



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:55 AM UTC

PDB ID : 1TLZ / pdb_00001tlz
Title : Tsx structure complexed with uridine
Authors : Ye, J.; van den Berg, B.
Deposited on : 2004-06-10
Resolution : 3.10 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

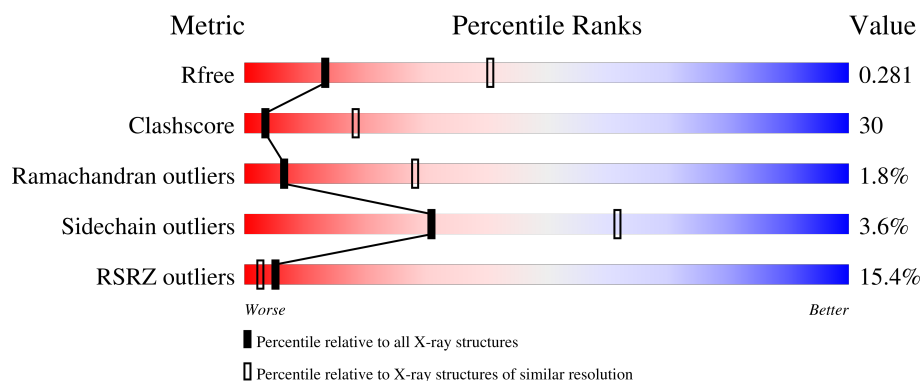
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	278	4% 44% 40% 6% 10%
1	B	278	24% 42% 43% • 11%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4176 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.

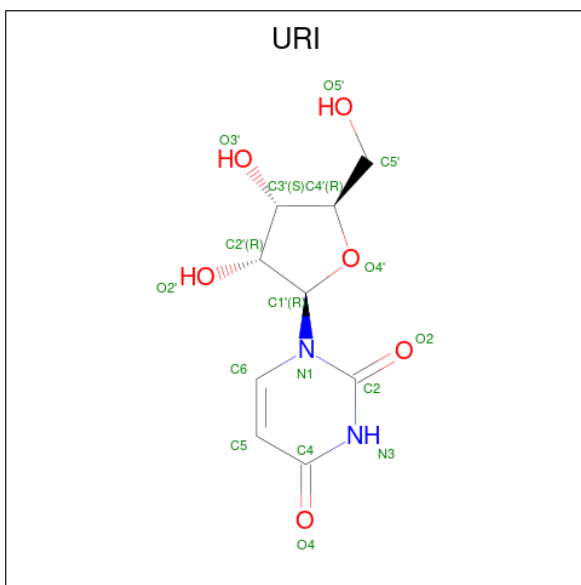
- Molecule 1 is a protein called Nucleoside-specific channel-forming protein tsx.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	251	Total	C	N	O	S	0	0	0
			2078	1340	347	386	5			
1	B	248	Total	C	N	O	S	0	0	0
			2048	1322	338	383	5			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	273	HIS	-	expression tag	UNP P0A927
A	274	HIS	-	expression tag	UNP P0A927
A	275	HIS	-	expression tag	UNP P0A927
A	276	HIS	-	expression tag	UNP P0A927
A	277	HIS	-	expression tag	UNP P0A927
A	278	HIS	-	expression tag	UNP P0A927
B	273	HIS	-	expression tag	UNP P0A927
B	274	HIS	-	expression tag	UNP P0A927
B	275	HIS	-	expression tag	UNP P0A927
B	276	HIS	-	expression tag	UNP P0A927
B	277	HIS	-	expression tag	UNP P0A927
B	278	HIS	-	expression tag	UNP P0A927

- Molecule 2 is URIDINE (CCD ID: URI) (formula: C₉H₁₂N₂O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			17	9	2	6		
2	B	1	Total	C	N	O	0	0
			17	9	2	6		

