



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:17 PM GMT

PDB ID : 3WS9
Title : Crystal structure of PDE10A in complex with a benzimidazole inhibitor
Authors : Amano, Y.; Honbou, K.
Deposited on : 2014-03-04
Resolution : 2.99 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

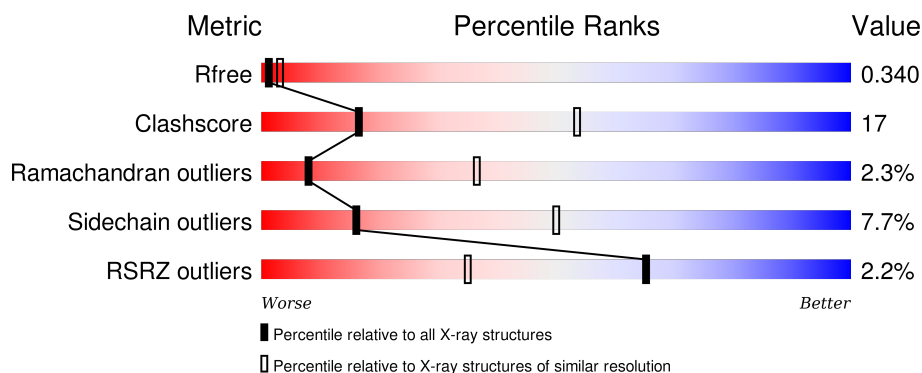
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.99 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}	91344	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

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. 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	345	<div> <div>3%</div> <div>57%</div> <div>34%</div> <div>6%</div> </div>
1	B	345	<div> <div>%</div> <div>58%</div> <div>32%</div> <div>6%</div> </div>

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	MG	A	802	-	-	-	X
3	MG	B	802	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5321 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 and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	324	Total	C	N	O	S	0	0	0
			2629	1678	447	479	25			
1	B	324	Total	C	N	O	S	0	0	0
			2629	1678	447	479	25			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	445	GLY	-	EXPRESSION TAG	UNP Q9Y233
A	446	SER	-	EXPRESSION TAG	UNP Q9Y233
A	447	HIS	-	EXPRESSION TAG	UNP Q9Y233
A	448	MET	-	EXPRESSION TAG	UNP Q9Y233
B	445	GLY	-	EXPRESSION TAG	UNP Q9Y233
B	446	SER	-	EXPRESSION TAG	UNP Q9Y233
B	447	HIS	-	EXPRESSION TAG	UNP Q9Y233
B	448	MET	-	EXPRESSION TAG	UNP Q9Y233

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

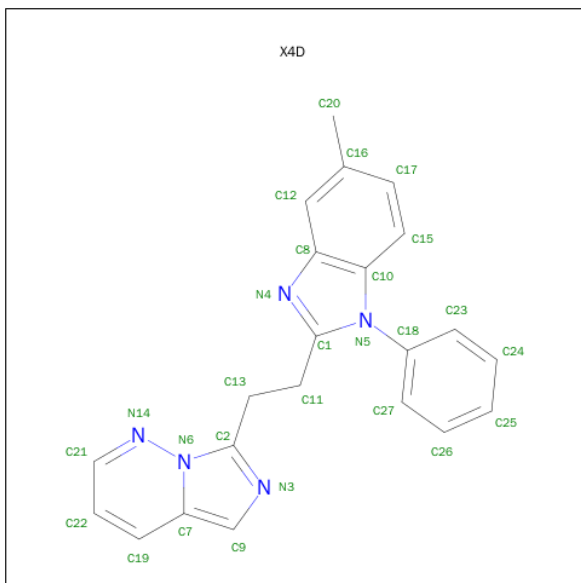
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		

- Molecule 4 is 7-[2-(5-METHYL-1-PHENYL-1H-BENZIMIDAZOL-2-YL)ETHYL]IMIDAZO [1,5-B]PYRIDAZINE (three-letter code: X4D) (formula: C₂₂H₁₉N₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	N	0	0
			27	22	5		
4	B	1	Total	C	N	0	0
			27	22	5		

