



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:18 PM GMT

PDB ID : 3TJD
Title : co-crystal structure of Jak2 with thienopyridine 19
Authors : Huang, X.
Deposited on : 2011-08-24
Resolution : 2.90 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

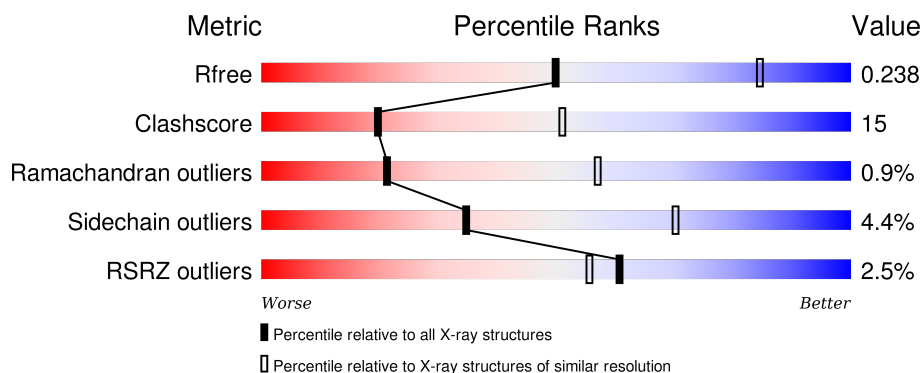
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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	298	<div> <div>2%</div> <div>63%</div> <div>29%</div> <div>• •</div> </div>
1	B	298	<div> <div>2%</div> <div>71%</div> <div>24%</div> <div>• •</div> </div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4886 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.

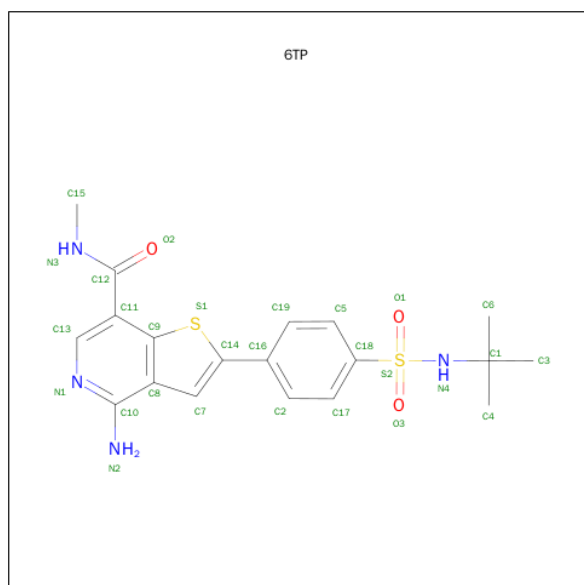
- Molecule 1 is a protein called Tyrosine-protein kinase JAK2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	286	Total	C	N	O	P	S	0	0	0
			2373	1506	408	443	2	14			
1	B	289	Total	C	N	O	P	S	0	0	0
			2395	1519	411	449	2	14			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	835	GLY	-	EXPRESSION TAG	UNP O60674
A	836	SER	-	EXPRESSION TAG	UNP O60674
B	835	GLY	-	EXPRESSION TAG	UNP O60674
B	836	SER	-	EXPRESSION TAG	UNP O60674

- Molecule 2 is 4-AMINO-2-[4-(TERT-BUTYLSULFAMOYL)PHENYL]-N-METHYLTHIEN O[3,2-C]PYRIDINE-7-CARBOXAMIDE (three-letter code: 6TP) (formula: C₁₉H₂₂N₄O₃S₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			28	19	4	3	2		
2	B	1	Total	C	N	O	S	0	0
			28	19	4	3	2		

