



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 08:48 AM UTC

PDB ID : 4F08 / pdb_00004f08
Title : Discovery and Optimization of C-2 Methyl Imidazo-pyrrolopyridines as Potent and Orally Bioavailable JAK1 Inhibitors with Selectivity over JAK2
Authors : Murray, J.M.
Deposited on : 2012-05-03
Resolution : 2.82 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

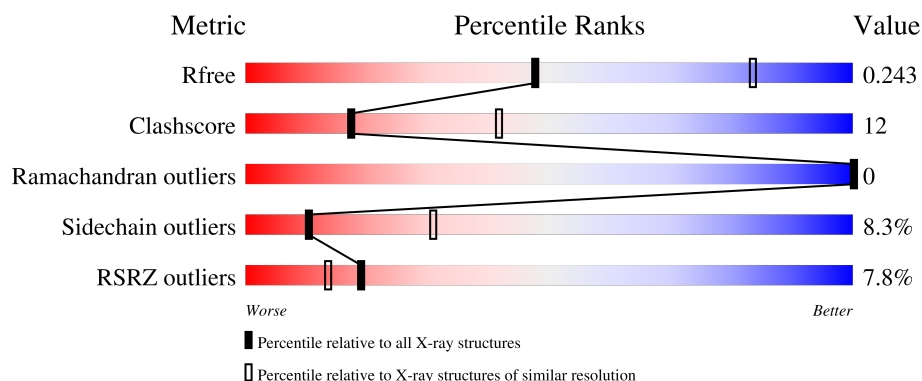
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.82 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4591 (2.84-2.80)
Clashscore	190562	5010 (2.84-2.80)
Ramachandran outliers	187476	4916 (2.84-2.80)
Sidechain outliers	187428	4918 (2.84-2.80)
RSRZ outliers	180081	4594 (2.84-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	302	
1	B	302	

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 4804 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.

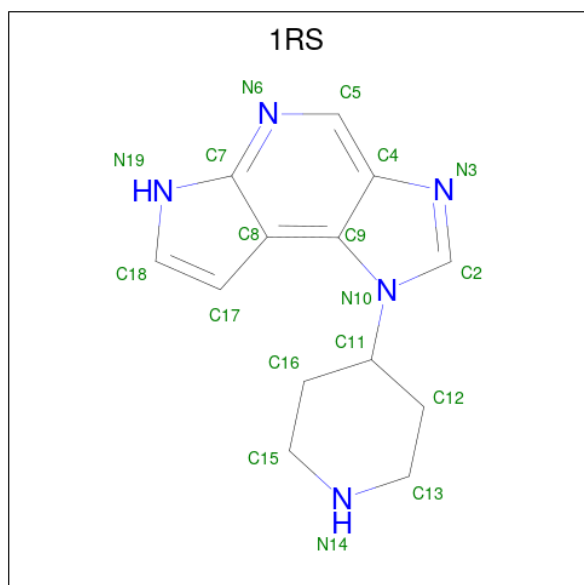
- Molecule 1 is a protein called Tyrosine-protein kinase JAK2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	290	Total	C	N	O	P	S	0	0	0
			2392	1517	412	447	2	14			
1	B	290	Total	C	N	O	P	S	0	0	0
			2376	1505	408	447	2	14			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	831	GLY	-	expression tag	UNP O60674
A	832	SER	-	expression tag	UNP O60674
B	831	GLY	-	expression tag	UNP O60674
B	832	SER	-	expression tag	UNP O60674

- Molecule 2 is 1-(piperidin-4-yl)-1,6-dihydroimidazo[4,5-d]pyrrolo[2,3-b]pyridine (CCD ID: 1RS) (formula: C₁₃H₁₅N₅).