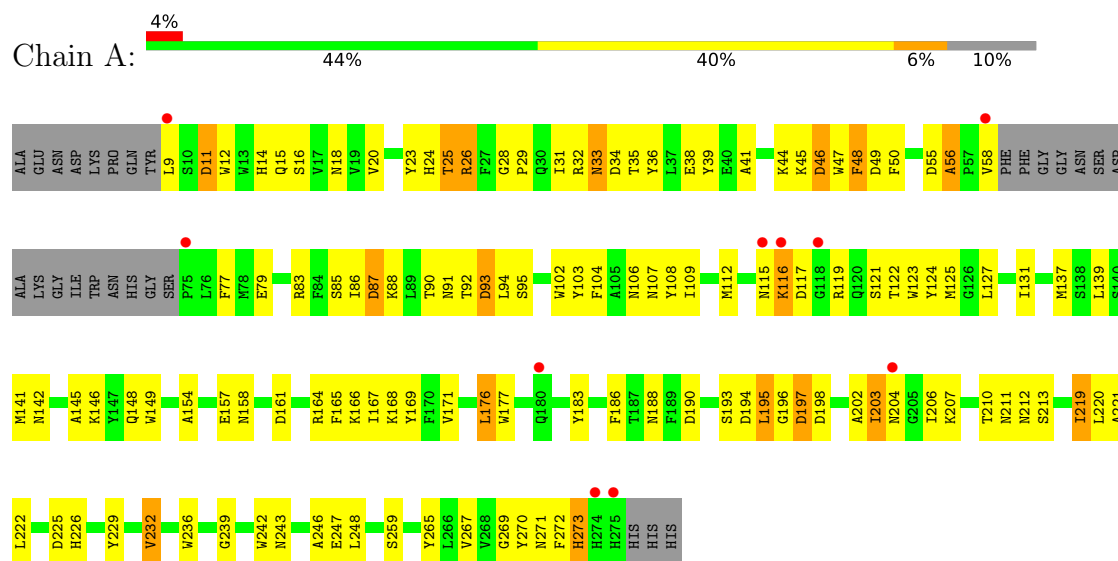
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	13	Total	O	0	0
			13	13		
3	B	3	Total	O	0	0
			3	3		

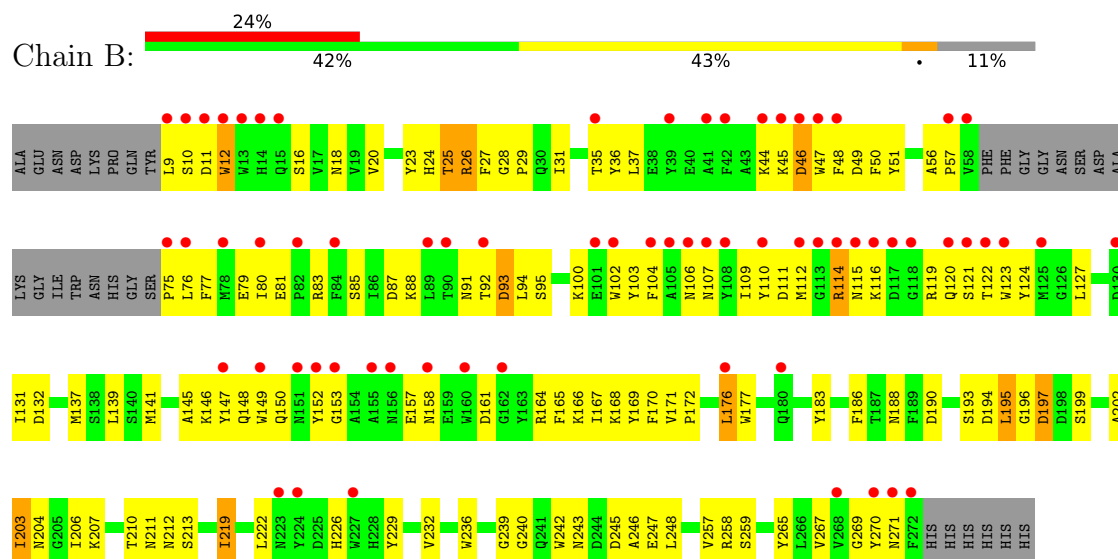
3 Residue-property plots

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: Nucleoside-specific channel-forming protein tsx



- Molecule 1: Nucleoside-specific channel-forming protein tsx



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	149.86Å 149.86Å 120.85Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.77 – 3.10 29.77 – 3.10	Depositor EDS
% Data completeness (in resolution range)	97.2 (29.77-3.10) 97.1 (29.77-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.08 (at 3.01Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.276 , 0.299 0.257 , 0.281	Depositor DCC
R_{free} test set	2085 reflections (6.58%)	wwPDB-VP
Wilson B-factor (Å ²)	56.2	Xtriage
Anisotropy	0.124	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 48.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	4176	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.84% 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:
URI

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	0.55	0/2154	1.07	16/2931 (0.5%)
1	B	0.41	0/2121	0.98	10/2886 (0.3%)
All	All	0.49	0/4275	1.03	26/5817 (0.4%)

There are no bond length outliers.

