



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 04:23 PM GMT

PDB ID : 3DGA
Title : Wild-type Plasmodium falciparum dihydrofolate reductase-thymidylatesynthase (PfDHFR-TS) complexed with RJF01302, NADPH, and dUMP
Authors : Dasgupta, T.; Chitnumsub, P.; Maneeruttanarungroj, C.; Kamchonwongpaisan, S.; Nichols, S.; Lyons, T.M.; Tirado-Rives, J.; Jorgensen, W.L.; Yuthavong, Y.; Anderson, K.S.
Deposited on : 2008-06-13
Resolution : 2.70 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

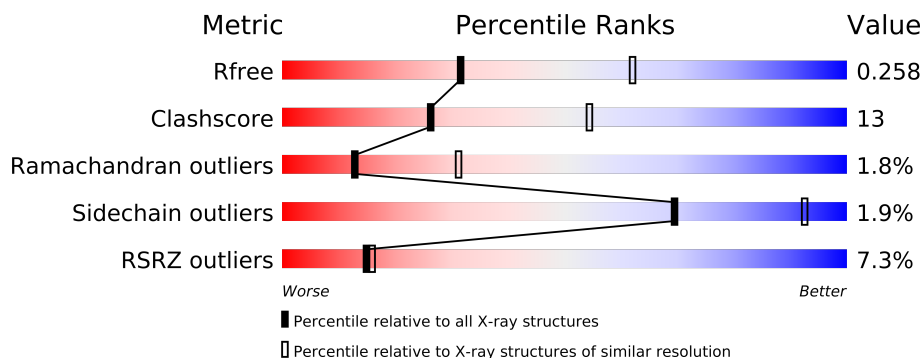
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.70 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}	66092	1557 (2.70-2.70)
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)
RSRZ outliers	66119	1559 (2.70-2.70)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	280	
1	B	280	
2	C	328	
2	D	328	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	RJ1	A	609	-	X
3	RJ1	B	609	-	X

2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 9510 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

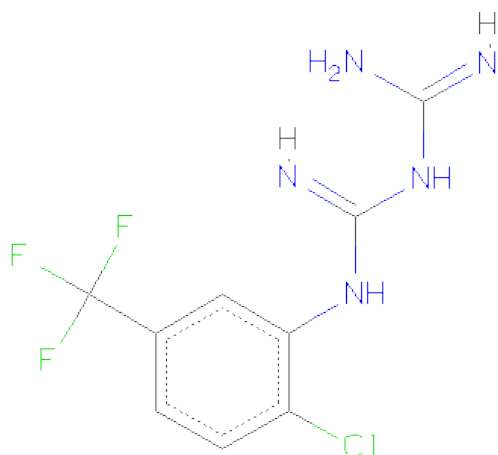
- Molecule 1 is a protein called Bifunctional dihydrofolate reductase-thymidylatesynthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	219	Total	C	N	O	S	0	0	0
			1815	1176	291	335	13			
1	B	224	Total	C	N	O	S	0	0	0
			1851	1195	298	346	12			

- Molecule 2 is a protein called Bifunctional dihydrofolate reductase-thymidylatesynthase.

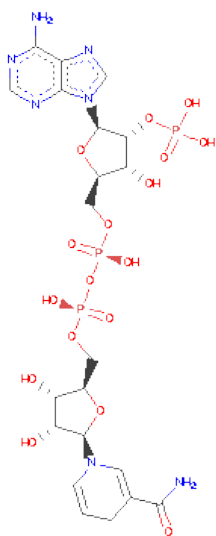
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	325	Total	C	N	O	S	0	0	0
			2705	1743	455	492	15			
2	D	326	Total	C	N	O	S	0	0	0
			2718	1749	456	498	15			

- Molecule 3 is N-[2-CHLORO-5-(TRIFLUOROMETHYL)PHENYL]IMIDODICARBONIMI DICDIAMIDE (three-letter code: RJ1) (formula: C₉H₉ClF₃N₅).



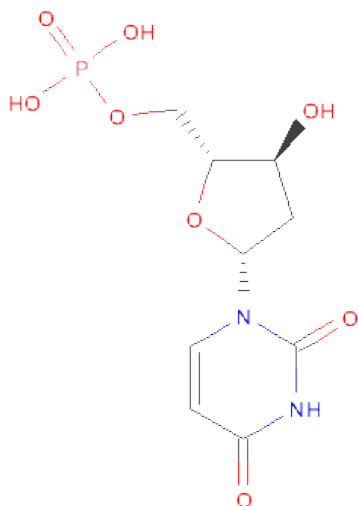
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Cl	F	N	0	0
			18	9	1	3	5		
3	B	1	Total	C	Cl	F	N	0	0
			18	9	1	3	5		

- Molecule 4 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDEPHOSPHATE (three-letter code: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
4	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 5 is 2'-DEOXYURIDINE 5'-MONOPHOSPHATE (three-letter code: UMP) (formula: $C_9H_{13}N_2O_8P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	C	1	Total	C	N	O	P	0	0
			20	9	2	8	1		
5	D	1	Total	C	N	O	P	0	0
			20	9	2	8	1		

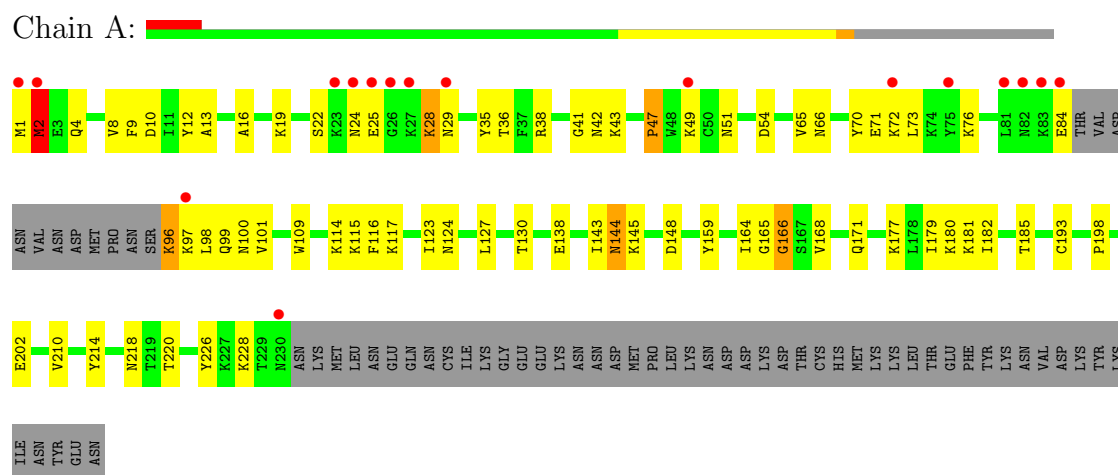
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	44	Total	O	0	0
			44	44		
6	B	18	Total	O	0	0
			18	18		
6	C	88	Total	O	0	0
			88	88		
6	D	99	Total	O	0	0
			99	99		

