



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2014 – 12:46 AM GMT

PDB ID : 1RL3
Title : Crystal structure of cAMP-free R1a subunit of PKA
Authors : Wu, J.; Brown, S.; Xuong, N.-H.; Taylor, S.S.
Deposited on : 2003-11-24
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>

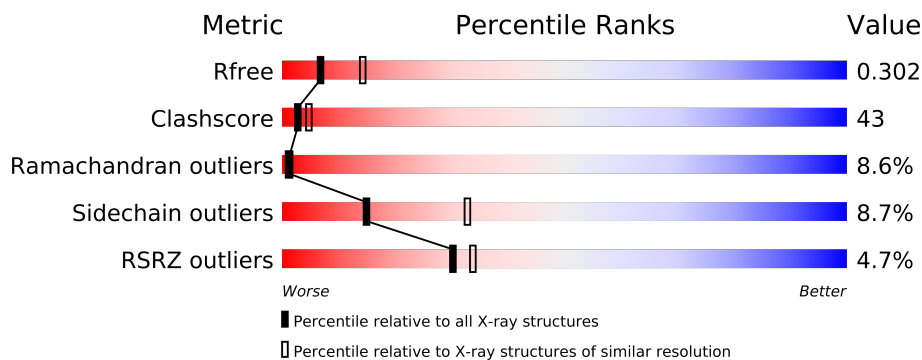
The following versions of software and data (see [references](#)) were used in the production of this report:

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	288	
1	B	288	

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	GOL	A	402	X	-
3	GOL	B	802	X	-
3	GOL	B	901	X	-

2 Entry composition i

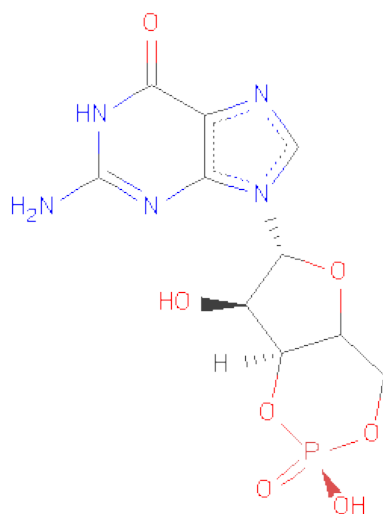
There are 4 unique types of molecules in this entry. The entry contains 4008 atoms, of which 0 are hydrogen and 0 are deuterium.

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 cAMP-dependent protein kinase type I-alpha regulatory chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	268	Total	C	N	O	S	0	0	0
			1991	1259	339	386	7			
1	B	259	Total	C	N	O	S	0	0	0
			1936	1232	325	372	7			

- Molecule 2 is CYCLIC GUANOSINE MONOPHOSPHATE (three-letter code: PCG) (formula: $C_{10}H_{12}N_5O_7P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 4 is water.

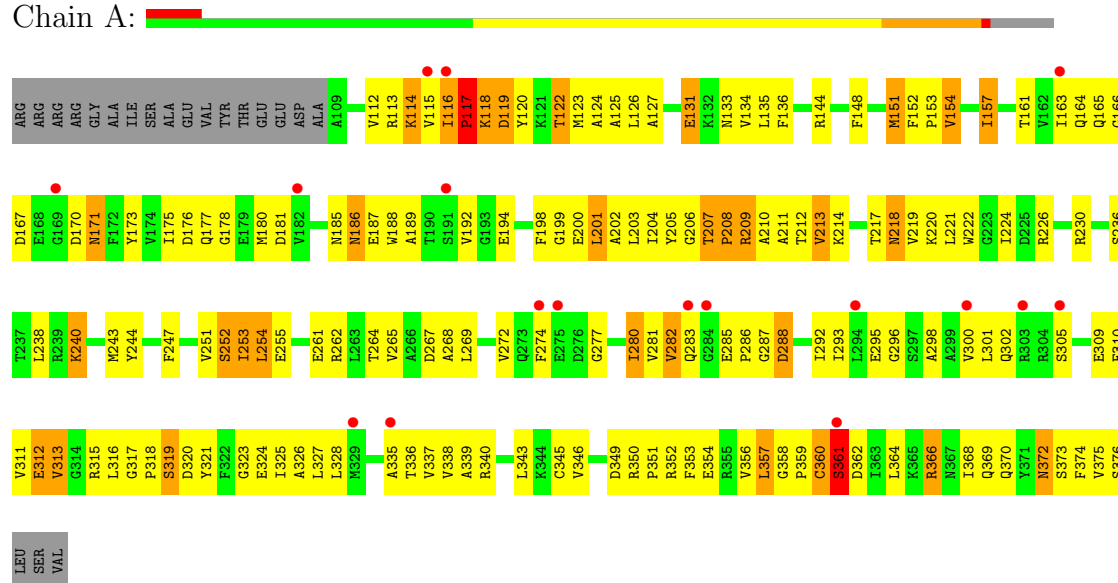
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	12	Total	O	0	0
			12	12		
4	B	5	Total	O	0	0
			5	5		

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.