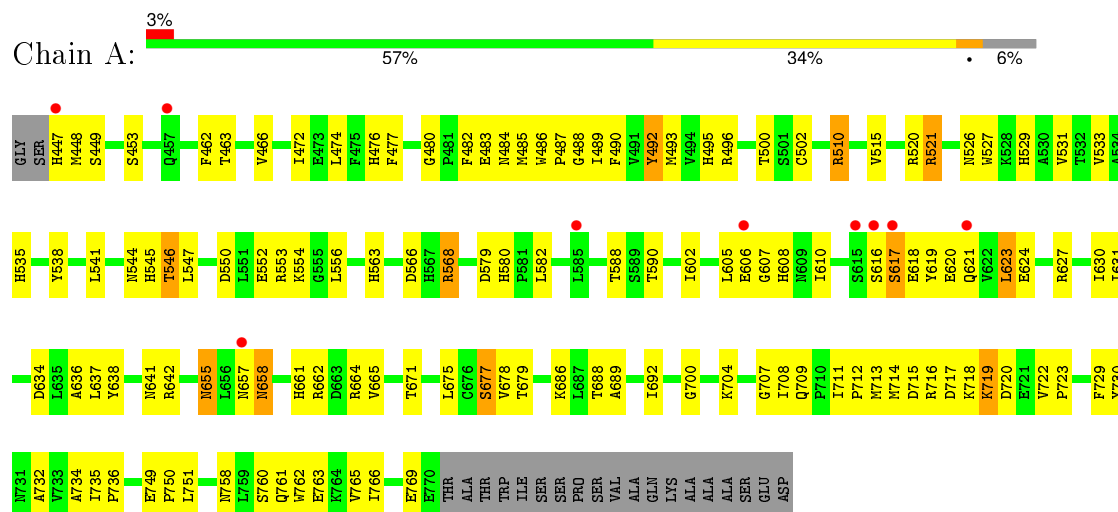
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	O	0	0
			2	2		
5	B	3	Total	O	0	0
			3	3		

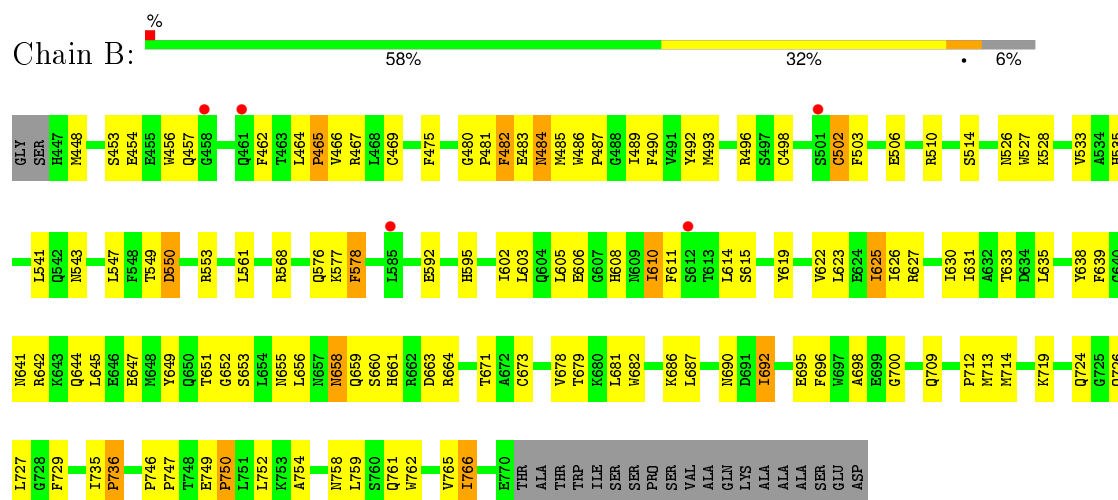
3 Residue-property plots

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 and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A



- Molecule 1: cAMP and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	50.42Å 81.54Å 161.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.50 – 2.99 29.50 – 2.99	Depositor EDS
% Data completeness (in resolution range)	99.1 (29.50-2.99) 99.3 (29.50-2.99)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.95 (at 3.00Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.223 , 0.346 0.221 , 0.340	Depositor DCC
R_{free} test set	695 reflections (5.24%)	DCC
Wilson B-factor (Å ²)	51.6	Xtriage
Anisotropy	0.101	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 39.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	2 of 13949 reflections (0.014%)	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	5321	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 4.20% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: ZN, X4D, MG

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	0.61	0/2694	0.78	0/3647
1	B	0.61	0/2694	0.79	1/3647 (0.0%)
All	All	0.61	0/5388	0.78	1/7294 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	656	LEU	CA-CB-CG	5.70	128.42	115.30

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	2629	0	2588	87	1
1	B	2629	0	2588	89	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	27	0	19	1	0
4	B	27	0	19	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	2	0	0	0	0
5	B	3	0	0	0	0
All	All	5321	0	5214	175	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 17.

