



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 30, 2017 – 08:31 PM EST

PDB ID : 5TQ6
Title : Design and Synthesis of a pan-JAK Kinase Inhibitor Clinical Candidate (PF-06263276) Suitable for Inhaled and Topical Delivery for the Treatment of Inflammatory Diseases of the Lungs and Skin
Authors : Chrencik, J.; Jones, P.
Deposited on : 2016-10-23
Resolution : 2.06 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	:	1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20028442
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)	:	rb-20028442

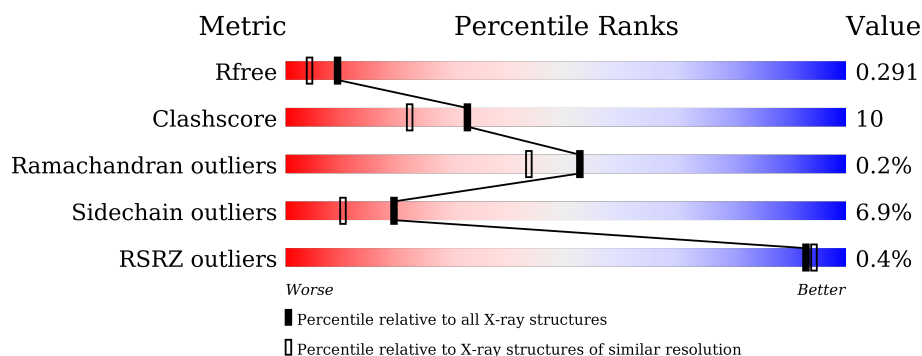
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.06 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	1799 (2.08-2.04)
Clashscore	102246	1910 (2.08-2.04)
Ramachandran outliers	100387	1893 (2.08-2.04)
Sidechain outliers	100360	1893 (2.08-2.04)
RSRZ outliers	91569	1802 (2.08-2.04)

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	304	<div> <div>59%</div> <div>29%</div> <div>• 9%</div> </div>
1	B	304	<div> <div>67%</div> <div>20%</div> <div>• 10%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4846 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.

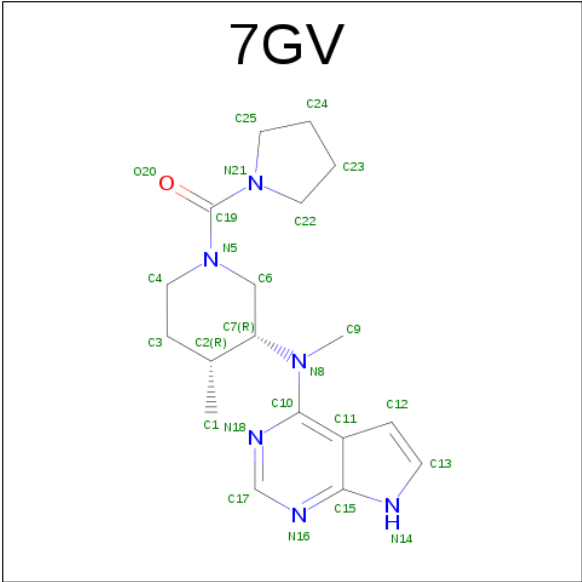
- Molecule 1 is a protein called Tyrosine-protein kinase JAK2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	276	Total	C	N	O	P	S	0	0	0
			2296	1459	394	428	2	13			
1	B	273	Total	C	N	O	P	S	0	0	0
			2277	1449	391	422	2	13			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	829	MET	-	initiating methionine	UNP O60674
A	830	GLY	-	expression tag	UNP O60674
A	831	HIS	-	expression tag	UNP O60674
A	832	HIS	-	expression tag	UNP O60674
A	833	HIS	-	expression tag	UNP O60674
A	834	HIS	-	expression tag	UNP O60674
A	835	HIS	-	expression tag	UNP O60674
A	836	HIS	-	expression tag	UNP O60674
A	1073	SER	MET	engineered mutation	UNP O60674
A	1076	THR	PHE	engineered mutation	UNP O60674
B	829	MET	-	initiating methionine	UNP O60674
B	830	GLY	-	expression tag	UNP O60674
B	831	HIS	-	expression tag	UNP O60674
B	832	HIS	-	expression tag	UNP O60674
B	833	HIS	-	expression tag	UNP O60674
B	834	HIS	-	expression tag	UNP O60674
B	835	HIS	-	expression tag	UNP O60674
B	836	HIS	-	expression tag	UNP O60674
B	1073	SER	MET	engineered mutation	UNP O60674
B	1076	THR	PHE	engineered mutation	UNP O60674