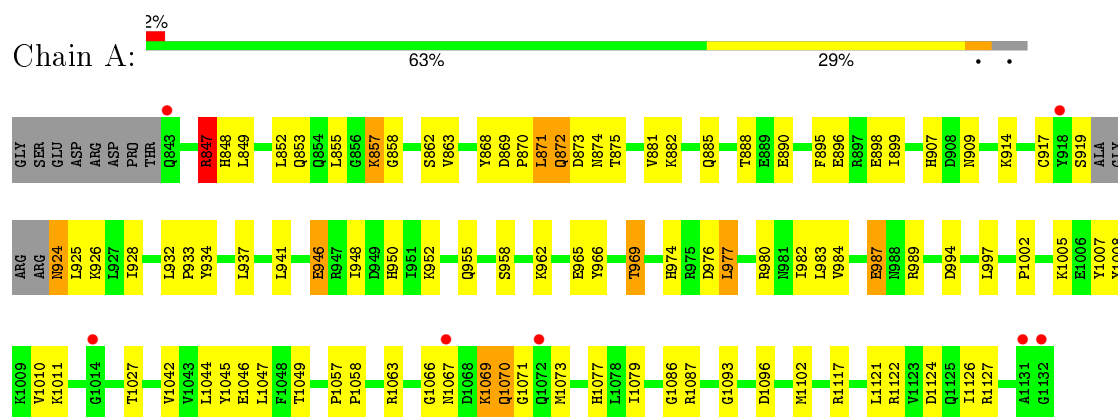
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	26	Total	O	0	0
			26	26		
3	B	36	Total	O	0	0
			36	36		

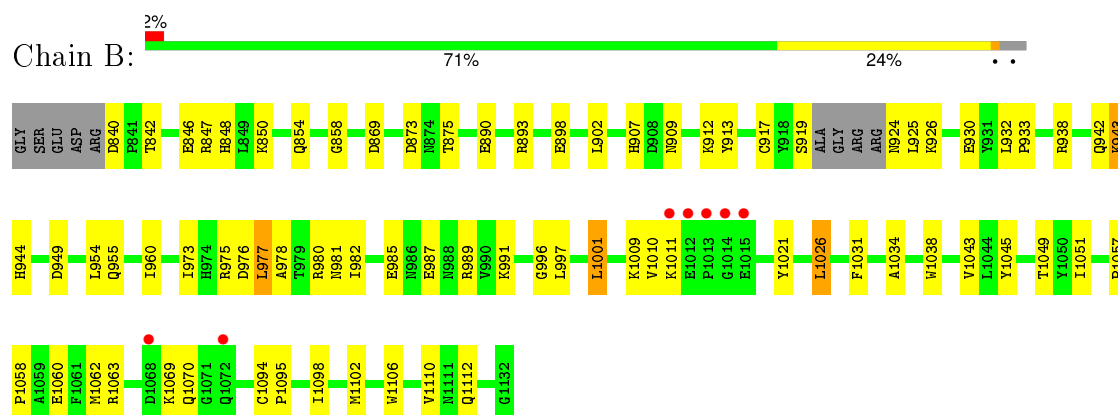
3 Residue-property plots [i](#)

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 41	Depositor
Cell constants a, b, c, α , β , γ	111.82Å 111.82Å 71.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.90 43.95 – 2.84	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-2.90) 94.9 (43.95-2.84)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 2.86Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.286 , 0.312 0.223 , 0.238	Depositor DCC
R_{free} test set	948 reflections (5.08%)	DCC
Wilson B-factor (Å ²)	52.6	Xtriage
Anisotropy	0.388	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 29.1	EDS
Estimated twinning fraction	0.028 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 19805 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4886	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.82% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: 6TP, 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	0.60	2/2389 (0.1%)	0.75	3/3211 (0.1%)
1	B	0.57	0/2412	0.71	0/3244
All	All	0.58	2/4801 (0.0%)	0.73	3/6455 (0.0%)

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	A	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	847	ARG	CZ-NH1	5.82	1.40	1.33
1	A	847	ARG	CZ-NH2	5.07	1.39	1.33

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	847	ARG	NE-CZ-NH1	-8.31	116.14	120.30
1	A	980	ARG	NE-CZ-NH2	-6.34	117.13	120.30
1	A	857	LYS	CD-CE-NZ	-5.02	100.16	111.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	847	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	2373	0	2343	81	0
1	B	2395	0	2362	63	0
2	A	28	0	22	4	0
2	B	28	0	22	2	0
3	A	26	0	0	2	0
3	B	36	0	0	7	0
All	All	4886	0	4749	143	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 15.

