



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 07:56 am BST

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

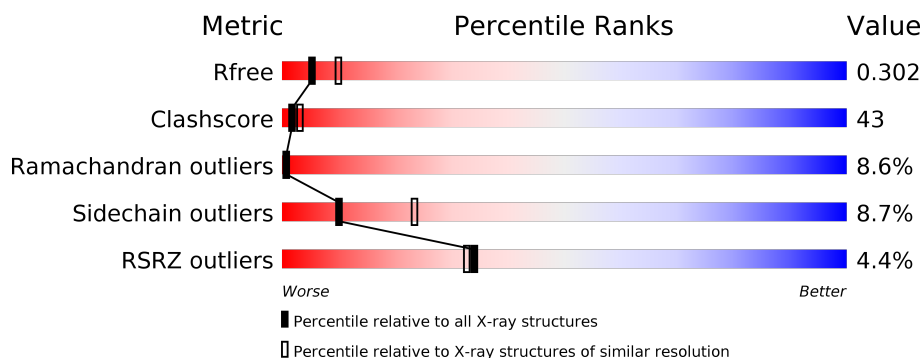
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (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 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	288	
1	B	288	

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

2 Entry composition [i](#)

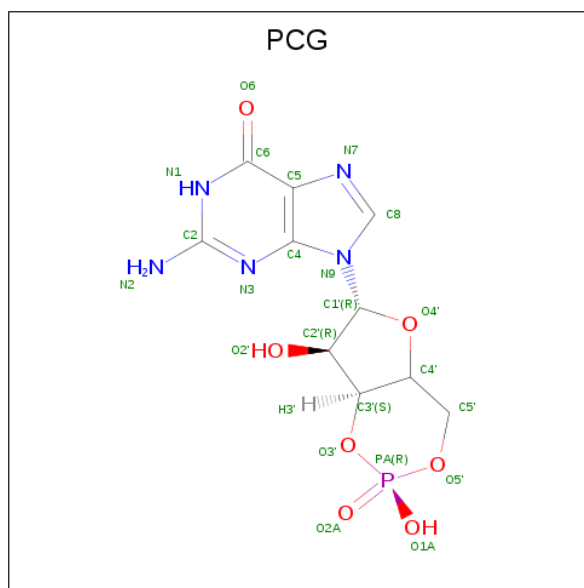
There are 4 unique types of molecules in this entry. The entry contains 4008 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 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		

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4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	90.30 Å 90.30 Å 177.60 Å 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 2.68 Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.240 , 0.285 0.263 , 0.302	Depositor DCC
R_{free} test set	2075 reflections (9.86%)	wwPDB-VP
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 , 43.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.094 for h,-h-k,-l	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.

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: 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 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	1991	0	1835	185	0
1	B	1936	0	1820	153	0
2	A	23	0	11	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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

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 43.

