



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:30 AM UTC

PDB ID : 2HRD / pdb_00002hrd
Title : Crystal structure of the uridine phosphorylase from Salmonella typhimurium in complex with thymine and phosphate ion at 1.70Å resolution
Authors : Timofeev, V.I.; Gabdulkhakov, A.G.; Mikhailov, A.M.
Deposited on : 2006-07-20
Resolution : 1.70 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

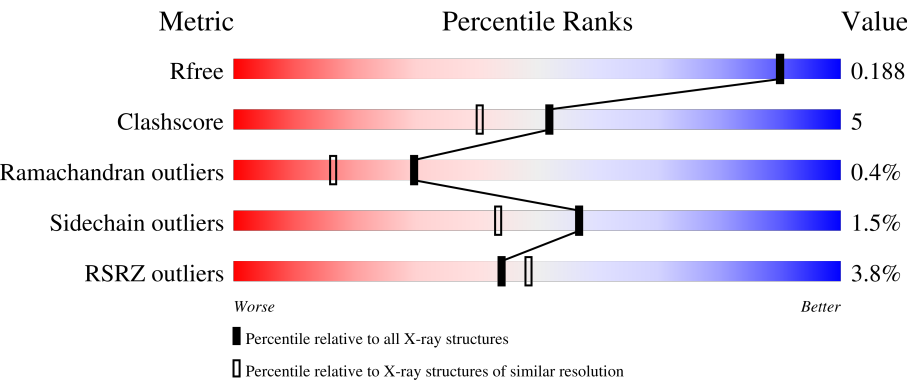
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	253	6% 86%12%..
1	B	253	2% 87%10%. .
1	C	253	4% 89%8%. .
1	D	253	3% 88%8%. .
1	E	253	4% 88%11%. .

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	253	

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	TDR	A	901[A]	-	X	-	-
3	TDR	A	901[B]	-	X	-	-
3	TDR	B	902	-	X	-	-
3	TDR	C	903[A]	-	X	-	-
3	TDR	C	903[B]	-	X	-	-
3	TDR	D	904[A]	-	X	-	-
3	TDR	D	904[B]	-	X	-	-
3	TDR	E	905[A]	-	X	-	-
3	TDR	E	905[B]	-	X	-	-
3	TDR	F	906[A]	-	X	-	-
3	TDR	F	906[B]	-	X	X	-
5	1PE	B	1001	-	-	-	X

2 Entry composition [i](#)

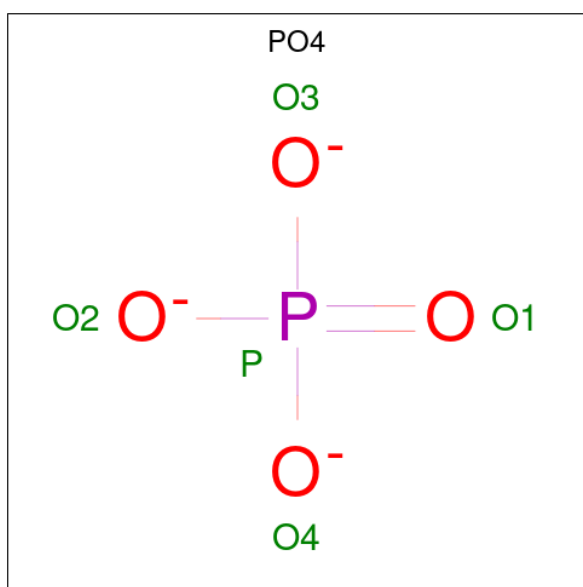
There are 6 unique types of molecules in this entry. The entry contains 12584 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Uridine phosphorylase.

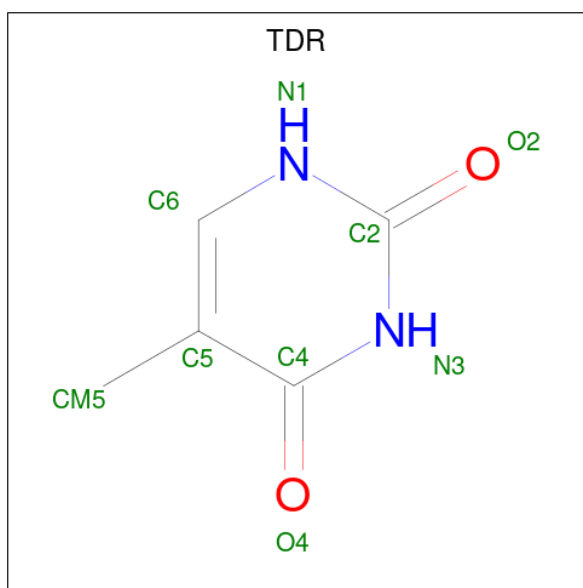
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	251	Total	C	N	O	S	0	16	0
			1973	1244	342	375	12			
1	B	246	Total	C	N	O	S	0	18	0
			1939	1223	337	368	11			
1	C	245	Total	C	N	O	S	0	17	0
			1947	1233	342	361	11			
1	D	244	Total	C	N	O	S	0	18	0
			1937	1219	335	372	11			
1	E	252	Total	C	N	O	S	0	18	0
			1991	1256	343	380	12			
1	F	245	Total	C	N	O	S	0	11	0
			1893	1195	328	359	11			

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O P 5 4 1	0	0
2	A	1	Total O P 10 8 2	0	1
2	B	1	Total O P 5 4 1	0	0
2	C	1	Total O P 5 4 1	0	0
2	D	1	Total O P 5 4 1	0	0
2	E	1	Total O P 5 4 1	0	0
2	F	1	Total O P 5 4 1	0	0
2	F	1	Total O P 10 8 2	0	1

- Molecule 3 is THYMINE (CCD ID: TDR) (formula: C₅H₆N₂O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O 18 10 4 4	0	1
3	B	1	Total C N O 9 5 2 2	0	0
3	C	1	Total C N O 18 10 4 4	0	1
3	D	1	Total C N O 18 10 4 4	0	1

Continued on next page...

Continued from previous page...