All (143) 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:1069:LYS:HA	1:A:1069:LYS:HE3	1.20	1.17
1:A:1069:LYS:HA	1:A:1069:LYS:CE	1.85	1.04
1:A:977:LEU:HD21	1:A:982:ILE:HD11	1.49	0.95
1:B:890:GLU:HG2	1:B:893:ARG:HH22	1.41	0.86
1:B:854:GLN:HG3	3:B:27:HOH:O	1.76	0.84
1:A:1008:PTR:HE1	1:A:1010:VAL:HB	1.59	0.83
1:A:873:ASP:O	1:A:875:THR:HG23	1.78	0.82
1:B:890:GLU:HG2	1:B:893:ARG:NH2	1.98	0.79
1:A:966:TYR:O	1:A:969:THR:HG22	1.83	0.79
1:A:977:LEU:CD2	1:A:982:ILE:HD11	2.17	0.75
1:B:976:ASP:O	1:B:981:ASN:ND2	2.21	0.74
1:A:1087:ARG:HH11	1:A:1087:ARG:HG3	1.58	0.69
1:A:1069:LYS:HE2	1:A:1070:GLN:HG3	1.74	0.69
1:A:952:LYS:HD2	1:A:955:GLN:HE21	1.58	0.68
1:A:1027:THR:HG22	1:A:1079:ILE:HD13	1.76	0.67
1:B:919:SER:HG	1:B:924:ASN:N	1.94	0.65
1:B:976:ASP:HB2	1:B:997:LEU:HD12	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:985:GLU:OE1	1:B:989:ARG:HD2	1.96	0.65
1:A:848:HIS:CG	1:A:870:PRO:HA	2.32	0.64
1:B:840:ASP:OD1	1:B:842:THR:HG22	1.97	0.64
1:A:1057:PRO:HB2	1:A:1058:PRO:HD3	1.80	0.64
1:A:1007:PTR:OH	3:A:13:HOH:O	2.15	0.63
1:B:975:ARG:HG2	1:B:1031:PHE:CD2	2.33	0.63
1:A:976:ASP:HB2	1:A:997:LEU:HD12	1.81	0.62
1:A:1069:LYS:HE2	1:A:1070:GLN:H	1.63	0.62
1:A:1069:LYS:CE	1:A:1069:LYS:CA	2.69	0.61
1:B:1057:PRO:HB2	1:B:1058:PRO:HD3	1.81	0.61
1:A:946:GLU:H	1:A:946:GLU:CD	2.00	0.61
1:B:942:GLN:HG2	1:B:1051:ILE:HB	1.82	0.61
1:B:917:CYS:SG	1:B:926:LYS:HE3	2.41	0.60
1:A:1069:LYS:HD3	1:A:1077:HIS:CD2	2.38	0.59
1:A:1044:LEU:HD23	1:A:1102:MET:HG3	1.84	0.59
1:A:857:LYS:HZ1	1:A:885:GLN:HG2	1.68	0.59
1:B:1045:TYR:O	1:B:1049:THR:HG23	2.04	0.58
1:A:1086:GLY:O	1:A:1087:ARG:HG3	2.03	0.58
1:A:952:LYS:HD2	1:A:955:GLN:NE2	2.18	0.58
1:A:977:LEU:HD21	1:A:982:ILE:CD1	2.28	0.57
1:B:917:CYS:HB3	1:B:926:LYS:HE3	1.86	0.56
1:B:873:ASP:OD1	1:B:875:THR:HG23	2.05	0.56
1:A:1087:ARG:HG3	1:A:1087:ARG:NH1	2.19	0.56
1:A:917:CYS:HB3	1:A:926:LYS:HB2	1.86	0.56
1:B:1038:TRP:CE3	1:B:1106:TRP:HA	2.40	0.56
1:A:994:ASP:HB2	2:A:1:6TP:H5	1.87	0.55
1:A:1045:TYR:O	1:A:1049:THR:HG23	2.07	0.55
1:B:917:CYS:HB3	1:B:926:LYS:HB2	1.89	0.54
1:B:846:GLU:CD	1:B:926:LYS:NZ	2.60	0.54
1:B:981:ASN:OD1	2:B:1:6TP:O3	2.27	0.53
1:B:858:GLY:HA3	2:B:1:6TP:H3B	1.89	0.53