All (334) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A: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	Interatomic distance (Å)	Clash overlap (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:A:318:PRO:O	1:A:319:SER:HB2	1.93	0.68
1:B:570:ASP:OD2	1:B:571:ASN:N	2.24	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:566:GLY:HA2	1:B:608:PRO:HB2	1.80	0.62
1:A:264:THR:HG22	1:A:356:VAL:HG12	1.82	0.62
1:B:630:ARG:HH11	1:B:630:ARG:CB	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:563:ILE:HD11	1:B:609:ARG:NE	2.17	0.59
1:B:660:TRP:HA	1:B:663:LEU:HD12	1.85	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: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:A:282:VAL:CB	1:A:336:THR:HA	2.34	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:OD2	1:B:546:ASP:N	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:A:176:ASP:OD2	1:A:222:TRP:NE1	2.36	0.56
1:B:752:ARG:O	1:B:756:VAL:HG22	2.05	0.56
1:B:533:ASN:OD1	1:B:535:LEU:HB2	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:588:TRP:C	1:B:588:TRP:CD1	2.80	0.54
1:B:571:ASN:C	1:B:603:LEU:HD11	2.29	0.54
1:B:674:PHE:HD1	1:B:678:GLN:HE21	1.56	0.54
1:B:766:ARG:O	1:B:767:ASN:ND2	2.40	0.54
1:A:116:ILE:O	1:A:116:ILE:HG12	2.08	0.54
1:B:669:LEU:HD22	1:B:746:VAL:HB	1.89	0.54
1:B:560:GLU:HA	3:B:901:GOL:H12	1.90	0.53
1:B:561:THR:OG1	1:B:614:LYS:HG2	2.08	0.53
1:B:511:TYR:CD1	1:B:512:VAL:N	2.77	0.53
1:B:657:LEU:HD23	1:B:661:GLU:HB3	1.89	0.53
1:A:163:ILE:CG2	1:A:209:ARG:HE	2.20	0.53
1:A:163:ILE:HG23	1:A:209:ARG:CD	2.38	0.53
1:B:533:ASN:HB3	1:B:536:PHE:HD2	1.74	0.53
1:B:571:ASN:HB2	1:B:573:TYR:CE2	2.43	0.53
1:A:282:VAL:CB	1:A:337:VAL:H	2.22	0.53
1:A:157:ILE:HG22	1:A:218:ASN:OD1	2.08	0.53
1:B:663:LEU:O	1:B:666:ALA:HB3	2.08	0.53
1:A:277:GLY:H	1:A:339:ALA:HB3	1.73	0.52
1:B:510:SER:O	1:B:511:TYR:CB	2.57	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: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:B:680:ILE:N	1:B:680:ILE:HD12	2.26	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:B:684:GLY:O	1:B:732:PRO:HB3	2.11	0.51
1:A:298:ALA:HB1	1:A:338:VAL:O	2.11	0.51
1:A:302:GLN:H	1:A:311:VAL:N	2.09	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:170:ASP:N	1:A:170:ASP:OD2	2.43	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: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:A:118:LYS:HB2	1:A:118:LYS:NZ	2.27	0.49
1:A:369:GLN:OE1	1:A:370:GLN:OE1	2.30	0.49
1:B:510:SER:O	1:B:511:TYR:HB3	2.12	0.49
1:B:565:GLN:HE22	1:B:585:ASN:N	2.10	0.49
1:B:727:LEU:HB3	1:B:753:PHE:CE2	2.46	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:119:ASP:C	1:A:119:ASP:OD2	2.49	0.49
1:A:144:ARG:HH21	1:A:144:ARG:HG3	1.77	0.49
1:A:268:ALA:HB1	1:A:352:ARG:HD2	1.93	0.49
1:B:627:ASP:O	1:B:631:ARG:HB2	2.13	0.49
1:A:126:LEU:HD11	1:A:151:MET:CE	2.43	0.49
1:B:688:ASP:HB3	1:B:750:ARG:CZ	2.42	0.49
1:A:247:PHE:HZ	1:A:293:ILE:O	1.95	0.49
1:A:356:VAL:HG23	1:A:357:LEU:HD13	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:ILE:HG21	1:A:194:GLU:HA	1.94	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:177:GLN:HA	1:A:194:GLU:OE1	2.13	0.48
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:615:ALA:HB1	1:B:617:THR:O	2.14	0.48
1:B:747:LYS:O	1:B:748:LEU:HD12	2.14	0.48
1:A:312:GLU:O	1:A:313:VAL:CB	2.62	0.47
1:B:588:TRP:CH2	1:B:591:SER:HB2	2.49	0.47
1:B:658:ASP:OD2	1:B:661:GLU:HG3	2.14	0.47
1:B:540:ASP:CG	1:B:543:GLU:HG3	2.35	0.47