All (175) 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:655:ASN:HD21	1:A:657:ASN:HB2	1.24	0.99
1:B:713:MET:CE	4:B:803:X4D:H7	1.96	0.95
1:A:521:ARG:HH11	1:A:521:ARG:HB3	1.31	0.93
1:B:655:ASN:H	1:B:661:HIS:HD2	1.25	0.84
1:A:641:ASN:HD22	1:A:664:ARG:HH11	1.22	0.84
1:A:655:ASN:ND2	1:A:657:ASN:HB2	1.92	0.83
1:B:533:VAL:HG13	1:B:673:CYS:HB3	1.60	0.82
1:A:521:ARG:NH1	1:A:521:ARG:HB3	2.00	0.76
1:A:483:GLU:HA	1:A:486:TRP:CE2	2.25	0.71
1:A:544:ASN:HB3	1:A:547:LEU:HD13	1.73	0.70
1:A:462:PHE:HZ	1:A:485:MET:HG2	1.56	0.70
1:A:462:PHE:CZ	1:A:485:MET:HG2	2.25	0.69
1:B:484:ASN:HD22	1:B:484:ASN:N	1.90	0.69
1:A:553:ARG:HG3	1:A:553:ARG:HH11	1.57	0.68
1:A:708:ILE:HG22	1:A:709:GLN:N	2.09	0.68
1:A:618:GLU:HA	1:A:621:GLN:HE21	1.59	0.68
1:B:713:MET:HE2	4:B:803:X4D:H7	1.73	0.67
1:A:732:ALA:HB2	1:B:687:LEU:HD21	1.75	0.67
1:B:642:ARG:HE	1:B:736:PRO:HB3	1.62	0.65
1:B:576:GLN:O	1:B:578:PHE:N	2.27	0.65
1:B:655:ASN:H	1:B:661:HIS:CD2	2.13	0.64
1:A:708:ILE:CG2	1:A:709:GLN:N	2.61	0.64
1:B:453:SER:HA	1:B:456:TRP:CD1	2.33	0.63
1:A:689:ALA:O	1:A:692:ILE:HG22	1.98	0.62
1:B:481:PRO:HG2	1:B:482:PHE:CE1	2.35	0.62
1:B:735:ILE:HB	1:B:736:PRO:HD3	1.82	0.62
1:A:708:ILE:CG2	1:A:709:GLN:H	2.12	0.62
1:B:498:CYS:HB3	1:B:502:CYS:SG	2.41	0.61
1:A:563:HIS:HA	1:A:630:ILE:HG12	1.81	0.61
1:A:704:LYS:O	1:A:707:GLY:N	2.33	0.61
1:A:580:HIS:ND1	1:A:582:LEU:HB2	2.16	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:641:ASN:HD22	1:B:664:ARG:HH11	1.48	0.60
1:A:492:TYR:C	1:A:492:TYR:CD2	2.75	0.59
1:A:655:ASN:HD22	1:A:657:ASN:H	1.50	0.59
1:B:641:ASN:HD22	1:B:664:ARG:NH1	1.99	0.59
1:B:692:ILE:HD11	1:B:696:PHE:CZ	2.38	0.59
1:A:677:SER:HB3	1:A:688:THR:HG21	1.85	0.58
1:B:735:ILE:HB	1:B:736:PRO:CD	2.33	0.58
1:B:465:PRO:O	1:B:467:ARG:N	2.37	0.58
1:A:711:ILE:HD12	1:A:712:PRO:HD2	1.85	0.58
1:A:490:PHE:HA	1:A:493:MET:HE2	1.86	0.58
1:B:498:CYS:CB	1:B:502:CYS:SG	2.92	0.57
1:A:713:MET:HG3	4:A:803:X4D:C8	2.34	0.57
1:A:760:SER:HA	1:A:763:GLU:HB2	1.87	0.57
1:A:624:GLU:HA	1:A:624:GLU:OE2	2.05	0.56
1:A:568:ARG:HG2	1:A:568:ARG:HH11	1.71	0.56
1:B:695:GLU:O	1:B:698:ALA:HB3	2.05	0.55
1:B:605:LEU:HB3	1:B:608:HIS:ND1	2.21	0.55
1:A:535:HIS:O	1:A:538:TYR:HB3	2.07	0.55
1:A:520:ARG:NE	1:A:566:ASP:OD1	2.38	0.55
1:A:553:ARG:HG3	1:A:553:ARG:NH1	2.21	0.55
1:A:485:MET:HB3	1:A:489:ILE:HD12	1.88	0.54
1:A:486:TRP:HB2	1:A:487:PRO:HD3	1.89	0.54
1:A:492:TYR:C	1:A:492:TYR:HD2	2.12	0.54
1:B:541:LEU:HD22	1:B:553:ARG:HG2	1.91	0.53
1:B:611:PHE:O	1:B:614:LEU:HB2	2.08	0.52
1:B:592:GLU:HA	1:B:595:HIS:HD2	1.74	0.52
1:B:713:MET:CE	4:B:803:X4D:C27	2.82	0.52