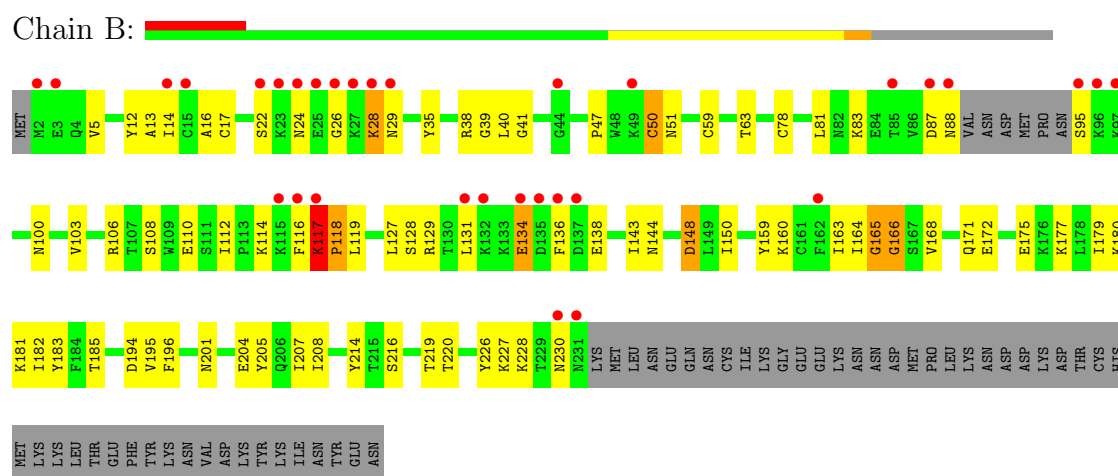
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

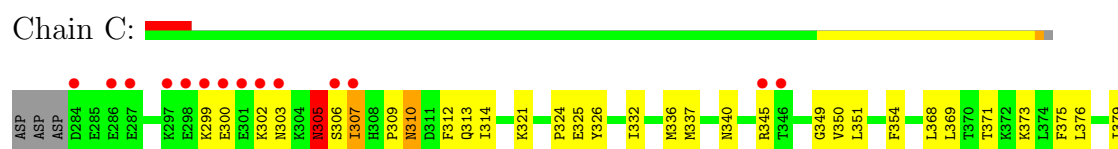
- Molecule 1: Bifunctional dihydrofolate reductase-thymidylatesynthase

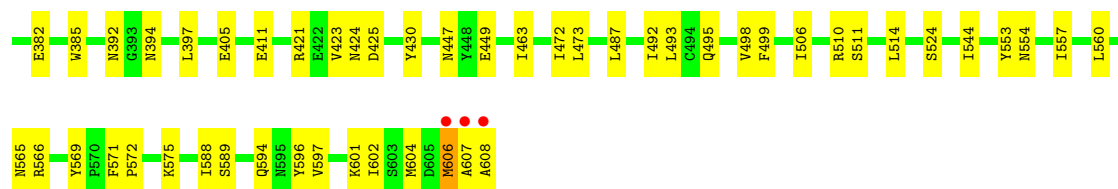


- Molecule 1: Bifunctional dihydrofolate reductase-thymidylatesynthase



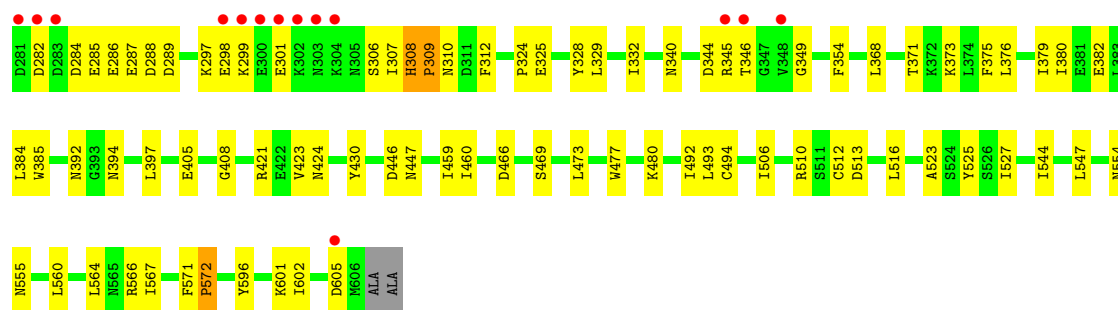
- Molecule 2: Bifunctional dihydrofolate reductase-thymidylatesynthase





● Molecule 2: Bifunctional dihydrofolate reductase-thymidylatesynthase

Chain D:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	57.38Å 156.01Å 164.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.18 – 2.70 29.18 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.18-2.70) 99.6 (29.18-2.70)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.40 (at 2.68Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.205 , 0.256 0.206 , 0.258	Depositor DCC
R_{free} test set	2084 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	39.3	Xtriage
Anisotropy	0.746	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 27.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 41425 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9510	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.59% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: NDP, UMP, RJ1

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	0.41	0/1849	0.71	3/2485 (0.1%)
1	B	0.39	0/1885	0.70	4/2536 (0.2%)
2	C	0.37	0/2776	0.64	0/3755
2	D	0.38	0/2789	0.63	0/3774
All	All	0.39	0/9299	0.67	7/12550 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	166	GLY	N-CA-C	-11.51	84.34	113.10
1	A	166	GLY	N-CA-C	-10.36	87.21	113.10
1	B	117	LYS	C-N-CD	8.61	146.47	128.40
1	A	165	GLY	N-CA-C	7.54	131.94	113.10
1	B	134	GLU	N-CA-C	7.13	130.27	111.00
1	A	25	GLU	N-CA-C	-6.35	93.86	111.00
1	B	165	GLY	N-CA-C	6.18	128.54	113.10

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1815	0	1846	63	0
1	B	1851	0	1871	69	0
2	C	2705	0	2634	64	0
2	D	2718	0	2636	58	0
3	A	18	0	6	3	0
3	B	18	0	5	3	0
4	A	48	0	26	5	0
4	B	48	0	26	6	0
5	C	20	0	11	0	0
5	D	20	0	11	0	0
6	A	44	0	0	0	0
6	B	18	0	0	0	0
6	C	88	0	0	0	0
6	D	99	0	0	2	0
All	All	9510	0	9072	244	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 13.