1:A:919:SER:HB3	1:A:924:ASN:ND2	2.24	0.53
1:A:1093:GLY:HA3	1:B:1112:GLN:O	2.10	0.52
1:A:932:LEU:HD12	1:A:983:LEU:CB	2.40	0.52
1:B:919:SER:HA	3:B:119:HOH:O	2.10	0.52
1:B:960:ILE:HD13	1:B:982:ILE:HD13	1.90	0.52
1:B:1095:PRO:HB2	1:B:1098:ILE:HG12	1.92	0.52
1:A:952:LYS:HE3	1:A:955:GLN:HE22	1.75	0.52
1:B:943:LYS:HD3	1:B:944:HIS:CD2	2.45	0.52
1:A:919:SER:HB3	1:A:924:ASN:HD21	1.75	0.52
1:A:870:PRO:HG2	1:A:871:LEU:HD13	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:932:LEU:HD12	1:A:983:LEU:HB2	1.92	0.52
1:A:1070:GLN:OE1	1:A:1073:MET:HG3	2.10	0.51
1:A:857:LYS:HZ1	1:A:885:GLN:CG	2.23	0.51
1:B:1034:ALA:CB	1:B:1110:VAL:HG13	2.41	0.51
1:B:938:ARG:HD2	3:B:48:HOH:O	2.10	0.50
1:B:976:ASP:OD1	1:B:980:ARG:NH2	2.45	0.50
1:A:974:HIS:HE1	1:A:994:ASP:O	1.95	0.50
1:B:975:ARG:NH2	3:B:18:HOH:O	2.45	0.50
1:A:849:LEU:HG	1:A:868:TYR:HD1	1.77	0.49
1:A:855:LEU:HD22	2:A:1:6TP:O2	2.12	0.49
1:B:1060:GLU:OE1	1:B:1063:ARG:NH2	2.46	0.48
1:A:862:SER:O	1:A:882:LYS:HA	2.13	0.48
1:B:907:HIS:HE1	1:B:909:ASN:HD22	1.61	0.48
1:A:863:VAL:HG23	2:A:1:6TP:C6	2.44	0.48
1:A:1063:ARG:HG2	1:A:1067:ASN:HD21	1.79	0.48
1:A:1069:LYS:HE2	1:A:1070:GLN:N	2.29	0.48
1:B:973:ILE:CG1	1:B:1001:LEU:HD13	2.44	0.48
1:B:1102:MET:HE3	1:B:1106:TRP:HH2	1.79	0.47
1:B:912:LYS:HB2	1:B:930:GLU:HG2	1.96	0.47
1:B:1021:TYR:HB3	1:B:1026:LEU:HD13	1.97	0.47
1:A:873:ASP:O	1:A:874:ASN:C	2.53	0.47
1:A:937:LEU:HD21	1:A:1047:LEU:HD21	1.97	0.47
1:A:974:HIS:HD2	1:A:977:LEU:N	2.13	0.47
1:B:846:GLU:CD	1:B:926:LYS:HZ3	2.18	0.46
1:A:1042:VAL:O	1:A:1046:GLU:HG3	2.16	0.46
1:B:960:ILE:CD1	1:B:982:ILE:HD13	2.45	0.46
1:A:977:LEU:HA	1:A:977:LEU:HD23	1.47	0.46
1:B:930:GLU:OE1	1:B:991:LYS:NZ	2.40	0.46
1:B:955:GLN:NE2	3:B:265:HOH:O	2.49	0.46
1:A:907:HIS:HE1	1:A:909:ASN:HD22	1.63	0.46
1:A:857:LYS:HZ1	1:A:885:GLN:CD	2.19	0.46
1:B:977:LEU:HD22	1:B:978:ALA:H	1.80	0.45
1:B:980:ARG:HG2	1:B:981:ASN:OD1	2.16	0.45
1:B:917:CYS:CB	1:B:926:LYS:HE3	2.46	0.45
1:B:847:ARG:HG3	1:B:848:HIS:CD2	2.51	0.45
1:A:965:GLU:OE1	1:A:965:GLU:HA	2.16	0.45
1:A:974:HIS:CE1	1:A:994:ASP:O	2.69	0.45
1:A:1066:GLY:O	1:A:1069:LYS:HB2	2.17	0.45
1:B:840:ASP:CG	1:B:842:THR:HG22	2.35	0.45
1:B:1045:TYR:HB2	1:B:1102:MET:HE3	1.99	0.45