- Molecule 2 is {(3R,4R)-4-methyl-3-[methyl(7H-pyrrolo[2,3-d]pyrimidin-4-yl)amino]piperidin-1-yl}(pyrrolidin-1-yl)methanone (three-letter code: 7GV) (formula: C₁₈H₂₆N₆O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			25	18	6	1		
2	B	1	Total	C	N	O	0	0
			25	18	6	1		

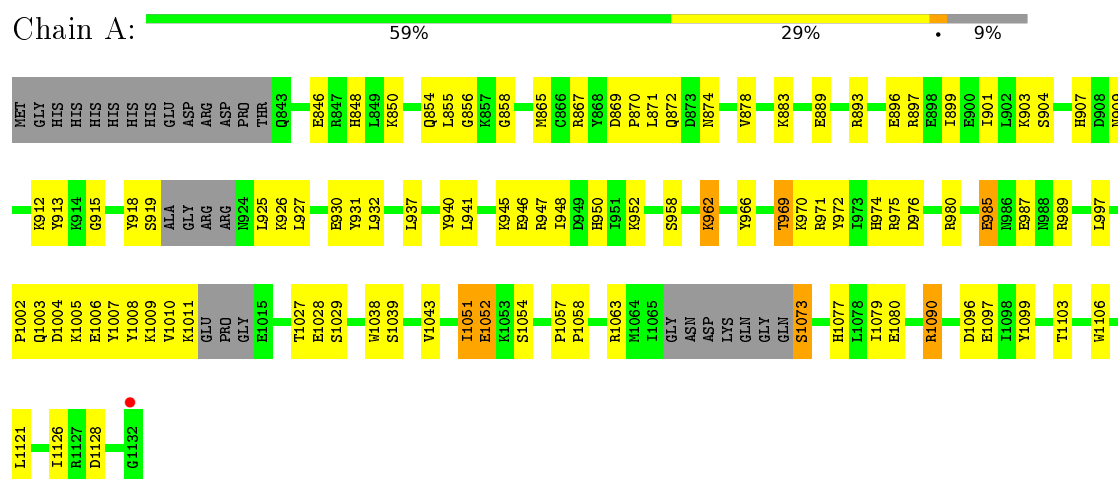
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	109	Total	O	0	0
			109	109		
3	B	114	Total	O	0	0
			114	114		

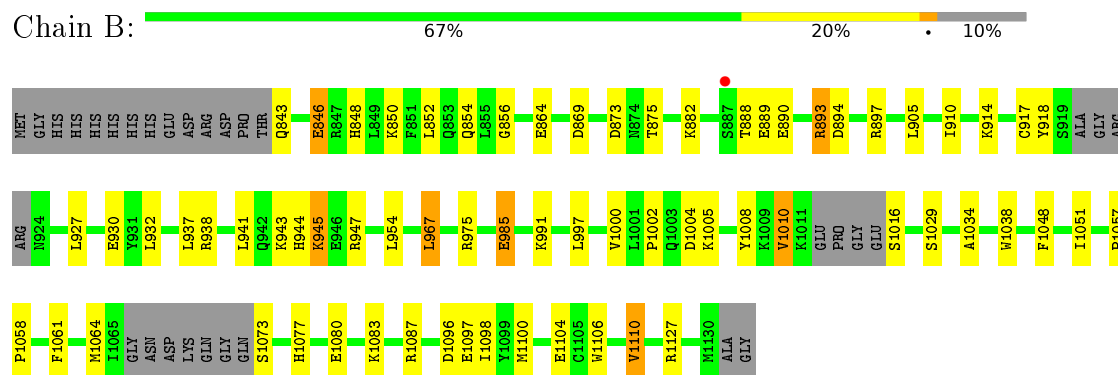
3 Residue-property plots

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: Tyrosine-protein kinase JAK2



• Molecule 1: Tyrosine-protein kinase JAK2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	36.98Å 183.40Å 52.58Å 90.00° 110.59° 90.00°	Depositor
Resolution (Å)	47.54 – 2.06 47.54 – 2.06	Depositor EDS
% Data completeness (in resolution range)	92.6 (47.54-2.06) 92.6 (47.54-2.06)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.54 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.257 , 0.289 0.259 , 0.291	Depositor DCC
R_{free} test set	1900 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	18.6	Xtriage
Anisotropy	0.103	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 29.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	0.486 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4846	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.64% 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: 7GV, PTR

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	0.58	0/2308	0.72	1/3101 (0.0%)
1	B	0.57	0/2289	0.73	1/3077 (0.0%)
All	All	0.58	0/4597	0.73	2/6178 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1051	ILE	CG1-CB-CG2	-5.24	99.88	111.40
1	B	967	LEU	CA-CB-CG	5.12	127.08	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	2296	0	2274	55	0
1	B	2277	0	2261	41	0
2	A	25	0	0	1	0
2	B	25	0	0	0	0
3	A	109	0	0	2	0
3	B	114	0	0	5	0
All	All	4846	0	4535	96	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 10.

