Full wwPDB X-ray Structure Validation Report ⓘ

Jan 4, 2024 – 05:57 am GMT

PDB ID	:	4YYY
Title	:	X-ray structure of the thymidine phosphorylase from Salmonella typhimurium in complex with uridine
Authors	:	Balaev, V.V.; Lashkov, A.A.; Gabdulkhakov, A.G.; Betzel, C.; Mikhailov, A.M.
Deposited on	:	2015-03-24
Resolution	:	2.43 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:

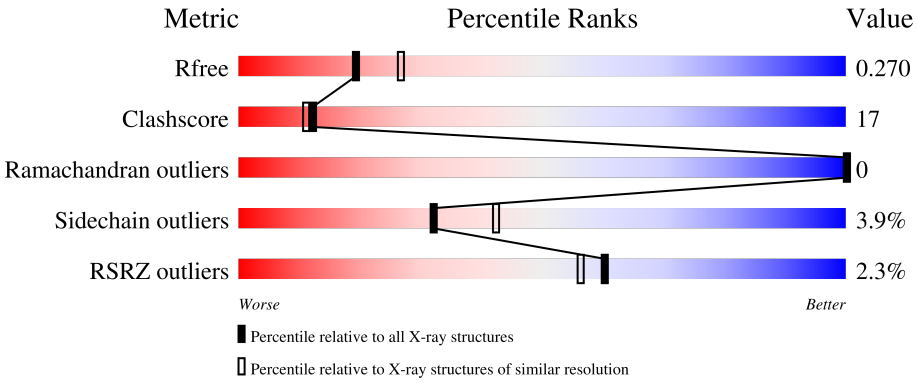
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

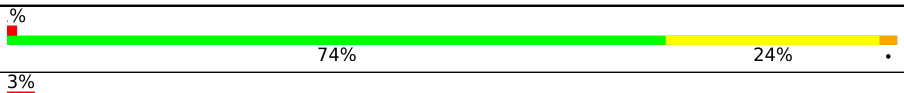
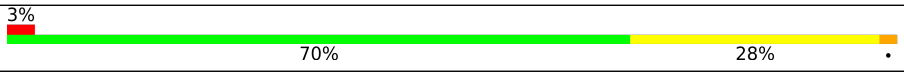
The reported resolution of this entry is 2.43 Å.

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}	130704	1564 (2.46-2.42)
Clashscore	141614	1631 (2.46-2.42)
Ramachandran outliers	138981	1617 (2.46-2.42)
Sidechain outliers	138945	1617 (2.46-2.42)
RSRZ outliers	127900	1547 (2.46-2.42)

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	440	
1	B	440	

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
2	CIT	A	502	-	X	-	-

2 Entry composition (i)

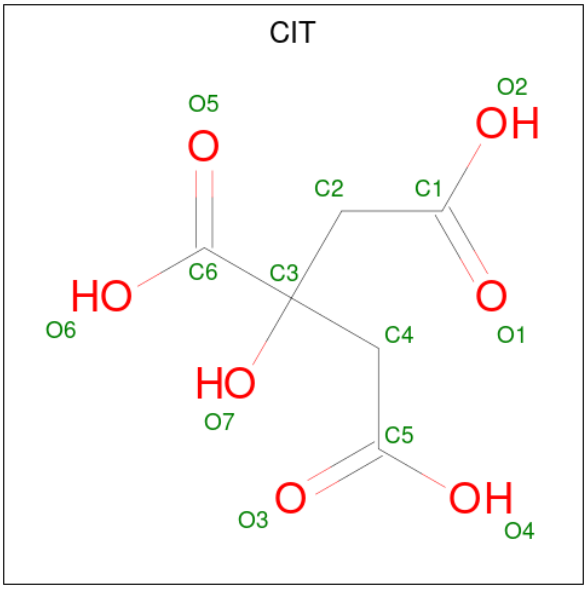
There are 5 unique types of molecules in this entry. The entry contains 6722 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Thymidine phosphorylase.

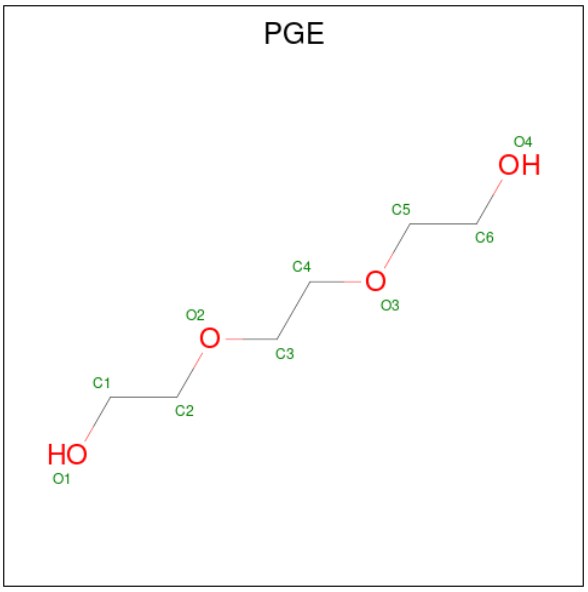
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	440	Total	C	N	O	S	1	2	0
			3305	2070	568	645	22			
1	B	440	Total	C	N	O	S	2	0	0
			3290	2062	566	640	22			

- Molecule 2 is CITRIC ACID (three-letter code: CIT) (formula: $C_6H_8O_7$).