All (244) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:201:ASN:HB3	1:B:204:GLU:HG3	1.42	0.98
1:A:127:LEU:HD23	1:A:143:ILE:HG13	1.53	0.87
2:C:300:GLU:HG2	2:C:300:GLU:O	1.72	0.87
2:D:299:LYS:HG2	2:D:299:LYS:O	1.77	0.82
2:C:376:LEU:HD22	2:C:379:ILE:HD11	1.60	0.81
1:B:182:ILE:HB	1:B:226:TYR:HB2	1.63	0.80
1:A:115:LYS:HE3	1:A:116:PHE:CZ	2.18	0.79
1:B:12:TYR:CE1	1:B:180:LYS:HD2	2.19	0.77
1:A:4:GLN:HE22	1:A:228:LYS:NZ	1.83	0.76
1:B:164:ILE:HB	3:B:609:RJ1:CL7	2.24	0.75
1:B:87:ASP:OD1	1:B:88:ASN:N	2.18	0.74
1:A:8:VAL:HA	1:A:76:LYS:HD3	1.67	0.73
2:C:305:ASN:O	2:C:307:ILE:HG13	1.89	0.73
1:B:129:ARG:NH1	1:B:129:ARG:HB2	2.04	0.72
1:B:117:LYS:O	1:B:118:PRO:C	2.20	0.70
1:B:117:LYS:O	1:B:119:LEU:N	2.27	0.68
1:B:41:GLY:HA2	1:B:47:PRO:HD3	1.75	0.67
2:D:494:CYS:SG	2:D:525:TYR:HE2	2.18	0.67
1:B:14:ILE:HG13	1:B:183:TYR:HB2	1.77	0.67
1:A:4:GLN:HE22	1:A:228:LYS:HZ1	1.42	0.66
2:C:506:ILE:HG12	2:C:544:ILE:HB	1.78	0.66
2:C:307:ILE:HG12	2:C:337:MET:SD	2.35	0.66