All (96) 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:985:GLU:HG3	1:B:991:LYS:HE3	1.65	0.78
1:B:941:LEU:O	1:B:945:LYS:HB3	1.85	0.75
1:B:894:ASP:OD1	1:B:897:ARG:NH2	2.21	0.73
1:B:1096:ASP:O	1:B:1100:MET:HG3	1.88	0.73
1:A:1052:GLU:HG3	3:B:4104:HOH:O	1.89	0.73
1:B:854:GLN:NE2	1:B:856:GLY:O	2.22	0.72
1:A:904:SER:O	1:A:970:LYS:NZ	2.24	0.70
1:A:883:LYS:HD3	1:A:926:LYS:HD3	1.73	0.68
1:B:846:GLU:HB2	1:B:917:CYS:SG	2.34	0.68
1:A:1054:SER:O	1:A:1063:ARG:NH2	2.27	0.67
1:A:1051:ILE:O	1:A:1051:ILE:HG22	1.95	0.66
1:B:905:LEU:HD13	1:B:910:ILE:HD13	1.78	0.66
1:B:938:ARG:HA	1:B:1051:ILE:HD13	1.78	0.66
1:A:947:ARG:HH11	1:A:987:GLU:HG3	1.61	0.65
1:B:985:GLU:HG3	1:B:991:LYS:CE	2.26	0.65
1:A:850:LYS:HE3	1:A:869:ASP:HB3	1.80	0.64
1:A:1097:GLU:HG3	1:A:1126:ILE:HG12	1.81	0.62
1:B:914:LYS:HE3	1:B:930:GLU:HA	1.79	0.62
1:B:1061:PHE:HA	1:B:1064:MET:HE3	1.85	0.58
1:A:1009:LYS:HE3	1:A:1028:GLU:O	2.04	0.58
1:B:1004:ASP:HB3	1:B:1005:LYS:HE3	1.86	0.58
1:B:1080:GLU:HG2	3:B:4109:HOH:O	2.03	0.58
1:A:1073:SER:O	1:A:1077:HIS:HD2	1.87	0.58
1:A:907:HIS:HE1	1:A:909:ASN:HD22	1.53	0.56
1:A:1011:LYS:HD2	1:A:1029:SER:OG	2.05	0.56
1:A:941:LEU:HD23	1:A:1051:ILE:CD1	2.36	0.56
1:B:893:ARG:HD3	3:B:4129:HOH:O	2.07	0.55
1:B:1010:VAL:O	1:B:1029:SER:HB3	2.06	0.55
1:A:940:TYR:OH	1:A:947:ARG:NH1	2.39	0.55
1:A:985:GLU:HB3	1:A:989:ARG:O	2.07	0.55
1:A:1028:GLU:HG2	3:A:4145:HOH:O	2.07	0.55
1:B:1034:ALA:HB3	1:B:1110:VAL:HG22	1.87	0.54
1:B:846:GLU:O	1:B:846:GLU:HG3	2.07	0.54
1:A:941:LEU:HD23	1:A:1051:ILE:HD13	1.91	0.53
1:A:850:LYS:NZ	1:A:872:GLN:OE1	2.35	0.52
1:B:1048:PHE:CZ	1:B:1098:ILE:HD13	2.44	0.52
1:B:854:GLN:OE1	1:B:864:GLU:HG2	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:850:LYS:HD3	1:B:869:ASP:HB2	1.92	0.52
1:A:1057:PRO:HB2	1:A:1058:PRO:HD3	1.91	0.51
1:A:1002:PRO:HD3	1:A:1008:PTR:HD2	1.92	0.51
1:A:1005:LYS:HE2	1:A:1007:PTR:O3P	2.11	0.51
1:A:865:MET:HE2	1:A:878:VAL:HG11	1.93	0.51
1:A:932:LEU:HD22	1:A:985:GLU:HA	1.92	0.50
1:B:890:GLU:HB3	3:B:4190:HOH:O	2.11	0.50
1:A:1005:LYS:NZ	1:A:1007:PTR:O3P	2.45	0.49
1:A:855:LEU:HD11	1:A:931:TYR:CE1	2.47	0.49
1:B:914:LYS:HE3	1:B:930:GLU:CA	2.43	0.48