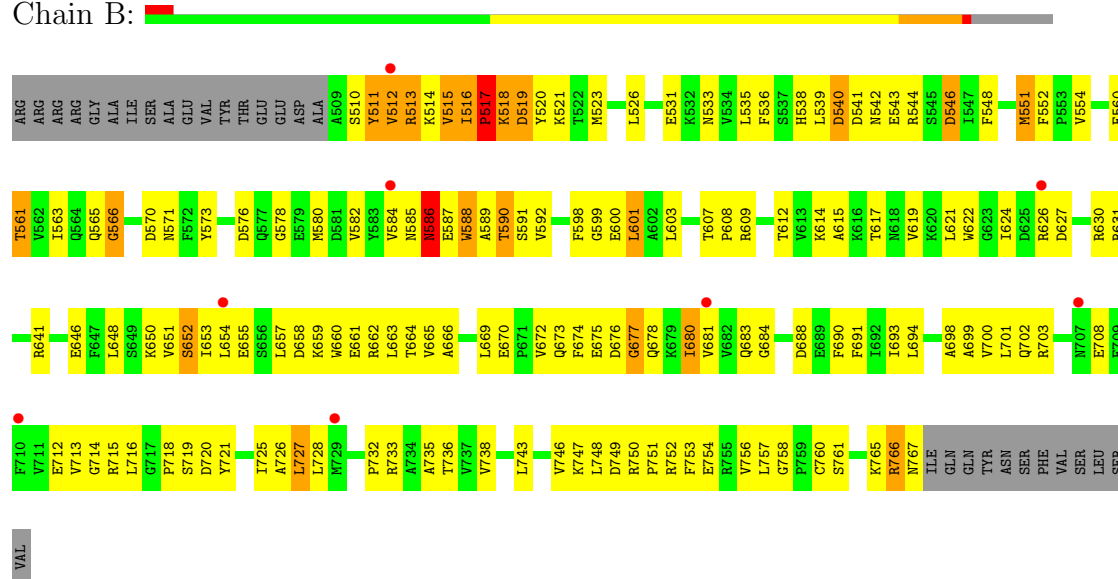
- Molecule 1: cAMP-dependent protein kinase type I-alpha regulatory chain

Chain A:



- Molecule 1: cAMP-dependent protein kinase type I-alpha regulatory chain

Chain B:



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	90.30Å 90.30Å 177.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.70 35.90 – 2.70	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.70) 87.5 (35.90-2.70)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.42 (at 2.68Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.240 , 0.285 0.263 , 0.302	Depositor DCC
R_{free} test set	1932 reflections (10.95%)	DCC
Wilson B-factor (Å ²)	58.6	Xtriage
Anisotropy	0.303	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 41.7	EDS
Estimated twinning fraction	0.094 for h,-h-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 21043 reflections	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	4008	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.04% 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: GOL, PCG

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	0.46	0/2029	0.78	2/2753 (0.1%)
1	B	0.45	0/1973	0.77	4/2674 (0.1%)
All	All	0.45	0/4002	0.77	6/5427 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	587	GLU	N-CA-CB	8.06	125.11	110.60
1	A	361	SER	CB-CA-C	-7.12	96.57	110.10
1	B	586	ASN	CB-CA-C	-7.07	96.27	110.40
1	B	587	GLU	CA-C-O	-6.93	105.54	120.10
1	B	587	GLU	CA-C-N	6.42	131.31	117.20
1	A	362	ASP	N-CA-C	-5.67	95.68	111.00

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	1991	0	1835	185	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1936	0	1820	153	0
2	A	23	0	11	3	0
2	B	23	0	11	2	0
3	A	6	0	4	2	0
3	B	12	0	8	2	0
4	A	12	0	0	1	0
4	B	5	0	0	0	0
All	All	4008	0	3689	334	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 43.

All (334) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:165:GLN:NE2	1:A:185:ASN:H	1.56	1.01
1:A:163:ILE:HB	1:A:213:VAL:HG23	1.43	1.00
1:A:316:LEU:HD22	1:A:320:ASP:HB3	1.46	0.97
1:B:756:VAL:HG23	1:B:757:LEU:HD13	1.49	0.95
1:A:165:GLN:HE22	1:A:185:ASN:N	1.65	0.94
1:A:300:VAL:HG13	1:A:335:ALA:HB1	1.49	0.93
1:B:585:ASN:O	1:B:586:ASN:HB2	1.67	0.92
1:A:133:ASN:OD1	1:A:135:LEU:HB2	1.71	0.91
1:A:350:ARG:HB3	1:A:351:PRO:HD3	1.54	0.89
1:B:630:ARG:HB3	1:B:630:ARG:NH1	1.88	0.88
1:A:199:GLY:HA2	2:A:401:PCG:O1A	1.73	0.88
1:A:117:PRO:HB3	1:B:517:PRO:HB3	1.54	0.87
1:A:163:ILE:CD1	1:A:209:ARG:HH21	1.92	0.82
1:A:171:ASN:CB	1:A:224:ILE:O	2.27	0.82
1:B:694:LEU:HD11	1:B:746:VAL:HG13	1.61	0.82
1:B:680:ILE:HG22	1:B:681:VAL:HG13	1.62	0.82
1:A:203:LEU:HD22	1:A:226:ARG:HB3	1.62	0.81
1:A:204:ILE:HG21	1:A:238:LEU:HD21	1.60	0.81
1:A:327:LEU:HA	1:A:350:ARG:HH22	1.45	0.81
1:A:217:THR:O	1:A:219:VAL:HG13	1.82	0.79
1:A:302:GLN:O	1:A:310:PHE:HA	1.82	0.79
1:A:114:LYS:O	1:A:116:ILE:HG22	1.80	0.79
1:B:630:ARG:HB3	1:B:630:ARG:HH11	1.45	0.78
1:A:165:GLN:HG3	1:A:211:ALA:HA	1.67	0.75
1:A:268:ALA:O	1:A:352:ARG:NH1	2.19	0.74
1:A:300:VAL:CG1	1:A:335:ALA:HB1	2.18	0.73
1:A:165:GLN:HE22	1:A:185:ASN:H	0.82	0.73
1:A:163:ILE:HG23	1:A:209:ARG:HE	1.55	0.72