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	116	LYS	N-CA-C	-9.42	101.71	113.01
1	B	203	ILE	N-CA-C	-9.14	100.19	110.62
1	A	203	ILE	N-CA-C	-8.28	101.18	110.62
1	A	145	ALA	N-CA-C	-7.26	98.89	110.14
1	B	145	ALA	N-CA-C	-7.16	98.57	109.95
1	B	197	ASP	N-CA-C	-6.95	103.63	111.14
1	A	104	PHE	N-CA-C	-6.91	95.83	107.99
1	A	56	ALA	CA-C-N	6.83	126.87	119.90
1	A	56	ALA	C-N-CA	6.83	126.87	119.90
1	A	26	ARG	N-CA-C	6.71	120.83	111.56
1	A	219	ILE	N-CA-C	6.53	117.52	108.12
1	B	104	PHE	N-CA-C	-6.46	95.73	107.75
1	A	28	GLY	CA-C-N	6.33	126.08	119.05
1	A	28	GLY	C-N-CA	6.33	126.08	119.05
1	A	86	ILE	N-CA-C	-5.76	104.90	110.72
1	B	26	ARG	N-CA-C	5.55	119.22	111.56
1	A	232	VAL	N-CA-C	5.52	115.94	107.77
1	B	28	GLY	CA-C-N	5.47	125.12	119.05
1	B	28	GLY	C-N-CA	5.47	125.12	119.05
1	A	197	ASP	N-CA-C	-5.44	105.25	111.07
1	A	103	TYR	N-CA-C	5.41	117.41	109.24
1	A	221	ALA	N-CA-C	5.33	117.59	108.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	219	ILE	N-CA-C	5.22	116.00	108.48
1	B	203	ILE	CA-C-N	5.05	127.76	120.38
1	B	203	ILE	C-N-CA	5.05	127.76	120.38
1	A	33	ASN	N-CA-C	-5.04	99.89	108.26

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	2078	0	1867	118	0
1	B	2048	0	1846	115	0
2	A	17	0	12	3	0
2	B	17	0	12	4	0
3	A	13	0	0	1	0
3	B	3	0	0	0	0
All	All	4176	0	3737	236	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 30.