1:B:563:ILE:CD1	1:B:609:ARG:NE	2.78	0.47
1:B:570:ASP:H	1:B:609:ARG:NH2	2.09	0.47
1:B:681:VAL:O	1:B:736:THR:HG23	2.14	0.47
1:B:516:ILE:O	1:B:516:ILE:HG12	2.14	0.47
1:B:661:GLU:O	1:B:664:THR:HB	2.15	0.47
1:A:126:LEU:HD13	1:A:222:TRP:CE3	2.49	0.47
1:B:542:ASN:N	1:B:542:ASN:ND2	2.62	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:A:127:ALA:O	1:A:131:GLU:HB3	2.15	0.46
1:B:641:ARG:HD3	1:B:663:LEU:HD22	1.97	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:A:240:LYS:HD2	1:A:244:TYR:HE2	1.81	0.46
1:B:672:VAL:HG11	1:B:747:LYS:HE2	1.97	0.46
1:A:220:LYS:O	1:A:221:LEU:HD23	2.16	0.46
1:A:188:TRP:CG	1:A:189:ALA:N	2.82	0.46
1:B:646:GLU:HB3	1:B:650:LYS:HZ1	1.80	0.46
1:B:727:LEU:HD13	1:B:753:PHE:CD2	2.50	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:262:ARG:HH11	1:A:262:ARG:CB	2.30	0.45
1:A:315:ARG:O	1:A:316:LEU:HG	2.17	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:A:369:GLN:NE2	1:A:373:SER:HB2	2.32	0.45
1:B:542:ASN:O	1:B:546:ASP:OD2	2.34	0.45
1:B:716:LEU:HA	1:B:720:ASP:OD2	2.16	0.45
1:B:727:LEU:HD23	1:B:750:ARG:HD3	1.98	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:A:120:TYR:OH	1:B:544:ARG:CZ	2.65	0.45
1:B:601:LEU:HB2	2:B:801:PCG:O3'	2.16	0.45
1:B:674:PHE:CZ	1:B:680:ILE:HG13	2.52	0.45
1:A:288:ASP:HB3	1:A:350:ARG:HE	1.82	0.44
1:B:758:GLY:C	1:B:760:CYS:H	2.21	0.44
1:A:144:ARG:HG3	1:A:144:ARG:NH2	2.32	0.44
1:A:157:ILE:HG22	1:A:218:ASN:HB3	1.98	0.44
1:B:565:GLN:NE2	1:B:585:ASN:H	2.14	0.44
1:B:580:MET:HB2	1:B:592:VAL:HB	2.00	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:B:657:LEU:HD11	1:B:760:CYS:SG	2.57	0.44
1:A:134:VAL:HG12	1:A:134:VAL:O	2.17	0.44
1:A:200:GLU:OE1	1:A:200:GLU:N	2.47	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:352:ARG:O	1:A:356:VAL:HG22	2.18	0.43
1:B:538:HIS:CD2	1:B:538:HIS:N	2.85	0.43
1:B:675:GLU:O	1:B:677:GLY:N	2.44	0.43
1:A:262:ARG:HH11	1:A:262:ARG:HB3	1.82	0.43
1:A:163:ILE:HG21	2:A:401:PCG:H5'2	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:A:214:LYS:NZ	4:A:906:HOH:O	2.52	0.43
1:B:512:VAL:O	1:B:513:ARG:CB	2.66	0.43
1:B:599:GLY:HA2	2:B:801:PCG:O1A	2.19	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:240:LYS:HD2	1:A:244:TYR:CE2	2.54	0.43
1:A:328:LEU:HD11	1:A:364:LEU:CD1	2.49	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:A:369:GLN:HE21	1:A:373:SER:HB2	1.84	0.42
1:B:693:ILE:HG23	1:B:743:LEU:HD11	2.01	0.42
1:A:207:THR:HA	1:A:208:PRO:HD2	1.76	0.42
1:B:563:ILE:HG12	1:B:609:ARG:HD2	2.02	0.42
1:B:664:THR:CG2	1:B:756:VAL:HG12	2.46	0.42
1:A:120:TYR:HA	1:B:523:MET:CE	2.49	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:253:ILE:C	1:A:255:GLU:H	2.23	0.42
1:B:749:ASP:O	1:B:750:ARG:C	2.58	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:116:ILE:O	1:A:116:ILE:CG1	2.67	0.42
1:A:173:TYR:HB2	1:A:198:PHE:CE1	2.54	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: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:B:757:LEU:O	1:B:760:CYS:HB2	2.20	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:526:LEU:HD11	1:B:551:MET:CE	2.47	0.40
1:B:688:ASP:HB2	1:B:727:LEU:CD2	2.51	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:A:188:TRP:CD1	1:A:188:TRP:C	2.93	0.40
1:B:552:PHE:CD1	1:B:552:PHE:O	2.75	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%)	0	0
1	B	257/288 (89%)	210 (82%)	30 (12%)	17 (7%)	1	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
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