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:163:ILE:HG23	1:A:209:ARG:NE	2.05	0.72
1:B:578:GLY:HA3	1:B:619:VAL:HG12	1.72	0.72
1:B:688:ASP:HB3	1:B:750:ARG:NH1	2.05	0.72
1:A:119:ASP:OD2	1:A:122:THR:HB	1.89	0.71
1:A:324:GLU:HB2	1:A:328:LEU:HD12	1.72	0.71
1:B:700:VAL:O	1:B:713:VAL:HG22	1.92	0.70
1:B:563:ILE:HD11	1:B:609:ARG:HE	1.57	0.70
1:B:540:ASP:OD1	1:B:543:GLU:HG3	1.92	0.69
1:A:171:ASN:HB2	1:A:224:ILE:O	1.92	0.69
1:A:366:ARG:HG2	1:A:366:ARG:HH11	1.56	0.69
1:B:750:ARG:HB3	1:B:751:PRO:HD3	1.74	0.69
1:A:361:SER:H	1:A:364:LEU:HG	1.58	0.69
1:A:272:VAL:HG13	1:A:272:VAL:O	1.93	0.69
1:B:603:LEU:HD22	1:B:626:ARG:HB3	1.76	0.68
1:A:178:GLY:HA3	1:A:219:VAL:HG12	1.76	0.68
1:A:205:TYR:HB2	1:A:207:THR:HG22	1.74	0.68
1:B:570:ASP:OD2	1:B:571:ASN:N	2.24	0.68
1:A:318:PRO:O	1:A:319:SER:HB2	1.93	0.68
1:B:646:GLU:HB3	1:B:650:LYS:NZ	2.10	0.67
1:B:570:ASP:H	1:B:609:ARG:HH21	1.43	0.67
1:B:571:ASN:HB3	1:B:624:ILE:O	1.95	0.67
1:A:180:MET:HG3	1:A:198:PHE:CE2	2.30	0.67
1:A:292:ILE:O	1:A:345:CYS:HB2	1.94	0.67
1:B:580:MET:HG3	1:B:598:PHE:CE2	2.30	0.66
1:A:171:ASN:HB3	1:A:224:ILE:O	1.94	0.66
1:B:526:LEU:HD11	1:B:551:MET:HE3	1.76	0.66
1:A:205:TYR:HB2	1:A:207:THR:CG2	2.25	0.65
1:B:590:THR:HG22	1:B:591:SER:H	1.60	0.65
1:A:315:ARG:HD3	1:A:316:LEU:H	1.62	0.65
1:A:366:ARG:HG2	1:A:366:ARG:NH1	2.11	0.65
1:A:163:ILE:HG12	1:A:209:ARG:HE	1.62	0.64
1:A:180:MET:HB2	1:A:192:VAL:HB	1.78	0.64
1:A:165:GLN:HG3	1:A:211:ALA:CA	2.27	0.64
1:B:518:LYS:O	1:B:519:ASP:O	2.15	0.64
1:B:600:GLU:OE2	1:B:601:LEU:HD13	1.98	0.64
1:A:372:ASN:N	1:A:372:ASN:HD22	1.95	0.64
1:B:580:MET:HG3	1:B:598:PHE:HE2	1.63	0.63
1:A:154:VAL:HG23	1:A:221:LEU:HB2	1.80	0.63
1:B:664:THR:HG22	1:B:756:VAL:HG12	1.81	0.63
1:B:670:GLU:OE2	1:B:747:LYS:HE3	1.99	0.63
1:A:163:ILE:HD11	1:A:209:ARG:HH21	1.63	0.62
1:B:566:GLY:HA2	1:B:608:PRO:HB2	1.80	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:264:THR:HG22	1:A:356:VAL:HG12	1.82	0.62
1:B:630:ARG:CB	1:B:630:ARG:HH11	2.11	0.62
1:B:765:LYS:O	1:B:767:ASN:N	2.31	0.62
1:A:164:GLN:HB3	1:A:167:ASP:OD2	1.99	0.61
1:A:349:ASP:OD2	1:A:351:PRO:HD2	2.00	0.61
1:A:315:ARG:HD3	1:A:316:LEU:N	2.16	0.60
2:A:401:PCG:H3'	2:A:401:PCG:N3	2.17	0.60
1:A:122:THR:HG22	1:A:123:MET:N	2.16	0.60
1:A:316:LEU:HD22	1:A:320:ASP:CB	2.28	0.60
1:B:652:SER:O	1:B:655:GLU:HG2	2.02	0.60
1:B:702:GLN:HB3	1:B:713:VAL:HG11	1.82	0.60
1:A:175:ILE:CG2	1:A:194:GLU:HA	2.31	0.60
1:B:698:ALA:HB1	1:B:738:VAL:O	2.02	0.60
1:B:565:GLN:HE22	1:B:585:ASN:H	1.50	0.59
1:B:702:GLN:HG3	1:B:703:ARG:N	2.17	0.59
1:A:157:ILE:HA	1:A:218:ASN:HA	1.84	0.59
1:A:358:GLY:N	1:A:359:PRO:CD	2.65	0.59
1:B:749:ASP:OD2	1:B:752:ARG:HD3	2.02	0.59
1:B:584:VAL:HG23	1:B:589:ALA:HB2	1.85	0.59
1:B:660:TRP:HA	1:B:663:LEU:HD12	1.85	0.59
1:B:563:ILE:HD11	1:B:609:ARG:NE	2.17	0.59
1:A:360:CYS:O	1:A:361:SER:HB3	2.02	0.59
1:B:516:ILE:HD13	1:B:552:PHE:HB3	1.85	0.58
1:A:163:ILE:HD12	1:A:213:VAL:HG21	1.84	0.58
1:B:761:SER:O	1:B:765:LYS:HG3	2.03	0.58
1:A:282:VAL:CB	1:A:336:THR:HA	2.34	0.58
1:A:133:ASN:HB3	1:A:136:PHE:CD2	2.39	0.58
1:A:180:MET:HG3	1:A:198:PHE:HE2	1.68	0.58
1:B:680:ILE:H	1:B:680:ILE:HD12	1.69	0.57
1:B:651:VAL:O	1:B:653:ILE:N	2.38	0.57
1:B:607:THR:HB	1:B:608:PRO:HD2	1.85	0.57
1:A:300:VAL:HA	1:A:336:THR:O	2.04	0.57
1:A:364:LEU:O	1:A:368:ILE:HG12	2.05	0.57
1:B:542:ASN:HD22	1:B:542:ASN:N	2.02	0.57
1:B:546:ASP:N	1:B:546:ASP:OD2	2.38	0.57
1:B:750:ARG:O	1:B:754:GLU:HG3	2.05	0.57
1:B:565:GLN:HB2	1:B:612:THR:OG1	2.05	0.56
1:B:752:ARG:O	1:B:756:VAL:HG22	2.05	0.56
1:A:176:ASP:OD2	1:A:222:TRP:NE1	2.36	0.56
1:B:533:ASN:OD1	1:B:535:LEU:HB2	2.05	0.56
1:B:698:ALA:HB2	1:B:743:LEU:HD22	1.87	0.56
1:B:699:ALA:HB1	1:B:712:GLU:HG2	1.88	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:327:LEU:HA	1:A:350:ARG:NH2	2.18	0.56
1:A:261:GLU:O	1:A:264:THR:HB	2.06	0.55
1:B:582:VAL:HB	1:B:590:THR:H	1.71	0.55
1:B:761:SER:OG	1:B:765:LYS:HE3	2.07	0.55
1:B:683:GLN:HA	1:B:733:ARG:O	2.06	0.55
1:A:324:GLU:O	1:A:328:LEU:HB2	2.07	0.54
1:A:366:ARG:CG	1:A:366:ARG:HH11	2.19	0.54
1:B:718:PRO:O	1:B:719:SER:HB3	2.06	0.54
1:A:202:ALA:O	1:A:206:GLY:N	2.39	0.54
1:A:350:ARG:HB3	1:A:351:PRO:CD	2.34	0.54
1:B:674:PHE:HD1	1:B:678:GLN:HE21	1.56	0.54
1:B:571:ASN:C	1:B:603:LEU:HD11	2.29	0.54
1:B:588:TRP:C	1:B:588:TRP:CD1	2.80	0.54
1:B:766:ARG:O	1:B:767:ASN:ND2	2.40	0.54
1:B:669:LEU:HD22	1:B:746:VAL:HB	1.89	0.54
1:A:116:ILE:O	1:A:116:ILE:HG12	2.08	0.54
1:B:561:THR:OG1	1:B:614:LYS:HG2	2.08	0.53
1:B:560:GLU:HA	3:B:901:GOL:H12	1.90	0.53