All (236) 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:210:THR:HG22	1:B:212:ASN:H	1.31	0.95
1:B:146:LYS:HE2	1:B:161:ASP:O	1.66	0.94
1:B:9:LEU:HB3	1:B:12:TRP:HE1	1.34	0.91
1:A:210:THR:HG22	1:A:212:ASN:H	1.41	0.86
1:A:146:LYS:HE2	1:A:161:ASP:O	1.77	0.84
1:B:29:PRO:HG3	1:B:242:TRP:CZ2	2.12	0.83
2:B:801:URI:O5'	2:B:801:URI:H6	1.80	0.81
1:B:204:ASN:HB2	1:B:247:GLU:H	1.47	0.78
1:A:29:PRO:HG3	1:A:242:TRP:CZ2	2.19	0.78
1:A:106:ASN:HD22	1:A:123:TRP:HE1	1.32	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:107:ASN:ND2	1:B:164:ARG:HH11	1.83	0.77
1:A:116:LYS:HA	1:A:119:ARG:HH11	1.50	0.75
1:B:122:THR:HG23	1:B:146:LYS:HD2	1.67	0.74
1:A:204:ASN:HB2	1:A:247:GLU:H	1.53	0.72
1:B:9:LEU:HB3	1:B:12:TRP:NE1	2.05	0.71
1:A:107:ASN:ND2	1:A:164:ARG:HH11	1.91	0.69
1:A:193:SER:O	1:A:211:ASN:ND2	2.26	0.69
1:B:29:PRO:HG3	1:B:242:TRP:CE2	2.29	0.67
1:A:131:ILE:HB	1:A:139:LEU:HD12	1.77	0.66
1:A:204:ASN:HB2	1:A:247:GLU:HB3	1.76	0.66
1:A:219:ILE:HG12	1:A:232:VAL:HG22	1.78	0.66
1:B:219:ILE:HG12	1:B:232:VAL:HG22	1.78	0.64
1:A:122:THR:HG23	1:A:146:LYS:HD2	1.79	0.64
1:B:193:SER:O	1:B:211:ASN:ND2	2.31	0.64
1:B:100:LYS:HD2	1:B:132:ASP:HB2	1.79	0.63
1:B:169:TYR:HE1	1:B:183:TYR:HE2	1.47	0.63
1:B:131:ILE:HB	1:B:139:LEU:HD12	1.81	0.62
1:B:46:ASP:O	1:B:88:LYS:HG2	1.99	0.62
1:B:45:LYS:O	1:B:47:TRP:N	2.33	0.62
1:B:122:THR:CG2	1:B:146:LYS:HD2	2.30	0.61
1:A:115:ASN:C	1:A:117:ASP:N	2.58	0.61
1:A:169:TYR:HE1	1:A:183:TYR:CE2	2.18	0.61
1:B:36:TYR:HB2	1:B:56:ALA:O	1.98	0.61
1:A:46:ASP:O	1:A:88:LYS:HG2	2.01	0.61
1:A:109:ILE:N	1:A:109:ILE:HD12	2.16	0.61
1:A:115:ASN:C	1:A:117:ASP:H	2.08	0.60
1:B:169:TYR:HE1	1:B:183:TYR:CE2	2.20	0.60
1:A:32:ARG:NH2	1:A:34:ASP:OD2	2.35	0.60
1:A:210:THR:HG22	1:A:212:ASN:N	2.14	0.60
1:B:210:THR:HG22	1:B:212:ASN:N	2.10	0.60
1:A:210:THR:HG22	1:A:212:ASN:O	2.01	0.60
1:A:112:MET:HG3	1:A:119:ARG:NH2	2.18	0.59
1:A:45:LYS:O	1:A:47:TRP:N	2.35	0.59
1:A:169:TYR:HE1	1:A:183:TYR:HE2	1.50	0.59
1:B:100:LYS:CG	1:B:132:ASP:HB2	2.32	0.59
1:A:20:VAL:O	1:A:35:THR:HA	2.02	0.59
1:A:36:TYR:HB2	1:A:56:ALA:O	2.03	0.59
1:A:44:LYS:HD2	1:A:49:ASP:HB3	1.85	0.59
1:A:204:ASN:HB2	1:A:247:GLU:CB	2.33	0.59
1:B:49:ASP:OD2	1:B:83:ARG:HB2	2.03	0.59
1:B:226:HIS:HB3	1:B:270:TYR:CE1	2.37	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:SER:HB2	1:A:271:ASN:OD1	2.04	0.58
1:B:148:GLN:O	1:B:158:ASN:HA	2.04	0.58
1:B:111:ASP:O	1:B:119:ARG:HB3	2.03	0.58
1:A:102:TRP:CZ3	1:A:127:LEU:HD11	2.39	0.58
1:B:102:TRP:CE3	1:B:127:LEU:HD11	2.40	0.57
1:B:204:ASN:HB2	1:B:247:GLU:HB3	1.87	0.57
1:B:16:SER:HB2	1:B:271:ASN:OD1	2.05	0.57