1:A:658:ASN:OD1	1:A:661:HIS:CD2	2.63	0.52
1:A:627:ARG:O	1:A:631:ILE:HG12	2.10	0.52
1:B:475:PHE:CD1	1:B:750:PRO:HB2	2.44	0.52
1:B:713:MET:HE1	4:B:803:X4D:H7	1.90	0.52
1:A:510:ARG:HD2	1:A:607:GLY:O	2.10	0.52
1:B:619:TYR:CE2	1:B:623:LEU:CD1	2.92	0.52
1:A:711:ILE:HB	1:A:712:PRO:CD	2.40	0.52
1:A:493:MET:SD	1:A:535:HIS:HA	2.50	0.52
1:A:521:ARG:CB	1:A:521:ARG:HH11	2.14	0.51
1:B:619:TYR:CE2	1:B:623:LEU:HD12	2.44	0.51
1:B:627:ARG:O	1:B:631:ILE:HG12	2.10	0.51
1:B:480:GLY:N	1:B:481:PRO:HD2	2.24	0.51
1:B:682:TRP:CZ2	1:B:686:LYS:HG3	2.45	0.51
1:B:483:GLU:HA	1:B:486:TRP:CE2	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:644:GLN:O	1:B:647:GLU:HB3	2.11	0.51
1:A:486:TRP:HE3	1:A:527:TRP:CD1	2.30	0.50
1:B:681:LEU:O	1:B:682:TRP:C	2.49	0.50
1:B:503:PHE:HD1	1:B:610:ILE:HB	1.76	0.50
1:A:474:LEU:HD13	1:A:476:HIS:CE1	2.46	0.50
1:A:617:SER:HA	1:A:620:GLU:HB2	1.93	0.50
1:B:533:VAL:HG13	1:B:673:CYS:CB	2.38	0.49
1:B:464:LEU:HD23	1:B:469:CYS:SG	2.52	0.49
1:B:484:ASN:N	1:B:484:ASN:ND2	2.60	0.49
1:B:526:ASN:OD1	1:B:528:LYS:HB3	2.12	0.49
1:B:484:ASN:HD22	1:B:484:ASN:H	1.59	0.49
1:B:700:GLY:HA3	1:B:714:MET:O	2.12	0.49
1:A:529:HIS:O	1:A:533:VAL:HG23	2.13	0.49
1:B:678:VAL:O	1:B:758:ASN:ND2	2.43	0.49
1:A:602:ILE:HA	1:A:605:LEU:HD12	1.95	0.48
1:A:634:ASP:OD1	1:A:636:ALA:HB3	2.12	0.48
1:A:679:THR:HG21	1:A:751:LEU:HD22	1.94	0.48
1:B:658:ASN:ND2	1:B:658:ASN:C	2.67	0.48
1:B:641:ASN:ND2	1:B:664:ARG:HH11	2.10	0.48
1:A:515:VAL:HG12	1:A:527:TRP:HZ3	1.78	0.48
1:B:746:PRO:N	1:B:747:PRO:CD	2.77	0.48
1:B:749:GLU:O	1:B:752:LEU:N	2.46	0.48
1:A:716:ARG:C	1:A:718:LYS:H	2.16	0.48
1:A:638:TYR:CD1	1:A:671:THR:HG21	2.48	0.48
1:A:761:GLN:O	1:A:765:VAL:HG23	2.14	0.48
1:B:622:VAL:O	1:B:625:ILE:HB	2.14	0.47
1:B:561:LEU:HG	1:B:561:LEU:O	2.13	0.47
1:A:722:VAL:O	1:A:723:PRO:C	2.53	0.47
1:B:486:TRP:HE3	1:B:527:TRP:CD1	2.31	0.47
1:A:658:ASN:O	1:A:661:HIS:HB2	2.14	0.47
1:A:678:VAL:O	1:A:758:ASN:ND2	2.48	0.47
1:B:610:ILE:HG13	1:B:610:ILE:H	1.53	0.47
1:A:495:HIS:CG	1:A:500:THR:HG22	2.49	0.47
1:A:735:ILE:HB	1:A:736:PRO:HD3	1.97	0.47
1:A:655:ASN:ND2	1:A:657:ASN:H	2.12	0.46
1:B:727:LEU:HD22	1:B:766:ILE:HD13	1.97	0.46
1:A:619:TYR:C	1:A:621:GLN:H	2.19	0.46
1:A:483:GLU:HA	1:A:486:TRP:CD2	2.49	0.46
1:A:711:ILE:HB	1:A:712:PRO:HD3	1.96	0.46
1:B:506:GLU:OE1	1:B:510:ARG:NH2	2.49	0.46
1:B:514:SER:HG	1:B:608:HIS:HE2	1.62	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:602:ILE:HA	1:A:605:LEU:CD1	2.46	0.46
1:A:675:LEU:O	1:A:678:VAL:HG22	2.16	0.45
1:B:679:THR:O	1:B:754:ALA:HB1	2.17	0.45
1:A:492:TYR:CE2	1:A:496:ARG:HG3	2.52	0.45