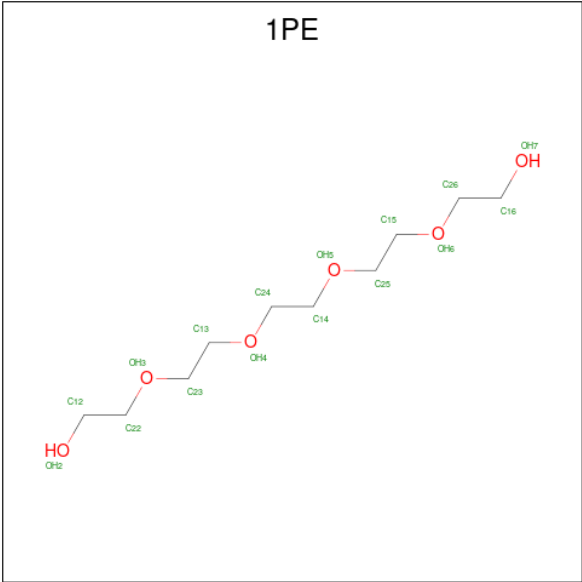
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	E	1	Total	C	N	O	0	1
			18	10	4	4		
3	F	1	Total	C	N	O	0	1
			18	10	4	4		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	E	1	Total	C	O	0	0
			6	3	3		
4	F	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: $C_{10}H_{22}O_6$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	C	O	0
			16	10	6	

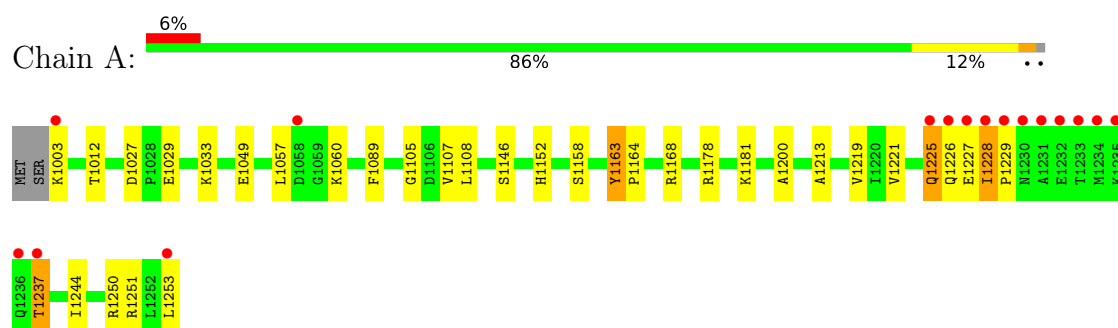
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	126	Total	O	0	1
			127	127		
6	B	136	Total	O	0	0
			136	136		
6	C	91	Total	O	0	0
			91	91		
6	D	111	Total	O	0	1
			112	112		
6	E	123	Total	O	0	2
			125	125		
6	F	118	Total	O	0	0
			118	118		

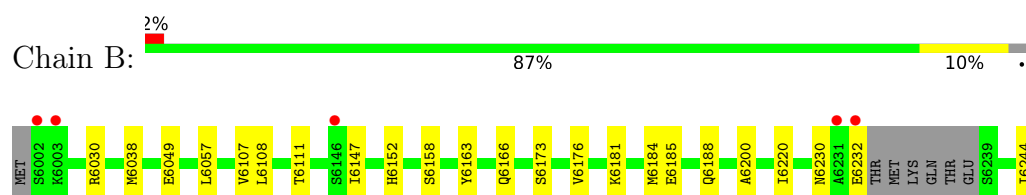
3 Residue-property plots

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.