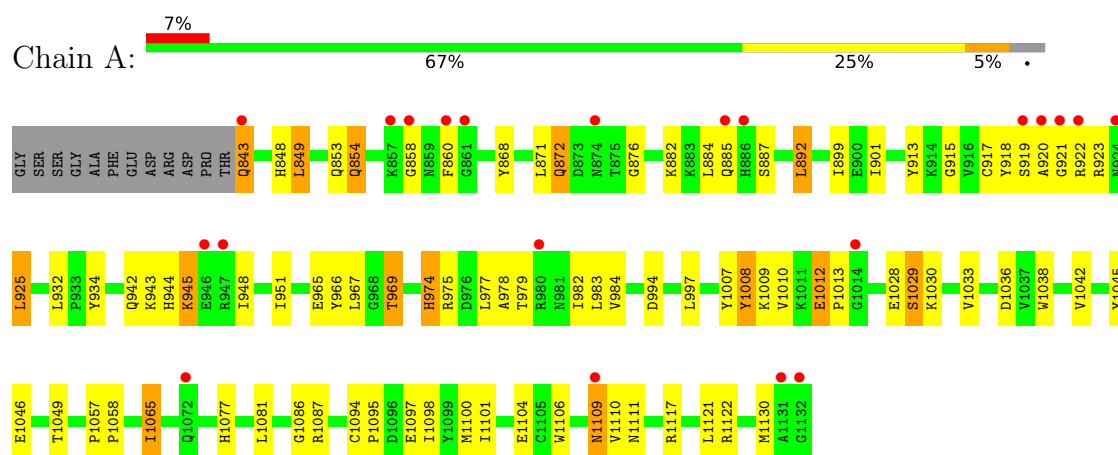


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			18	13	5		
2	B	1	Total	C	N	0	0
			18	13	5		

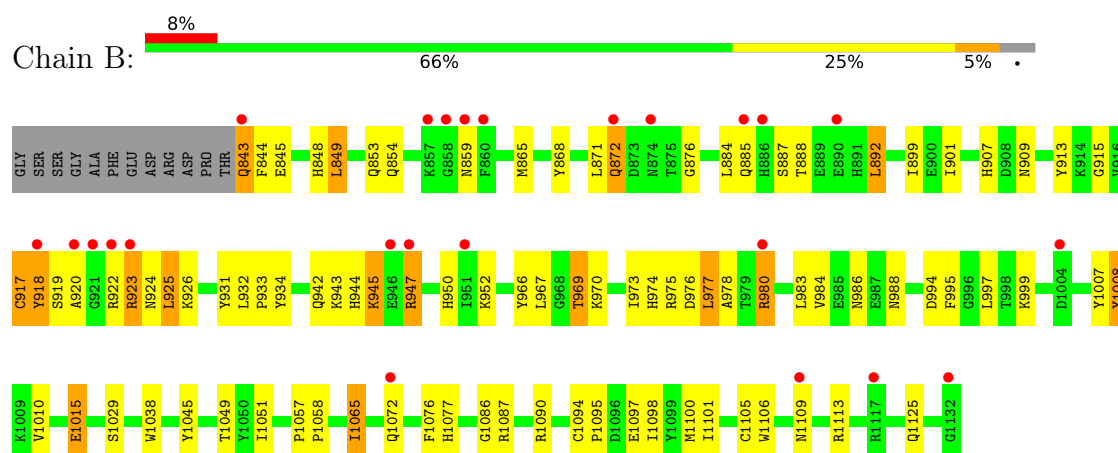
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Tyrosine-protein kinase JAK2



• Molecule 1: Tyrosine-protein kinase JAK2



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	110.92Å 110.92Å 70.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.61 – 2.82 49.61 – 2.82	Depositor EDS
% Data completeness (in resolution range)	99.6 (49.61-2.82) 99.7 (49.61-2.82)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.17 (at 2.81Å)	Xtriage
Refinement program	BUSTER-TNT, BUSTER 2.11.2	Depositor
R, R_{free}	0.213 , 0.241 0.210 , 0.243	Depositor DCC
R_{free} test set	1034 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	54.7	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.031 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4804	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PTR, 1RS

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	0.73	0/2409	1.00	6/3240 (0.2%)
1	B	0.76	3/2392 (0.1%)	0.98	4/3220 (0.1%)
All	All	0.75	3/4801 (0.1%)	0.99	10/6460 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	978	ALA	C-O	-5.36	1.18	1.23
1	B	950	HIS	CG-ND1	-5.26	1.32	1.38
1	B	974	HIS	CG-ND1	-5.04	1.32	1.38