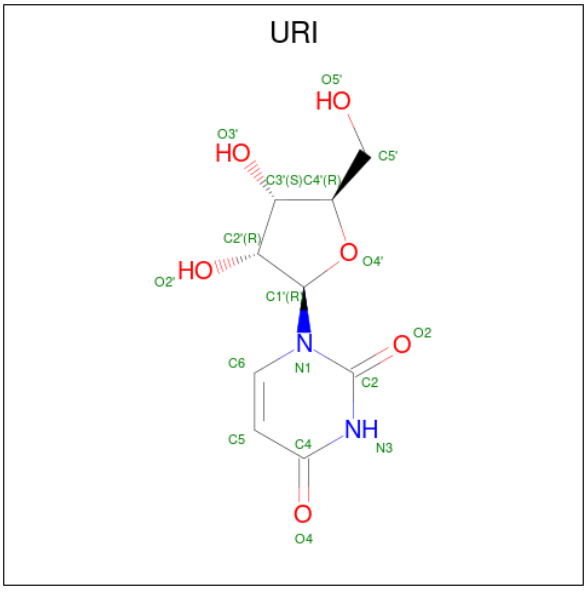
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	6	7		
2	A	1	Total	C	O	0	0
			13	6	7		
2	B	1	Total	C	O	0	0
			13	6	7		

- Molecule 3 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 10 6 4	0	0

- Molecule 4 is URIDINE (three-letter code: URI) (formula: $C_9H_{12}N_2O_6$).



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

- Molecule 5 is water.

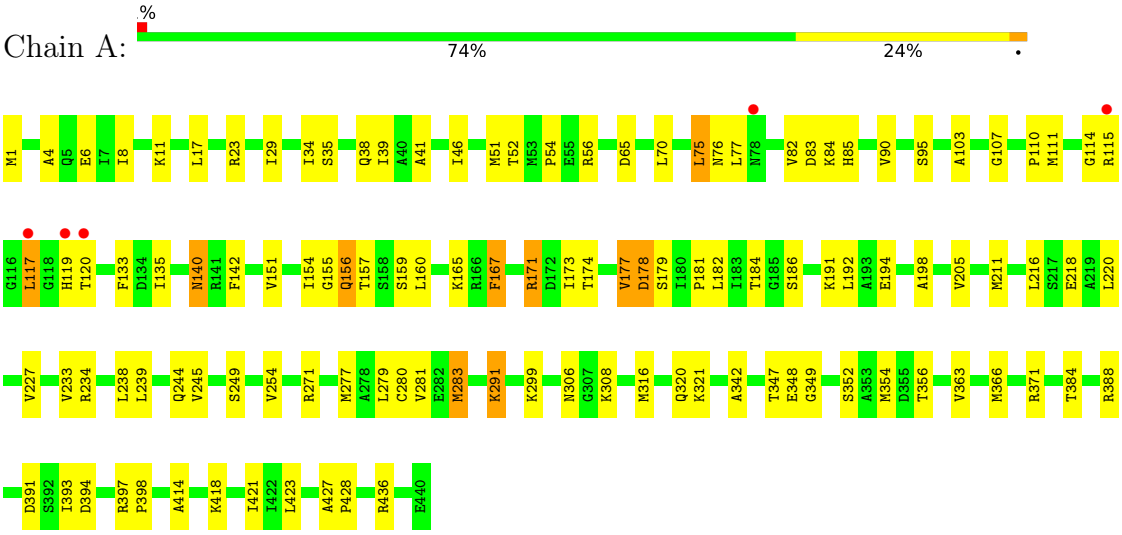
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	29	Total O 29 29	0	0
5	B	15	Total O 15 15	0	0

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.