1:B:657:LEU:HD23	1:B:661:GLU:HB3	1.89	0.53
1:B:511:TYR:CD1	1:B:512:VAL:N	2.77	0.53
1:A:163:ILE:HG23	1:A:209:ARG:CD	2.38	0.53
1:A:163:ILE:CG2	1:A:209:ARG:HE	2.20	0.53
1:B:571:ASN:HB2	1:B:573:TYR:CE2	2.43	0.53
1:B:533:ASN:HB3	1:B:536:PHE:HD2	1.74	0.53
1:A:282:VAL:CB	1:A:337:VAL:H	2.22	0.53
1:B:663:LEU:O	1:B:666:ALA:HB3	2.08	0.53
1:A:157:ILE:HG22	1:A:218:ASN:OD1	2.08	0.53
1:B:510:SER:O	1:B:511:TYR:CB	2.57	0.52
1:A:277:GLY:H	1:A:339:ALA:HB3	1.73	0.52
1:B:749:ASP:CG	1:B:751:PRO:HD2	2.30	0.52
1:B:516:ILE:HD13	1:B:552:PHE:CB	2.38	0.52
1:A:209:ARG:O	1:A:211:ALA:N	2.42	0.52
1:A:163:ILE:O	1:A:212:THR:HG23	2.10	0.52
1:B:512:VAL:CB	1:B:631:ARG:CZ	2.88	0.52
1:B:680:ILE:N	1:B:680:ILE:HD12	2.26	0.51
1:A:116:ILE:HD13	1:A:152:PHE:HB3	1.93	0.51
1:A:200:GLU:HG2	1:A:201:LEU:N	2.23	0.51
1:A:209:ARG:HH11	1:A:209:ARG:HG2	1.75	0.51
1:A:350:ARG:HD2	1:A:354:GLU:OE2	2.11	0.51
1:A:353:PHE:CE1	1:A:357:LEU:HD23	2.46	0.51
1:B:691:PHE:CD1	1:B:691:PHE:N	2.79	0.51
1:A:252:SER:HB2	1:A:321:TYR:HE2	1.75	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:684:GLY:O	1:B:732:PRO:HB3	2.11	0.51
1:A:302:GLN:H	1:A:311:VAL:N	2.09	0.51
1:A:298:ALA:HB1	1:A:338:VAL:O	2.11	0.51
1:B:665:VAL:O	1:B:669:LEU:HG	2.11	0.51
1:B:565:GLN:N	1:B:612:THR:OG1	2.33	0.50
1:A:277:GLY:H	1:A:339:ALA:CB	2.23	0.50
1:A:118:LYS:HE2	1:A:148:PHE:O	2.12	0.50
1:A:181:ASP:O	1:A:213:VAL:HA	2.12	0.50
1:B:614:LYS:HE2	3:B:901:GOL:O3	2.12	0.50
1:A:170:ASP:N	1:A:170:ASP:OD2	2.43	0.50
1:A:122:THR:O	1:A:125:ALA:HB3	2.12	0.50
1:A:204:ILE:CG2	1:A:238:LEU:HD21	2.37	0.50
1:B:727:LEU:HB3	1:B:753:PHE:CE2	2.46	0.49
1:B:565:GLN:HE22	1:B:585:ASN:N	2.10	0.49
1:A:118:LYS:HB2	1:A:118:LYS:NZ	2.27	0.49
1:B:510:SER:O	1:B:511:TYR:HB3	2.12	0.49
1:A:369:GLN:OE1	1:A:370:GLN:OE1	2.30	0.49
1:A:205:TYR:CB	1:A:207:THR:HG22	2.41	0.49
1:A:209:ARG:NH1	1:A:209:ARG:HG2	2.27	0.49
1:B:688:ASP:C	1:B:727:LEU:HD21	2.32	0.49
1:A:117:PRO:O	1:A:118:LYS:HB2	2.13	0.49
1:B:680:ILE:H	1:B:680:ILE:CD1	2.24	0.49
1:A:268:ALA:HB1	1:A:352:ARG:HD2	1.93	0.49
1:A:119:ASP:C	1:A:119:ASP:OD2	2.49	0.49
1:A:144:ARG:HG3	1:A:144:ARG:HH21	1.77	0.49
1:B:627:ASP:O	1:B:631:ARG:HB2	2.13	0.49
1:B:688:ASP:HB3	1:B:750:ARG:CZ	2.42	0.49
1:A:126:LEU:HD11	1:A:151:MET:CE	2.43	0.49
1:A:356:VAL:HG23	1:A:357:LEU:HD13	1.95	0.49
1:A:247:PHE:HZ	1:A:293:ILE:O	1.95	0.49
1:B:519:ASP:CG	1:B:520:TYR:H	2.15	0.49
1:A:163:ILE:CD1	1:A:209:ARG:NH2	2.69	0.48
1:A:185:ASN:O	1:A:187:GLU:N	2.46	0.48
1:A:207:THR:O	1:A:208:PRO:O	2.31	0.48
1:A:265:VAL:O	1:A:269:LEU:HG	2.13	0.48
1:A:133:ASN:HB3	1:A:136:PHE:HD2	1.78	0.48
1:B:666:ALA:HA	1:B:669:LEU:HD12	1.94	0.48
1:B:726:ALA:C	1:B:728:LEU:H	2.16	0.48
1:A:288:ASP:HB3	1:A:350:ARG:HH11	1.79	0.48
1:B:746:VAL:O	1:B:746:VAL:HG23	2.13	0.48
1:A:175:ILE:HG21	1:A:194:GLU:HA	1.94	0.48
1:A:177:GLN:HA	1:A:194:GLU:OE1	2.13	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:153:PRO:O	1:A:154:VAL:HG13	2.13	0.48
1:A:243:MET:HG2	1:A:243:MET:O	2.13	0.48
1:B:563:ILE:HG13	1:B:609:ARG:HG3	1.94	0.48
1:B:571:ASN:CB	1:B:624:ILE:O	2.61	0.48
1:A:317:GLY:N	1:A:320:ASP:OD2	2.30	0.48
1:B:747:LYS:O	1:B:748:LEU:HD12	2.14	0.48
1:B:615:ALA:HB1	1:B:617:THR:O	2.14	0.48
1:B:658:ASP:OD2	1:B:661:GLU:HG3	2.14	0.47
1:B:588:TRP:CH2	1:B:591:SER:HB2	2.49	0.47
1:A:312:GLU:O	1:A:313:VAL:CB	2.62	0.47
1:B:681:VAL:O	1:B:736:THR:HG23	2.14	0.47
1:B:570:ASP:H	1:B:609:ARG:NH2	2.09	0.47
1:B:563:ILE:CD1	1:B:609:ARG:NE	2.78	0.47
1:B:540:ASP:CG	1:B:543:GLU:HG3	2.35	0.47
1:B:661:GLU:O	1:B:664:THR:HB	2.15	0.47
1:B:516:ILE:O	1:B:516:ILE:HG12	2.14	0.47
1:B:542:ASN:N	1:B:542:ASN:ND2	2.62	0.47
1:A:126:LEU:HD13	1:A:222:TRP:CE3	2.49	0.47
1:A:119:ASP:O	1:A:123:MET:HB2	2.15	0.46
1:A:171:ASN:HD22	1:A:173:TYR:HE2	1.61	0.46
1:A:226:ARG:O	1:A:230:ARG:HG3	2.15	0.46
1:A:293:ILE:HA	1:A:345:CYS:HB3	1.96	0.46
1:A:301:LEU:HA	1:A:312:GLU:N	2.29	0.46
1:B:514:LYS:O	1:B:515:VAL:C	2.54	0.46
1:B:521:LYS:HD2	1:B:521:LYS:HA	1.60	0.46
1:B:641:ARG:HD3	1:B:663:LEU:HD22	1.97	0.46
1:A:127:ALA:O	1:A:131:GLU:HB3	2.15	0.46
1:A:171:ASN:ND2	1:A:173:TYR:CE2	2.83	0.46
1:A:323:GLY:O	1:A:324:GLU:C	2.53	0.46
1:B:672:VAL:HG11	1:B:747:LYS:HE2	1.97	0.46
1:A:240:LYS:HD2	1:A:244:TYR:HE2	1.81	0.46
1:A:220:LYS:O	1:A:221:LEU:HD23	2.16	0.46
1:B:727:LEU:HD13	1:B:753:PHE:CD2	2.50	0.46
1:B:646:GLU:HB3	1:B:650:LYS:HZ1	1.80	0.46
1:A:188:TRP:CG	1:A:189:ALA:N	2.82	0.46
1:A:280:ILE:C	1:A:282:VAL:H	2.20	0.46
1:A:262:ARG:O	1:A:265:VAL:HB	2.16	0.46
1:B:690:PHE:CE1	1:B:721:TYR:HB2	2.50	0.46
1:A:269:LEU:HD22	1:A:346:VAL:HB	1.98	0.46