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	974	HIS	N-CA-C	6.79	118.48	111.14
1	A	922	ARG	CB-CA-C	-6.01	100.09	110.72
1	B	872	GLN	N-CA-C	-5.95	104.05	111.90
1	A	1110	VAL	N-CA-C	5.58	116.97	110.62
1	A	948	ILE	N-CA-C	5.57	116.42	107.73
1	B	923	ARG	N-CA-C	-5.41	99.27	110.80
1	B	980	ARG	N-CA-C	-5.39	105.40	111.28
1	A	872	GLN	N-CA-C	-5.33	105.42	112.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	858	GLY	N-CA-C	-5.24	103.32	111.12
1	B	950	HIS	N-CA-C	5.07	116.80	111.28

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	947	ARG	Sidechain

5.2 Too-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 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	2392	0	2356	60	3
1	B	2376	0	2328	59	3
2	A	18	0	15	1	0
2	B	18	0	15	2	0
All	All	4804	0	4714	119	3

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 12.

All (119) 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:B:920:ALA:H	1:B:923:ARG:HA	1.26	0.99
1:B:907:HIS:CD2	1:B:909:ASN:H	1.81	0.98
1:A:951:ILE:HD13	1:A:1130:MET:SD	2.05	0.96
1:A:1104:GLU:OE2	1:A:1122:ARG:NH1	2.01	0.93
1:B:907:HIS:HD2	1:B:909:ASN:H	1.10	0.93
1:B:999:LYS:HD2	1:B:1008:PTR:HE2	1.51	0.91
1:A:951:ILE:HD11	1:A:1130:MET:HB3	1.56	0.86
1:A:1104:GLU:CD	1:A:1122:ARG:HH12	1.86	0.84
1:A:951:ILE:CD1	1:A:1130:MET:SD	2.72	0.78
1:A:1117:ARG:O	1:A:1121:LEU:HD13	1.84	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:951:ILE:CD1	1:A:1130:MET:HB3	2.17	0.75
1:B:1086:GLY:O	1:B:1087:ARG:HD3	1.88	0.74
1:B:907:HIS:HD2	1:B:909:ASN:N	1.85	0.72
1:A:965:GLU:OE1	1:A:1117:ARG:HG3	1.90	0.72
1:B:920:ALA:H	1:B:923:ARG:CA	2.04	0.68
1:A:943:LYS:HG2	1:A:944:HIS:CE1	2.31	0.66
1:A:1008:PTR:HE1	1:A:1010:VAL:HB	1.77	0.66
1:A:975:ARG:NH2	1:A:1012:GLU:OE1	2.21	0.66
1:A:966:TYR:O	1:A:969:THR:HB	1.96	0.65
1:B:843:GLN:NE2	1:B:918:TYR:OH	2.29	0.65
1:A:1121:LEU:HD21	1:B:888:THR:HG22	1.79	0.65
1:B:983:LEU:HD21	2:B:1201:1RS:H14	1.79	0.63
1:B:966:TYR:O	1:B:969:THR:HB	1.99	0.62
1:A:843:GLN:N	1:A:843:GLN:CD	2.58	0.61
1:A:1081:LEU:HD11	1:A:1086:GLY:HA3	1.82	0.61
1:B:1045:TYR:O	1:B:1049:THR:HG23	2.02	0.60
1:B:859:ASN:ND2	1:B:1015:GLU:HG3	2.17	0.59
1:A:1081:LEU:CD1	1:A:1086:GLY:HA3	2.33	0.59
1:B:942:GLN:O	1:B:945:LYS:HD2	2.01	0.59
1:A:942:GLN:O	1:A:945:LYS:HD2	2.02	0.59
1:A:951:ILE:CD1	1:A:1130:MET:CB	2.80	0.59
1:B:844:PHE:O	1:B:917:CYS:HB2	2.03	0.58
1:B:920:ALA:HA	1:B:923:ARG:CB	2.34	0.58
2:B:1201:1RS:H15	2:B:1201:1RS:H13	1.85	0.57
1:A:983:LEU:HD21	2:A:1201:1RS:H14	1.86	0.56
1:B:943:LYS:HG2	1:B:944:HIS:CE1	2.41	0.55
1:B:913:TYR:CZ	1:B:915:GLY:HA2	2.42	0.55
1:B:999:LYS:HD2	1:B:1008:PTR:CE2	2.32	0.55
1:A:868:TYR:O	1:A:876:GLY:HA3	2.07	0.54
1:B:868:TYR:O	1:B:876:GLY:HA3	2.08	0.54
1:B:919:SER:O	1:B:920:ALA:HB3	2.08	0.53
1:B:919:SER:HB2	1:B:922:ARG:O	2.09	0.53
1:B:865:MET:HE3	1:B:931:TYR:CZ	2.43	0.52
1:B:918:TYR:CD1	1:B:918:TYR:N	2.77	0.52
1:A:1086:GLY:O	1:A:1087:ARG:HD3	2.10	0.52
1:B:849:LEU:HD13	1:B:868:TYR:HD1	1.75	0.51
1:A:1038:TRP:CE3	1:A:1106:TRP:HA	2.46	0.51
1:B:907:HIS:CD2	1:B:909:ASN:N	2.64	0.51
1:A:978:ALA:O	1:A:982:ILE:HG12	2.10	0.50
1:A:932:LEU:HD12	1:A:983:LEU:HB2	1.94	0.50