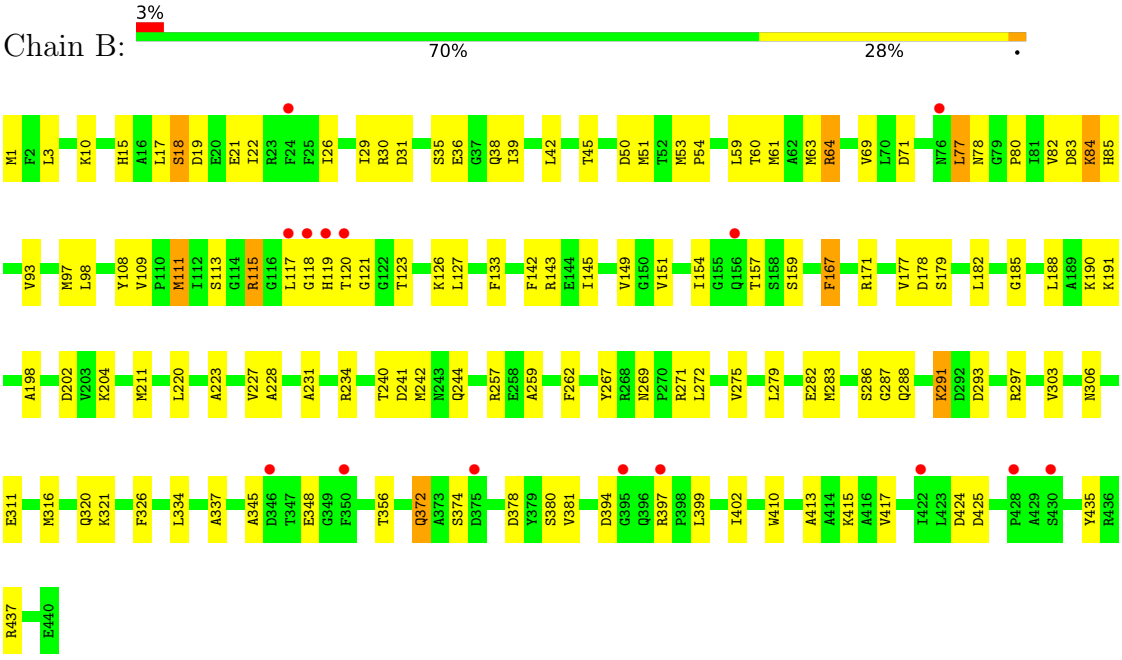
• Molecule 1: Thymidine phosphorylase

Chain A:



• Molecule 1: Thymidine phosphorylase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α, β, γ	193.93Å 193.93Å 57.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.06 – 2.43 48.06 – 2.43	Depositor EDS
% Data completeness (in resolution range)	92.5 (48.06-2.43) 92.8 (48.06-2.43)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.32 (at 2.42Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.204 , 0.272 0.213 , 0.270	Depositor DCC
R_{free} test set	1904 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	54.6	Xtriage
Anisotropy	0.690	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 50.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	0.029 for -k,-h,-l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	6722	wwPDB-VP
Average B, all atoms (Å^2)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.83% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: URI, PGE, CIT

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	0.58	0/3353	0.65	0/4535
1	B	0.53	0/3338	0.60	0/4514
All	All	0.55	0/6691	0.63	0/9049

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	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	171	ARG	Sidechain
1	A	177	VAL	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	3305	0	3316	106	1
1	B	3290	0	3306	118	2
2	A	26	0	10	3	0
2	B	13	0	5	3	0
3	A	10	0	14	1	0
4	B	34	0	24	0	0
5	A	29	0	0	2	0
5	B	15	0	0	0	0
All	All	6722	0	6675	227	2

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 17.