1:A:210:THR:CG2	1:A:212:ASN:O	2.53	0.56
1:B:109:ILE:N	1:B:109:ILE:HD12	2.20	0.55
1:A:112:MET:HG3	1:A:119:ARG:HH22	1.72	0.54
1:B:20:VAL:O	1:B:35:THR:HA	2.07	0.54
1:A:139:LEU:HD23	1:A:169:TYR:HB3	1.90	0.54
1:A:29:PRO:HG3	1:A:242:TRP:CE2	2.42	0.54
1:B:195:LEU:HD22	1:B:196:GLY:H	1.71	0.54
1:A:115:ASN:HB2	1:A:117:ASP:HB3	1.90	0.54
1:A:148:GLN:O	1:A:158:ASN:HA	2.07	0.53
1:A:77:PHE:CZ	2:A:800:URI:H2'	2.44	0.53
1:B:107:ASN:HB3	1:B:124:TYR:HB2	1.91	0.53
1:A:206:ILE:HD12	1:A:243:ASN:ND2	2.24	0.53
1:B:169:TYR:CE1	1:B:183:TYR:HE2	2.26	0.53
1:A:77:PHE:CE2	2:A:800:URI:H2'	2.44	0.53
1:B:93:ASP:OD1	1:B:95:SER:HB3	2.08	0.53
1:B:100:LYS:CD	1:B:132:ASP:HB2	2.38	0.53
1:A:38:GLU:HG3	1:A:55:ASP:OD1	2.09	0.52
1:B:239:GLY:N	1:B:259:SER:HB2	2.24	0.52
1:A:239:GLY:N	1:A:259:SER:HB2	2.24	0.52
1:A:210:THR:CG2	1:A:212:ASN:H	2.17	0.52
1:B:176:LEU:HD22	1:B:177:TRP:N	2.25	0.52
1:B:204:ASN:HB2	1:B:247:GLU:N	2.23	0.52
1:B:141:MET:HE3	1:B:165:PHE:HZ	1.74	0.52
1:B:139:LEU:HD23	1:B:169:TYR:HB3	1.91	0.52
1:B:196:GLY:O	1:B:207:LYS:HD3	2.10	0.52
1:A:226:HIS:HB3	1:A:270:TYR:CE1	2.44	0.52
1:A:102:TRP:CE3	1:A:127:LEU:HD11	2.46	0.52
1:B:50:PHE:HE1	1:B:80:ILE:HB	1.75	0.51
1:A:166:LYS:HG3	1:A:188:ASN:HD21	1.75	0.51
1:A:87:ASP:O	1:A:91:ASN:N	2.43	0.51
1:B:44:LYS:HD2	1:B:49:ASP:HB3	1.91	0.51
1:A:122:THR:CG2	1:A:146:LYS:HD2	2.40	0.51
1:A:137:MET:CB	1:A:171:VAL:HG22	2.41	0.51
1:A:204:ASN:HD22	1:A:206:ILE:HG12	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:TYR:CE1	1:A:183:TYR:HE2	2.28	0.51
1:B:137:MET:CB	1:B:171:VAL:HG22	2.41	0.50
1:A:137:MET:HB2	1:A:171:VAL:HG22	1.91	0.50
1:A:167:ILE:O	1:A:186:PHE:HA	2.11	0.50
1:B:120:GLN:HA	1:B:150:GLN:O	2.11	0.50
1:A:93:ASP:OD1	1:A:95:SER:HB3	2.11	0.50
1:A:195:LEU:O	1:A:211:ASN:ND2	2.45	0.50
1:B:166:LYS:HD3	1:B:168:LYS:HE3	1.94	0.50
1:B:92:THR:O	1:B:94:LEU:HG	2.11	0.50
1:B:206:ILE:HD12	1:B:243:ASN:ND2	2.27	0.49
1:A:116:LYS:HA	1:A:119:ARG:HD2	1.93	0.49
1:B:79:GLU:HG3	1:B:109:ILE:HG13	1.93	0.49
1:A:92:THR:O	1:A:94:LEU:HG	2.12	0.49
1:B:115:ASN:C	1:B:115:ASN:HD22	2.19	0.49
1:A:206:ILE:HD11	1:A:246:ALA:HA	1.95	0.49
1:B:26:ARG:O	1:B:239:GLY:HA3	2.13	0.49
1:B:107:ASN:ND2	1:B:164:ARG:NH1	2.59	0.49
1:B:112:MET:HG3	1:B:119:ARG:HH11	1.77	0.49
1:A:12:TRP:CD1	1:A:12:TRP:H	2.31	0.49
1:A:85:SER:HB3	1:A:88:LYS:HB2	1.95	0.49
1:A:9:LEU:HG	1:A:11:ASP:H	1.78	0.49
1:B:9:LEU:HG	1:B:11:ASP:HB2	1.95	0.48
1:B:229:TYR:CD1	1:B:229:TYR:N	2.81	0.48
1:A:90:THR:HG22	1:A:92:THR:OG1	2.12	0.48
1:B:102:TRP:CZ3	1:B:127:LEU:HD11	2.48	0.48
1:A:203:ILE:HD12	1:A:247:GLU:HG2	1.96	0.48
1:A:204:ASN:HB2	1:A:247:GLU:N	2.23	0.48