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:164:ILE:HB	3:A:609:RJ1:CL7	2.34	0.65
2:C:307:ILE:H	2:C:307:ILE:HD12	1.59	0.65
2:D:513:ASP:OD2	2:D:516:LEU:HB2	1.96	0.65
1:B:168:VAL:HG23	4:B:610:NDP:O2N	1.97	0.63
1:B:22:SER:HB2	1:B:24:ASN:ND2	2.13	0.63
1:A:12:TYR:HD1	1:A:181:LYS:HB2	1.65	0.62
1:A:71:GLU:CD	1:A:71:GLU:H	2.04	0.61
2:D:332:ILE:HD13	2:D:560:LEU:HD22	1.83	0.61
1:B:13:ALA:HB2	1:B:179:ILE:HD12	1.82	0.61
2:D:298:GLU:HA	2:D:298:GLU:OE1	2.02	0.60
2:C:332:ILE:CD1	2:C:514:LEU:HB3	2.32	0.60
1:B:129:ARG:HH11	1:B:129:ARG:HB2	1.67	0.59
1:B:208:ILE:HD13	1:B:227:LYS:HD2	1.83	0.59
2:D:329:LEU:HD22	2:D:564:LEU:HD12	1.85	0.59
2:C:463:ILE:HG23	2:C:498:VAL:HG21	1.85	0.59
1:B:28:LYS:O	1:B:29:ASN:C	2.40	0.59
2:C:492:ILE:HD11	2:C:510:ARG:HD3	1.84	0.59
1:A:166:GLY:HA3	4:A:610:NDP:PA	2.43	0.59
2:D:373:LYS:HE2	2:D:375:PHE:CE1	2.38	0.58
2:D:309:PRO:HA	2:D:312:PHE:HD2	1.67	0.58
1:B:108:SER:O	1:B:112:ILE:HG13	2.04	0.58
1:B:116:PHE:O	1:B:117:LYS:O	2.21	0.58
1:A:182:ILE:HB	1:A:226:TYR:HB2	1.86	0.58
2:D:494:CYS:SG	2:D:525:TYR:CE2	2.97	0.58
1:A:168:VAL:HG23	4:A:610:NDP:O2N	2.03	0.58
2:C:324:PRO:HB2	2:C:571:PHE:HE2	1.69	0.58
2:D:459:ILE:HG13	2:D:460:ILE:N	2.19	0.57
2:C:313:GLN:HG3	2:C:314:ILE:H	1.70	0.57
1:A:96:LYS:HE3	1:A:97:LYS:H	1.68	0.57
1:A:70:TYR:O	1:A:73:LEU:N	2.38	0.57
2:D:284:ASP:O	2:D:287:GLU:HG3	2.06	0.56
1:A:41:GLY:HA2	1:A:47:PRO:HD3	1.88	0.56
1:B:143:ILE:HG22	1:B:144:ASN:N	2.21	0.56
1:B:166:GLY:HA3	4:B:610:NDP:PA	2.46	0.56
2:D:506:ILE:HG12	2:D:544:ILE:HB	1.88	0.56
1:B:127:LEU:HD12	1:B:143:ILE:HB	1.87	0.55
1:B:81:LEU:HD12	1:B:83:LYS:HE2	1.89	0.55
1:A:1:MET:HG2	1:A:2:MET:N	2.22	0.55
1:B:35:TYR:CZ	1:B:38:ARG:HD3	2.42	0.55
1:B:205:TYR:CD1	1:B:228:LYS:HA	2.42	0.54
1:A:42:ASN:ND2	1:A:43:LYS:HG3	2.23	0.54
1:B:143:ILE:HG23	1:B:148:ASP:HB3	1.90	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:309:PRO:HA	2:D:312:PHE:CD2	2.42	0.54
2:C:423:VAL:O	2:C:424:ASN:HB2	2.07	0.54
2:D:307:ILE:O	2:D:309:PRO:HD3	2.07	0.54
1:A:114:LYS:NZ	1:A:114:LYS:HB3	2.23	0.54
2:C:473:LEU:CD1	2:C:495:GLN:HG3	2.38	0.54
1:B:129:ARG:CB	1:B:129:ARG:HH11	2.20	0.54
2:C:321:LYS:HD2	2:C:326:TYR:CE1	2.42	0.54
1:B:12:TYR:HE1	1:B:180:LYS:HD2	1.70	0.53
2:D:566:ARG:NH1	2:D:602:ILE:HD11	2.23	0.53
1:B:207:ILE:HB	2:D:567:ILE:HD13	1.91	0.52
2:C:345:ARG:HG2	2:C:345:ARG:HH11	1.74	0.52
2:C:325:GLU:HG3	2:C:369:LEU:HD22	1.91	0.52
1:A:19:LYS:HG2	1:A:36:THR:HG22	1.92	0.52
2:D:555:ASN:ND2	2:D:605:ASP:OD1	2.43	0.52
1:B:12:TYR:CD1	1:B:181:LYS:HB2	2.45	0.52
1:B:164:ILE:O	3:B:609:RJ1:CL7	2.65	0.52
2:C:506:ILE:HG13	2:D:354:PHE:CE2	2.44	0.52
2:D:423:VAL:O	2:D:424:ASN:HB2	2.08	0.52
1:B:216:SER:O	1:B:219:THR:HG22	2.10	0.52
2:D:421:ARG:HH11	2:D:421:ARG:HG2	1.75	0.52
2:D:397:LEU:HD21	2:D:405:GLU:HB2	1.91	0.52
2:C:376:LEU:HD22	2:C:379:ILE:CD1	2.35	0.51
1:A:35:TYR:CZ	1:A:38:ARG:HD3	2.45	0.51
2:D:324:PRO:HB2	2:D:571:PHE:HE2	1.74	0.51
1:A:28:LYS:O	1:A:29:ASN:HB2	2.10	0.51
1:B:131:LEU:HD22	1:B:136:PHE:CE2	2.46	0.51
2:C:309:PRO:HG2	2:C:310:ASN:H	1.75	0.51
1:B:194:ASP:OD1	1:B:195:VAL:HG13	2.11	0.51
2:C:447:ASN:OD1	2:C:449:GLU:HG2	2.10	0.51
1:B:28:LYS:HD2	1:B:28:LYS:C	2.32	0.50
1:A:177:LYS:HB3	1:A:228:LYS:NZ	2.27	0.50
1:A:171:GLN:HA	1:A:198:PRO:HG2	1.92	0.50
2:C:354:PHE:CE2	2:D:506:ILE:HG13	2.45	0.50
1:A:109:TRP:CE2	1:A:117:LYS:HD2	2.46	0.50
2:C:299:LYS:NZ	2:C:302:LYS:HB2	2.25	0.50
2:D:308:HIS:O	2:D:310:ASN:N	2.44	0.50
2:C:307:ILE:CD1	2:C:307:ILE:H	2.18	0.50
2:C:392:ASN:OD1	2:C:394:ASN:HB2	2.11	0.50
2:C:493:LEU:CD2	2:D:492:ILE:HG21	2.42	0.50
1:A:164:ILE:O	3:A:609:RJ1:CL7	2.67	0.49
2:D:376:LEU:HD22	2:D:379:ILE:HD11	1.94	0.49
2:C:560:LEU:HD21	2:C:604:MET:HE1	1.94	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:C:382:GLU:O	2:C:385:TRP:HB3	2.12	0.49
2:D:493:LEU:C	2:D:493:LEU:HD12	2.32	0.49
1:B:78:CYS:HB3	1:B:83:LYS:O	2.13	0.49
2:D:446:ASP:OD1	2:D:447:ASN:N	2.44	0.49
1:A:8:VAL:HA	1:A:76:LYS:CD	2.42	0.49
2:C:307:ILE:N	2:C:307:ILE:HD12	2.27	0.49
1:B:214:TYR:O	1:B:220:THR:HA	2.13	0.48
2:D:473:LEU:HA	2:D:494:CYS:O	2.14	0.48
2:C:313:GLN:HG3	2:C:314:ILE:N	2.27	0.48
1:A:24:ASN:O	1:A:24:ASN:CG	2.51	0.48
1:A:144:ASN:HD22	1:A:145:LYS:N	2.11	0.48
1:B:171:GLN:O	1:B:175:GLU:HG2	2.14	0.48
2:C:312:PHE:HA	2:C:565:ASN:OD1	2.13	0.48
1:B:134:GLU:C	1:B:136:PHE:H	2.17	0.48
2:C:493:LEU:HD22	2:D:492:ILE:HG21	1.95	0.48
2:C:332:ILE:HD11	2:C:514:LEU:HB3	1.96	0.47
1:A:180:LYS:HE2	2:D:285:GLU:OE1	2.13	0.47
1:A:16:ALA:HA	1:A:185:THR:HB	1.97	0.47
2:D:572:PRO:HB3	2:D:596:TYR:HA	1.95	0.47
2:C:566:ARG:CZ	2:C:602:ILE:HD11	2.44	0.47
2:D:512:CYS:SG	2:D:547:LEU:HD22	2.54	0.47
1:B:168:VAL:O	1:B:172:GLU:HG2	2.15	0.47
1:A:214:TYR:O	1:A:220:THR:HA	2.15	0.47
1:A:84:GLU:CD	1:A:84:GLU:N	2.68	0.47
2:C:575:LYS:HE3	2:C:594:GLN:OE1	2.14	0.47
2:C:411:GLU:H	2:C:411:GLU:CD	2.18	0.47
2:C:492:ILE:HG21	2:D:493:LEU:CD2	2.45	0.46
1:B:106:ARG:NH1	1:B:110:GLU:OE2	2.43	0.46
2:D:328:TYR:CZ	2:D:332:ILE:HD11	2.50	0.46
1:A:4:GLN:NE2	1:A:228:LYS:NZ	2.58	0.46