All (227) 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:114:GLY:HA2	1:A:156:GLN:OE1	1.50	1.08
1:A:157:THR:HG22	1:A:159:SER:N	1.67	1.08
1:A:157:THR:HG22	1:A:159:SER:H	0.90	1.06
1:A:4:ALA:HB3	1:A:38:GLN:HE22	1.28	0.97
1:A:4:ALA:HB3	1:A:38:GLN:NE2	1.84	0.93
1:A:178:ASP:OD2	5:A:601:HOH:O	1.90	0.89
1:B:372:GLN:O	1:B:372:GLN:HG2	1.71	0.89
1:B:287:GLY:C	1:B:288:GLN:HG2	1.94	0.88
1:B:242:MET:CE	1:B:356:THR:HG21	2.04	0.88
1:A:244:GLN:HE21	1:A:271:ARG:CZ	1.86	0.87
1:A:171:ARG:NH1	1:A:177:VAL:O	2.08	0.86
1:A:157:THR:HG21	1:A:159:SER:HB3	1.54	0.86
1:A:4:ALA:CB	1:A:38:GLN:NE2	2.39	0.85
1:A:4:ALA:CB	1:A:38:GLN:HE22	1.89	0.84
1:B:17:LEU:O	1:B:21:GLU:HG3	1.81	0.81
1:B:120:THR:HG22	1:B:121:GLY:N	1.95	0.81
1:A:352:SER:HA	1:A:428:PRO:HG3	1.63	0.81
1:A:34:ILE:HG23	1:A:38:GLN:HE21	1.45	0.80
1:A:299:LYS:NZ	5:A:602:HOH:O	2.15	0.79
1:A:157:THR:CG2	1:A:159:SER:HB3	2.13	0.77
1:B:109:VAL:HG12	1:B:111:MET:HG2	1.66	0.77
1:B:291:LYS:HD3	1:B:291:LYS:N	2.00	0.77
1:A:157:THR:CG2	1:A:159:SER:H	1.85	0.76
1:A:133:PHE:CZ	1:A:320:GLN:NE2	2.54	0.75
1:A:117:LEU:HD12	1:A:117:LEU:H	1.52	0.74
1:B:120:THR:HG22	1:B:121:GLY:H	1.52	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:SER:O	1:B:19:ASP:HB3	1.89	0.73
1:A:119:HIS:CD2	1:A:120:THR:H	2.07	0.72
1:B:71:ASP:O	1:B:143:ARG:NH2	2.22	0.72
2:B:503:CIT:O7	2:B:503:CIT:O3	2.07	0.70
1:B:19:ASP:H	1:B:22:ILE:HG13	1.56	0.69
1:A:95:SER:HB3	1:A:111:MET:HE1	1.73	0.69
1:A:177:VAL:HB	1:A:178:ASP:OD1	1.93	0.69
1:B:287:GLY:O	1:B:288:GLN:CG	2.41	0.69
1:A:115:ARG:HB2	1:A:156:GLN:HE22	1.57	0.68
1:B:287:GLY:O	1:B:288:GLN:HG2	1.92	0.68
1:A:179:SER:OG	1:A:182:LEU:HB2	1.93	0.68
1:B:10:LYS:NZ	1:B:21:GLU:OE1	2.27	0.68
1:B:133:PHE:CZ	1:B:320:GLN:NE2	2.62	0.68
1:B:171:ARG:HD3	1:B:182:LEU:HD13	1.76	0.67
1:B:119:HIS:O	1:B:120:THR:HB	1.94	0.67
1:A:245:VAL:HG23	1:A:354:MET:HE1	1.77	0.67
1:A:354:MET:HG2	1:A:421:ILE:HG12	1.77	0.66
1:B:242:MET:HE1	1:B:356:THR:HG21	1.78	0.66
1:A:114:GLY:CA	1:A:156:GLN:OE1	2.37	0.66
1:A:211:MET:HG3	1:A:216:LEU:HB3	1.78	0.66
1:B:287:GLY:C	1:B:288:GLN:CG	2.64	0.66
1:B:282:GLU:OE1	1:B:437:ARG:NH2	2.29	0.65
1:B:17:LEU:O	1:B:21:GLU:CG	2.44	0.65
1:A:117:LEU:C	1:A:117:LEU:HD13	2.16	0.65
1:B:177:VAL:HG23	1:B:178:ASP:H	1.61	0.64
1:A:177:VAL:N	1:A:178:ASP:OD1	2.30	0.63
1:A:178:ASP:OD1	1:A:178:ASP:N	2.28	0.63
1:A:320:GLN:O	1:A:321:LYS:HB2	1.97	0.63
1:B:293:ASP:OD1	1:B:293:ASP:C	2.36	0.63
1:B:117:LEU:N	1:B:117:LEU:HD22	2.12	0.63
1:B:157:THR:HG23	1:B:159:SER:H	1.63	0.62
2:A:501:CIT:O4	2:A:501:CIT:O7	2.15	0.62
1:B:50:ASP:OD1	1:B:51:MET:N	2.31	0.62
1:B:109:VAL:HG12	1:B:109:VAL:O	1.99	0.61
1:A:85:HIS:HB2	1:A:191:LYS:HG2	1.82	0.61
1:A:186:SER:OG	2:A:502:CIT:O3	2.07	0.60
1:A:82:VAL:HG12	1:A:283:MET:HE3	1.83	0.59
1:B:119:HIS:O	1:B:120:THR:CB	2.50	0.59
1:B:17:LEU:HB2	1:B:22:ILE:HD11	1.84	0.58
1:B:83:ASP:OD1	1:B:84:LYS:N	2.32	0.58
1:B:204:LYS:HD2	1:B:242:MET:HG2	1.84	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:17:LEU:O	1:B:18:SER:CB	2.50	0.58
1:A:119:HIS:CD2	1:A:120:THR:N	2.72	0.57
1:B:35:SER:HB3	1:B:38:GLN:HG3	1.86	0.57
1:B:320:GLN:C	1:B:321:LYS:HG2	2.23	0.57
1:B:424:ASP:OD1	1:B:425:ASP:N	2.36	0.57
1:A:117:LEU:C	1:A:117:LEU:CD1	2.73	0.57
1:A:347:THR:OG1	1:A:348:GLU:N	2.38	0.57
1:A:29:ILE:CD1	1:A:39:ILE:HG12	2.34	0.56
1:A:115:ARG:CB	1:A:156:GLN:HE22	2.18	0.56
