211:C:H2'	2:F:212:U:C6	2.47	0.50
1:A:451:ARG:HH22	1:B:553:ASP:CG	2.19	0.49
1:A:477:TRP:HZ3	1:A:508:LEU:HD13	1.76	0.49
1:B:399:HIS:ND1	3:B:603:SO4:O2	2.44	0.49
2:E:210:U:H2'	2:E:211:C:C6	2.48	0.49
1:B:405:VAL:O	1:B:407:LEU:HG	2.13	0.49
1:A:472:VAL:HG22	1:A:475:LEU:HB2	1.95	0.49
2:E:210:U:O2'	2:E:211:C:OP1	2.28	0.49
2:E:176:C:C2'	2:E:177:C:H5'	2.43	0.48
2:E:213:G:N2	2:E:214:U:O2	2.46	0.48
2:E:188:C:OP2	2:E:188:C:H4'	2.12	0.48
1:A:445:TRP:CE2	1:A:451:ARG:HG3	2.48	0.48
1:B:463:LYS:HB2	1:B:463:LYS:HE3	1.72	0.48
1:B:388:ARG:HB3	1:B:439:VAL:HA	1.96	0.47
1:A:335:ARG:HG2	1:A:342:ARG:O	2.14	0.47
1:A:349:LEU:HD23	1:A:382:PHE:HB3	1.97	0.47
1:A:457:ARG:NH2	1:A:475:LEU:O	2.48	0.47
1:B:373:LEU:HD23	1:B:489:LEU:HD11	1.97	0.47
1:B:374:ARG:HH21	1:B:504:ARG:HD3	1.79	0.47
1:A:435:CYS:O	1:A:439:VAL:HG13	2.15	0.46
1:B:370:LYS:O	1:B:372:PRO:HD3	2.15	0.46
1:A:351:ARG:HA	1:A:355:PHE:HD2	1.81	0.46
1:B:516:LEU:O	1:B:521:VAL:HG13	2.15	0.45
2:E:217:U:OP1	2:E:217:U:H3'	2.17	0.45
1:A:515:LEU:HD12	1:A:515:LEU:HA	1.83	0.45
2:E:175:G:O2'	2:E:176:C:H6	1.99	0.45
2:E:184:A:H2'	2:E:185:G:C8	2.51	0.45
1:A:329:PHE:CE2	1:A:331:LEU:HB2	2.52	0.45
1:B:454:PHE:O	1:B:458:VAL:HG23	2.15	0.45
1:B:482:MET:HB2	2:F:213:G:OP1	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:429:TYR:OH	3:A:607:SO4:O4	2.27	0.45
1:A:569:SER:HA	1:A:570:PRO:HD3	1.73	0.45
1:B:429:TYR:CE1	1:B:459:ARG:HB2	2.52	0.45
1:A:459:ARG:HD3	1:B:324:ARG:NH1	2.32	0.45
2:E:184:A:H2'	2:E:185:G:H8	1.82	0.45
1:A:375:TYR:O	1:A:378:MET:HG2	2.17	0.45
1:A:444:LEU:HD12	1:A:444:LEU:HA	1.84	0.45
1:B:523:LEU:HD23	1:B:523:LEU:HA	1.84	0.45
1:A:342:ARG:HH11	1:A:342:ARG:HA	1.82	0.44
1:A:554:LEU:HD23	1:A:557:ARG:HH12	1.81	0.44
2:E:192:U:HO2'	2:E:193:G:H8	1.65	0.44
1:B:342:ARG:HD3	1:B:343:ARG:H	1.82	0.44
1:B:393:ARG:O	1:B:397:VAL:HG23	2.18	0.44
2:E:193:G:H2'	2:E:194:C:O4'	2.17	0.44
1:B:323:GLY:O	1:B:545:GLN:CD	2.60	0.44
1:A:477:TRP:CE3	1:A:479:ILE:HB	2.52	0.44
1:B:436:LEU:HD12	1:B:436:LEU:HA	1.84	0.44
1:A:354:PHE:CE1	1:A:507:ILE:HG23	2.53	0.44
1:B:512:LEU:HD12	1:B:512:LEU:HA	1.81	0.44
2:F:191:C:N4	2:F:192:U:O4	2.51	0.43
1:A:401:MET:C	1:A:403:PRO:HD3	2.44	0.42
1:A:419:ILE:HA	1:A:420:PRO:HD3	1.85	0.42
1:A:477:TRP:HZ2	1:A:509:SER:HB3	1.83	0.42
1:A:335:ARG:NH1	1:A:343:ARG:O	2.52	0.42
1:B:462:LEU:HD12	1:B:462:LEU:HA	1.85	0.42
1:A:472:VAL:CG2	1:A:475:LEU:HB2	2.49	0.42
1:B:414:GLU:CD	1:B:414:GLU:H	2.28	0.42
1:A:336:ARG:O	1:A:338:ALA:N	2.53	0.42
2:E:213:G:C2	2:E:214:U:O2	2.72	0.42
2:E:215:U:O2'	2:E:216:G:H5'	2.20	0.41
1:B:335:ARG:HH22	1:B:345:GLN:NE2	2.19	0.41
1:B:374:ARG:NH2	1:B:504:ARG:HH11	2.17	0.41
2:E:176:C:H2'	2:E:177:C:H5'	2.03	0.41
1:A:476:MET:C	1:A:477:TRP:CG	2.99	0.41
1:A:409:ARG:O	1:A:410:ALA:HB3	2.21	0.41
2:E:193:G:H2'	2:E:194:C:C6	2.56	0.40
1:A:477:TRP:HH2	1:A:508:LEU:HB3	1.85	0.40
1:B:331:LEU:HB3	1:B:389:HIS:NE2	2.36	0.40
1:B:375:TYR:HA	1:B:378:MET:HG3	2.02	0.40
2:E:209:C:O2'	2:E:210:U:P	2.80	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:220:C:O2'	2:F:188:C:OP1[8_565]	2.12	0.08