1:A:1109:ASN:ND2	1:A:1111:ASN:OD1	2.45	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:920:ALA:N	1:B:923:ARG:HA	2.10	0.50
1:B:924:ASN:O	1:B:926:LYS:HG3	2.13	0.49
1:B:1098:ILE:HA	1:B:1101:ILE:HD12	1.95	0.49
1:B:1105:CYS:O	1:B:1113:ARG:HD3	2.13	0.49
1:A:974:HIS:O	1:A:975:ARG:HB2	2.13	0.49
1:B:976:ASP:O	1:B:977:LEU:C	2.55	0.49
1:B:871:LEU:O	1:B:872:GLN:C	2.56	0.48
1:B:907:HIS:CD2	1:B:909:ASN:HB2	2.49	0.48
1:B:934:TYR:CE2	1:B:986:ASN:HA	2.48	0.48
1:B:884:LEU:HD12	1:B:892:LEU:HD13	1.96	0.47
1:A:899:ILE:CD1	1:A:925:LEU:HD13	2.43	0.47
1:B:934:TYR:HB2	1:B:984:VAL:O	2.14	0.47
1:B:1090:ARG:HD2	1:B:1094:CYS:O	2.14	0.46
1:B:1057:PRO:HB2	1:B:1058:PRO:HD3	1.98	0.46
1:A:975:ARG:HD3	1:A:997:LEU:O	2.15	0.46
1:A:913:TYR:CZ	1:A:915:GLY:HA2	2.51	0.46
1:A:1065:ILE:HD13	1:A:1077:HIS:HB2	1.98	0.45
1:B:859:ASN:ND2	1:B:1015:GLU:CG	2.79	0.45
1:A:918:TYR:CD1	1:A:918:TYR:O	2.69	0.45
1:B:973:ILE:HG22	1:B:975:ARG:HG3	1.97	0.45
1:B:1100:MET:HE3	1:B:1100:MET:HB2	1.80	0.45
1:A:977:LEU:HD12	1:A:977:LEU:HA	1.79	0.45
1:A:871:LEU:O	1:A:872:GLN:C	2.60	0.45
1:A:849:LEU:HD13	1:A:868:TYR:HD1	1.80	0.44
1:A:1045:TYR:O	1:A:1049:THR:HG23	2.17	0.44
1:A:1094:CYS:HA	1:A:1095:PRO:HD3	1.89	0.44
1:B:925:LEU:HD23	1:B:925:LEU:HA	1.69	0.44
1:A:1100:MET:HE3	1:A:1100:MET:HB2	1.82	0.44
1:B:859:ASN:CG	1:B:1015:GLU:HG3	2.42	0.44
1:B:899:ILE:CD1	1:B:925:LEU:HD13	2.47	0.44
1:B:1065:ILE:HD13	1:B:1077:HIS:HB2	1.99	0.44
1:A:848:HIS:HB2	1:A:868:TYR:CE1	2.53	0.43
1:A:1057:PRO:HB2	1:A:1058:PRO:HD3	1.99	0.43
1:B:969:THR:HG22	1:B:970:LYS:CG	2.49	0.43
1:B:994:ASP:C	1:B:995:PHE:CD1	2.97	0.43
1:A:943:LYS:CG	1:A:944:HIS:CE1	3.00	0.43
1:A:1098:ILE:HA	1:A:1101:ILE:HD12	2.00	0.43
1:A:943:LYS:CD	1:A:944:HIS:CE1	3.02	0.43
1:A:934:TYR:HB2	1:A:984:VAL:O	2.19	0.42
1:B:865:MET:HE3	1:B:931:TYR:CE2	2.54	0.42
1:A:951:ILE:HD11	1:A:1130:MET:SD	2.58	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:920:ALA:N	1:A:921:GLY:HA3	2.34	0.42
1:A:1009:LYS:HA	1:A:1029:SER:O	2.19	0.42
1:B:1051:ILE:HG22	1:B:1051:ILE:O	2.20	0.42
1:B:1094:CYS:HA	1:B:1095:PRO:HD3	1.93	0.42
1:A:1012:GLU:HA	1:A:1013:PRO:HD3	1.92	0.42
1:A:1042:VAL:O	1:A:1045:TYR:HB3	2.20	0.42
1:B:932:LEU:HA	1:B:933:PRO:HD3	1.70	0.42
1:B:1072:GLN:HB3	1:B:1076:PHE:CE2	2.55	0.42
1:B:848:HIS:HB2	1:B:868:TYR:CE1	2.55	0.42
1:B:976:ASP:OD1	1:B:980:ARG:NH2	2.53	0.42
1:B:1125:GLN:HA	1:B:1125:GLN:OE1	2.19	0.42
1:A:965:GLU:OE1	1:A:1117:ARG:CG	2.65	0.41
1:A:1087:ARG:HD3	1:A:1087:ARG:HA	1.87	0.41
1:A:854:GLN:HE21	1:A:854:GLN:HB3	1.72	0.41
1:B:952:LYS:NZ	1:B:988:ASN:OD1	2.49	0.41
1:A:1033:VAL:O	1:A:1036:ASP:HB2	2.21	0.41
1:B:1038:TRP:CE3	1:B:1106:TRP:HA	2.56	0.41
1:A:884:LEU:HD12	1:A:892:LEU:HD13	2.01	0.41
1:B:975:ARG:NH1	1:B:997:LEU:O	2.52	0.41
1:A:921:GLY:C	1:A:923:ARG:H	2.28	0.41
1:A:974:HIS:HE1	1:A:994:ASP:O	2.04	0.41
1:A:932:LEU:HD12	1:A:983:LEU:CB	2.50	0.41
1:A:849:LEU:HD12	1:A:849:LEU:HA	1.76	0.41
1:A:925:LEU:HA	1:A:925:LEU:HD23	1.80	0.40
1:A:979:THR:HG23	1:A:1046:GLU:OE2	2.21	0.40
1:A:882:LYS:HE3	1:A:994:ASP:OD1	2.22	0.40
1:B:849:LEU:HD12	1:B:849:LEU:HA	1.94	0.40