1:A:869:ASP:OD2	1:A:872:GLN:HA	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:958:SER:O	1:A:962:LYS:HG3	2.17	0.44
1:B:1058:PRO:O	1:B:1062:MET:HG2	2.18	0.44
1:B:850:LYS:HE2	1:B:869:ASP:HB3	1.99	0.44
1:B:1102:MET:HE3	1:B:1106:TRP:CH2	2.52	0.44
1:A:950:HIS:HE1	1:A:1093:GLY:O	2.00	0.44
1:A:1122:ARG:O	1:A:1126:ILE:HG13	2.17	0.44
1:B:1009:LYS:NZ	3:B:192:HOH:O	2.50	0.44
1:A:1045:TYR:HB2	1:A:1102:MET:HE3	1.99	0.44
1:A:857:LYS:HG3	1:A:858:GLY:N	2.33	0.44
1:B:977:LEU:HD23	1:B:977:LEU:HA	1.46	0.44
1:B:902:LEU:HG	1:B:913:TYR:HB2	1.99	0.44
1:A:941:LEU:HD12	1:A:948:ILE:CD1	2.48	0.44
1:B:976:ASP:CB	1:B:997:LEU:HD12	2.47	0.43
1:A:1007:PTR:O2P	1:A:1007:PTR:HE1	2.18	0.43
1:A:896:GLU:CG	3:A:28:HOH:O	2.65	0.43
1:B:917:CYS:SG	1:B:926:LYS:CE	3.06	0.43
1:B:846:GLU:OE1	1:B:926:LYS:NZ	2.50	0.43
1:A:888:THR:HG22	1:A:890:GLU:H	1.84	0.43
1:A:871:LEU:CD1	1:A:871:LEU:N	2.81	0.43
1:A:1124:ASP:HA	1:A:1127:ARG:HD3	2.01	0.43
1:B:1010:VAL:HG23	1:B:1010:VAL:O	2.19	0.42
1:B:917:CYS:SG	1:B:926:LYS:NZ	2.92	0.42
1:A:882:LYS:NZ	1:A:898:GLU:OE2	2.40	0.42
1:B:932:LEU:HA	1:B:933:PRO:HD2	1.77	0.42
1:A:934:TYR:HB2	1:A:984:VAL:HB	2.02	0.42
1:B:989:ARG:NH2	3:B:71:HOH:O	2.50	0.42
1:A:932:LEU:HA	1:A:933:PRO:HD2	1.90	0.42
1:A:941:LEU:HD12	1:A:948:ILE:HD12	2.02	0.42
1:A:914:LYS:HB2	1:A:928:ILE:O	2.20	0.42
1:B:898:GLU:OE1	1:B:996:GLY:HA2	2.19	0.42
1:A:1002:PRO:HG2	1:A:1005:LYS:HB2	2.02	0.42
1:B:978:ALA:HA	1:B:1043:VAL:HG22	2.01	0.42
1:A:974:HIS:HD2	1:A:976:ASP:H	1.68	0.41
1:A:895:PHE:O	1:A:899:ILE:HG13	2.21	0.41
1:A:881:VAL:HG22	1:A:928:ILE:HD12	2.03	0.41
1:A:873:ASP:OD1	1:A:875:THR:OG1	2.26	0.41
1:A:847:ARG:H	1:A:847:ARG:HG2	1.61	0.41
1:A:872:GLN:HE21	1:A:872:GLN:HB3	1.60	0.41
1:A:852:LEU:O	1:A:853:GLN:HB3	2.20	0.41
1:B:973:ILE:HG13	1:B:1001:LEU:HD13	2.02	0.40
1:B:954:LEU:HD23	1:B:954:LEU:HA	1.85	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:849:LEU:HG	1:A:868:TYR:CD1	2.57	0.40
1:B:1094:CYS:HA	1:B:1095:PRO:HD3	1.95	0.40
1:A:994:ASP:CB	2:A:1:6TP:H5	2.51	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	280/298 (94%)	263 (94%)	14 (5%)	3 (1%)	17	51
1	B	283/298 (95%)	270 (95%)	11 (4%)	2 (1%)	26	63
All	All	563/596 (94%)	533 (95%)	25 (4%)	5 (1%)	21	57