1:B:448:MET:HE3	1:B:602:ILE:HG13	1.97	0.45
1:B:729:PHE:CZ	4:B:803:X4D:C7	2.99	0.45
1:B:659:GLN:NE2	1:B:663:ASP:OD1	2.50	0.45
1:A:715:ASP:OD1	1:A:717:ASP:N	2.47	0.45
1:B:503:PHE:CD1	1:B:610:ILE:HB	2.52	0.45
1:B:498:CYS:HB2	1:B:502:CYS:SG	2.57	0.44
1:A:735:ILE:HD13	1:A:735:ILE:HA	1.75	0.44
1:B:576:GLN:C	1:B:578:PHE:H	2.15	0.44
1:A:718:LYS:O	1:A:719:LYS:C	2.55	0.44
1:A:762:TRP:O	1:A:766:ILE:HG13	2.18	0.44
1:B:642:ARG:NE	1:B:736:PRO:HB3	2.28	0.44
1:B:490:PHE:CE2	1:B:561:LEU:HD22	2.52	0.44
1:A:655:ASN:C	1:A:655:ASN:HD22	2.21	0.43
1:A:485:MET:HB3	1:A:489:ILE:CD1	2.48	0.43
1:B:649:TYR:O	1:B:651:THR:N	2.51	0.43
1:A:527:TRP:O	1:A:531:VAL:HG23	2.19	0.43
1:B:493:MET:SD	1:B:535:HIS:HA	2.58	0.43
1:A:526:ASN:OD1	1:A:526:ASN:C	2.55	0.43
1:B:484:ASN:ND2	1:B:485:MET:HG3	2.34	0.43
1:B:635:LEU:HD13	1:B:639:PHE:HE1	1.84	0.43
1:A:482:PHE:HB3	1:A:485:MET:HB2	1.99	0.43
1:B:758:ASN:O	1:B:759:LEU:C	2.57	0.43
1:B:549:THR:O	1:B:550:ASP:C	2.57	0.43
1:B:724:GLN:HA	1:B:766:ILE:HD11	2.00	0.42
1:A:641:ASN:HA	1:A:664:ARG:NH1	2.33	0.42
1:A:472:ILE:HD11	1:A:489:ILE:HG23	2.01	0.42
1:A:541:LEU:HD22	1:A:553:ARG:HG2	2.01	0.42
1:B:644:GLN:HG3	1:B:664:ARG:CZ	2.49	0.42
1:B:543:ASN:HD22	1:B:543:ASN:N	2.16	0.42
1:B:533:VAL:CG1	1:B:673:CYS:HB3	2.42	0.42
1:A:730:TYR:HA	1:A:734:ALA:HB3	2.02	0.42
1:A:550:ASP:O	1:A:554:LYS:HG2	2.20	0.42
1:B:603:LEU:C	1:B:605:LEU:H	2.23	0.42
1:A:477:PHE:HB3	1:A:535:HIS:CE1	2.55	0.42
1:B:746:PRO:N	1:B:747:PRO:HD3	2.34	0.42
1:B:645:LEU:HD11	1:B:664:ARG:O	2.20	0.41
1:B:486:TRP:HB2	1:B:487:PRO:CD	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:635:LEU:HD13	1:B:639:PHE:CE1	2.55	0.41
1:A:556:LEU:HA	1:A:556:LEU:HD23	1.89	0.41
1:B:492:TYR:CZ	1:B:496:ARG:HG3	2.54	0.41
1:A:623:LEU:HD23	1:A:623:LEU:HA	1.88	0.41
1:B:615:SER:O	1:B:619:TYR:N	2.42	0.41
1:A:608:HIS:O	1:A:610:ILE:HG23	2.19	0.41
1:B:462:PHE:HE1	1:B:485:MET:HG2	1.86	0.41
1:B:626:ILE:O	1:B:627:ARG:C	2.59	0.41
1:A:700:GLY:HA3	1:A:714:MET:O	2.20	0.41
1:B:726:GLN:HG3	1:B:762:TRP:CH2	2.56	0.41
1:B:761:GLN:O	1:B:765:VAL:HG23	2.20	0.41
1:B:484:ASN:ND2	1:B:484:ASN:H	2.19	0.41
1:B:453:SER:HA	1:B:456:TRP:NE1	2.35	0.41
1:B:489:ILE:O	1:B:493:MET:HG3	2.21	0.41
1:A:552:GLU:CD	1:A:662:ARG:HH21	2.24	0.40
1:B:651:THR:O	1:B:653:SER:N	2.54	0.40
1:B:630:ILE:O	1:B:633:THR:OG1	2.33	0.40
1:A:749:GLU:N	1:A:750:PRO:HD2	2.36	0.40
1:A:580:HIS:CE1	1:A:582:LEU:HB2	2.56	0.40
1:A:462:PHE:CD2	1:A:488:GLY:HA3	2.56	0.40
1:A:729:PHE:O	1:A:730:TYR:C	2.59	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:449:SER:OG	1:B:614:LEU:O[1_545]	2.16	0.04