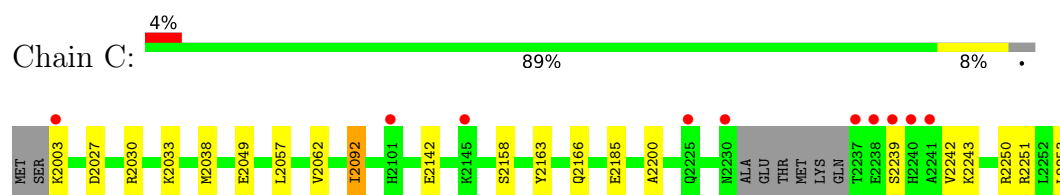
- Molecule 1: Uridine phosphorylase



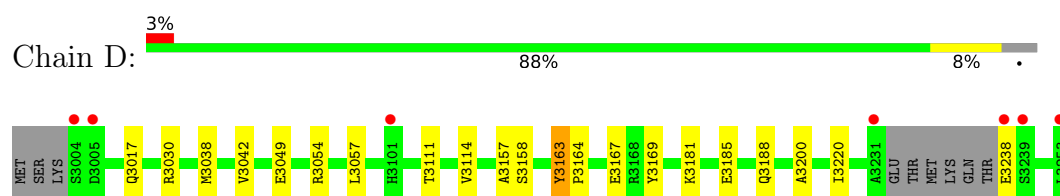
- Molecule 1: Uridine phosphorylase



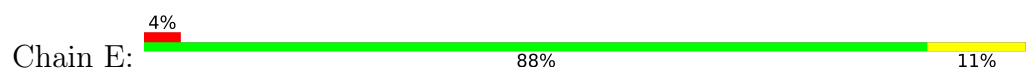
- Molecule 1: Uridine phosphorylase

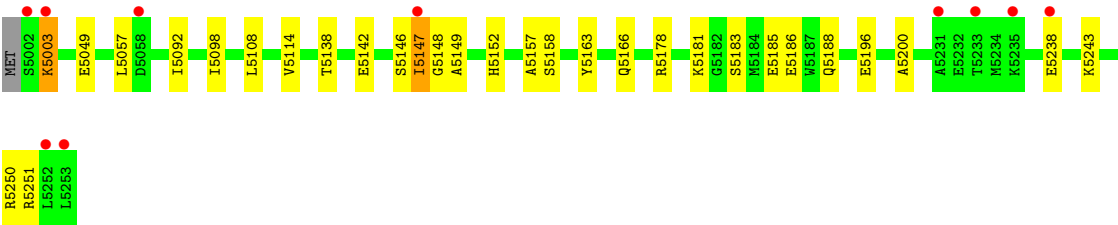


- Molecule 1: Uridine phosphorylase

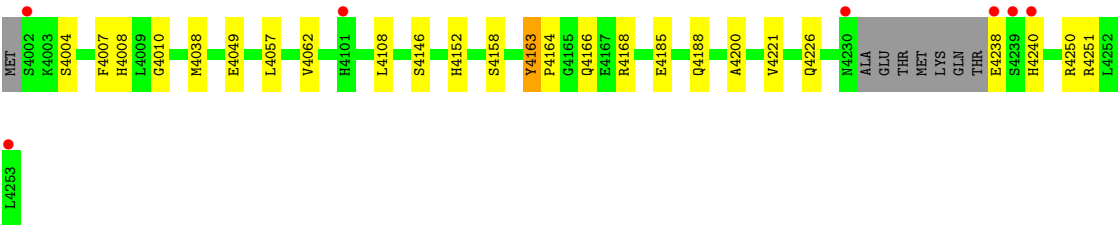
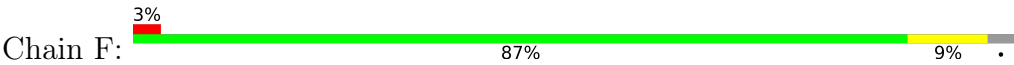


- Molecule 1: Uridine phosphorylase





● Molecule 1: Uridine phosphorylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	88.20Å 123.40Å 133.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.70 30.00 – 1.70	Depositor EDS
% Data completeness (in resolution range)	99.9 (30.00-1.70) 99.9 (30.00-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.17 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.166 , 0.189 0.165 , 0.188	Depositor DCC
R_{free} test set	7980 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	7.7	Xtriage
Anisotropy	0.190	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 37.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12584	wwPDB-VP
Average B, all atoms (Å ²)	8.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, GOL, 1PE, TDR

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	0.40	0/2051	0.72	0/2772
1	B	0.39	0/2019	0.70	0/2728
1	C	0.40	0/2023	0.72	0/2736
1	D	0.40	0/2018	0.72	0/2731
1	E	0.40	0/2066	0.73	0/2793
1	F	0.41	0/1953	0.73	0/2644
All	All	0.40	0/12130	0.72	0/16404

There are no bond length outliers.

There are no bond angle outliers.

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	1973	0	2035	36	0
1	B	1939	0	1999	17	0
1	C	1947	0	2014	15	0
1	D	1937	0	1971	11	0
1	E	1991	0	2054	26	0
1	F	1893	0	1926	18	0
2	A	15	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	5	0	0	1	0
2	C	5	0	0	0	0
2	D	5	0	0	1	0
2	E	5	0	0	0	0
2	F	15	0	0	0	0
3	A	18	0	12	0	0
3	B	9	0	6	3	0
3	C	18	0	12	1	0
3	D	18	0	12	1	0
3	E	18	0	12	2	0
3	F	18	0	12	5	0
4	A	6	0	8	0	0
4	B	6	0	8	1	0
4	C	6	0	8	1	0
4	E	6	0	8	0	0
4	F	6	0	8	1	0
5	B	16	0	22	1	0
6	A	127	0	0	0	0
6	B	136	0	0	0	0
6	C	91	0	0	1	0
6	D	112	0	0	0	0
6	E	125	0	0	1	0
6	F	118	0	0	3	0
All	All	12584	0	12127	122	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 5.

All (122) 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:E:5147[A]:ILE:CD1	1:E:5243[A]:LYS:HD2	1.83	1.08
1:E:5147[A]:ILE:HD12	1:E:5243[A]:LYS:HD2	1.36	1.06
1:E:5098[B]:ILE:HG22	1:E:5188[B]:GLN:HG2	1.32	1.06
1:A:1057[A]:LEU:HD21	1:A:1250[A]:ARG:CG	1.86	1.05
1:E:5098[B]:ILE:HG22	1:E:5188[B]:GLN:CG	1.93	0.98
1:E:5057[B]:LEU:HD21	1:E:5250[B]:ARG:HB2	1.47	0.97
1:A:1221[B]:VAL:CG1	1:A:1229:PRO:HA	1.94	0.96
1:A:1057[A]:LEU:CD2	1:A:1250[A]:ARG:HG2	1.98	0.93
1:A:1057[A]:LEU:HD21	1:A:1250[A]:ARG:HG3	1.51	0.92
1:E:5098[B]:ILE:CG2	1:E:5188[B]:GLN:HG3	2.01	0.91