1:A:326:ALA:N	3:A:402:GOL:O3	2.41	0.46
1:B:648:LEU:HD22	1:B:654:LEU:CD1	2.45	0.46
1:A:113:ARG:O	1:A:114:LYS:HB2	2.15	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:552:PHE:HE1	1:B:554:VAL:HG11	1.81	0.45
1:B:701:LEU:O	1:B:735:ALA:HB1	2.16	0.45
1:A:209:ARG:C	1:A:211:ALA:H	2.19	0.45
1:A:315:ARG:O	1:A:316:LEU:HG	2.17	0.45
1:A:262:ARG:HH11	1:A:262:ARG:CB	2.30	0.45
1:A:375:VAL:O	1:A:376:SER:HB3	2.16	0.45
1:B:748:LEU:HD23	1:B:753:PHE:HA	1.98	0.45
1:B:519:ASP:CG	1:B:520:TYR:N	2.70	0.45
1:B:511:TYR:CG	1:B:512:VAL:N	2.83	0.45
1:B:727:LEU:HD23	1:B:750:ARG:HD3	1.98	0.45
1:B:542:ASN:O	1:B:546:ASP:OD2	2.34	0.45
1:A:369:GLN:NE2	1:A:373:SER:HB2	2.32	0.45
1:B:716:LEU:HA	1:B:720:ASP:OD2	2.16	0.45
1:B:674:PHE:CZ	1:B:680:ILE:HG13	2.52	0.45
1:A:272:VAL:HG22	1:A:274:PHE:CE1	2.52	0.45
1:A:247:PHE:CZ	1:A:293:ILE:O	2.70	0.45
1:B:601:LEU:HB2	2:B:801:PCG:O3'	2.16	0.45
1:A:120:TYR:OH	1:B:544:ARG:CZ	2.65	0.45
1:B:758:GLY:C	1:B:760:CYS:H	2.21	0.44
1:A:288:ASP:HB3	1:A:350:ARG:HE	1.82	0.44
1:B:565:GLN:NE2	1:B:585:ASN:H	2.14	0.44
1:A:157:ILE:HG22	1:A:218:ASN:HB3	1.98	0.44
1:A:144:ARG:HG3	1:A:144:ARG:NH2	2.32	0.44
1:B:580:MET:HB2	1:B:592:VAL:HB	2.00	0.44
1:B:657:LEU:HD11	1:B:760:CYS:SG	2.57	0.44
1:A:353:PHE:O	1:A:354:GLU:C	2.56	0.44
1:A:117:PRO:CB	1:B:517:PRO:HB3	2.35	0.44
1:A:134:VAL:HG12	1:A:134:VAL:O	2.17	0.44
1:B:584:VAL:C	1:B:586:ASN:H	2.20	0.44
1:B:563:ILE:HG12	1:B:609:ARG:CD	2.47	0.44
1:A:200:GLU:OE1	1:A:200:GLU:N	2.47	0.44
1:A:352:ARG:O	1:A:356:VAL:HG22	2.18	0.43
1:B:675:GLU:O	1:B:677:GLY:N	2.44	0.43
1:B:538:HIS:CD2	1:B:538:HIS:N	2.85	0.43
1:A:163:ILE:HG21	2:A:401:PCG:H5'2	1.99	0.43
1:A:262:ARG:HH11	1:A:262:ARG:HB3	1.82	0.43
1:A:311:VAL:HG13	1:A:311:VAL:O	2.19	0.43
1:A:318:PRO:O	1:A:319:SER:CB	2.62	0.43
1:A:277:GLY:N	1:A:339:ALA:HB3	2.33	0.43
1:A:135:LEU:HD21	1:A:267:ASP:HB3	1.99	0.43
1:A:163:ILE:HG23	1:A:209:ARG:HD3	2.00	0.43
1:B:599:GLY:HA2	2:B:801:PCG:O1A	2.19	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:512:VAL:O	1:B:513:ARG:CB	2.66	0.43
1:A:214:LYS:NZ	4:A:906:HOH:O	2.52	0.43
1:A:122:THR:CG2	1:A:123:MET:N	2.81	0.43
1:A:247:PHE:O	1:A:251:VAL:HG23	2.19	0.43
1:B:539:LEU:HD13	1:B:544:ARG:HA	2.01	0.43
1:A:328:LEU:HD11	1:A:364:LEU:CD1	2.49	0.43
1:A:240:LYS:HD2	1:A:244:TYR:CE2	2.54	0.43
1:A:126:LEU:HD11	1:A:151:MET:HE3	2.00	0.42
1:A:285:GLU:HA	1:A:286:PRO:HD3	1.81	0.42
1:B:693:ILE:HG23	1:B:743:LEU:HD11	2.01	0.42
1:A:369:GLN:HE21	1:A:373:SER:HB2	1.84	0.42
1:B:664:THR:CG2	1:B:756:VAL:HG12	2.46	0.42
1:B:563:ILE:HG12	1:B:609:ARG:HD2	2.02	0.42
1:A:207:THR:HA	1:A:208:PRO:HD2	1.76	0.42
1:A:118:LYS:HB3	1:A:122:THR:HG22	2.00	0.42
1:B:675:GLU:H	1:B:678:GLN:CD	2.22	0.42
1:A:120:TYR:HA	1:B:523:MET:CE	2.49	0.42
1:B:749:ASP:O	1:B:750:ARG:C	2.58	0.42
1:A:253:ILE:C	1:A:255:GLU:H	2.23	0.42
1:A:200:GLU:CD	1:A:201:LEU:HD13	2.39	0.42
1:B:563:ILE:O	1:B:563:ILE:HG23	2.19	0.42
1:A:173:TYR:HB2	1:A:198:PHE:CE1	2.54	0.42
1:A:116:ILE:O	1:A:116:ILE:CG1	2.67	0.42
1:B:715:ARG:C	1:B:716:LEU:HG	2.40	0.42
1:A:288:ASP:OD1	1:A:288:ASP:N	2.52	0.42
1:A:288:ASP:HB3	1:A:350:ARG:NH1	2.34	0.42
1:B:700:VAL:HG23	1:B:714:GLY:O	2.20	0.42
1:B:757:LEU:O	1:B:760:CYS:HB2	2.20	0.41
1:A:277:GLY:CA	1:A:339:ALA:HB3	2.50	0.41
1:B:659:LYS:O	1:B:662:ARG:HB3	2.19	0.41
1:A:118:LYS:HB3	1:A:122:THR:CG2	2.49	0.41
1:A:185:ASN:O	1:A:186:ASN:C	2.58	0.41
1:B:554:VAL:HG23	1:B:621:LEU:HB2	2.02	0.41
1:A:327:LEU:CA	1:A:350:ARG:HH22	2.23	0.41
1:A:372:ASN:ND2	1:A:372:ASN:N	2.67	0.41
1:B:585:ASN:O	1:B:586:ASN:CB	2.50	0.41
1:B:526:LEU:HD21	1:B:548:PHE:HA	2.03	0.41
1:B:576:ASP:HB2	1:B:622:TRP:CD1	2.55	0.41
1:A:253:ILE:HG13	1:A:254:LEU:N	2.36	0.41
1:A:254:LEU:HD22	1:A:357:LEU:HD11	2.02	0.41
1:B:684:GLY:C	1:B:732:PRO:HB3	2.41	0.41
1:A:188:TRP:CD1	1:A:189:ALA:N	2.89	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:326:ALA:HB2	3:A:402:GOL:O3	2.21	0.41
1:A:152:PHE:HE1	1:A:154:VAL:CG1	2.33	0.40
1:B:688:ASP:HB2	1:B:727:LEU:CD2	2.51	0.40
1:B:526:LEU:HD11	1:B:551:MET:CE	2.47	0.40
1:A:163:ILE:HD13	1:A:209:ARG:NH2	2.36	0.40
1:A:324:GLU:HG2	1:A:325:ILE:H	1.86	0.40
1:A:296:GLY:HA3	1:A:343:LEU:HA	2.04	0.40
1:A:117:PRO:HB3	1:B:517:PRO:CB	2.40	0.40
1:A:124:ALA:O	1:A:125:ALA:C	2.60	0.40
1:B:552:PHE:CD1	1:B:552:PHE:O	2.75	0.40
1:A:188:TRP:CD1	1:A:188:TRP:C	2.93	0.40
1:B:721:TYR:C	1:B:721:TYR:CD1	2.95	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	266/288 (92%)	196 (74%)	42 (16%)	28 (10%)	1	0
1	B	257/288 (89%)	210 (82%)	30 (12%)	17 (7%)	2	2
All	All	523/576 (91%)	406 (78%)	72 (14%)	45 (9%)	1	1