2:D:349:GLY:C	2:D:554:ASN:ND2	2.69	0.46
1:B:12:TYR:HD1	1:B:181:LYS:HB2	1.81	0.46
1:A:144:ASN:C	1:A:144:ASN:HD22	2.18	0.46
2:C:300:GLU:CG	2:C:300:GLU:O	2.50	0.46
2:D:408:GLY:O	2:D:423:VAL:HG13	2.16	0.46
1:B:100:ASN:OD1	1:B:159:TYR:HB3	2.16	0.46
1:B:128:SER:CB	1:B:131:LEU:HD12	2.45	0.46
2:D:477:TRP:HB2	2:D:492:ILE:HG23	1.98	0.46
2:D:368:LEU:HD13	2:D:376:LEU:HD11	1.97	0.46
1:A:51:ASN:HB3	1:A:54:ASP:HB3	1.98	0.46
1:B:177:LYS:HE2	1:B:204:GLU:OE1	2.16	0.45
1:A:42:ASN:HB2	1:A:193:CYS:HA	1.98	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:210:VAL:HG23	1:A:210:VAL:O	2.15	0.45
1:A:210:VAL:HG21	2:C:326:TYR:HE2	1.82	0.45
1:A:101:VAL:HG22	1:A:123:ILE:HB	1.96	0.45
1:B:22:SER:C	1:B:24:ASN:H	2.20	0.45
2:C:350:VAL:HG12	2:C:553:TYR:CD1	2.51	0.45
2:C:368:LEU:CD1	2:C:376:LEU:HD21	2.46	0.45
2:C:493:LEU:C	2:C:493:LEU:HD12	2.37	0.45
1:B:41:GLY:O	1:B:195:VAL:HG22	2.17	0.45
2:D:382:GLU:O	2:D:385:TRP:HB3	2.17	0.45
1:A:166:GLY:HA3	4:A:610:NDP:O1A	2.17	0.45
1:A:4:GLN:NE2	1:A:228:LYS:HZ3	2.15	0.45
1:A:4:GLN:HE22	1:A:228:LYS:HZ3	1.62	0.45
1:A:144:ASN:HD21	1:A:145:LYS:NZ	2.15	0.45
2:C:607:ALA:O	2:C:608:ALA:OXT	2.35	0.45
2:C:472:ILE:C	2:C:473:LEU:HD12	2.37	0.45
3:A:609:RJ1:N17	3:A:609:RJ1:CL7	2.87	0.45
2:C:510:ARG:HG3	2:C:511:SER:N	2.32	0.44
2:C:499:PHE:CZ	2:D:340:ASN:HB3	2.52	0.44
1:B:201:ASN:HB3	1:B:204:GLU:CG	2.30	0.44
2:C:492:ILE:HG21	2:D:493:LEU:HD22	1.98	0.44
1:B:131:LEU:HD22	1:B:136:PHE:CZ	2.53	0.44
2:D:523:ALA:O	2:D:527:ILE:HG13	2.17	0.44
3:B:609:RJ1:N17	3:B:609:RJ1:CL7	2.88	0.44
1:B:38:ARG:O	1:B:196:PHE:HB3	2.18	0.43
1:A:65:VAL:HG12	1:A:66:ASN:N	2.33	0.43
2:D:480:LYS:HD3	6:D:1151:HOH:O	2.17	0.43
2:C:421:ARG:HD2	2:C:425:ASP:HB3	1.99	0.43
2:D:492:ILE:HD11	2:D:510:ARG:HD3	2.00	0.43
2:D:368:LEU:HA	6:D:1036:HOH:O	2.17	0.43
1:A:100:ASN:OD1	1:A:159:TYR:HB3	2.19	0.43
2:C:397:LEU:HD21	2:C:405:GLU:HB2	2.00	0.43
2:C:305:ASN:O	2:C:307:ILE:N	2.51	0.43
1:B:129:ARG:HG3	4:B:610:NDP:C2A	2.48	0.43
1:B:16:ALA:HA	1:B:185:THR:HB	1.99	0.43
1:B:59:CYS:O	1:B:63:THR:HG23	2.19	0.43
2:C:572:PRO:HB3	2:C:596:TYR:HA	2.00	0.43
1:B:114:LYS:O	1:B:117:LYS:N	2.47	0.43
1:B:13:ALA:O	1:B:182:ILE:HA	2.19	0.42
2:D:421:ARG:NH1	2:D:421:ARG:HG2	2.34	0.42
2:C:325:GLU:OE2	2:C:371:THR:HG23	2.19	0.42
2:D:392:ASN:OD1	2:D:394:ASN:HB2	2.20	0.42
2:D:380:ILE:O	2:D:384:LEU:HG	2.19	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:C:336:MET:HE2	2:C:557:ILE:HG23	2.02	0.42
2:C:368:LEU:HD23	2:C:596:TYR:CE1	2.55	0.42
1:B:40:LEU:O	4:B:610:NDP:H2N	2.20	0.42
1:A:65:VAL:HG11	1:A:98:LEU:HB3	2.02	0.42
2:C:321:LYS:HD2	2:C:326:TYR:CD1	2.55	0.42
1:A:115:LYS:HE3	1:A:116:PHE:CE1	2.53	0.42
1:A:72:LYS:HB3	1:A:72:LYS:HE2	1.88	0.42
2:D:325:GLU:OE2	2:D:371:THR:HG23	2.20	0.42
1:B:163:ILE:C	1:B:165:GLY:H	2.22	0.42
1:A:65:VAL:CG1	1:A:66:ASN:N	2.83	0.42
2:C:314:ILE:HD12	2:C:565:ASN:HB3	2.03	0.41
1:A:109:TRP:CZ2	1:A:117:LYS:HD2	2.55	0.41
1:A:124:ASN:N	1:A:124:ASN:HD22	2.18	0.41
1:A:127:LEU:O	4:A:610:NDP:H1B	2.19	0.41
1:A:28:LYS:HD3	2:C:373:LYS:NZ	2.35	0.41
1:B:17:CYS:HA	1:B:39:GLY:O	2.19	0.41
1:B:12:TYR:CE2	1:B:160:LYS:HD3	2.55	0.41
2:C:373:LYS:HE2	2:C:375:PHE:CZ	2.55	0.41
1:A:177:LYS:HB3	1:A:228:LYS:HZ1	1.86	0.41
1:B:114:LYS:O	1:B:117:LYS:HB2	2.21	0.41
1:A:71:GLU:N	1:A:71:GLU:CD	2.73	0.41
2:D:466:ASP:OD2	2:D:469:SER:HB2	2.21	0.41
1:B:41:GLY:H	1:B:195:VAL:HG23	1.85	0.41
1:A:99:GLN:HE21	1:A:123:ILE:HG13	1.86	0.41
2:C:606:MET:O	2:C:606:MET:HG2	2.21	0.41
1:B:5:VAL:HG11	1:B:150:ILE:HD12	2.01	0.41
1:A:127:LEU:CD2	1:A:143:ILE:HG13	2.38	0.41
1:B:127:LEU:O	4:B:610:NDP:H1B	2.20	0.41
2:D:344:ASP:C	2:D:346:THR:H	2.23	0.41
2:D:601:LYS:HB2	2:D:601:LYS:HE3	1.89	0.41
2:D:285:GLU:O	2:D:288:ASP:HB2	2.20	0.41
2:C:349:GLY:C	2:C:554:ASN:HD22	2.24	0.41
1:B:103:VAL:O	1:B:164:ILE:HG12	2.20	0.41
1:B:166:GLY:HA3	4:B:610:NDP:O2A	2.21	0.41
1:B:22:SER:HB2	1:B:24:ASN:HD22	1.85	0.41
1:A:181:LYS:NZ	2:D:289:ASP:OD2	2.52	0.41
1:A:9:PHE:O	1:A:10:ASP:C	2.57	0.41
2:C:340:ASN:O	2:C:351:LEU:HA	2.21	0.41
1:A:138:GLU:H	1:A:138:GLU:CD	2.24	0.41
2:C:569:TYR:CD2	2:C:597:VAL:HG12	2.56	0.41
1:B:196:PHE:CD1	1:B:196:PHE:N	2.89	0.41
1:A:145:LYS:HE2	1:A:148:ASP:OD2	2.20	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:50:CYS:SG	1:B:51:ASN:N	2.94	0.41
1:A:218:ASN:CG	1:A:218:ASN:O	2.58	0.41
2:C:601:LYS:HB2	2:C:601:LYS:HE3	1.86	0.41
2:D:328:TYR:OH	2:D:332:ILE:HD11	2.20	0.40
2:D:284:ASP:O	2:D:285:GLU:C	2.60	0.40
1:A:13:ALA:HB2	1:A:179:ILE:HD12	2.02	0.40
1:A:130:THR:HG23	4:A:610:NDP:P2B	2.61	0.40
2:C:588:ILE:HG23	2:C:589:SER:N	2.36	0.40
2:D:308:HIS:C	2:D:310:ASN:H	2.24	0.40
1:B:128:SER:HB3	1:B:131:LEU:HB2	2.03	0.40
1:B:143:ILE:CG2	1:B:144:ASN:N	2.83	0.40
2:C:345:ARG:NH1	2:C:345:ARG:HG2	2.36	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	215/280 (77%)	200 (93%)	10 (5%)	5 (2%)	10	24
1	B	220/280 (79%)	189 (86%)	28 (13%)	3 (1%)	16	41
2	C	323/328 (98%)	301 (93%)	18 (6%)	4 (1%)	19	45
2	D	324/328 (99%)	291 (90%)	26 (8%)	7 (2%)	10	25
All	All	1082/1216 (89%)	981 (91%)	82 (8%)	19 (2%)	13	31