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1057[A]:LEU:HD21	1:A:1250[A]:ARG:HG2	1.51	0.90
1:A:1221[B]:VAL:HG11	1:A:1229:PRO:HA	1.52	0.90
1:E:5098[B]:ILE:CG2	1:E:5188[B]:GLN:CG	2.50	0.90
1:A:1057[A]:LEU:CG	1:A:1250[A]:ARG:HG2	2.01	0.89
1:E:5057[B]:LEU:HD23	1:E:5250[B]:ARG:HA	1.63	0.78
1:E:5147[A]:ILE:HD13	1:E:5243[A]:LYS:HD2	1.64	0.78
1:A:1057[A]:LEU:HD11	1:A:1250[A]:ARG:NE	1.98	0.77
1:A:1057[A]:LEU:HG	1:A:1250[A]:ARG:HG2	1.68	0.76
1:E:5057[B]:LEU:CD2	1:E:5250[B]:ARG:HB2	2.17	0.75
1:A:1178:ARG:HA	1:A:1181:LYS:HE2	1.71	0.73
1:F:4168:ARG:HH12	3:F:906[B]:TDR:HN3	1.40	0.68
1:F:4166:GLN:NE2	3:F:906[B]:TDR:O4	2.26	0.67
1:A:1221[B]:VAL:HG13	1:A:1229:PRO:HA	1.75	0.67
1:F:4038:MET:HG2	1:F:4057:LEU:HD13	1.76	0.67
1:E:5057[B]:LEU:HD23	1:E:5250[B]:ARG:CA	2.26	0.66
1:A:1221[B]:VAL:HG11	1:A:1229:PRO:CA	2.25	0.64
1:E:5183:SER:O	1:E:5186[B]:GLU:HG2	1.96	0.64
1:F:4057:LEU:HD11	1:F:4250[B]:ARG:HE	1.61	0.64
1:B:6185[B]:GLU:CD	1:B:6188[B]:GLN:OE1	2.41	0.64
1:E:5158:SER:HB3	1:E:5200:ALA:HB2	1.79	0.64
1:E:5147[A]:ILE:CD1	1:E:5243[A]:LYS:CD	2.71	0.63
1:F:4038:MET:SD	1:F:4062[B]:VAL:HG21	2.38	0.63
1:A:1057[A]:LEU:HD11	1:A:1250[A]:ARG:HE	1.63	0.63
1:A:1158:SER:HB3	1:A:1200:ALA:HB2	1.83	0.61
1:A:1060[A]:LYS:HD2	1:A:1253:LEU:HB3	1.83	0.60
1:E:5147[A]:ILE:HD13	1:E:5243[A]:LYS:CD	2.30	0.60
1:A:1146[A]:SER:OG	1:A:1251:ARG:NH2	2.35	0.59
1:A:1029[B]:GLU:OE2	1:A:1033[B]:LYS:NZ	2.33	0.59
1:B:6158:SER:HB3	1:B:6200:ALA:HB2	1.83	0.59
1:C:2158:SER:HB3	1:C:2200:ALA:HB2	1.85	0.59
1:A:1057[A]:LEU:CD2	1:A:1250[A]:ARG:CG	2.62	0.58
1:F:4158:SER:HB3	1:F:4200:ALA:HB2	1.86	0.57
4:C:1013:GOL:H2	1:D:3111:THR:HB	1.87	0.56
1:A:1226:GLN:HE21	1:A:1228:ILE:HG13	1.70	0.56
1:B:6038:MET:HG2	1:B:6057:LEU:HD13	1.88	0.55
1:D:3042[B]:VAL:HG11	1:D:3054:ARG:NH2	2.22	0.55
1:D:3158:SER:HB3	1:D:3200:ALA:HB2	1.87	0.55
1:E:5003:LYS:H	1:E:5003:LYS:HD2	1.72	0.55
1:C:2142:GLU:OE2	1:C:2251[B]:ARG:NH1	2.38	0.54
1:E:5178:ARG:HA	1:E:5181[B]:LYS:HE2	1.89	0.54
1:B:6166:GLN:HE22	3:B:902:TDR:HN3	1.55	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:2166:GLN:HE22	3:C:903[A]:TDR:HN3	1.56	0.54
1:E:5148[A]:GLY:O	1:E:5149[A]:ALA:HB2	2.08	0.54
1:C:2030[B]:ARG:HD2	1:C:2033:LYS:HD2	1.90	0.53
1:A:1049:GLU:HB3	1:F:4049:GLU:HB3	1.90	0.53
1:C:2027:ASP:HB3	1:C:2030[A]:ARG:HG3	1.91	0.53
1:B:6107:VAL:HG21	1:B:6244:ILE:HD12	1.91	0.52
1:C:2185[B]:GLU:HG3	6:C:99:HOH:O	2.09	0.52
1:A:1027:ASP:CG	1:A:1029[B]:GLU:HG3	2.35	0.52
1:C:2038:MET:HG2	1:C:2057[A]:LEU:HD13	1.92	0.51
1:B:6049:GLU:HB3	1:D:3049:GLU:HB3	1.93	0.50
1:B:6184:MET:HB3	5:B:1001:1PE:H162	1.92	0.50
1:C:2239:SER:HB3	1:C:2243:LYS:HE3	1.94	0.50
1:F:4166:GLN:HE22	3:F:906[A]:TDR:HN3	1.60	0.50
1:C:2092[A]:ILE:HD12	1:C:2242:VAL:HA	1.93	0.50
1:B:6111:THR:HB	4:F:1011:GOL:H2	1.94	0.49
1:E:5166:GLN:HE22	3:E:905[A]:TDR:HN3	1.61	0.49
1:C:2038:MET:SD	1:C:2062[B]:VAL:HG21	2.52	0.49
1:C:2057[B]:LEU:HD22	1:C:2250:ARG:HG3	1.94	0.49
1:B:6185[B]:GLU:HA	1:B:6188[B]:GLN:OE1	2.13	0.48
1:F:4146[B]:SER:OG	1:F:4251:ARG:NH2	2.46	0.48
1:B:6057:LEU:HG	1:B:6250:ARG:HG2	1.95	0.48
1:B:6166:GLN:OE1	3:B:902:TDR:O4	2.31	0.48
3:F:906[B]:TDR:CM5	6:F:197:HOH:O	2.62	0.48
1:F:4007[A]:PHE:HD2	1:F:4008:HIS:CE1	2.32	0.47
1:A:1228:ILE:HG13	1:A:1228:ILE:O	2.14	0.47
1:B:6030[A]:ARG:HH22	2:B:802:PO4:P	2.38	0.47
3:E:905[B]:TDR:CM5	6:E:243:HOH:O	2.62	0.47
1:C:2092[A]:ILE:CD1	1:C:2242:VAL:HG22	2.45	0.47
1:F:4108:LEU:HD22	1:F:4152:HIS:HB2	1.96	0.47
1:A:1029[B]:GLU:OE2	1:A:1033[B]:LYS:HE3	2.15	0.46
1:B:6176:VAL:O	1:B:6181[A]:LYS:HE2	2.15	0.46
1:F:4221:VAL:CG2	1:F:4226:GLN:HG3	2.45	0.46
1:E:5108:LEU:HD22	1:E:5152:HIS:HB2	1.98	0.45
1:D:3038:MET:HG2	1:D:3057:LEU:HD13	1.99	0.45
1:A:1219:VAL:O	1:A:1237:THR:HG21	2.16	0.45
1:A:1168:ARG:HG2	1:A:1227:GLU:HG2	1.99	0.44
1:E:5146[B]:SER:OG	1:E:5251:ARG:NH2	2.51	0.44
1:A:1029[B]:GLU:OE2	1:A:1033[B]:LYS:CE	2.66	0.44
1:F:4185:GLU:HG3	6:F:33:HOH:O	2.18	0.44
1:A:1003:LYS:HE2	1:A:1012:THR:HG21	1.99	0.44
1:D:3030[A]:ARG:HH22	2:D:804:PO4:P	2.41	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:4108:LEU:CD2	1:F:4152:HIS:HB2	2.48	0.44
1:D:3114:VAL:HB	1:D:3157:ALA:HA	1.99	0.43
1:B:6057:LEU:HB3	1:B:6253:LEU:HD11	2.01	0.43
1:A:1163:TYR:HB2	1:A:1164:PRO:CD	2.49	0.43
1:A:1226:GLN:NE2	1:A:1228:ILE:O	2.52	0.43
1:A:1225:GLN:HE21	1:A:1225:GLN:HB2	1.64	0.43
1:A:1228:ILE:HB	1:F:4007[B]:PHE:CE2	2.54	0.43
1:E:5185[B]:GLU:HA	1:E:5188[B]:GLN:OE1	2.19	0.42
1:F:4004:SER:HB2	1:F:4010:GLY:HA2	2.02	0.42
1:E:5057[B]:LEU:CD2	1:E:5250[B]:ARG:CB	2.91	0.42
1:A:1105:GLY:HA2	1:A:1237:THR:HG23	2.01	0.42
1:C:2049:GLU:HB3	1:E:5049:GLU:HB3	2.00	0.42
1:F:4185:GLU:HA	1:F:4188[A]:GLN:OE1	2.19	0.42
1:A:1107:VAL:HG21	1:A:1244:ILE:HD12	2.01	0.41
1:F:4163:TYR:HB2	1:F:4164:PRO:CD	2.49	0.41
3:F:906[B]:TDR:H5M2	6:F:197:HOH:O	2.20	0.41
1:D:3163:TYR:HB2	1:D:3164:PRO:CD	2.51	0.41
1:E:5138:THR:O	1:E:5142[B]:GLU:HG2	2.20	0.41
1:C:2030[B]:ARG:CD	1:C:2033:LYS:HD2	2.50	0.41
1:A:1089:PHE:O	1:A:1213:ALA:HA	2.21	0.41
1:B:6173:SER:HA	4:B:1014:GOL:H2	2.01	0.41
1:B:6220:ILE:HD11	3:B:902:TDR:H6	2.02	0.41
1:D:3220:ILE:HD11	3:D:904[A]:TDR:H6	2.03	0.41
1:B:6108:LEU:HD22	1:B:6152:HIS:HB2	2.02	0.41
1:D:3167:GLU:HG2	1:D:3169:TYR:CE1	2.56	0.41
1:D:3185[A]:GLU:HA	1:D:3188[A]:GLN:OE1	2.21	0.41
1:A:1178:ARG:CA	1:A:1181:LYS:HE2	2.47	0.40
1:A:1108:LEU:HD22	1:A:1152:HIS:HB2	2.03	0.40
1:E:5114:VAL:HB	1:E:5157:ALA:HA	2.03	0.40
1:C:2057[B]:LEU:HB3	1:C:2253:LEU:HD11	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	265/253 (105%)	262 (99%)	2 (1%)	1 (0%)	30	16
1	B	259/253 (102%)	256 (99%)	2 (1%)	1 (0%)	30	16
1	C	259/253 (102%)	255 (98%)	3 (1%)	1 (0%)	30	16
1	D	258/253 (102%)	253 (98%)	4 (2%)	1 (0%)	30	16
1	E	268/253 (106%)	263 (98%)	4 (2%)	1 (0%)	30	16
1	F	251/253 (99%)	246 (98%)	4 (2%)	1 (0%)	30	16
All	All	1560/1518 (103%)	1535 (98%)	19 (1%)	6 (0%)	30	16