1:B:118:GLY:O	1:B:119:HIS:HB2	2.06	0.56
1:A:349:GLY:O	1:A:393:ILE:N	2.34	0.56
1:B:61:MET:SD	1:B:64:ARG:NH1	2.78	0.56
1:B:177:VAL:HG23	1:B:178:ASP:N	2.20	0.56
1:A:244:GLN:HE21	1:A:271:ARG:NE	2.02	0.55
1:A:8:ILE:HD13	1:A:41:ALA:HB3	1.88	0.55
1:A:157:THR:CG2	1:A:159:SER:CB	2.84	0.55
1:A:119:HIS:CG	1:A:120:THR:N	2.75	0.55
1:B:241:ASP:OD1	1:B:435:TYR:OH	2.13	0.55
1:A:184:THR:HG21	1:A:227:VAL:HG21	1.89	0.55
1:B:18:SER:O	1:B:19:ASP:CB	2.50	0.55
1:A:133:PHE:CE1	1:A:320:GLN:NE2	2.75	0.54
1:A:192:LEU:HD22	1:A:233:VAL:HG21	1.88	0.54
1:B:188:LEU:HD21	1:B:228:ALA:HB2	1.88	0.54
1:A:82:VAL:HG22	1:A:198:ALA:HB3	1.89	0.53
1:B:26:ILE:HA	1:B:29:ILE:HD12	1.90	0.53
1:B:117:LEU:HD22	1:B:117:LEU:H	1.72	0.53
1:B:17:LEU:HB3	1:B:22:ILE:HG12	1.90	0.53
1:B:97:MET:HE1	1:B:272:LEU:HD21	1.90	0.53
1:B:120:THR:CG2	1:B:121:GLY:H	2.19	0.53
1:B:120:THR:CG2	1:B:121:GLY:N	2.65	0.52
1:A:167:PHE:O	1:A:171:ARG:HG3	2.09	0.52
1:B:109:VAL:CG1	1:B:111:MET:HG2	2.39	0.52
1:A:95:SER:HB3	1:A:111:MET:CE	2.40	0.52
1:A:245:VAL:H	1:A:354:MET:HE1	1.74	0.52
1:B:303:VAL:HA	1:B:306:ASN:OD1	2.10	0.52
1:B:348:GLU:HG2	1:B:394:ASP:HA	1.92	0.51
1:B:149:VAL:HG13	1:B:151:VAL:H	1.76	0.51
1:A:83:ASP:OD1	1:A:84:LYS:N	2.38	0.51
1:A:347:THR:HG23	1:A:423:LEU:HD12	1.93	0.51
1:B:78:ASN:OD1	1:B:108:TYR:OH	2.18	0.51
1:B:244:GLN:HE21	1:B:271:ARG:CZ	2.23	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:363:VAL:HA	1:A:366:MET:HE2	1.93	0.51
1:B:202:ASP:OD1	1:B:240:THR:OG1	2.16	0.51
1:A:244:GLN:NE2	1:A:271:ARG:NE	2.59	0.51
1:B:211:MET:HE2	1:B:220:LEU:HD22	1.92	0.51
1:A:414:ALA:O	1:A:418:LYS:HG3	2.11	0.50
1:A:46:ILE:HG23	1:A:51:MET:HG3	1.93	0.50
1:A:70:LEU:HD12	1:A:194:GLU:HA	1.93	0.50
1:A:11:LYS:HD3	1:A:17:LEU:HG	1.93	0.50
1:A:103:ALA:HA	1:A:107:GLY:O	2.12	0.49
1:A:394:ASP:OD1	1:A:397:ARG:N	2.45	0.49
1:A:388:ARG:N	1:A:391:ASP:OD2	2.40	0.49
1:B:53:MET:HB3	1:B:54:PRO:HD3	1.95	0.49
1:B:293:ASP:OD1	1:B:293:ASP:O	2.30	0.49
1:B:80:PRO:HG3	1:B:287:GLY:O	2.12	0.49
1:A:142:PHE:CD1	1:A:154:ILE:HD13	2.47	0.49
1:A:177:VAL:CB	1:A:178:ASP:OD1	2.61	0.49
1:B:227:VAL:O	1:B:231:ALA:N	2.42	0.49
1:B:402:ILE:HB	1:B:410:TRP:NE1	2.28	0.49
1:A:157:THR:HG22	1:A:159:SER:CB	2.43	0.49
1:A:238:LEU:HD12	1:A:279:LEU:HD23	1.95	0.49
1:B:234:ARG:NH1	1:B:288:GLN:OE1	2.45	0.49
1:A:95:SER:CB	1:A:111:MET:HE1	2.41	0.48
1:B:35:SER:O	1:B:39:ILE:HG13	2.12	0.48
1:B:97:MET:CE	1:B:259:ALA:HA	2.43	0.48
1:B:204:LYS:HD2	1:B:242:MET:CG	2.42	0.48
1:B:179:SER:HB3	1:B:182:LEU:HB2	1.96	0.48
1:B:223:ALA:O	1:B:227:VAL:HG23	2.14	0.48
1:A:157:THR:HG22	1:A:159:SER:CA	2.40	0.48
1:B:326:PHE:HZ	1:B:334:LEU:HD21	1.79	0.48
1:A:254:VAL:HG13	1:A:384[B]:THR:HG22	1.95	0.47
1:B:345:ALA:HB2	1:B:399:LEU:HD11	1.96	0.47
1:B:17:LEU:O	1:B:18:SER:HB2	2.13	0.47
1:B:77:LEU:HD12	1:B:77:LEU:HA	1.67	0.47
1:A:52:THR:HB	1:A:54:PRO:HD2	1.97	0.47
1:B:42:LEU:O	1:B:45:THR:HG22	2.15	0.47
1:B:64:ARG:HE	1:B:69:VAL:HG11	1.79	0.47
1:B:31:ASP:OD1	1:B:31:ASP:N	2.47	0.47
1:A:291:LYS:HD2	1:A:291:LYS:HA	1.60	0.47
3:A:503:PGE:H6	3:A:503:PGE:H42	1.71	0.47
1:B:372:GLN:O	1:B:374:SER:N	2.48	0.47
1:A:115:ARG:N	1:A:156:GLN:OE1	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:113:SER:HA	2:B:503:CIT:O1	2.15	0.47
1:B:97:MET:HE2	1:B:259:ALA:HA	1.97	0.46
1:A:84:LYS:NZ	2:A:501:CIT:O3	2.40	0.46
1:B:142:PHE:CD1	1:B:154:ILE:HD13	2.51	0.46