1:B:848:HIS:HB3	1:B:869:ASP:C	2.34	0.48
1:B:938:ARG:HA	1:B:1051:ILE:CD1	2.43	0.48
1:A:854:GLN:NE2	1:A:856:GLY:O	2.46	0.47
1:A:974:HIS:O	1:A:975:ARG:HB2	2.14	0.47
1:B:1034:ALA:CB	1:B:1110:VAL:HG22	2.45	0.47
1:B:1073:SER:O	1:B:1077:HIS:ND1	2.37	0.47
1:A:958:SER:O	1:A:962:LYS:HG3	2.15	0.47
1:A:1099:TYR:O	1:A:1103:THR:HG23	2.15	0.47
1:B:932:LEU:HD13	1:B:985:GLU:HG2	1.96	0.47
1:A:1004:ASP:OD1	1:A:1005:LYS:HG3	2.16	0.46
1:A:1027:THR:HG22	1:A:1079:ILE:HD13	1.96	0.46
1:A:901:ILE:HD12	1:A:972:TYR:CZ	2.50	0.46
1:B:873:ASP:OD1	1:B:875:THR:OG1	2.27	0.46
1:A:966:TYR:O	1:A:969:THR:HG22	2.16	0.45
1:B:944:HIS:CD2	1:B:947:ARG:NH1	2.84	0.45
1:A:846:GLU:OE2	1:A:919:SER:HB3	2.16	0.45
1:B:850:LYS:HD3	1:B:869:ASP:CB	2.46	0.45
1:A:901:ILE:HD12	1:A:972:TYR:CE1	2.52	0.45
1:A:855:LEU:HD11	1:A:931:TYR:HE1	1.82	0.44
1:A:1038:TRP:CE3	1:A:1106:TRP:HA	2.51	0.44
1:A:952:LYS:NZ	1:A:987:GLU:OE2	2.51	0.44
1:A:867:ARG:HD3	1:A:874:ASN:HA	2.00	0.44
1:A:913:TYR:CZ	1:A:915:GLY:HA2	2.53	0.44
1:A:858:GLY:HA3	2:A:4000:7GV:C23	2.48	0.44
1:B:954:LEU:HB2	1:B:1127:ARG:HD2	1.98	0.44
1:A:1090:ARG:NH2	1:A:1096:ASP:CG	2.72	0.43
1:B:1100:MET:O	1:B:1104:GLU:HG3	2.18	0.43
1:B:975:ARG:HD3	1:B:997:LEU:O	2.17	0.43
1:B:1008:PTR:HD1	1:B:1008:PTR:C	2.47	0.43
1:B:1038:TRP:CE3	1:B:1106:TRP:HA	2.53	0.43
1:B:882:LYS:HB3	1:B:927:LEU:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1005:LYS:CE	1:A:1007:PTR:O3P	2.66	0.43
1:A:976:ASP:OD1	1:A:980:ARG:NH2	2.51	0.43
1:A:899:ILE:HG12	1:A:927:LEU:HD13	2.01	0.43
1:A:1039:SER:O	1:A:1043:VAL:HG23	2.19	0.42
1:A:903:LYS:HD2	3:A:4115:HOH:O	2.19	0.42
1:A:1051:ILE:O	1:A:1051:ILE:CG2	2.64	0.42
1:A:945:LYS:HA	1:A:948:ILE:HD12	2.01	0.42
1:B:848:HIS:O	1:B:869:ASP:N	2.39	0.42
1:A:975:ARG:HD3	1:A:997:LEU:O	2.19	0.42
1:A:893:ARG:HG2	1:A:897:ARG:CZ	2.49	0.42
1:A:1010:VAL:O	1:A:1029:SER:HB3	2.20	0.42
1:B:1000:VAL:HG22	3:B:4125:HOH:O	2.19	0.41
1:B:1057:PRO:HB2	1:B:1058:PRO:HD3	2.02	0.41
1:A:848:HIS:CG	1:A:870:PRO:HA	2.55	0.41
1:B:1051:ILE:O	1:B:1051:ILE:HG22	2.20	0.41
1:A:912:LYS:HD3	1:A:930:GLU:CD	2.41	0.41
1:B:888:THR:HG22	1:B:889:GLU:N	2.35	0.40
1:A:971:ARG:NH2	1:A:1005:LYS:O	2.54	0.40