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1163	TYR
1	E	5163	TYR
1	F	4163	TYR
1	B	6163	TYR
1	C	2163	TYR
1	D	3163	TYR

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	215/202 (106%)	212 (99%)	3 (1%)	59	46
1	B	212/202 (105%)	207 (98%)	5 (2%)	43	26
1	C	211/202 (104%)	208 (99%)	3 (1%)	59	46
1	D	211/202 (104%)	207 (98%)	4 (2%)	50	34
1	E	217/202 (107%)	211 (97%)	6 (3%)	38	21
1	F	204/202 (101%)	202 (99%)	2 (1%)	68	58
All	All	1270/1212 (105%)	1247 (98%)	23 (2%)	57	36

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1225	GLN
1	A	1228	ILE
1	A	1237	THR
1	B	6147[A]	ILE
1	B	6147[B]	ILE
1	B	6230[A]	ASN
1	B	6230[B]	ASN
1	B	6232	GLU
1	C	2003	LYS
1	C	2092[A]	ILE
1	C	2092[B]	ILE
1	D	3017[A]	GLN
1	D	3017[B]	GLN
1	D	3181	LYS
1	D	3238	GLU
1	E	5003	LYS
1	E	5092	ILE
1	E	5147[A]	ILE
1	E	5147[B]	ILE
1	E	5196	GLU
1	E	5238	GLU
1	F	4238	GLU
1	F	4240	HIS

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	1014	ASN
1	A	1225	GLN
1	A	1240	HIS
1	B	6226	GLN
1	C	2226	GLN
1	D	3226	GLN
1	E	5101	HIS
1	E	5103	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