1:A:277:MET:O	1:A:281:VAL:HG23	2.16	0.46
1:B:190:LYS:H	1:B:190:LYS:HG3	1.50	0.46
1:B:337:ALA:HB2	1:B:380:SER:O	2.16	0.45
1:A:75:LEU:HD22	1:A:76:ASN:N	2.32	0.45
1:A:306:ASN:OD1	1:A:308:LYS:HB2	2.17	0.45
1:B:60:THR:OG1	1:B:185:GLY:HA2	2.17	0.45
1:B:10:LYS:O	1:B:15:HIS:HB2	2.16	0.45
1:B:198:ALA:HB1	1:B:286:SER:HB3	1.99	0.45
1:A:135:ILE:O	1:A:155:GLY:N	2.49	0.44
1:A:110:PRO:O	1:A:110:PRO:HG2	2.17	0.44
1:A:363:VAL:HA	1:A:366:MET:CE	2.46	0.44
1:B:118:GLY:O	1:B:119:HIS:CB	2.64	0.44
1:B:123:THR:HA	1:B:126:LYS:HD2	1.98	0.44
1:A:244:GLN:HA	1:A:354:MET:HE2	1.97	0.44
1:A:280:CYS:HA	1:A:283:MET:HG3	1.99	0.44
1:A:342:ALA:HB1	1:A:398:PRO:HB3	1.99	0.44
1:B:145:ILE:O	1:B:149:VAL:HG12	2.17	0.44
1:B:269:ASN:HB3	1:B:272:LEU:HB3	1.98	0.44
1:A:84:LYS:HD2	1:A:279:LEU:HD13	2.00	0.44
1:B:30:ARG:NE	1:B:31:ASP:OD1	2.48	0.44
1:B:115:ARG:HD3	1:B:115:ARG:HA	1.30	0.44
1:A:157:THR:HG22	1:A:159:SER:HB3	1.98	0.44
1:B:287:GLY:O	1:B:288:GLN:HG3	2.17	0.44
1:B:111:MET:HE3	1:B:111:MET:HB3	1.59	0.44
1:B:167:PHE:O	1:B:171:ARG:HG3	2.17	0.44
1:A:140:ASN:OD1	1:A:140:ASN:N	2.52	0.43
1:B:77:LEU:HG	1:B:108:TYR:CZ	2.54	0.43
1:A:29:ILE:HD11	1:A:39:ILE:HG12	2.00	0.43
1:A:90:VAL:HG21	1:A:356:THR:O	2.17	0.43
1:A:142:PHE:CE1	1:A:154:ILE:HD13	2.53	0.43
1:B:151:VAL:O	1:B:316:MET:HG3	2.18	0.43
1:A:11:LYS:HB2	1:A:17:LEU:HD21	2.00	0.43
1:B:293:ASP:OD1	1:B:297:ARG:HG3	2.19	0.43
1:B:191:LYS:HA	1:B:191:LYS:HD3	1.85	0.43
1:B:97:MET:HG2	1:B:262:PHE:CD1	2.54	0.43
1:B:240:THR:HG21	1:B:275:VAL:HG13	2.01	0.43
1:B:85:HIS:HD2	2:B:503:CIT:C1	2.32	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:59:LEU:O	1:B:63:MET:HG3	2.19	0.42
1:B:142:PHE:CE1	1:B:154:ILE:HD13	2.54	0.42
1:A:191:LYS:HD2	1:A:191:LYS:HA	1.90	0.42
1:A:211:MET:HE2	1:A:220:LEU:HD22	2.01	0.42
1:A:321:LYS:N	1:A:321:LYS:CD	2.82	0.42
1:B:394:ASP:OD1	1:B:397:ARG:HD3	2.19	0.42
1:B:93:VAL:HG22	1:B:93:VAL:O	2.19	0.42
1:B:117:LEU:HA	1:B:117:LEU:HD13	1.59	0.42
1:A:427:ALA:HA	1:A:428:PRO:HD2	1.85	0.42
1:B:378:ASP:O	1:B:381:VAL:HG22	2.19	0.42
1:B:262:PHE:HA	1:B:267:TYR:O	2.20	0.42
1:A:211:MET:CE	1:A:220:LEU:HD22	2.50	0.42
1:B:127:LEU:HD13	1:B:133:PHE:CZ	2.55	0.42
1:A:110:PRO:HB2	1:A:142:PHE:HZ	1.84	0.42
1:A:151:VAL:O	1:A:316:MET:HG3	2.20	0.42
1:A:173:ILE:HG13	1:A:174:THR:HG23	2.02	0.42
1:B:98:LEU:HD21	1:B:283:MET:SD	2.59	0.42
1:B:257:ARG:HE	1:B:257:ARG:HB2	1.42	0.42
1:A:23:ARG:NH2	1:A:65:ASP:OD2	2.53	0.41
1:A:218:GLU:OE1	1:A:436:ARG:NH1	2.53	0.41
1:B:133:PHE:CE1	1:B:320:GLN:NE2	2.87	0.41
1:B:157:THR:HG23	1:B:159:SER:N	2.31	0.41
1:A:249:SER:HB2	1:A:254:VAL:CG1	2.50	0.41
1:A:205:VAL:CG2	1:A:239:LEU:HB3	2.50	0.41
1:A:244:GLN:NE2	1:A:271:ARG:CD	2.84	0.41
1:A:244:GLN:NE2	1:A:271:ARG:CZ	2.67	0.41
1:B:337:ALA:HA	1:B:380:SER:OG	2.21	0.41
1:A:321:LYS:N	1:A:321:LYS:HD3	2.36	0.41
1:B:117:LEU:H	1:B:117:LEU:CD2	2.33	0.41
1:B:202:ASP:HB2	1:B:279:LEU:HD21	2.03	0.41
1:B:241:ASP:HB2	1:B:435:TYR:HE1	1.86	0.41
1:B:415:LYS:HD3	1:B:415:LYS:HA	1.65	0.41
1:A:56:ARG:HE	1:A:181:PRO:HB3	1.86	0.41
1:B:244:GLN:NE2	1:B:271:ARG:CZ	2.84	0.40
1:B:413:ALA:O	1:B:417:VAL:HG23	2.22	0.40
1:A:245:VAL:HG23	1:A:354:MET:CE	2.47	0.40
1:B:17:LEU:CB	1:B:22:ILE:HG12	2.51	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:SER:OG	1:B:311:GLU:OE2[7_444]	2.03	0.17
1:A:35:SER:OG	1:B:36:GLU:OE2[1_556]	2.17	0.03