1:B:167:ILE:O	1:B:186:PHE:HA	2.13	0.48
1:B:131:ILE:HB	1:B:139:LEU:CD1	2.43	0.48
1:B:121:SER:HB3	1:B:149:TRP:HB2	1.96	0.48
1:B:110:TYR:CZ	1:B:112:MET:HB2	2.48	0.48
1:A:26:ARG:HB2	1:A:236:TRP:CD2	2.49	0.48
1:A:18:ASN:OD1	1:A:269:GLY:HA3	2.14	0.48
1:B:210:THR:HG22	1:B:212:ASN:O	2.13	0.48
1:A:219:ILE:HG23	1:A:232:VAL:HG22	1.95	0.48
1:B:212:ASN:O	1:B:213:SER:HB2	2.13	0.48
1:A:131:ILE:HB	1:A:139:LEU:CD1	2.44	0.47
1:A:176:LEU:HD22	1:A:177:TRP:CD1	2.49	0.47
1:B:57:PRO:HD3	1:B:75:PRO:O	2.14	0.47
1:A:23:TYR:O	1:A:24:HIS:HB2	2.13	0.47
1:A:107:ASN:HB3	1:A:124:TYR:HB2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:801:URI:H6	2:B:801:URI:HO5'	1.77	0.47
1:B:137:MET:HB2	1:B:171:VAL:HG22	1.96	0.47
1:A:194:ASP:O	1:A:197:ASP:HB2	2.15	0.47
1:A:212:ASN:O	1:A:213:SER:HB2	2.14	0.47
1:B:107:ASN:HD22	1:B:164:ARG:HH11	1.60	0.47
1:A:49:ASP:OD2	1:A:49:ASP:C	2.58	0.47
1:B:164:ARG:HB2	1:B:190:ASP:OD2	2.15	0.47
1:B:18:ASN:OD1	1:B:269:GLY:HA3	2.15	0.46
1:B:202:ALA:C	1:B:204:ASN:N	2.69	0.46
1:A:107:ASN:HD22	1:A:164:ARG:HH11	1.61	0.46
1:B:168:LYS:HB3	1:B:186:PHE:HB3	1.96	0.46
1:B:87:ASP:O	1:B:91:ASN:N	2.47	0.46
1:A:229:TYR:CD1	1:A:229:TYR:N	2.83	0.46
1:B:10:SER:C	1:B:12:TRP:H	2.24	0.46
1:A:195:LEU:HD22	1:A:196:GLY:H	1.80	0.46
1:B:232:VAL:HB	1:B:265:TYR:HB2	1.98	0.46
1:A:121:SER:HB3	1:A:149:TRP:HB2	1.99	0.45
1:B:204:ASN:HB2	1:B:247:GLU:CB	2.46	0.45
1:A:154:ALA:O	1:A:157:GLU:HG3	2.16	0.45
1:B:77:PHE:CE2	2:B:801:URI:H3'	2.51	0.45
1:A:14:HIS:O	1:A:41:ALA:HA	2.16	0.45
1:A:26:ARG:O	1:A:239:GLY:HA3	2.16	0.45
1:A:92:THR:O	1:A:94:LEU:N	2.50	0.45
1:A:167:ILE:O	1:A:167:ILE:HG23	2.17	0.45
1:B:20:VAL:HG13	1:B:267:VAL:HG22	1.99	0.45
1:A:202:ALA:HB1	1:A:247:GLU:O	2.17	0.45
1:A:24:HIS:HA	1:A:31:ILE:HG23	1.98	0.45
1:A:198:ASP:N	1:A:198:ASP:OD2	2.50	0.45
1:A:220:LEU:HD12	1:A:220:LEU:HA	1.69	0.45
1:A:15:GLN:NE2	1:A:39:TYR:OH	2.36	0.44
1:A:196:GLY:O	1:A:207:LYS:HD3	2.18	0.44
1:A:225:ASP:O	1:A:273:HIS:HE1	2.00	0.44
1:B:210:THR:CG2	1:B:212:ASN:O	2.66	0.44
1:A:9:LEU:HD21	1:A:11:ASP:OD1	2.18	0.44
1:B:85:SER:OG	1:B:88:LYS:HD3	2.18	0.44
1:B:166:LYS:HG3	1:B:188:ASN:HD21	1.82	0.44
1:A:204:ASN:HD22	1:A:206:ILE:CD1	2.31	0.44
1:B:92:THR:O	1:B:94:LEU:N	2.51	0.43
1:B:195:LEU:CD2	1:B:199:SER:OG	2.66	0.43
1:B:27:PHE:O	1:B:240:GLY:HA3	2.18	0.43
1:B:51:TYR:O	1:B:81:GLU:N	2.45	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:50:PHE:CD2	1:A:50:PHE:C	2.94	0.43
1:A:164:ARG:HB2	1:A:190:ASP:OD2	2.18	0.43
1:A:90:THR:HG22	1:A:90:THR:O	2.19	0.43
1:B:106:ASN:HD22	1:B:123:TRP:HE1	1.65	0.43
1:A:141:MET:HE3	1:A:165:PHE:HZ	1.84	0.43