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Mol	Chain	Res	Type
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%)	7	16
1	B	188/244 (77%)	175 (93%)	13 (7%)	15	35
All	All	379/488 (78%)	346 (91%)	33 (9%)	10	23

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

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Mol	Chain	Res	Type
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 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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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	-	22,26,26	2.77	7 (31%)	26,41,41	4.41	14 (53%)
3	GOL	B	901	-	5,5,5	4.42	5 (100%)	5,5,5	5.79	3 (60%)
2	PCG	B	801	-	22,26,26	3.08	8 (36%)	26,41,41	4.18	14 (53%)
3	GOL	A	402	-	5,5,5	4.70	5 (100%)	5,5,5	5.76	3 (60%)
3	GOL	B	802	-	5,5,5	4.47	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	-	-	0/0/31/31	1/4/4/4
3	GOL	B	901	-	-	3/4/4/4	-
2	PCG	B	801	-	-	0/0/31/31	0/4/4/4
3	GOL	A	402	-	-	2/4/4/4	-
3	GOL	B	802	-	-	2/4/4/4	-

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	GOL	C3-C2	-7.79	1.19	1.51
3	B	802	GOL	C3-C2	-7.31	1.21	1.51
3	B	901	GOL	C3-C2	-7.23	1.22	1.51
2	B	801	PCG	PA-O5'	-7.06	1.49	1.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	PCG	PA-O5'	-6.55	1.50	1.57
2	B	801	PCG	C6-N1	6.11	1.43	1.33
2	B	801	PCG	C4-N3	6.06	1.45	1.35
2	A	401	PCG	C6-N1	5.61	1.42	1.33
2	B	801	PCG	O3'-C3'	5.13	1.52	1.44
2	A	401	PCG	C4-N3	4.92	1.43	1.35
3	A	402	GOL	O1-C1	4.77	1.62	1.42
3	B	802	GOL	O1-C1	4.63	1.62	1.42
2	A	401	PCG	PA-O3'	-4.61	1.50	1.57
3	B	901	GOL	O1-C1	4.48	1.61	1.42
2	B	801	PCG	C6-C5	3.94	1.48	1.41
2	A	401	PCG	O3'-C3'	3.86	1.50	1.44
2	A	401	PCG	C2-N1	3.67	1.41	1.35
3	B	901	GOL	O3-C3	3.43	1.56	1.42
3	B	802	GOL	O3-C3	3.38	1.56	1.42
3	A	402	GOL	O2-C2	-3.30	1.33	1.43
2	B	801	PCG	C2-N1	3.27	1.41	1.35
2	B	801	PCG	C5'-C4'	-3.05	1.46	1.51
3	A	402	GOL	O3-C3	2.95	1.54	1.42
2	A	401	PCG	C6-C5	2.92	1.46	1.41
3	A	402	GOL	C1-C2	-2.75	1.40	1.51
3	B	901	GOL	C1-C2	-2.66	1.40	1.51
3	B	802	GOL	O2-C2	-2.65	1.35	1.43
3	B	901	GOL	O2-C2	-2.59	1.35	1.43
3	B	802	GOL	C1-C2	-2.55	1.41	1.51
2	B	801	PCG	PA-O3'	-2.43	1.53	1.57