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	440/440 (100%)	421 (96%)	19 (4%)	0	100	100
1	B	438/440 (100%)	420 (96%)	18 (4%)	0	100	100
All	All	878/880 (100%)	841 (96%)	37 (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	346/344 (101%)	331 (96%)	15 (4%)	29	38
1	B	344/344 (100%)	332 (96%)	12 (4%)	36	47
All	All	690/688 (100%)	663 (96%)	27 (4%)	32	42

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	6	GLU

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Mol	Chain	Res	Type
1	A	75	LEU
1	A	77	LEU
1	A	117	LEU
1	A	140	ASN
1	A	156	GLN
1	A	160	LEU
1	A	165	LYS
1	A	167	PHE
1	A	178	ASP
1	A	234	ARG
1	A	283	MET
1	A	291	LYS
1	A	371	ARG
1	B	1	MET
1	B	3	LEU
1	B	18	SER
1	B	64	ARG
1	B	77	LEU
1	B	82	VAL
1	B	84	LYS
1	B	111	MET
1	B	115	ARG
1	B	167	PHE
1	B	291	LYS
1	B	372	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	ASN
1	A	38	GLN
1	A	119	HIS
1	A	244	GLN
1	A	320	GLN
1	B	32	ASN
1	B	119	HIS
1	B	320	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

6 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	CIT	A	501	-	12,12,12	1.09	0	17,17,17	1.86	4 (23%)
4	URI	B	502	-	18,18,18	0.35	0	26,26,26	0.69	0
4	URI	B	501	-	18,18,18	0.31	0	26,26,26	0.51	0
3	PGE	A	503	-	9,9,9	0.32	0	8,8,8	0.54	0
2	CIT	B	503	-	12,12,12	1.13	0	17,17,17	1.86	6 (35%)
2	CIT	A	502	-	12,12,12	1.66	3 (25%)	17,17,17	1.91	5 (29%)

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	CIT	A	501	-	-	12/16/16/16	-
4	URI	B	502	-	-	2/6/22/22	0/2/2/2
4	URI	B	501	-	-	6/6/22/22	0/2/2/2
3	PGE	A	503	-	-	4/7/7/7	-
2	CIT	B	503	-	-	9/16/16/16	-
2	CIT	A	502	-	-	11/16/16/16	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	502	CIT	C3-C6	-3.64	1.49	1.53
2	A	502	CIT	C4-C3	-2.06	1.51	1.53
2	A	502	CIT	O4-C5	-2.01	1.24	1.30