1:B:100:LYS:HG3	1:B:132:ASP:HB2	2.01	0.43
1:A:272:PHE:CD1	1:A:272:PHE:N	2.86	0.43
1:B:194:ASP:O	1:B:197:ASP:HB2	2.18	0.43
1:A:123:TRP:CD2	1:A:125:MET:HE2	2.54	0.43
1:B:85:SER:HB3	1:B:88:LYS:HB2	1.99	0.43
1:B:195:LEU:O	1:B:211:ASN:ND2	2.52	0.43
1:A:202:ALA:O	1:A:203:ILE:C	2.62	0.42
1:B:25:THR:HA	1:B:236:TRP:CH2	2.53	0.42
1:B:76:LEU:HD23	1:B:112:MET:HE2	2.00	0.42
1:B:161:ASP:OD2	1:B:193:SER:HA	2.19	0.42
2:B:801:URI:H3'	2:B:801:URI:C6	2.49	0.42
1:A:25:THR:HA	1:A:236:TRP:CH2	2.54	0.42
1:A:49:ASP:OD2	1:A:49:ASP:O	2.38	0.42
1:B:24:HIS:HA	1:B:31:ILE:HG23	2.01	0.42
1:B:153:GLY:HA2	1:B:157:GLU:OE1	2.20	0.42
1:B:203:ILE:HD12	1:B:247:GLU:HG2	2.00	0.42
1:A:107:ASN:ND2	1:A:164:ARG:NH1	2.63	0.42
1:B:147:TYR:HE2	1:B:149:TRP:CE2	2.38	0.42
1:A:202:ALA:C	1:A:204:ASN:N	2.74	0.42
1:B:115:ASN:OD1	1:B:152:TYR:CE2	2.73	0.42
1:A:272:PHE:O	1:A:273:HIS:HB3	2.20	0.41
1:B:26:ARG:HB2	1:B:236:TRP:CD2	2.55	0.41
1:A:20:VAL:HG13	1:A:267:VAL:HG22	2.03	0.41
1:B:204:ASN:ND2	1:B:245:ASP:O	2.53	0.41
1:A:168:LYS:CB	1:A:186:PHE:HB3	2.50	0.41
1:B:167:ILE:O	1:B:167:ILE:HG23	2.20	0.41
1:A:58:VAL:O	1:A:58:VAL:HG12	2.20	0.41
1:A:176:LEU:HD22	1:A:177:TRP:N	2.35	0.41
1:A:232:VAL:HB	1:A:265:TYR:HB2	2.01	0.41
1:B:204:ASN:HD22	1:B:206:ILE:CD1	2.34	0.41
1:A:25:THR:HG22	3:A:812:HOH:O	2.21	0.41
1:A:33:ASN:C	1:A:34:ASP:OD1	2.64	0.41
1:B:202:ALA:O	1:B:203:ILE:C	2.62	0.41
1:A:9:LEU:HB3	1:A:12:TRP:HE1	1.85	0.41
1:B:23:TYR:CZ	1:B:24:HIS:CE1	3.09	0.41
1:B:103:TYR:O	1:B:127:LEU:HD22	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:ASP:OD2	1:A:83:ARG:HB2	2.20	0.41
1:B:9:LEU:HG	1:B:11:ASP:H	1.86	0.41
1:B:195:LEU:O	1:B:197:ASP:N	2.51	0.41
1:B:206:ILE:HD11	1:B:246:ALA:HA	2.02	0.41
1:B:219:ILE:HG23	1:B:232:VAL:HG22	2.03	0.41
1:B:170:PHE:CE2	1:B:172:PRO:HG3	2.56	0.41
1:B:257:VAL:O	1:B:258:ARG:HD3	2.20	0.41
1:A:33:ASN:O	1:A:34:ASP:OD1	2.39	0.40
1:A:79:GLU:HG3	1:A:109:ILE:HG13	2.02	0.40
1:A:123:TRP:CG	1:A:125:MET:HE2	2.56	0.40
1:B:20:VAL:HG22	1:B:267:VAL:HG22	2.02	0.40
1:B:37:LEU:HD13	1:B:37:LEU:HA	1.90	0.40
1:A:55:ASP:OD2	2:A:800:URI:O3'	2.22	0.40
1:B:88:LYS:N	1:B:88:LYS:HD2	2.36	0.40
1:A:106:ASN:HD22	1:A:123:TRP:NE1	2.10	0.40
1:A:108:TYR:HA	1:A:122:THR:O	2.22	0.40
1:B:168:LYS:HB3	1:B:186:PHE:CB	2.51	0.40
1:B:203:ILE:H	1:B:203:ILE:HG13	1.74	0.40
1:A:48:PHE:CD2	1:A:48:PHE:O	2.74	0.40
1:B:23:TYR:O	1:B:24:HIS:HB2	2.21	0.40
1:B:137:MET:HB3	1:B:171:VAL:HG22	2.04	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	247/278 (89%)	224 (91%)	19 (8%)	4 (2%)	7	30
1	B	244/278 (88%)	219 (90%)	20 (8%)	5 (2%)	6	25
All	All	491/556 (88%)	443 (90%)	39 (8%)	9 (2%)	6	28