27 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	TDR	E	905[B]	-	9,9,9	1.79	3 (33%)	12,12,12	3.47	8 (66%)
2	PO4	A	808[B]	-	4,4,4	1.03	0	6,6,6	0.45	0
2	PO4	F	807[A]	-	4,4,4	1.02	0	6,6,6	0.38	0
3	TDR	D	904[B]	-	9,9,9	1.79	3 (33%)	12,12,12	3.45	8 (66%)
2	PO4	F	806	-	4,4,4	1.01	0	6,6,6	0.45	0
3	TDR	C	903[A]	-	9,9,9	1.75	3 (33%)	12,12,12	3.40	6 (50%)
3	TDR	B	902	-	9,9,9	1.51	2 (22%)	12,12,12	4.02	8 (66%)
4	GOL	F	1011	-	5,5,5	0.38	0	5,5,5	0.20	0
2	PO4	B	802	-	4,4,4	1.03	0	6,6,6	0.60	0
5	1PE	B	1001	-	15,15,15	0.46	0	14,14,14	0.26	0
2	PO4	D	804	-	4,4,4	1.09	0	6,6,6	0.53	0
2	PO4	F	807[B]	-	4,4,4	0.96	0	6,6,6	0.44	0
2	PO4	E	805	-	4,4,4	1.03	0	6,6,6	0.46	0
3	TDR	C	903[B]	-	9,9,9	1.83	3 (33%)	12,12,12	3.39	8 (66%)
2	PO4	A	801	-	4,4,4	1.05	0	6,6,6	0.27	0
4	GOL	E	1015	-	5,5,5	0.36	0	5,5,5	0.29	0
2	PO4	C	803	-	4,4,4	1.01	0	6,6,6	0.56	0
3	TDR	A	901[A]	-	9,9,9	1.72	3 (33%)	12,12,12	3.46	8 (66%)
3	TDR	F	906[A]	-	9,9,9	1.66	3 (33%)	12,12,12	3.57	6 (50%)
4	GOL	B	1014	-	5,5,5	0.41	0	5,5,5	0.28	0
2	PO4	A	808[A]	-	4,4,4	0.99	0	6,6,6	0.49	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	TDR	E	905[A]	-	9,9,9	1.65	3 (33%)	12,12,12	3.62	6 (50%)
3	TDR	D	904[A]	-	9,9,9	1.63	3 (33%)	12,12,12	3.56	6 (50%)
4	GOL	C	1013	-	5,5,5	0.40	0	5,5,5	0.29	0
3	TDR	A	901[B]	-	9,9,9	1.80	3 (33%)	12,12,12	3.47	7 (58%)
3	TDR	F	906[B]	-	9,9,9	1.67	3 (33%)	12,12,12	3.43	6 (50%)
4	GOL	A	1254	-	5,5,5	0.40	0	5,5,5	0.35	0

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
3	TDR	E	905[A]	-	-	-	0/1/1/1
3	TDR	E	905[B]	-	-	-	0/1/1/1
5	1PE	B	1001	-	-	6/13/13/13	-
3	TDR	D	904[A]	-	-	-	0/1/1/1
3	TDR	D	904[B]	-	-	-	0/1/1/1
4	GOL	C	1013	-	-	2/4/4/4	-
3	TDR	C	903[A]	-	-	-	0/1/1/1
3	TDR	C	903[B]	-	-	-	0/1/1/1
3	TDR	B	902	-	-	-	0/1/1/1
3	TDR	A	901[A]	-	-	-	0/1/1/1
3	TDR	A	901[B]	-	-	-	0/1/1/1
4	GOL	A	1254	-	-	4/4/4/4	-
4	GOL	B	1014	-	-	1/4/4/4	-
4	GOL	F	1011	-	-	1/4/4/4	-
3	TDR	F	906[A]	-	-	-	0/1/1/1
4	GOL	E	1015	-	-	2/4/4/4	-
3	TDR	F	906[B]	-	-	-	0/1/1/1

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	903[A]	TDR	C6-C5	3.32	1.39	1.34
3	E	905[B]	TDR	C6-C5	3.29	1.39	1.34
3	C	903[B]	TDR	C6-C5	3.29	1.39	1.34
3	D	904[B]	TDR	C6-C5	3.29	1.39	1.34
3	A	901[B]	TDR	C6-C5	3.27	1.39	1.34
3	A	901[A]	TDR	C6-C5	3.26	1.39	1.34

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	905[A]	TDR	C6-C5	3.22	1.38	1.34
3	F	906[A]	TDR	C6-C5	3.20	1.38	1.34
3	D	904[A]	TDR	C6-C5	3.11	1.38	1.34
3	F	906[B]	TDR	C6-C5	3.09	1.38	1.34
3	B	902	TDR	C6-C5	2.71	1.38	1.34
3	C	903[B]	TDR	C4-C5	2.61	1.49	1.44
3	D	904[B]	TDR	C4-C5	2.47	1.48	1.44
3	A	901[B]	TDR	C4-C5	2.46	1.48	1.44
3	E	905[B]	TDR	C4-C5	2.39	1.48	1.44
3	A	901[A]	TDR	C4-C5	2.37	1.48	1.44
3	A	901[B]	TDR	C4-N3	-2.34	1.34	1.38
3	C	903[A]	TDR	C4-C5	2.29	1.48	1.44
3	C	903[A]	TDR	C4-N3	-2.28	1.34	1.38
3	C	903[B]	TDR	C4-N3	-2.27	1.34	1.38
3	E	905[B]	TDR	C4-N3	-2.26	1.34	1.38
3	D	904[B]	TDR	C4-N3	-2.26	1.34	1.38
3	F	906[B]	TDR	C4-N3	-2.22	1.34	1.38
3	F	906[A]	TDR	C4-C5	2.19	1.48	1.44
3	A	901[A]	TDR	C4-N3	-2.16	1.34	1.38
3	D	904[A]	TDR	C4-C5	2.15	1.48	1.44
3	E	905[A]	TDR	C4-C5	2.08	1.48	1.44
3	E	905[A]	TDR	C4-N3	-2.06	1.35	1.38
3	D	904[A]	TDR	C4-N3	-2.03	1.35	1.38
3	F	906[A]	TDR	C4-N3	-2.03	1.35	1.38
3	B	902	TDR	C4-C5	2.03	1.48	1.44
3	F	906[B]	TDR	C4-C5	2.02	1.48	1.44

All (77) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	905[B]	TDR	N1-C2-N3	6.88	122.43	115.17
3	A	901[B]	TDR	N1-C2-N3	6.88	122.42	115.17
3	D	904[B]	TDR	N1-C2-N3	6.84	122.38	115.17
3	C	903[B]	TDR	N1-C2-N3	6.75	122.29	115.17
3	E	905[A]	TDR	N1-C2-N3	6.70	122.24	115.17
3	D	904[A]	TDR	N1-C2-N3	6.65	122.19	115.17
3	F	906[B]	TDR	N1-C2-N3	6.59	122.12	115.17
3	F	906[A]	TDR	N1-C2-N3	6.58	122.11	115.17
3	C	903[A]	TDR	N1-C2-N3	6.52	122.05	115.17
3	A	901[A]	TDR	N1-C2-N3	6.51	122.03	115.17
3	B	902	TDR	N1-C2-N3	6.38	121.90	115.17
3	B	902	TDR	O4-C4-C5	-6.29	117.72	124.92