All (3) 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:853:GLN:OE1	1:B:947:ARG:NH2[1_554]	1.90	0.30
1:A:853:GLN:OE1	1:B:947:ARG:CZ[1_554]	2.05	0.15
1:A:853:GLN:OE1	1:B:947:ARG:NH1[1_554]	2.16	0.04

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	286/302 (95%)	275 (96%)	11 (4%)	0	100	100
1	B	286/302 (95%)	278 (97%)	8 (3%)	0	100	100
All	All	572/604 (95%)	553 (97%)	19 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	260/271 (96%)	239 (92%)	21 (8%)	11	32
1	B	257/271 (95%)	235 (91%)	22 (9%)	10	29
All	All	517/542 (95%)	474 (92%)	43 (8%)	10	31

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	843	GLN
1	A	849	LEU
1	A	854	GLN
1	A	860	PHE
1	A	885	GLN
1	A	887	SER
1	A	892	LEU
1	A	901	ILE

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Mol	Chain	Res	Type
1	A	917	CYS
1	A	919	SER
1	A	925	LEU
1	A	945	LYS
1	A	967	LEU
1	A	969	THR
1	A	1012	GLU
1	A	1028	GLU
1	A	1029	SER
1	A	1030	LYS
1	A	1065	ILE
1	A	1097	GLU
1	A	1109	ASN
1	B	843	GLN
1	B	845	GLU
1	B	849	LEU
1	B	853	GLN
1	B	854	GLN
1	B	885	GLN
1	B	887	SER
1	B	892	LEU
1	B	901	ILE
1	B	917	CYS
1	B	918	TYR
1	B	925	LEU
1	B	945	LYS
1	B	967	LEU
1	B	969	THR
1	B	977	LEU
1	B	1010	VAL
1	B	1015	GLU
1	B	1029	SER
1	B	1065	ILE
1	B	1097	GLU
1	B	1109	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	848	HIS
1	A	854	GLN
1	A	859	ASN