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	232/257 (90%)	213 (92%)	18 (8%)	1 (0%)	30	65
1	B	220/257 (86%)	198 (90%)	21 (10%)	1 (0%)	24	60
All	All	452/514 (88%)	411 (91%)	39 (9%)	2 (0%)	30	65

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	337	THR
1	B	328	GLY

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	207/222 (93%)	176 (85%)	31 (15%)	3	14
1	B	200/222 (90%)	159 (80%)	41 (20%)	1	7
All	All	407/444 (92%)	335 (82%)	72 (18%)	2	10

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	317	SER
1	A	320	LEU
1	A	333	ARG
1	A	337	THR
1	A	342	ARG
1	A	343	ARG
1	A	352	LEU
1	A	356	GLU
1	A	381	VAL
1	A	386	LEU
1	A	396	SER
1	A	398	LEU
1	A	412	GLN
1	A	414	GLU
1	A	415	LEU
1	A	418	LEU
1	A	428	VAL
1	A	436	LEU
1	A	439	VAL
1	A	444	LEU
1	A	453	GLN
1	A	462	LEU
1	A	474	GLU
1	A	504	ARG
1	A	508	LEU
1	A	512	LEU
1	A	521	VAL
1	A	524	VAL
1	A	552	GLN
1	A	560	ILE
1	A	567	GLU
1	B	317	SER
1	B	320	LEU
1	B	324	ARG
1	B	327	HIS
1	B	331	LEU
1	B	342	ARG
1	B	349	LEU
1	B	352	LEU
1	B	378	MET
1	B	398	LEU
1	B	401	MET

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Mol	Chain	Res	Type
1	B	405	VAL
1	B	407	LEU
1	B	408	SER
1	B	414	GLU
1	B	415	LEU
1	B	418	LEU
1	B	419	ILE
1	B	428	VAL
1	B	437	THR
1	B	439	VAL
1	B	444	LEU
1	B	452	LEU
1	B	453	GLN
1	B	456	SER
1	B	457	ARG
1	B	462	LEU
1	B	466	LYS
1	B	468	GLU
1	B	478	LYS
1	B	488	LYS
1	B	504	ARG
1	B	512	LEU
1	B	521	VAL
1	B	524	VAL
1	B	534	VAL
1	B	537	LYS
1	B	541	ARG
1	B	544	ARG
1	B	558	ARG
1	B	562	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	453	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	E	46/51 (90%)	24 (52%)	3 (6%)

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Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	F	46/51 (90%)	25 (54%)	2 (4%)
All	All	92/102 (90%)	49 (53%)	5 (5%)

All (49) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	E	171	G
2	E	172	A
2	E	175	G
2	E	176	C
2	E	177	C
2	E	178	G
2	E	182	U
2	E	183	C
2	E	184	A
2	E	185	G
2	E	187	U
2	E	188	C
2	E	189	G
2	E	190	G
2	E	193	G
2	E	198	G
2	E	200	A
2	E	204	U
2	E	210	U
2	E	211	C
2	E	214	U
2	E	215	U
2	E	216	G
2	E	217	U
2	F	175	G
2	F	176	C
2	F	177	C
2	F	178	G
2	F	182	U
2	F	183	C
2	F	184	A
2	F	185	G
2	F	187	U
2	F	188	C
2	F	189	G
2	F	190	G