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	46	ASP
1	A	93	ASP
1	B	46	ASP
1	B	93	ASP
1	B	12	TRP
1	B	114	ARG
1	A	273	HIS
1	B	116	LYS
1	A	87	ASP

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	212/233 (91%)	204 (96%)	8 (4%)	29	60
1	B	209/233 (90%)	202 (97%)	7 (3%)	33	63
All	All	421/466 (90%)	406 (96%)	15 (4%)	31	62

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ASP
1	A	25	THR
1	A	48	PHE
1	A	142	ASN
1	A	176	LEU
1	A	195	LEU
1	A	222	LEU
1	A	248	LEU
1	B	25	THR
1	B	48	PHE
1	B	114	ARG
1	B	176	LEU
1	B	195	LEU
1	B	222	LEU

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Mol	Chain	Res	Type
1	B	248	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	A	24	HIS
1	A	91	ASN
1	A	106	ASN
1	A	107	ASN
1	A	151	ASN
1	A	158	ASN
1	A	180	GLN
1	A	188	ASN
1	A	204	ASN
1	A	211	ASN
1	A	223	ASN
1	A	243	ASN
1	A	256	ASN
1	A	273	HIS
1	B	91	ASN
1	B	106	ASN
1	B	107	ASN
1	B	115	ASN
1	B	120	GLN
1	B	151	ASN
1	B	158	ASN
1	B	180	GLN
1	B	188	ASN
1	B	204	ASN
1	B	223	ASN
1	B	237	HIS
1	B	256	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	$\# Z > 2$
2	URI	A	800	-	18,18,18	0.60	0	26,26,26	1.10	1 (3%)
2	URI	B	801	-	18,18,18	0.66	0	26,26,26	0.97	1 (3%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	URI	A	800	-	-	2/6/22/22	0/2/2/2
2	URI	B	801	-	-	0/6/22/22	0/2/2/2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	800	URI	C2'-C1'-N1	-2.74	105.63	113.25
2	B	801	URI	C2'-C1'-N1	-2.71	105.71	113.25

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	800	URI	O4'-C4'-C5'-O5'
2	A	800	URI	C3'-C4'-C5'-O5'

There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	800	URI	3	0
2	B	801	URI	4	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	251/278 (90%)	0.10	10 (3%) 42 23	28, 41, 68, 90	0
1	B	248/278 (89%)	1.40	67 (27%) 1 1	37, 77, 124, 164	0
All	All	499/556 (89%)	0.75	77 (15%) 5 2	28, 52, 110, 164	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	115	ASN	9.7
1	B	116	LYS	6.7
1	B	75	PRO	6.4
1	B	114	ARG	6.2
1	B	10	SER	5.8
1	B	9	LEU	5.7
1	B	272	PHE	5.7
1	B	13	TRP	5.4
1	B	15	GLN	5.0
1	B	14	HIS	4.8
1	B	44	LYS	4.7
1	A	116	LYS	4.6
1	B	47	TRP	4.4
1	B	120	GLN	4.2
1	B	117	ASP	4.2
1	B	118	GLY	4.0
1	B	48	PHE	3.9
1	B	45	LYS	3.8
1	A	9	LEU	3.7
1	B	39	TYR	3.6
1	B	113	GLY	3.5
1	B	46	ASP	3.4
1	B	108	TYR	3.4
1	B	160	TRP	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	89	LEU	3.3
1	A	115	ASN	3.3
1	B	84	PHE	3.2
1	B	106	ASN	3.2
1	B	92	THR	3.2
1	A	275	HIS	3.2
1	B	12	TRP	3.1
1	A	58	VAL	3.1
1	B	123	TRP	3.1
1	B	80	ILE	3.0
1	B	112	MET	3.0
1	B	42	PHE	3.0
1	B	180	GLN	3.0
1	B	152	TYR	2.9
1	B	125	MET	2.9
1	B	155	ALA	2.9
1	B	35	THR	2.9
1	B	156	ASN	2.8
1	B	162	GLY	2.8
1	B	110	TYR	2.8
1	B	227	TRP	2.7
1	A	274	HIS	2.7
1	A	118	GLY	2.7
1	B	90	THR	2.7
1	B	149	TRP	2.6
1	B	82	PRO	2.5
1	B	271	ASN	2.5
1	B	104	PHE	2.5
1	B	147	TYR	2.5
1	B	158	ASN	2.5
1	B	76	LEU	2.5
1	B	268	VAL	2.5
1	B	121	SER	2.4
1	B	107	ASN	2.4
1	A	75	PRO	2.4
1	B	57	PRO	2.4
1	B	58	VAL	2.4
1	B	151	ASN	2.3
1	B	11	ASP	2.3
1	B	122	THR	2.3
1	B	176	LEU	2.3
1	A	204	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	270	TYR	2.2
1	B	223	ASN	2.2
1	B	105	ALA	2.2
1	B	102	TRP	2.2
1	A	180	GLN	2.2
1	B	153	GLY	2.2
1	B	130	ASP	2.1
1	B	224	TYR	2.1
1	B	78	MET	2.1
1	B	101	GLU	2.1
1	B	41	ALA	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
2	URI	B	801	17/17	0.88	0.11	39,43,50,52	0
2	URI	A	800	17/17	0.96	0.10	12,17,21,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.