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Mol	Chain	Res	Type
1	A	874	ASN
1	A	906	GLN
1	A	909	ASN
1	A	944	HIS
1	A	955	GLN
1	A	974	HIS
1	A	1067	ASN
1	A	1077	HIS
1	A	1109	ASN
1	A	1112	GLN
1	B	843	GLN
1	B	848	HIS
1	B	854	GLN
1	B	872	GLN
1	B	874	ASN
1	B	906	GLN
1	B	907	HIS
1	B	909	ASN
1	B	924	ASN
1	B	944	HIS
1	B	955	GLN
1	B	1067	ASN
1	B	1077	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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
1	PTR	B	1008	1	15,16,17	1.66	1 (6%)	17,22,24	0.71	1 (5%)
1	PTR	A	1008	1	15,16,17	1.72	1 (6%)	17,22,24	1.24	3 (17%)
1	PTR	B	1007	1	15,16,17	1.48	1 (6%)	17,22,24	0.91	1 (5%)
1	PTR	A	1007	1	15,16,17	1.64	1 (6%)	17,22,24	0.48	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	PTR	B	1008	1	-	3/10/11/13	0/1/1/1
1	PTR	A	1008	1	-	1/10/11/13	0/1/1/1
1	PTR	B	1007	1	-	0/10/11/13	0/1/1/1
1	PTR	A	1007	1	-	0/10/11/13	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1008	PTR	OH-CZ	-5.98	1.27	1.40
1	A	1008	PTR	OH-CZ	-5.94	1.27	1.40
1	A	1007	PTR	OH-CZ	-5.87	1.27	1.40
1	B	1007	PTR	OH-CZ	-5.05	1.29	1.40

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1007	PTR	O3P-P-OH	2.85	113.73	105.32
1	A	1008	PTR	OH-CZ-CE2	2.69	127.27	119.22
1	A	1008	PTR	O3P-P-OH	2.36	112.28	105.32
1	B	1008	PTR	O3P-P-OH	2.26	112.00	105.32
1	A	1008	PTR	OH-CZ-CE1	-2.16	112.75	119.22

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	1008	PTR	CZ-OH-P-O1P
1	B	1008	PTR	CA-CB-CG-CD1
1	B	1008	PTR	CA-CB-CG-CD2
1	B	1008	PTR	CZ-OH-P-O1P

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	1008	PTR	2	0
1	A	1008	PTR	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	1RS	A	1201	-	21,21,21	1.31	3 (14%)	24,30,30	2.92	14 (58%)
2	1RS	B	1201	-	21,21,21	1.21	3 (14%)	24,30,30	2.83	13 (54%)

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	1RS	A	1201	-	-	1/4/12/12	0/4/4/4
2	1RS	B	1201	-	-	1/4/12/12	0/4/4/4

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1201	1RS	C4-N3	-3.24	1.33	1.39
2	B	1201	1RS	C4-N3	-2.80	1.34	1.39
2	A	1201	1RS	C5-N6	-2.31	1.29	1.34
2	B	1201	1RS	C5-N6	-2.25	1.29	1.34
2	B	1201	1RS	C7-N6	-2.21	1.31	1.35
2	A	1201	1RS	C7-N6	-2.15	1.31	1.35