All (45) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	114	LYS
1	A	117	PRO
1	A	186	ASN
1	A	208	PRO
1	A	210	ALA
1	A	280	ILE
1	A	309	GLU
1	A	312	GLU

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Mol	Chain	Res	Type
1	A	313	VAL
1	A	340	ARG
1	B	511	TYR
1	B	512	VAL
1	B	513	ARG
1	B	515	VAL
1	B	517	PRO
1	B	519	ASP
1	B	586	ASN
1	B	652	SER
1	A	115	VAL
1	A	166	GLY
1	A	252	SER
1	A	282	VAL
1	A	305	SER
1	B	676	ASP
1	B	766	ARG
1	A	112	VAL
1	A	118	LYS
1	A	209	ARG
1	A	319	SER
1	A	360	CYS
1	B	708	GLU
1	A	119	ASP
1	A	281	VAL
1	A	283	GLN
1	A	295	GLU
1	A	374	PHE
1	B	518	LYS
1	B	566	GLY
1	A	254	LEU
1	A	287	GLY
1	A	361	SER
1	B	680	ILE
1	B	727	LEU
1	B	677	GLY
1	B	725	ILE

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	191/244 (78%)	171 (90%)	20 (10%)	10	22
1	B	188/244 (77%)	175 (93%)	13 (7%)	22	48
All	All	379/488 (78%)	346 (91%)	33 (9%)	15	33