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1070	GLN
1	B	987	GLU
1	B	1070	GLN
1	A	987	GLU
1	A	1071	GLY

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	260/269 (97%)	245 (94%)	15 (6%)	25	58
1	B	263/269 (98%)	255 (97%)	8 (3%)	48	83
All	All	523/538 (97%)	500 (96%)	23 (4%)	35	70

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	847	ARG
1	A	871	LEU
1	A	872	GLN
1	A	924	ASN
1	A	925	LEU
1	A	946	GLU
1	A	969	THR
1	A	977	LEU
1	A	987	GLU
1	A	989	ARG
1	A	1011	LYS
1	A	1069	LYS
1	A	1096	ASP
1	A	1117	ARG
1	A	1121	LEU
1	B	925	LEU
1	B	943	LYS
1	B	949	ASP
1	B	977	LEU
1	B	1001	LEU
1	B	1011	LYS
1	B	1026	LEU
1	B	1069	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	848	HIS
1	A	872	GLN
1	A	874	ASN
1	A	909	ASN
1	A	924	ASN
1	A	950	HIS
1	A	955	GLN

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Mol	Chain	Res	Type
1	A	974	HIS
1	A	1067	ASN
1	A	1077	HIS
1	B	843	GLN
1	B	909	ASN
1	B	924	ASN
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	14,16,17	0.55	0	18,22,24	0.77	0
1	PTR	A	1008	1	14,16,17	0.67	0	18,22,24	0.96	2 (11%)
1	PTR	B	1007	1	14,16,17	0.48	0	18,22,24	0.84	0
1	PTR	B	1008	1	14,16,17	0.77	0	18,22,24	0.87	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

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
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 (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1008	PTR	O-C-CA	-2.31	119.48	125.49
1	B	1008	PTR	O-C-CA	-2.24	119.66	125.49
1	A	1008	PTR	O3P-P-O1P	2.07	117.24	110.58

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	1007	PTR	2	0
1	A	1008	PTR	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates 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	6TP	A	1	-	24,30,30	2.90	9 (37%)	32,46,46	2.64	13 (40%)
2	6TP	B	1	-	24,30,30	3.24	12 (50%)	32,46,46	2.51	10 (31%)

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	6TP	A	1	-	-	0/22/22/22	0/3/3/3
2	6TP	B	1	-	-	0/22/22/22	0/3/3/3