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	322/345 (93%)	272 (84%)	44 (14%)	6 (2%)	10	43
1	B	322/345 (93%)	271 (84%)	42 (13%)	9 (3%)	6	30
All	All	644/690 (93%)	543 (84%)	86 (13%)	15 (2%)	8	36

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	484	ASN
1	B	466	VAL
1	B	577	LYS
1	A	719	LYS
1	B	625	ILE
1	B	719	LYS
1	A	546	THR
1	B	502	CYS
1	B	578	PHE
1	A	590	THR
1	B	652	GLY
1	B	750	PRO
1	A	545	HIS
1	A	480	GLY
1	B	465	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	293/308 (95%)	267 (91%)	26 (9%)	12	42
1	B	293/308 (95%)	274 (94%)	19 (6%)	21	58
All	All	586/616 (95%)	541 (92%)	45 (8%)	16	50

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	447	HIS

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Mol	Chain	Res	Type
1	A	448	MET
1	A	453	SER
1	A	463	THR
1	A	466	VAL
1	A	492	TYR
1	A	502	CYS
1	A	510	ARG
1	A	521	ARG
1	A	546	THR
1	A	568	ARG
1	A	579	ASP
1	A	588	THR
1	A	606	GLU
1	A	616	SER
1	A	617	SER
1	A	623	LEU
1	A	637	LEU
1	A	642	ARG
1	A	655	ASN
1	A	658	ASN
1	A	665	VAL
1	A	677	SER
1	A	686	LYS
1	A	720	ASP
1	A	769	GLU
1	B	454	GLU
1	B	457	GLN
1	B	482	PHE
1	B	484	ASN
1	B	547	LEU
1	B	550	ASP
1	B	568	ARG
1	B	606	GLU
1	B	610	ILE
1	B	638	TYR
1	B	658	ASN
1	B	660	SER
1	B	671	THR
1	B	690	ASN
1	B	692	ILE
1	B	709	GLN
1	B	712	PRO