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Mol	Chain	Res	Type
2	F	191	C
2	F	193	G
2	F	199	A
2	F	200	A
2	F	203	G
2	F	204	U
2	F	210	U
2	F	211	C
2	F	213	G
2	F	214	U
2	F	215	U
2	F	216	G
2	F	217	U

All (5) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	E	187	U
2	E	209	C
2	E	210	U
2	F	209	C
2	F	210	U

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](#)

12 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	SO4	A	606	-	4,4,4	0.24	0	6,6,6	0.17	0
3	SO4	A	602	-	4,4,4	0.24	0	6,6,6	0.14	0
3	SO4	A	601	-	4,4,4	0.18	0	6,6,6	0.21	0
3	SO4	A	608	-	4,4,4	0.22	0	6,6,6	0.26	0
3	SO4	B	604	-	4,4,4	0.24	0	6,6,6	0.07	0
3	SO4	A	605	-	4,4,4	0.22	0	6,6,6	0.09	0
3	SO4	A	604	-	4,4,4	0.26	0	6,6,6	0.24	0
3	SO4	B	603	-	4,4,4	0.21	0	6,6,6	0.23	0
3	SO4	B	601	-	4,4,4	0.22	0	6,6,6	0.19	0
3	SO4	A	603	-	4,4,4	0.22	0	6,6,6	0.21	0
3	SO4	B	602	-	4,4,4	0.23	0	6,6,6	0.22	0
3	SO4	A	607	-	4,4,4	0.27	0	6,6,6	0.19	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	601	SO4	3	0
3	B	603	SO4	1	0
3	A	607	SO4	1	0

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	240/257 (93%)	-0.02	8 (3%) 49 28	35, 57, 127, 157	0
1	B	228/257 (88%)	0.35	14 (6%) 27 14	48, 84, 139, 187	0
2	E	47/51 (92%)	0.64	7 (14%) 5 3	80, 111, 200, 228	0
2	F	47/51 (92%)	1.37	11 (23%) 2 1	114, 157, 275, 331	0
All	All	562/616 (91%)	0.30	40 (7%) 22 11	35, 79, 164, 331	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	497	PRO	6.4
1	B	474	GLU	5.3
1	A	358	LEU	4.7
1	A	494	GLY	4.2
1	A	465	GLY	3.6
2	E	216	G	3.6
1	B	340	GLY	3.5
1	B	407	LEU	3.5
1	A	338	ALA	3.3
2	F	200	A	3.3
2	E	204	U	3.3
1	A	533	SER	3.2
2	F	187	U	3.1
2	E	187	U	3.0
2	F	179	C	3.0
2	F	204	U	2.9
1	B	411	ALA	2.9
1	B	475	LEU	2.9
1	A	368	PRO	2.8
1	B	477	TRP	2.8
2	F	209	C	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	538	ASN	2.8
2	F	172	A	2.7
2	F	188	C	2.6
2	E	215	U	2.6
2	F	171	G	2.5
1	B	471	SER	2.4
1	B	323	GLY	2.4
2	E	175	G	2.4
2	F	220	C	2.3
2	F	219	C	2.3
1	B	472	VAL	2.3
1	B	413	GLY	2.2
1	B	498	PRO	2.2
2	F	195	U	2.2
2	E	188	C	2.1
1	B	478	LYS	2.1
1	B	563	GLY	2.1
2	E	209	C	2.1
1	A	316	GLY	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	SO4	B	602	5/5	0.81	0.18	113,129,136,141	0
3	SO4	A	602	5/5	0.84	0.21	115,120,122,130	0
3	SO4	B	604	5/5	0.87	0.08	138,138,146,147	0
3	SO4	A	605	5/5	0.89	0.08	139,146,154,155	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	A	608	5/5	0.90	0.37	97,117,121,138	0
3	SO4	A	606	5/5	0.94	0.17	80,86,92,97	0
3	SO4	B	603	5/5	0.95	0.08	70,84,93,97	0
3	SO4	A	607	5/5	0.95	0.09	56,65,72,75	0
3	SO4	B	601	5/5	0.96	0.14	55,69,84,90	0
3	SO4	A	603	5/5	0.96	0.21	95,96,109,135	0
3	SO4	A	601	5/5	0.98	0.14	42,43,65,76	5
3	SO4	A	604	5/5	0.98	0.10	61,75,81,83	0

6.5 Other polymers [i](#)

There are no such residues in this entry.