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	116	ILE
1	A	117	PRO
1	A	122	THR
1	A	131	GLU
1	A	151	MET
1	A	154	VAL
1	A	157	ILE
1	A	161	THR
1	A	171	ASN
1	A	201	LEU
1	A	207	THR
1	A	213	VAL
1	A	218	ASN
1	A	236	SER
1	A	240	LYS
1	A	253	ILE
1	A	288	ASP
1	A	357	LEU
1	A	366	ARG
1	A	372	ASN
1	B	516	ILE
1	B	517	PRO
1	B	531	GLU
1	B	540	ASP
1	B	541	ASP
1	B	546	ASP
1	B	551	MET
1	B	561	THR
1	B	586	ASN
1	B	588	TRP
1	B	590	THR
1	B	601	LEU
1	B	673	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	142	ASN
1	A	165	GLN
1	A	171	ASN
1	A	177	GLN
1	A	369	GLN
1	A	372	ASN
1	B	538	HIS
1	B	542	ASN
1	B	564	GLN
1	B	565	GLN
1	B	577	GLN
1	B	585	ASN
1	B	678	GLN
1	B	767	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 ⓘ

5 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	PCG	A	401	-	26,26,26	2.49	12 (46%)	38,41,41	3.70	14 (36%)
3	GOL	A	402	-	5,5,5	4.71	5 (100%)	5,5,5	5.73	3 (60%)
2	PCG	B	801	-	26,26,26	2.82	13 (50%)	38,41,41	3.33	16 (42%)
3	GOL	B	802	-	5,5,5	4.48	5 (100%)	5,5,5	5.74	3 (60%)
3	GOL	B	901	-	5,5,5	4.44	5 (100%)	5,5,5	5.77	3 (60%)

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	PCG	A	401	-	-	0/4/31/31	0/0/4/4
3	GOL	A	402	-	-	0/4/4/4	0/0/0/0
2	PCG	B	801	-	-	0/4/31/31	0/0/4/4
3	GOL	B	802	-	-	0/4/4/4	0/0/0/0
3	GOL	B	901	-	-	0/4/4/4	0/0/0/0

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	GOL	C3-C2	-7.91	1.19	1.52
3	B	802	GOL	C3-C2	-7.44	1.21	1.52
3	B	901	GOL	C3-C2	-7.35	1.22	1.52
2	B	801	PCG	C4-N9	5.99	1.46	1.37
2	A	401	PCG	C4-N9	5.64	1.46	1.37
2	B	801	PCG	C4-N3	5.59	1.45	1.35
2	B	801	PCG	O3'-C3'	4.98	1.52	1.44
3	A	402	GOL	O1-C1	4.64	1.62	1.42
2	A	401	PCG	C4-N3	4.53	1.43	1.35
2	B	801	PCG	PA-O5'	-4.52	1.49	1.58
3	B	802	GOL	O1-C1	4.51	1.62	1.42
3	B	901	GOL	O1-C1	4.36	1.61	1.42
2	B	801	PCG	C6-C5	4.24	1.48	1.41
2	A	401	PCG	PA-O5'	-4.20	1.50	1.58
2	A	401	PCG	PA-O3'	-4.13	1.50	1.58
2	B	801	PCG	C6-N1	3.86	1.43	1.37
2	A	401	PCG	O3'-C3'	3.72	1.50	1.44
2	B	801	PCG	C2-N3	3.60	1.38	1.33
3	B	901	GOL	O3-C3	3.34	1.56	1.42
2	A	401	PCG	C6-N1	3.34	1.42	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	802	GOL	O3-C3	3.29	1.56	1.42
2	A	401	PCG	C2-N1	3.24	1.41	1.36
3	A	402	GOL	O2-C2	-3.21	1.33	1.43
2	A	401	PCG	C6-C5	3.16	1.46	1.41
2	B	801	PCG	C5'-C4'	-2.93	1.46	1.51
3	A	402	GOL	C1-C2	-2.90	1.40	1.52
3	A	402	GOL	O3-C3	2.87	1.54	1.42
3	B	901	GOL	C1-C2	-2.81	1.40	1.52
2	B	801	PCG	C2-N1	2.80	1.41	1.36
2	A	401	PCG	C2-N3	2.74	1.37	1.33
3	B	802	GOL	C1-C2	-2.69	1.41	1.52
3	B	802	GOL	O2-C2	-2.59	1.35	1.43
3	B	901	GOL	O2-C2	-2.53	1.35	1.43
2	B	801	PCG	C5-C4	2.28	1.45	1.40
2	B	801	PCG	C2-N2	2.25	1.35	1.32
2	B	801	PCG	PA-O3'	-2.23	1.53	1.58
2	A	401	PCG	C8-N9	2.18	1.39	1.36
2	A	401	PCG	C5-C4	2.16	1.45	1.40
2	A	401	PCG	C2-N2	2.09	1.35	1.32
2	B	801	PCG	C8-N9	2.02	1.39	1.36