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	902	TDR	C4-N3-C2	-5.63	118.62	126.37
3	F	906[A]	TDR	O4-C4-C5	-5.43	118.70	124.92
3	B	902	TDR	C5-C4-N3	5.38	120.00	115.32
3	B	902	TDR	O2-C2-N1	-5.35	117.27	122.79
3	E	905[A]	TDR	O4-C4-C5	-5.33	118.82	124.92
3	D	904[A]	TDR	O4-C4-C5	-5.16	119.01	124.92
3	E	905[A]	TDR	C4-N3-C2	-5.04	119.43	126.37
3	F	906[A]	TDR	C4-N3-C2	-4.95	119.56	126.37
3	D	904[A]	TDR	C4-N3-C2	-4.94	119.57	126.37
3	F	906[B]	TDR	C4-N3-C2	-4.78	119.78	126.37
3	A	901[B]	TDR	C4-N3-C2	-4.76	119.81	126.37
3	F	906[B]	TDR	O4-C4-C5	-4.76	119.47	124.92
3	A	901[A]	TDR	C4-N3-C2	-4.73	119.86	126.37
3	E	905[B]	TDR	C4-N3-C2	-4.70	119.90	126.37
3	A	901[A]	TDR	O4-C4-C5	-4.67	119.58	124.92
3	C	903[A]	TDR	C4-N3-C2	-4.65	119.96	126.37
3	D	904[B]	TDR	C4-N3-C2	-4.65	119.97	126.37
3	C	903[A]	TDR	O4-C4-C5	-4.63	119.62	124.92
3	D	904[A]	TDR	O2-C2-N1	-4.57	118.08	122.79
3	E	905[A]	TDR	O2-C2-N1	-4.46	118.19	122.79
3	C	903[B]	TDR	C4-N3-C2	-4.45	120.24	126.37
3	E	905[A]	TDR	C5-C4-N3	4.41	119.16	115.32
3	F	906[A]	TDR	C5-C4-N3	4.39	119.14	115.32
3	E	905[B]	TDR	O2-C2-N1	-4.38	118.27	122.79
3	D	904[B]	TDR	O2-C2-N1	-4.38	118.27	122.79
3	F	906[A]	TDR	O2-C2-N1	-4.32	118.33	122.79
3	A	901[A]	TDR	O2-C2-N1	-4.27	118.39	122.79
3	A	901[B]	TDR	O2-C2-N1	-4.22	118.43	122.79
3	A	901[B]	TDR	O4-C4-C5	-4.22	120.09	124.92
3	F	906[B]	TDR	C5-C4-N3	4.19	118.97	115.32
3	C	903[B]	TDR	O2-C2-N1	-4.19	118.47	122.79
3	D	904[A]	TDR	C5-C4-N3	4.14	118.92	115.32
3	E	905[B]	TDR	O4-C4-C5	-4.09	120.24	124.92
3	C	903[A]	TDR	O2-C2-N1	-4.08	118.58	122.79
3	A	901[A]	TDR	C5-C4-N3	4.08	118.87	115.32
3	D	904[B]	TDR	O4-C4-C5	-4.06	120.27	124.92
3	C	903[A]	TDR	C5-C4-N3	4.02	118.82	115.32
3	F	906[B]	TDR	O2-C2-N1	-3.98	118.69	122.79
3	A	901[B]	TDR	C5-C4-N3	3.80	118.63	115.32
3	E	905[B]	TDR	C5-C4-N3	3.62	118.47	115.32
3	D	904[B]	TDR	C5-C4-N3	3.60	118.45	115.32
3	E	905[B]	TDR	C5-C6-N1	-3.55	118.59	122.42

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	901[B]	TDR	C5-C6-N1	-3.55	118.60	122.42
3	C	903[B]	TDR	O4-C4-C5	-3.52	120.89	124.92
3	B	902	TDR	C5-C6-N1	-3.50	118.65	122.42
3	C	903[B]	TDR	C5-C4-N3	3.45	118.32	115.32
3	D	904[B]	TDR	C5-C6-N1	-3.43	118.73	122.42
3	C	903[A]	TDR	C5-C6-N1	-3.42	118.73	122.42
3	D	904[A]	TDR	C5-C6-N1	-3.39	118.77	122.42
3	F	906[A]	TDR	C5-C6-N1	-3.38	118.78	122.42
3	F	906[B]	TDR	C5-C6-N1	-3.38	118.78	122.42
3	E	905[A]	TDR	C5-C6-N1	-3.36	118.80	122.42
3	A	901[A]	TDR	C5-C6-N1	-3.33	118.84	122.42
3	C	903[B]	TDR	CM5-C5-C6	-3.07	120.07	123.23
3	C	903[B]	TDR	C5-C6-N1	-3.05	119.14	122.42
3	C	903[B]	TDR	CM5-C5-C4	3.02	122.00	118.78
3	D	904[B]	TDR	CM5-C5-C6	-2.51	120.65	123.23
3	E	905[B]	TDR	CM5-C5-C6	-2.47	120.69	123.23
3	B	902	TDR	CM5-C5-C4	2.41	121.36	118.78
3	A	901[B]	TDR	CM5-C5-C6	-2.26	120.91	123.23
3	D	904[B]	TDR	CM5-C5-C4	2.20	121.13	118.78
3	A	901[A]	TDR	CM5-C5-C6	-2.09	121.08	123.23
3	E	905[B]	TDR	CM5-C5-C4	2.08	121.00	118.78
3	A	901[A]	TDR	CM5-C5-C4	2.06	120.98	118.78
3	B	902	TDR	CM5-C5-C6	-2.05	121.12	123.23

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1254	GOL	O1-C1-C2-O2
4	E	1015	GOL	C1-C2-C3-O3
5	B	1001	1PE	OH4-C13-C23-OH3
5	B	1001	1PE	OH2-C12-C22-OH3
4	A	1254	GOL	O1-C1-C2-C3
4	A	1254	GOL	C1-C2-C3-O3
4	C	1013	GOL	O1-C1-C2-C3
5	B	1001	1PE	OH7-C16-C26-OH6
4	E	1015	GOL	O2-C2-C3-O3
4	C	1013	GOL	O1-C1-C2-O2
4	F	1011	GOL	O2-C2-C3-O3
5	B	1001	1PE	C16-C26-OH6-C15
5	B	1001	1PE	C13-C23-OH3-C22
4	A	1254	GOL	O2-C2-C3-O3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
5	B	1001	1PE	OH6-C15-C25-OH5
4	B	1014	GOL	O1-C1-C2-O2

There are no ring outliers.