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Mol	Chain	Res	Type
1	B	736	PRO
1	B	766	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	447	HIS
1	A	621	GLN
1	A	641	ASN
1	A	655	ASN
1	A	743	GLN
1	B	461	GLN
1	B	484	ASN
1	B	543	ASN
1	B	545	HIS
1	B	621	GLN
1	B	641	ASN
1	B	661	HIS
1	B	690	ASN
1	B	709	GLN
1	B	761	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

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$
4	X4D	A	803	-	27,31,31	1.46	3 (11%)	22,44,44	2.33	6 (27%)
4	X4D	B	803	-	27,31,31	1.48	4 (14%)	22,44,44	2.19	6 (27%)

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
4	X4D	A	803	-	-	0/9/9/9	0/5/5/5
4	X4D	B	803	-	-	0/9/9/9	0/5/5/5

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	803	X4D	C18-N5	-4.54	1.40	1.46
4	A	803	X4D	C18-N5	-3.55	1.42	1.46
4	A	803	X4D	C12-C8	-2.63	1.37	1.41
4	B	803	X4D	C12-C8	-2.59	1.37	1.41
4	B	803	X4D	C10-N5	-2.37	1.36	1.39
4	B	803	X4D	C22-C21	3.47	1.48	1.37
4	A	803	X4D	C22-C21	3.86	1.49	1.37

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	803	X4D	C16-C12-C8	-6.86	115.28	121.07
4	B	803	X4D	C16-C12-C8	-6.38	115.69	121.07
4	A	803	X4D	C22-C19-C7	-5.06	119.63	121.21
4	A	803	X4D	C17-C15-C10	-3.63	113.24	119.42
4	B	803	X4D	C19-C22-C21	-3.15	114.39	118.88
4	B	803	X4D	C17-C15-C10	-3.14	114.07	119.42
4	A	803	X4D	C19-C22-C21	-2.15	115.83	118.88
4	A	803	X4D	C15-C17-C16	2.15	124.73	121.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	803	X4D	C22-C21-N14	2.16	126.97	123.86
4	A	803	X4D	C22-C21-N14	2.24	127.08	123.86
4	B	803	X4D	C15-C10-C8	2.57	125.68	121.36
4	B	803	X4D	C17-C16-C12	3.87	123.29	118.49

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	803	X4D	1	0
4	B	803	X4D	5	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	324/345 (93%)	0.24	9 (2%) 56 27	21, 42, 78, 98	0
1	B	324/345 (93%)	0.04	5 (1%) 76 49	19, 40, 65, 85	0
All	All	648/690 (93%)	0.14	14 (2%) 65 35	19, 41, 74, 98	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	615	SER	3.3
1	A	616	SER	3.3
1	A	585	LEU	3.1
1	B	585	LEU	3.0
1	A	447	HIS	2.9
1	B	461	GLN	2.8
1	A	621	GLN	2.4
1	A	606	GLU	2.4
1	B	612	SER	2.3
1	A	457	GLN	2.2
1	A	657	ASN	2.1
1	A	617	SER	2.1
1	B	458	GLY	2.0
1	B	501	SER	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. 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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	MG	B	802	1/1	0.89	0.25	2.66	14,14,14,14	0
3	MG	A	802	1/1	0.95	0.26	2.60	19,19,19,19	0
4	X4D	A	803	27/27	0.91	0.22	1.10	30,38,44,44	0
4	X4D	B	803	27/27	0.92	0.24	0.51	42,47,52,56	0
2	ZN	A	801	1/1	0.97	0.14	-	36,36,36,36	0
2	ZN	B	801	1/1	0.98	0.12	-	34,34,34,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.