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2	MET
1	A	28	LYS
1	B	117	LYS
2	C	306	SER
2	D	430	TYR
1	A	49	LYS

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Mol	Chain	Res	Type
2	C	305	ASN
2	C	310	ASN
2	C	430	TYR
2	D	282	ASP
2	D	309	PRO
1	A	202	GLU
2	D	306	SER
2	D	345	ARG
2	D	297	LYS
1	B	118	PRO
1	A	47	PRO
1	B	26	GLY
2	D	572	PRO

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	208/268 (78%)	204 (98%)	4 (2%)	69	92
1	B	213/268 (80%)	207 (97%)	6 (3%)	56	86
2	C	299/302 (99%)	293 (98%)	6 (2%)	68	92
2	D	302/302 (100%)	299 (99%)	3 (1%)	85	97
All	All	1022/1140 (90%)	1003 (98%)	19 (2%)	69	92

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

Mol	Chain	Res	Type
1	A	2	MET
1	A	22	SER
1	A	96	LYS
1	A	144	ASN
1	B	28	LYS
1	B	50	CYS
1	B	95	SER
1	B	138	GLU
1	B	148	ASP

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Mol	Chain	Res	Type
1	B	230	ASN
2	C	303	ASN
2	C	305	ASN
2	C	307	ILE
2	C	487	LEU
2	C	524	SER
2	C	606	MET
2	D	286	GLU
2	D	301	GLU
2	D	308	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	GLN
1	A	42	ASN
1	A	99	GLN
1	A	144	ASN
1	B	24	ASN
1	B	29	ASN
1	B	99	GLN
1	B	121	ASN
1	B	144	ASN
1	B	203	ASN
2	C	303	ASN
2	C	316	ASN
2	C	394	ASN
2	C	407	ASN
2	C	415	ASN
2	C	424	ASN
2	C	554	ASN
2	D	305	ASN
2	D	316	ASN
2	D	394	ASN
2	D	424	ASN
2	D	554	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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$
3	RJ1	A	609	-	18,18,18	1.28	2 (11%)	26,26,26	2.52	5 (19%)
4	NDP	A	610	-	52,52,52	1.71	9 (17%)	80,80,80	1.87	17 (21%)
3	RJ1	B	609	-	18,18,18	1.27	3 (16%)	26,26,26	2.84	6 (23%)
4	NDP	B	610	-	52,52,52	1.67	7 (13%)	80,80,80	1.98	17 (21%)
5	UMP	C	611	-	21,21,21	1.99	7 (33%)	26,31,31	1.79	9 (34%)
5	UMP	D	611	-	21,21,21	2.01	8 (38%)	26,31,31	1.85	9 (34%)