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	PCG	O3'-C3'-C2'	12.01	127.39	115.61
3	B	901	GOL	O3-C3-C2	10.53	160.67	110.20
3	A	402	GOL	O3-C3-C2	10.52	160.64	110.20
3	B	802	GOL	O3-C3-C2	10.36	159.86	110.20
2	A	401	PCG	O3'-C3'-C2'	10.27	125.68	115.61
2	A	401	PCG	O5'-PA-O3'	-9.40	92.73	105.68
2	A	401	PCG	O1A-PA-O3'	8.74	127.41	107.04
2	A	401	PCG	C5-C6-N1	-7.86	112.68	123.43
2	B	801	PCG	C5-C6-N1	-7.69	112.91	123.43
2	B	801	PCG	O1A-PA-O3'	7.23	123.89	107.04
3	B	802	GOL	O2-C2-C3	6.81	139.12	109.12
2	B	801	PCG	C6-N1-C2	6.76	126.67	115.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	901	GOL	O2-C2-C3	6.73	138.76	109.12
3	A	402	GOL	O2-C2-C3	6.63	138.33	109.12
2	A	401	PCG	C6-N1-C2	6.57	126.37	115.93
2	B	801	PCG	C2'-C3'-C4'	-5.53	93.42	103.22
2	A	401	PCG	C2'-C3'-C4'	-5.26	93.90	103.22
2	A	401	PCG	C1'-N9-C4	-4.59	118.58	126.64
2	B	801	PCG	C1'-N9-C4	-4.43	118.87	126.64
2	A	401	PCG	O3'-C3'-C4'	4.15	113.84	110.71
2	B	801	PCG	O5'-PA-O3'	-3.95	100.23	105.68
2	B	801	PCG	O3'-C3'-C4'	3.76	113.55	110.71
3	B	802	GOL	O1-C1-C2	3.56	127.25	110.20
2	B	801	PCG	N3-C2-N1	-3.37	122.72	127.22
2	A	401	PCG	O2'-C2'-C3'	3.35	120.69	111.17
3	B	901	GOL	O1-C1-C2	3.34	126.24	110.20
2	B	801	PCG	O4'-C4'-C5'	3.25	122.08	112.37
2	A	401	PCG	N3-C2-N1	-3.25	122.89	127.22
3	A	402	GOL	O1-C1-C2	3.23	125.69	110.20
2	B	801	PCG	C6-C5-C4	-3.07	117.87	120.80
2	B	801	PCG	C5'-C4'-C3'	-2.89	106.64	112.49
2	B	801	PCG	O2'-C2'-C3'	2.71	118.85	111.17
2	A	401	PCG	C6-C5-C4	-2.69	118.22	120.80
2	B	801	PCG	C3'-C2'-C1'	2.64	105.75	99.89
2	A	401	PCG	O4'-C4'-C5'	2.57	120.05	112.37
2	A	401	PCG	C3'-C2'-C1'	2.41	105.22	99.89
2	A	401	PCG	C2-N3-C4	-2.11	112.95	115.36

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	901	GOL	C1-C2-C3-O3
3	A	402	GOL	C1-C2-C3-O3
3	B	802	GOL	O1-C1-C2-C3
3	B	802	GOL	C1-C2-C3-O3
3	B	901	GOL	O1-C1-C2-C3
3	B	901	GOL	O1-C1-C2-O2
3	A	402	GOL	O1-C1-C2-O2