There are no symmetry-related clashes.

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	266/304 (88%)	257 (97%)	9 (3%)	0	100	100
1	B	263/304 (86%)	250 (95%)	12 (5%)	1 (0%)	39	28
All	All	529/608 (87%)	507 (96%)	21 (4%)	1 (0%)	52	43

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1002	PRO

5.3.2 Protein sidechains [i](#)

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	253/275 (92%)	234 (92%)	19 (8%)	17	7
1	B	252/275 (92%)	236 (94%)	16 (6%)	22	12
All	All	505/550 (92%)	470 (93%)	35 (7%)	19	10

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	871	LEU
1	A	889	GLU
1	A	896	GLU
1	A	918	TYR
1	A	925	LEU
1	A	937	LEU
1	A	946	GLU
1	A	950	HIS
1	A	962	LYS
1	A	969	THR
1	A	985	GLU
1	A	1003	GLN
1	A	1006	GLU
1	A	1052	GLU
1	A	1073	SER
1	A	1080	GLU
1	A	1090	ARG
1	A	1121	LEU
1	A	1128	ASP
1	B	843	GLN
1	B	846	GLU
1	B	852	LEU
1	B	893	ARG

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Mol	Chain	Res	Type
1	B	918	TYR
1	B	937	LEU
1	B	943	LYS
1	B	945	LYS
1	B	967	LEU
1	B	985	GLU
1	B	1010	VAL
1	B	1016	SER
1	B	1083	LYS
1	B	1087	ARG
1	B	1097	GLU
1	B	1110	VAL

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	859	ASN
1	A	885	GLN
1	A	909	ASN
1	A	950	HIS
1	A	955	GLN
1	A	1003	GLN
1	A	1077	HIS
1	B	885	GLN
1	B	909	ASN
1	B	942	GLN
1	B	944	HIS
1	B	955	GLN
1	B	1111	ASN
1	B	1129	ASN

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	A	1007	1	13,16,17	0.63	0	19,22,24	0.84	0
1	PTR	A	1008	1	13,16,17	0.69	0	19,22,24	0.89	0
1	PTR	B	1007	1	13,16,17	0.77	0	19,22,24	0.96	1 (5%)
1	PTR	B	1008	1	13,16,17	0.68	0	19,22,24	0.85	1 (5%)

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	A	1007	1	-	0/9/11/13	0/1/1/1
1	PTR	A	1008	1	-	0/9/11/13	0/1/1/1
1	PTR	B	1007	1	-	0/9/11/13	0/1/1/1
1	PTR	B	1008	1	-	0/9/11/13	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	1007	PTR	P-OH-CZ	-2.06	117.99	123.85
1	B	1008	PTR	O3P-P-O2P	2.03	114.91	107.44

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	1007	PTR	3	0
1	A	1008	PTR	1	0
1	B	1008	PTR	1	0

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

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	7GV	A	4000	-	27,28,28	0.55	0	23,40,40	0.83	1 (4%)
2	7GV	B	4001	-	27,28,28	0.68	0	23,40,40	1.00	1 (4%)

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	7GV	A	4000	-	-	0/16/36/36	0/4/4/4
2	7GV	B	4001	-	-	0/16/36/36	0/4/4/4

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	4000	7GV	C17-N18-C10	2.51	117.55	111.64
2	B	4001	7GV	C17-N18-C10	2.68	117.96	111.64

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	4000	7GV	1	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	274/304 (90%)	-0.42	1 (0%) 93 94	6, 19, 34, 41	0
1	B	271/304 (89%)	-0.39	1 (0%) 93 94	8, 18, 37, 42	0
All	All	545/608 (89%)	-0.40	2 (0%) 93 94	6, 18, 36, 42	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1132	GLY	2.4
1	B	887	SER	2.3

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. 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
1	PTR	B	1007	16/17	0.92	0.10	-	25,29,30,31	4
1	PTR	B	1008	16/17	0.92	0.13	-	25,29,36,37	4
1	PTR	A	1008	16/17	0.90	0.14	-	26,32,36,36	5
1	PTR	A	1007	16/17	0.91	0.12	-	26,29,35,35	4

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(\AA^2)	Q<0.9
2	7GV	A	4000	25/25	0.94	0.12	0.45	18,20,22,23	0
2	7GV	B	4001	25/25	0.94	0.10	-0.19	16,17,18,18	0

6.5 Other polymers [i](#)

There are no such residues in this entry.