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	RJ1	A	609	-	-	0/14/14/14	0/1/1/1
4	NDP	A	610	-	-	0/35/77/77	0/3/5/5
3	RJ1	B	609	-	-	0/14/14/14	0/1/1/1
4	NDP	B	610	-	-	0/35/77/77	0/3/5/5
5	UMP	C	611	-	-	0/7/22/22	0/2/2/2
5	UMP	D	611	-	-	0/7/22/22	0/2/2/2

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	610	NDP	C4N-C5N	-5.05	1.38	1.49
4	A	610	NDP	C2N-C3N	4.98	1.44	1.34
4	B	610	NDP	C4N-C3N	-4.69	1.41	1.50
4	B	610	NDP	C2N-C3N	4.41	1.43	1.34
5	D	611	UMP	C2-N1	4.33	1.43	1.38
4	A	610	NDP	C4N-C5N	-4.29	1.39	1.49
4	A	610	NDP	C4N-C3N	-4.23	1.42	1.50
5	C	611	UMP	C2-N1	4.08	1.42	1.38
5	C	611	UMP	O4'-C1'	4.01	1.51	1.42
5	D	611	UMP	O4'-C1'	3.95	1.51	1.42
4	A	610	NDP	C8A-N9A	3.79	1.42	1.36
4	B	610	NDP	C8A-N9A	3.75	1.42	1.36
5	C	611	UMP	C4-N3	3.53	1.43	1.37
3	A	609	RJ1	C13-N17	-2.91	1.29	1.33
5	D	611	UMP	C6-C5	2.87	1.40	1.36
5	C	611	UMP	C6-C5	2.81	1.40	1.36
4	A	610	NDP	O4D-C4D	-2.78	1.38	1.45
5	D	611	UMP	P-OP3	-2.72	1.44	1.54
5	D	611	UMP	C4-N3	2.68	1.41	1.37
3	B	609	RJ1	C13-N17	-2.68	1.29	1.33
4	B	610	NDP	C4A-N3A	2.65	1.39	1.35
5	D	611	UMP	C5-C4	2.56	1.47	1.40
4	A	610	NDP	C2N-N1N	2.46	1.41	1.36
4	B	610	NDP	C3B-C2B	-2.36	1.47	1.53
4	A	610	NDP	C4A-N3A	2.35	1.39	1.35
5	C	611	UMP	P-OP3	-2.34	1.46	1.54
5	C	611	UMP	C5-C4	2.31	1.46	1.40
4	A	610	NDP	C3B-C2B	-2.25	1.47	1.53
4	A	610	NDP	PA-O1A	-2.24	1.42	1.51
4	B	610	NDP	PA-O1A	-2.23	1.42	1.51
3	A	609	RJ1	F10-C8	2.19	1.41	1.32
5	D	611	UMP	P-OP2	-2.14	1.46	1.54
5	D	611	UMP	O4'-C4'	2.07	1.49	1.45
3	B	609	RJ1	F10-C8	2.05	1.40	1.32
3	B	609	RJ1	F11-C8	2.04	1.40	1.32
5	C	611	UMP	P-OP2	-2.01	1.47	1.54

All (63) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	609	RJ1	C15-N14-C13	10.32	141.56	125.26
4	B	610	NDP	O4B-C1B-N9A	8.22	116.08	108.44
3	A	609	RJ1	C15-N14-C13	8.18	138.18	125.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	609	RJ1	C5-N12-C13	7.09	149.72	126.32
3	A	609	RJ1	C5-N12-C13	6.73	148.53	126.32
4	A	610	NDP	O4B-C1B-N9A	6.35	114.35	108.44
4	B	610	NDP	C4B-O4B-C1B	-5.23	104.06	109.75
4	A	610	NDP	C4B-O4B-C1B	-5.14	104.17	109.75
4	B	610	NDP	C1D-N1N-C2N	-4.82	112.81	121.02
5	C	611	UMP	O4'-C1'-N1	4.66	116.44	107.68
4	A	610	NDP	C3N-C2N-N1N	-4.66	116.44	123.05
5	D	611	UMP	O4'-C1'-N1	4.65	116.42	107.68
4	A	610	NDP	C1D-N1N-C2N	-4.50	113.35	121.02
4	B	610	NDP	N3A-C2A-N1A	-4.29	125.12	128.71
4	B	610	NDP	C3N-C2N-N1N	-4.28	116.98	123.05
4	A	610	NDP	C5N-C4N-C3N	4.23	124.07	112.60
4	B	610	NDP	C5N-C4N-C3N	4.17	123.89	112.60
3	B	609	RJ1	N12-C13-N14	3.87	126.48	117.24
4	A	610	NDP	N3A-C2A-N1A	-3.81	125.52	128.71
3	A	609	RJ1	N12-C13-N14	3.76	126.23	117.24
4	A	610	NDP	C3B-C2B-C1B	-3.73	95.47	102.73
4	B	610	NDP	C3B-C2B-C1B	-3.70	95.53	102.73
3	A	609	RJ1	C5-C4-CL7	3.66	124.32	119.45
3	B	609	RJ1	C5-C4-CL7	3.51	124.13	119.45
5	D	611	UMP	O4'-C1'-C2'	-3.08	100.17	106.25
4	A	610	NDP	C6N-N1N-C2N	3.06	123.44	119.44
4	A	610	NDP	C4N-C3N-C2N	-3.00	118.03	121.68
5	C	611	UMP	O4'-C1'-C2'	-2.98	100.38	106.25
4	B	610	NDP	C4N-C3N-C2N	-2.97	118.06	121.68
4	B	610	NDP	PN-O3-PA	2.97	140.39	131.68
4	A	610	NDP	PN-O3-PA	2.96	140.35	131.68
5	D	611	UMP	C6-N1-C2	2.86	123.50	119.51
4	A	610	NDP	C8A-N9A-C4A	-2.86	104.72	106.90
5	C	611	UMP	C6-N1-C2	2.78	123.38	119.51
4	A	610	NDP	C3D-C2D-C1D	-2.74	95.90	101.35
4	A	610	NDP	O7N-C7N-N7N	-2.70	116.12	122.93
4	B	610	NDP	O7N-C7N-N7N	-2.69	116.13	122.93
5	D	611	UMP	O4'-C4'-C3'	-2.68	98.85	105.66
4	B	610	NDP	C6N-N1N-C2N	2.66	122.92	119.44
5	D	611	UMP	C4'-O4'-C1'	2.65	116.15	109.44
4	B	610	NDP	C8A-N9A-C4A	-2.63	104.89	106.90
5	C	611	UMP	N3-C2-N1	-2.58	113.82	115.97
5	D	611	UMP	C2'-C1'-N1	2.56	120.73	114.08
4	B	610	NDP	O3B-C3B-C4B	2.52	118.49	111.08
5	D	611	UMP	C6-N1-C1'	-2.48	113.59	119.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	610	NDP	C3D-C2D-C1D	-2.44	96.50	101.35
5	C	611	UMP	C6-N1-C1'	-2.41	113.77	119.35
5	C	611	UMP	C4'-O4'-C1'	2.40	115.51	109.44
5	D	611	UMP	N3-C2-N1	-2.39	113.98	115.97
5	C	611	UMP	C2'-C1'-N1	2.36	120.23	114.08
4	A	610	NDP	O3B-C3B-C4B	2.35	118.01	111.08
4	A	610	NDP	O3B-C3B-C2B	2.35	118.11	111.20
4	A	610	NDP	C4N-C5N-C6N	-2.29	118.41	122.61
4	B	610	NDP	C4N-C5N-C6N	-2.28	118.42	122.61
3	A	609	RJ1	C6-C1-C8	2.25	122.70	119.61
4	B	610	NDP	O3B-C3B-C2B	2.23	117.77	111.20
3	B	609	RJ1	C6-C1-C8	2.21	122.66	119.61
3	B	609	RJ1	N14-C13-N17	-2.11	115.59	122.40
5	D	611	UMP	C2'-C3'-C4'	2.10	107.44	102.73
5	C	611	UMP	O4'-C4'-C3'	-2.08	100.40	105.66
4	A	610	NDP	C2D-C3D-C4D	2.07	106.77	102.65
4	B	610	NDP	C2D-C3D-C4D	2.05	106.75	102.65
5	C	611	UMP	C2'-C3'-C4'	2.01	107.23	102.73