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	502	CIT	C3-C4-C5	-4.67	102.51	113.81
2	B	503	CIT	C3-C4-C5	-4.57	102.74	113.81
2	A	501	CIT	O6-C6-C3	4.54	120.93	113.05
2	A	502	CIT	O6-C6-C3	3.06	118.37	113.05
2	A	501	CIT	C3-C4-C5	-3.02	106.50	113.81
2	B	503	CIT	C4-C3-C6	-2.82	104.05	110.11
2	B	503	CIT	O6-C6-C3	2.65	117.66	113.05
2	A	502	CIT	O2-C1-C2	2.25	121.59	114.35
2	A	502	CIT	C4-C3-C2	2.25	115.02	109.16
2	A	501	CIT	O5-C6-C3	-2.24	119.09	122.25
2	B	503	CIT	O2-C1-C2	2.22	121.48	114.35
2	B	503	CIT	O2-C1-O1	-2.21	117.79	123.30
2	A	501	CIT	O2-C1-C2	2.15	121.26	114.35
2	B	503	CIT	C3-C2-C1	-2.04	108.87	113.81
2	A	502	CIT	O7-C3-C4	-2.02	104.68	109.40

There are no chirality outliers.

All (44) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	CIT	C1-C2-C3-O7
2	A	501	CIT	C1-C2-C3-C4
2	A	501	CIT	O7-C3-C6-O5
2	A	501	CIT	O7-C3-C6-O6
2	A	501	CIT	C4-C3-C6-O5
2	A	501	CIT	C4-C3-C6-O6
2	A	502	CIT	C1-C2-C3-O7
2	A	502	CIT	C1-C2-C3-C4
2	A	502	CIT	C1-C2-C3-C6
4	B	501	URI	C2'-C1'-N1-C2
4	B	501	URI	C2'-C1'-N1-C6
3	A	503	PGE	C6-C5-O3-C4
3	A	503	PGE	O2-C3-C4-O3
2	A	501	CIT	C1-C2-C3-C6

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Mol	Chain	Res	Type	Atoms
2	A	501	CIT	C6-C3-C4-C5
2	A	502	CIT	O7-C3-C6-O5
2	A	502	CIT	O7-C3-C6-O6
2	B	503	CIT	O7-C3-C6-O5
2	A	502	CIT	C2-C3-C6-O5
2	A	502	CIT	C2-C3-C6-O6
2	B	503	CIT	C2-C3-C6-O5
2	B	503	CIT	C2-C3-C6-O6
2	A	501	CIT	C2-C3-C4-C5
4	B	502	URI	O4'-C4'-C5'-O5'
4	B	501	URI	O4'-C4'-C5'-O5'
4	B	501	URI	C3'-C4'-C5'-O5'
2	B	503	CIT	C4-C3-C6-O6
4	B	501	URI	O4'-C1'-N1-C6
4	B	501	URI	O4'-C1'-N1-C2
3	A	503	PGE	C1-C2-O2-C3
2	B	503	CIT	O7-C3-C6-O6
2	A	502	CIT	C4-C3-C6-O6
2	B	503	CIT	C4-C3-C6-O5
2	B	503	CIT	O2-C1-C2-C3
2	B	503	CIT	O1-C1-C2-C3
2	A	501	CIT	O7-C3-C4-C5
2	A	501	CIT	C3-C4-C5-O3
2	A	502	CIT	O1-C1-C2-C3
3	A	503	PGE	C3-C4-O3-C5
2	A	502	CIT	O2-C1-C2-C3
4	B	502	URI	C3'-C4'-C5'-O5'
2	A	501	CIT	C3-C4-C5-O4
2	A	502	CIT	O7-C3-C4-C5
2	B	503	CIT	O7-C3-C4-C5

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	CIT	2	0
3	A	503	PGE	1	0
2	B	503	CIT	3	0
2	A	502	CIT	1	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

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	440/440 (100%)	-0.21	5 (1%)	80	79	44, 63, 90, 141	1 (0%)
1	B	440/440 (100%)	0.05	15 (3%)	45	42	60, 79, 115, 132	1 (0%)
All	All	880/880 (100%)	-0.08	20 (2%)	60	56	44, 71, 108, 141	2 (0%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	120	THR	4.6
1	B	76	ASN	4.0
1	B	24	PHE	3.6
1	B	120	THR	3.2
1	B	118	GLY	3.1
1	B	430	SER	2.9
1	B	422	ILE	2.9
1	B	117	LEU	2.9
1	B	350	PHE	2.7
1	B	397	ARG	2.6
1	B	119	HIS	2.4
1	B	346	ASP	2.3
1	B	428	PRO	2.3
1	B	375	ASP	2.2
1	B	395	GLY	2.2
1	A	115	ARG	2.1
1	A	119	HIS	2.1
1	A	117	LEU	2.1
1	B	156	GLN	2.1
1	A	78	ASN	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 monosaccharides 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, 95^{th} 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	PGE	A	503	10/10	0.76	0.26	57,73,78,79	10
4	URI	B	502	17/17	0.84	0.18	81,87,92,96	17
2	CIT	A	501	13/13	0.86	0.18	60,66,71,71	13
2	CIT	A	502	13/13	0.86	0.17	75,80,85,86	13
4	URI	B	501	17/17	0.90	0.15	74,87,100,101	0
2	CIT	B	503	13/13	0.90	0.15	73,81,85,85	0

6.5 Other polymers (i)

There are no such residues in this entry.