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	901	GOL	O3-C3-C2	10.45	160.67	109.71
3	A	402	GOL	O3-C3-C2	10.44	160.64	109.71
3	B	802	GOL	O3-C3-C2	10.28	159.86	109.71
2	B	801	PCG	O3'-C3'-C2'	10.23	127.39	115.62
2	A	401	PCG	O1A-PA-O3'	9.76	127.41	107.39
2	A	401	PCG	O5'-PA-O3'	-9.53	92.73	105.81
2	A	401	PCG	C6-C5-N7	-9.15	132.91	134.14
2	A	401	PCG	O3'-C3'-C2'	8.74	125.68	115.62
2	B	801	PCG	O1A-PA-O3'	8.04	123.89	107.39
3	B	802	GOL	O2-C2-C3	6.78	139.12	108.22
3	B	901	GOL	O2-C2-C3	6.70	138.76	108.22
3	A	402	GOL	O2-C2-C3	6.61	138.33	108.22
2	B	801	PCG	C6-C5-N7	-6.40	133.28	134.14
2	B	801	PCG	C4'-O4'-C1'	-5.96	103.28	109.75
2	B	801	PCG	C2'-C3'-C4'	-5.01	93.42	103.16
2	A	401	PCG	C4'-O4'-C1'	-4.96	104.36	109.75
2	A	401	PCG	C2'-C3'-C4'	-4.76	93.90	103.16
2	A	401	PCG	C1'-N9-C4	-4.66	118.58	126.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	PCG	C1'-N9-C4	-4.50	118.87	126.64
2	B	801	PCG	C6-N1-C2	4.09	126.67	119.51
2	B	801	PCG	O5'-PA-O3'	-4.07	100.23	105.81
2	A	401	PCG	C8-N9-C1'	4.05	134.36	126.38
2	A	401	PCG	C6-N1-C2	3.92	126.37	119.51
2	A	401	PCG	O3'-C3'-C4'	3.88	113.84	110.73
2	B	801	PCG	C8-N9-C1'	3.82	133.91	126.38
3	B	802	GOL	O1-C1-C2	3.60	127.25	109.71
2	B	801	PCG	O3'-C3'-C4'	3.51	113.55	110.73
3	B	901	GOL	O1-C1-C2	3.39	126.24	109.71
3	A	402	GOL	O1-C1-C2	3.28	125.69	109.71
2	A	401	PCG	O2'-C2'-C3'	3.23	120.69	111.20
2	B	801	PCG	O4'-C4'-C5'	3.21	122.08	111.97
2	B	801	PCG	C5'-C4'-C3'	-2.83	106.64	112.62
2	B	801	PCG	O2'-C2'-C3'	2.60	118.85	111.20
2	A	401	PCG	O4'-C4'-C5'	2.56	120.05	111.97
2	B	801	PCG	C3'-C2'-C1'	2.43	105.75	99.98
2	B	801	PCG	C5-C4-N9	-2.34	103.78	107.16
2	A	401	PCG	C3'-C2'-C1'	2.21	105.22	99.98
2	A	401	PCG	C5-C4-N9	-2.16	104.04	107.16
2	B	801	PCG	PA-O5'-C5'	2.13	125.80	118.66

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	268/288 (93%)	0.57	17 (6%) 19 21	20, 51, 94, 101	0
1	B	259/288 (89%)	0.51	8 (3%) 47 52	26, 49, 78, 95	0
All	All	527/576 (91%)	0.54	25 (4%) 30 34	20, 50, 91, 101	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	300	VAL	6.1
1	A	294	LEU	5.1
1	A	335	ALA	3.9
1	B	729	MET	3.7
1	A	116	ILE	3.7
1	B	707	ASN	3.4
1	A	283	GLN	3.4
1	A	305	SER	3.2
1	A	115	VAL	3.0
1	A	303	ARG	3.0
1	A	182	VAL	2.9
1	A	274	PHE	2.7
1	B	681	VAL	2.7
1	A	329	MET	2.7
1	B	512	VAL	2.5
1	B	584	VAL	2.5
1	A	361	SER	2.5
1	A	163	ILE	2.5
1	B	710	PHE	2.3
1	A	191	SER	2.3
1	A	169	GLY	2.2
1	A	284	GLY	2.1
1	A	275	GLU	2.1
1	B	626	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	654	LEU	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(\AA^2)	Q<0.9
2	PCG	A	401	23/23	0.24	0.27	31,34,39,44	0
3	GOL	B	901	6/6	0.19	0.25	61,66,67,67	0
3	GOL	B	802	6/6	0.22	-0.35	37,38,39,41	0
2	PCG	B	801	23/23	0.20	-0.45	28,32,33,36	0
3	GOL	A	402	6/6	0.19	-0.46	53,55,56,58	0

6.5 Other polymers

There are no such residues in this entry.