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1201	1RS	C8-C7-N6	-7.27	120.71	126.40
2	A	1201	1RS	C8-C7-N6	-7.18	120.79	126.40
2	A	1201	1RS	C12-C11-N10	5.26	117.22	111.40
2	A	1201	1RS	C16-C11-N10	4.46	116.34	111.40
2	B	1201	1RS	C12-C11-N10	4.11	115.95	111.40
2	B	1201	1RS	C4-N3-C2	3.87	109.16	104.40
2	B	1201	1RS	C8-C9-C4	-3.85	116.89	121.19
2	A	1201	1RS	C4-N3-C2	3.68	108.92	104.40
2	B	1201	1RS	C16-C11-N10	3.66	115.46	111.40
2	A	1201	1RS	C8-C9-C4	-3.65	117.10	121.19
2	B	1201	1RS	C7-N19-C18	3.35	110.05	108.11
2	A	1201	1RS	C5-N6-C7	3.32	122.70	116.19
2	A	1201	1RS	C13-C12-C11	-3.27	104.82	110.77
2	B	1201	1RS	N19-C7-N6	3.16	131.81	126.69
2	B	1201	1RS	C9-N10-C2	3.06	107.17	105.33
2	B	1201	1RS	C15-C16-C11	-3.01	105.30	110.77
2	B	1201	1RS	C5-N6-C7	2.99	122.05	116.19
2	A	1201	1RS	N19-C7-N6	2.99	131.53	126.69
2	A	1201	1RS	C11-N10-C2	-2.98	122.02	128.01
2	B	1201	1RS	C11-N10-C2	-2.91	122.15	128.01
2	A	1201	1RS	C7-N19-C18	2.87	109.77	108.11
2	A	1201	1RS	C9-N10-C2	2.84	107.03	105.33
2	A	1201	1RS	C15-C16-C11	-2.73	105.81	110.77
2	B	1201	1RS	C13-C12-C11	-2.58	106.08	110.77
2	B	1201	1RS	N10-C2-N3	-2.27	110.72	113.94
2	A	1201	1RS	N10-C2-N3	-2.12	110.92	113.94
2	A	1201	1RS	C9-C8-C17	2.06	135.58	132.05

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1201	1RS	C16-C11-N10-C2
2	B	1201	1RS	C16-C11-N10-C2

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1201	1RS	1	0
2	B	1201	1RS	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	288/302 (95%)	0.17	21 (7%) 21 15	30, 51, 109, 167	0
1	B	288/302 (95%)	0.32	24 (8%) 17 12	32, 51, 107, 167	0
All	All	576/604 (95%)	0.24	45 (7%) 19 13	30, 51, 108, 167	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	859	ASN	6.7
1	A	860	PHE	5.4
1	B	886	HIS	5.3
1	A	921	GLY	5.0
1	B	843	GLN	4.5
1	A	922	ARG	4.4
1	A	920	ALA	4.1
1	B	860	PHE	4.1
1	B	885	GLN	4.0
1	A	886	HIS	3.8
1	B	890	GLU	3.6
1	A	924	ASN	3.5
1	B	918	TYR	3.3
1	B	920	ALA	3.1
1	B	921	GLY	3.1
1	A	947	ARG	3.0
1	B	1109	ASN	3.0
1	B	947	ARG	2.9
1	B	946	GLU	2.9
1	B	922	ARG	2.8
1	A	919	SER	2.7
1	B	858	GLY	2.6
1	B	857	LYS	2.6
1	A	1132	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	872	GLN	2.6
1	B	1004	ASP	2.6
1	B	1132	GLY	2.5
1	A	843	GLN	2.5
1	A	874	ASN	2.4
1	B	874	ASN	2.4
1	A	858	GLY	2.3
1	B	923	ARG	2.3
1	B	1072	GLN	2.3
1	B	980	ARG	2.3
1	A	1072	GLN	2.2
1	B	1117	ARG	2.2
1	A	861	GLY	2.2
1	A	1014	GLY	2.2
1	A	885	GLN	2.1
1	A	980	ARG	2.1
1	A	946	GLU	2.1
1	B	951	ILE	2.1
1	A	857	LYS	2.1
1	A	1109	ASN	2.1
1	A	1131	ALA	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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($\text{\AA}^2$)	Q<0.9
1	PTR	A	1007	16/17	0.90	0.10	41,57,103,109	0
1	PTR	B	1007	16/17	0.91	0.11	54,59,99,104	0
1	PTR	B	1008	16/17	0.92	0.11	51,82,101,103	0
1	PTR	A	1008	16/17	0.94	0.09	34,61,77,79	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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($\text{\AA}^2$)	Q<0.9
2	1RS	A	1201	18/18	0.93	0.10	20,20,20,20	0
2	1RS	B	1201	18/18	0.95	0.13	20,20,20,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.