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	6TP	C18-S2	-4.16	1.70	1.76
2	B	1	6TP	O3-S2	-3.14	1.40	1.43
2	B	1	6TP	C4-C1	-2.25	1.46	1.52
2	B	1	6TP	C18-S2	-2.24	1.73	1.76
2	A	1	6TP	C13-C11	2.23	1.43	1.39
2	A	1	6TP	C17-C2	2.50	1.43	1.38
2	B	1	6TP	C10-C8	2.53	1.47	1.45
2	A	1	6TP	C17-C18	2.64	1.43	1.38
2	B	1	6TP	C17-C2	2.75	1.43	1.38
2	A	1	6TP	C5-C18	2.76	1.43	1.38
2	A	1	6TP	C10-C8	2.85	1.47	1.45
2	B	1	6TP	C13-C11	3.18	1.44	1.39
2	B	1	6TP	C19-C16	3.18	1.46	1.39
2	A	1	6TP	C19-C16	3.59	1.47	1.39
2	B	1	6TP	C5-C18	3.69	1.44	1.38
2	B	1	6TP	C12-N3	3.76	1.37	1.33
2	B	1	6TP	C17-C18	3.79	1.45	1.38
2	B	1	6TP	C2-C16	4.52	1.49	1.39
2	A	1	6TP	C2-C16	4.62	1.49	1.39
2	A	1	6TP	S2-N4	10.19	1.76	1.61
2	B	1	6TP	S2-N4	11.00	1.77	1.61

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	6TP	C5-C18-S2	-6.14	112.68	119.78
2	B	1	6TP	C5-C18-S2	-6.03	112.81	119.78
2	B	1	6TP	C11-C13-N1	-5.82	118.23	125.06
2	A	1	6TP	C11-C13-N1	-5.26	118.89	125.06
2	B	1	6TP	C2-C17-C18	-5.11	113.83	119.48
2	A	1	6TP	C2-C17-C18	-4.04	115.01	119.48
2	B	1	6TP	O1-S2-N4	-3.99	98.43	107.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	6TP	C15-N3-C12	-3.96	117.08	121.80
2	A	1	6TP	O3-S2-O1	-3.57	114.81	119.54
2	A	1	6TP	C8-C10-N2	-3.09	117.55	121.11
2	B	1	6TP	C11-C12-N3	-3.01	112.99	117.50
2	A	1	6TP	C19-C5-C18	-2.89	116.28	119.48
2	B	1	6TP	C19-C5-C18	-2.63	116.57	119.48
2	A	1	6TP	C18-S2-N4	-2.60	104.47	108.11
2	B	1	6TP	C8-C10-N2	-2.53	118.19	121.11
2	A	1	6TP	C13-N1-C10	2.32	121.58	118.88
2	A	1	6TP	C3-C1-N4	2.41	115.54	109.00
2	B	1	6TP	C13-N1-C10	2.61	121.91	118.88
2	A	1	6TP	O1-S2-C18	2.80	111.50	107.96
2	B	1	6TP	O3-S2-O1	2.96	123.47	119.54
2	A	1	6TP	O2-C12-C11	3.02	126.04	121.59
2	B	1	6TP	C17-C18-C5	5.32	127.53	120.42
2	A	1	6TP	C17-C18-C5	5.80	128.17	120.42

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
2	A	1	6TP	4	0
2	B	1	6TP	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 [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	284/298 (95%)	-0.24	7 (2%) 61 55	13, 31, 64, 83	0
1	B	287/298 (96%)	-0.25	7 (2%) 62 57	12, 30, 62, 84	0
All	All	571/596 (95%)	-0.24	14 (2%) 61 55	12, 30, 64, 84	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1132	GLY	5.5
1	B	1013	PRO	4.8
1	B	1012	GLU	3.9
1	A	918	TYR	3.4
1	B	1014	GLY	3.3
1	B	1015	GLU	3.1
1	A	1067	ASN	3.1
1	A	1131	ALA	2.7
1	A	843	GLN	2.6
1	A	1014	GLY	2.5
1	B	1068	ASP	2.4
1	B	1011	LYS	2.4
1	A	1072	GLN	2.4
1	B	1072	GLN	2.0

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(\AA^2)	Q<0.9
1	PTR	B	1008	16/17	0.86	0.20	-	41,54,71,72	0
1	PTR	A	1008	16/17	0.90	0.17	-	34,38,51,52	0
1	PTR	A	1007	16/17	0.91	0.16	-	37,48,67,69	0
1	PTR	B	1007	16/17	0.92	0.15	-	42,47,52,52	0

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	6TP	B	1	28/28	0.95	0.18	-0.15	13,20,25,29	0
2	6TP	A	1	28/28	0.97	0.14	-0.31	7,16,26,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.