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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	219/280 (78%)	0.16	17 (7%) 13 13	22, 37, 81, 90	0
1	B	224/280 (80%)	0.75	32 (14%) 3 4	29, 65, 90, 90	0
2	C	325/328 (99%)	-0.14	17 (5%) 26 29	18, 32, 88, 90	0
2	D	326/328 (99%)	-0.27	14 (4%) 34 38	20, 31, 86, 90	0
All	All	1094/1216 (89%)	0.07	80 (7%) 15 16	18, 36, 88, 90	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	25	GLU	7.3
2	C	607	ALA	6.4
1	A	1	MET	6.4
2	D	281	ASP	6.1
2	D	301	GLU	6.1
1	A	2	MET	6.0
2	C	302	LYS	6.0
2	D	282	ASP	5.8
1	B	26	GLY	5.3
2	C	300	GLU	5.2
1	A	24	ASN	5.2
1	A	23	LYS	5.1
2	D	299	LYS	5.1
2	C	608	ALA	4.9
1	B	231	ASN	4.9
1	B	2	MET	4.9
2	D	300	GLU	4.5
2	C	299	LYS	4.5
1	B	95	SER	4.3
2	C	301	GLU	4.2
1	B	22	SER	4.1

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Mol	Chain	Res	Type	RSRZ
1	B	136	PHE	3.9
1	B	135	ASP	3.9
1	A	25	GLU	3.9
1	B	24	ASN	3.8
1	B	29	ASN	3.8
2	D	283	ASP	3.8
2	D	303	ASN	3.8
2	C	303	ASN	3.7
1	B	85	THR	3.7
1	B	23	LYS	3.7
1	A	27	LYS	3.6
1	B	230	ASN	3.6
2	C	287	GLU	3.6
1	A	29	ASN	3.5
1	B	28	LYS	3.2
1	A	75	TYR	3.2
2	D	346	THR	3.2
1	B	117	LYS	3.2
2	C	286	GLU	3.1
2	C	346	THR	3.1
1	B	131	LEU	3.1
1	A	82	ASN	3.1
1	B	116	PHE	3.1
1	A	230	ASN	3.0
2	D	605	ASP	2.9
2	C	298	GLU	2.9
1	B	27	LYS	2.9
1	B	3	GLU	2.8
1	B	115	LYS	2.8
1	A	49	LYS	2.7
1	B	14	ILE	2.7
2	C	606	MET	2.7
1	B	88	ASN	2.7
2	C	306	SER	2.7
1	B	44	GLY	2.7
2	D	345	ARG	2.6
1	B	15	CYS	2.6
2	D	304	LYS	2.6
1	A	26	GLY	2.6
2	C	307	ILE	2.6
1	B	137	ASP	2.5
1	B	96	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
2	D	348	VAL	2.5
2	D	298	GLU	2.4
1	A	83	LYS	2.4
2	C	297	LYS	2.4
1	B	132	LYS	2.3
2	C	345	ARG	2.3
1	A	81	LEU	2.3
1	A	97	LYS	2.3
1	B	162	PHE	2.3
1	B	87	ASP	2.2
2	D	302	LYS	2.2
1	B	97	LYS	2.2
1	A	72	LYS	2.1
1	B	134	GLU	2.1
1	B	49	LYS	2.1
2	C	284	ASP	2.1
1	A	84	GLU	2.1

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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	RJ1	A	609	18/18	0.40	4.69	68,80,85,87	0
3	RJ1	B	609	18/18	0.47	3.93	86,89,90,90	0
5	UMP	D	611	20/20	0.14	-0.28	36,44,46,47	0
4	NDP	B	610	48/48	0.17	-0.57	64,72,85,86	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	UMP	C	611	20/20	0.12	-0.69	40,46,52,53	0
4	NDP	A	610	48/48	0.11	-1.11	28,33,39,40	0

6.5 Other polymers ⓘ

There are no such residues in this entry.