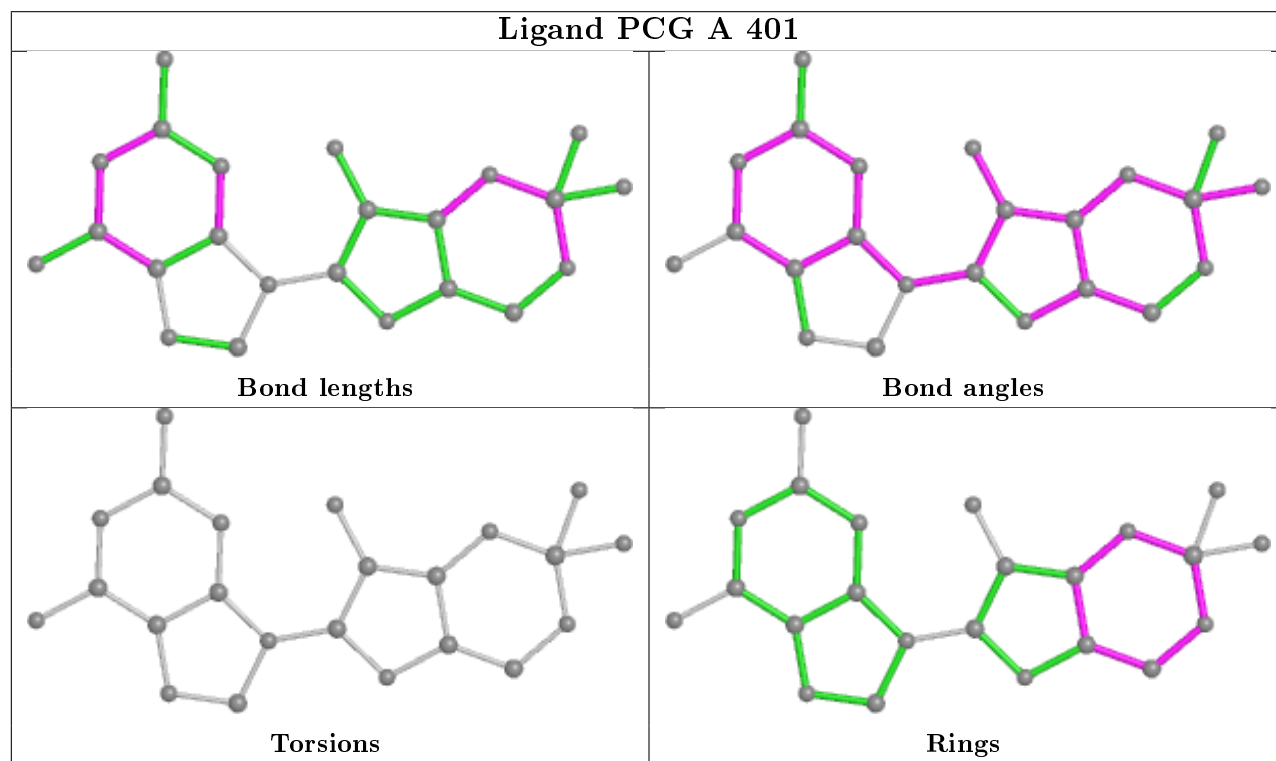
All (1) ring outliers are listed below:

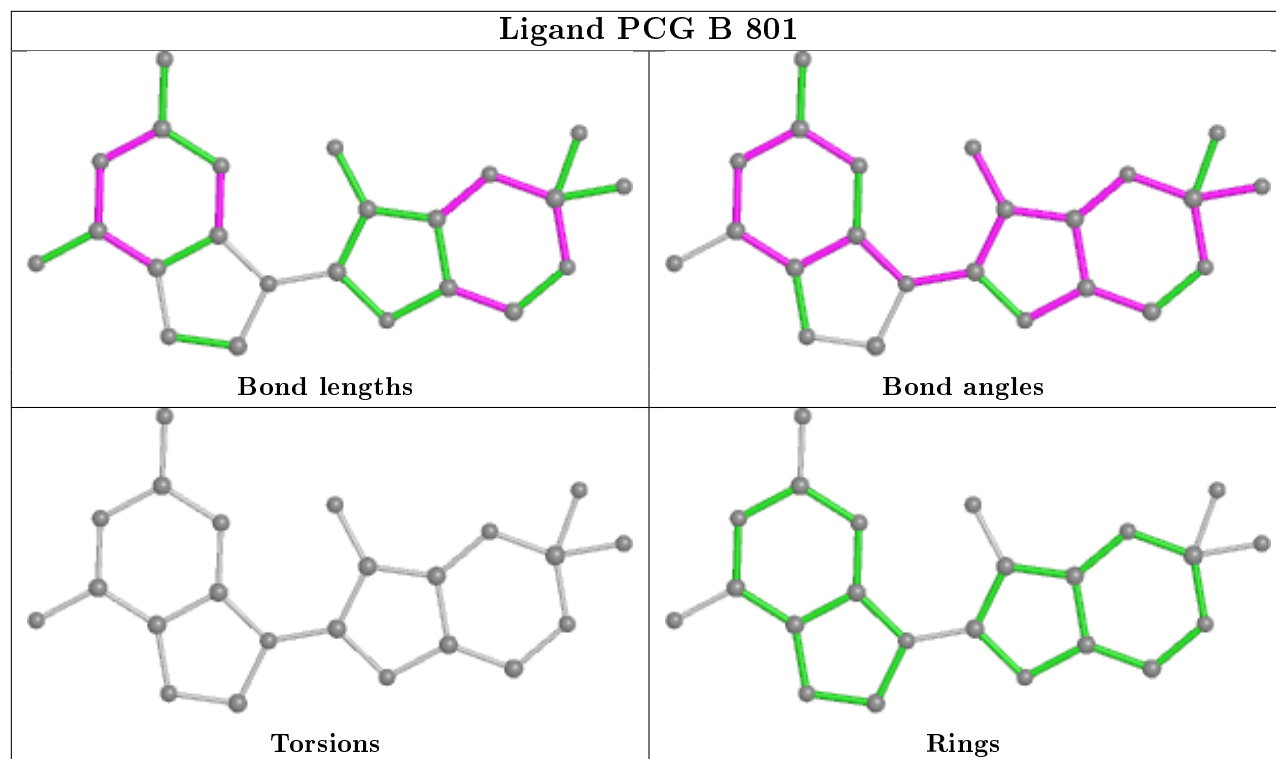
Mol	Chain	Res	Type	Atoms
2	A	401	PCG	C3'-C4'-C5'-O3'-O5'-PA

4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	PCG	3	0
3	B	901	GOL	2	0
2	B	801	PCG	2	0
3	A	402	GOL	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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	268/288 (93%)	0.51	15 (5%) 24 23	20, 51, 94, 101	0
1	B	259/288 (89%)	0.44	8 (3%) 49 49	26, 49, 78, 95	0
All	All	527/576 (91%)	0.48	23 (4%) 34 33	20, 50, 91, 101	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	300	VAL	6.1
1	A	294	LEU	5.0
1	A	335	ALA	3.8
1	A	116	ILE	3.6
1	B	729	MET	3.4
1	A	283	GLN	3.3
1	B	707	ASN	3.3
1	A	305	SER	3.1
1	A	115	VAL	2.9
1	A	303	ARG	2.9
1	A	182	VAL	2.9
1	A	274	PHE	2.6
1	B	681	VAL	2.6
1	A	329	MET	2.5
1	B	512	VAL	2.5
1	A	163	ILE	2.4
1	B	584	VAL	2.4
1	A	361	SER	2.4
1	B	710	PHE	2.3
1	A	191	SER	2.3
1	A	169	GLY	2.1
1	B	626	ARG	2.1
1	B	654	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