13 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	905[B]	TDR	1	0
3	C	903[A]	TDR	1	0
3	B	902	TDR	3	0
4	F	1011	GOL	1	0
2	B	802	PO4	1	0
5	B	1001	1PE	1	0
2	D	804	PO4	1	0
3	F	906[A]	TDR	1	0
4	B	1014	GOL	1	0
3	E	905[A]	TDR	1	0
3	D	904[A]	TDR	1	0
4	C	1013	GOL	1	0
3	F	906[B]	TDR	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

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	251/253 (99%)	-0.24	16 (6%) 25 27	2, 5, 21, 39	20 (7%)
1	B	246/253 (97%)	-0.39	6 (2%) 59 64	2, 5, 17, 31	17 (6%)
1	C	245/253 (96%)	-0.14	10 (4%) 41 45	2, 7, 23, 39	18 (7%)
1	D	244/253 (96%)	-0.31	7 (2%) 53 58	2, 6, 18, 33	18 (7%)
1	E	252/253 (99%)	-0.40	10 (3%) 42 46	2, 5, 20, 27	18 (7%)
1	F	245/253 (96%)	-0.45	7 (2%) 53 58	2, 5, 16, 31	10 (4%)
All	All	1483/1518 (97%)	-0.32	56 (3%) 44 48	2, 6, 19, 39	101 (6%)

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1228	ILE	10.7
1	A	1233	THR	9.3
1	C	2237	THR	8.5
1	D	3004	SER	6.8
1	E	5002	SER	5.7
1	A	1230	ASN	5.5
1	F	4230	ASN	5.1
1	A	1237	THR	4.8
1	B	6002	SER	4.7
1	E	5253	LEU	4.5
1	D	3231	ALA	4.5
1	A	1225	GLN	4.2
1	B	6253	LEU	4.2
1	B	6232	GLU	4.1
1	B	6231	ALA	4.0
1	C	2238	GLU	3.9
1	A	1234	MET	3.8
1	D	3238	GLU	3.8
1	A	1058	ASP	3.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	1231	ALA	3.7
1	F	4002	SER	3.4
1	E	5003	LYS	3.4
1	A	1253	LEU	3.4
1	F	4238	GLU	3.3
1	D	3005[A]	ASP	3.3
1	E	5147[A]	ILE	3.2
1	C	2101[A]	HIS	3.1
1	A	1229	PRO	3.1
1	A	1226	GLN	3.0
1	F	4240	HIS	3.0
1	C	2003	LYS	2.9
1	E	5233	THR	2.9
1	C	2230	ASN	2.9
1	A	1232	GLU	2.8
1	C	2239	SER	2.8
1	B	6003	LYS	2.7
1	C	2240	HIS	2.7
1	E	5238	GLU	2.7
1	E	5231	ALA	2.6
1	A	1227	GLU	2.5
1	E	5235	LYS	2.4
1	D	3101[A]	HIS	2.4
1	D	3239	SER	2.3
1	A	1003	LYS	2.3
1	F	4253	LEU	2.3
1	F	4101	HIS	2.3
1	A	1236	GLN	2.3
1	C	2225	GLN	2.2
1	B	6146	SER	2.2
1	F	4239	SER	2.1
1	C	2241	ALA	2.1
1	C	2145[A]	LYS	2.1
1	D	3253	LEU	2.1
1	A	1235	LYS	2.1
1	E	5058	ASP	2.0
1	E	5252	LEU	2.0

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 oligosaccharides in this entry.

6.4 Ligands ⓘ

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
5	1PE	B	1001	16/16	0.64	0.41	9,27,33,33	16
3	TDR	B	902	9/9	0.65	0.20	12,12,13,15	0
3	TDR	C	903[B]	9/9	0.68	0.23	21,21,21,21	9
3	TDR	C	903[A]	9/9	0.68	0.23	19,20,20,20	9
4	GOL	E	1015	6/6	0.69	0.20	41,42,42,42	0
4	GOL	F	1011	6/6	0.72	0.22	33,33,33,34	0
4	GOL	B	1014	6/6	0.73	0.19	41,42,42,42	0
4	GOL	A	1254	6/6	0.73	0.25	35,35,35,35	0
3	TDR	F	906[B]	9/9	0.74	0.18	8,9,9,10	9
3	TDR	F	906[A]	9/9	0.74	0.18	11,11,11,12	9
3	TDR	A	901[A]	9/9	0.79	0.18	14,15,15,15	9
3	TDR	D	904[A]	9/9	0.79	0.16	8,8,9,9	9
3	TDR	D	904[B]	9/9	0.79	0.16	17,17,17,17	9
3	TDR	E	905[A]	9/9	0.79	0.17	10,10,10,10	9
3	TDR	E	905[B]	9/9	0.79	0.17	17,17,17,17	9
3	TDR	A	901[B]	9/9	0.79	0.18	17,17,17,17	9
2	PO4	F	807[A]	5/5	0.83	0.13	15,15,16,16	5
2	PO4	F	807[B]	5/5	0.83	0.13	17,17,18,18	5
4	GOL	C	1013	6/6	0.84	0.15	33,33,33,34	0
2	PO4	A	808[B]	5/5	0.95	0.09	23,23,23,23	5
2	PO4	A	808[A]	5/5	0.95	0.09	15,15,15,15	5
2	PO4	B	802	5/5	0.99	0.04	7,7,8,8	0
2	PO4	C	803	5/5	0.99	0.03	8,8,9,9	0
2	PO4	D	804	5/5	0.99	0.04	7,7,7,7	0
2	PO4	E	805	5/5	0.99	0.02	5,5,6,7	0
2	PO4	F	806	5/5	0.99	0.03	5,6,7,7	0
2	PO4	A	801	5/5	0.99	0.03	4,4,4,5	0

6.5 Other polymers ⓘ

There are no such residues in this entry.