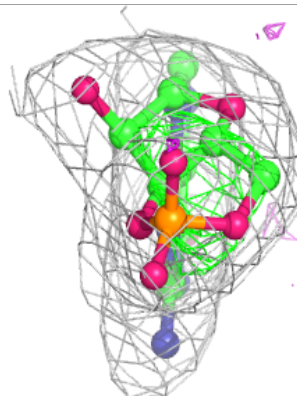
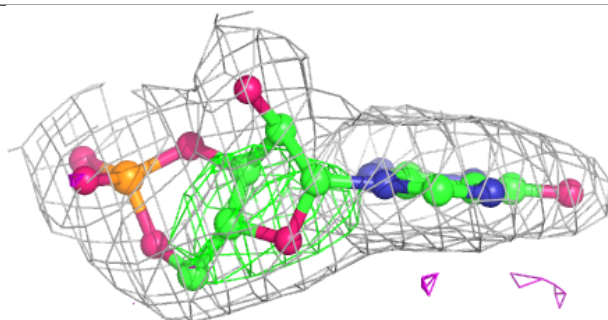
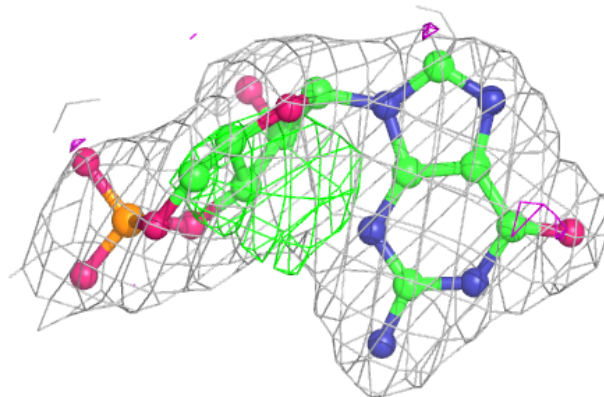
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	GOL	B	901	6/6	0.70	0.19	61,66,67,67	0
3	GOL	A	402	6/6	0.91	0.19	53,55,56,58	0
2	PCG	A	401	23/23	0.93	0.24	31,34,39,44	0
3	GOL	B	802	6/6	0.93	0.22	37,38,39,41	0
2	PCG	B	801	23/23	0.95	0.20	28,32,33,36	0

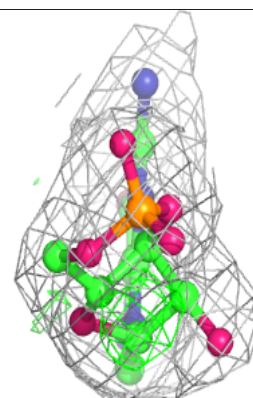
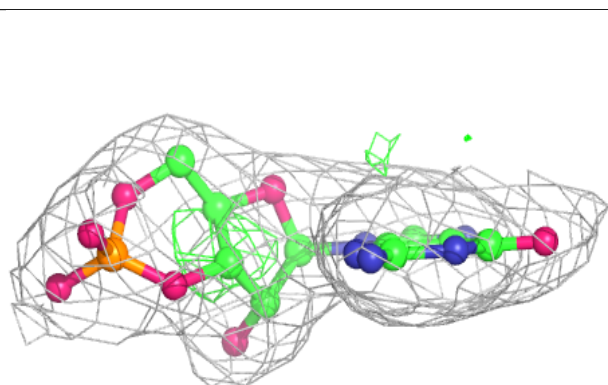
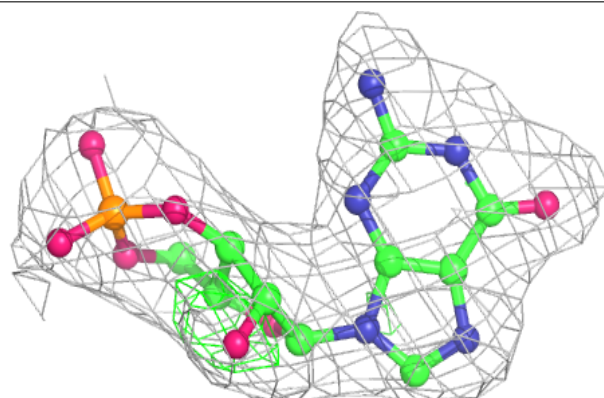
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around PCG A 401:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around PCG B 801:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.