Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 04:02 PM GMT

PDB ID	:	4E6Q
Title	:	JAK2 kinase (JH1 domain) triple mutant in complex with compound 12
Authors	:	Murray, J.M.
Deposited on	:	2012-03-15
Resolution	:	1.95 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

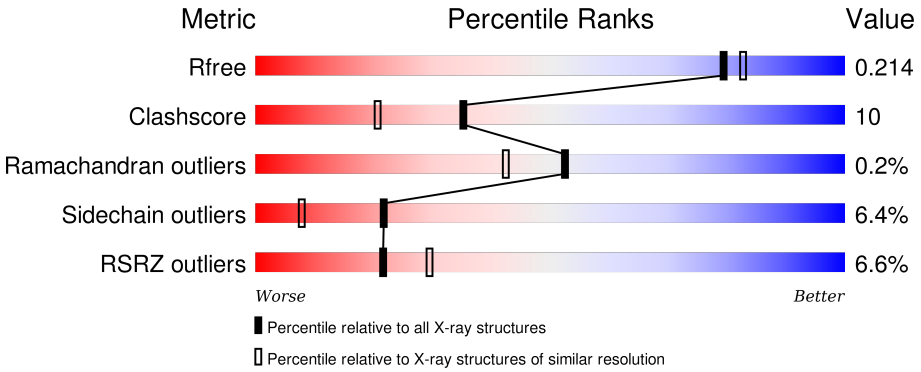
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	2910 (1.96-1.92)
Clashscore	102246	3095 (1.96-1.92)
Ramachandran outliers	100387	3062 (1.96-1.92)
Sidechain outliers	100360	3062 (1.96-1.92)
RSRZ outliers	91569	2915 (1.96-1.92)

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 <=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	6% 73% 24%
1	B	298	7% 76% 20%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5148 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.

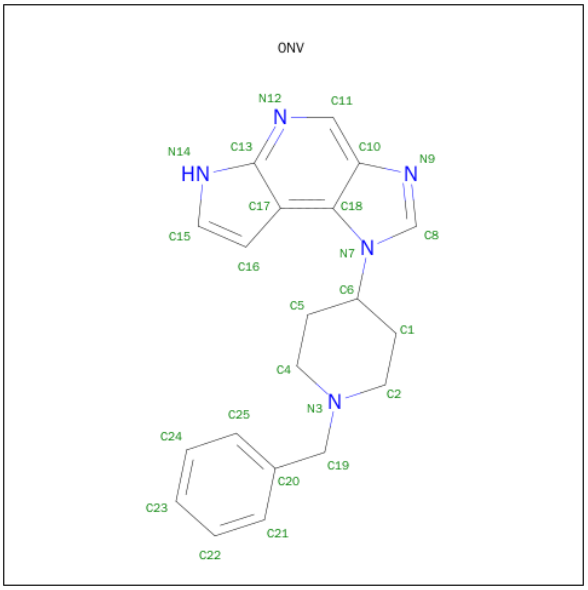
- Molecule 1 is a protein called Tyrosine-protein kinase JAK2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	298	Total	C	N	O	P	S	20	0	0
			2459	1558	426	459	2	14			
1	B	298	Total	C	N	O	P	S	0	0	0
			2456	1556	424	460	2	14			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	853	ARG	GLN	ENGINEERED MUTATION	UNP O60674
A	931	PHE	TYR	ENGINEERED MUTATION	UNP O60674
A	939	GLU	ASP	ENGINEERED MUTATION	UNP O60674
B	853	ARG	GLN	ENGINEERED MUTATION	UNP O60674
B	931	PHE	TYR	ENGINEERED MUTATION	UNP O60674
B	939	GLU	ASP	ENGINEERED MUTATION	UNP O60674

- Molecule 2 is 1-(1-BENZYLPIPERIDIN-4-YL)-1,6-DIHYDROIMIDAZO[4,5-D]PYRROLO [2,3-B]PYRIDINE (three-letter code: 0NV) (formula: $C_{20}H_{21}N_5$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C N 25 20 5	0	0
2	B	1	Total C N 25 20 5	0	0

- Molecule 3 is water.

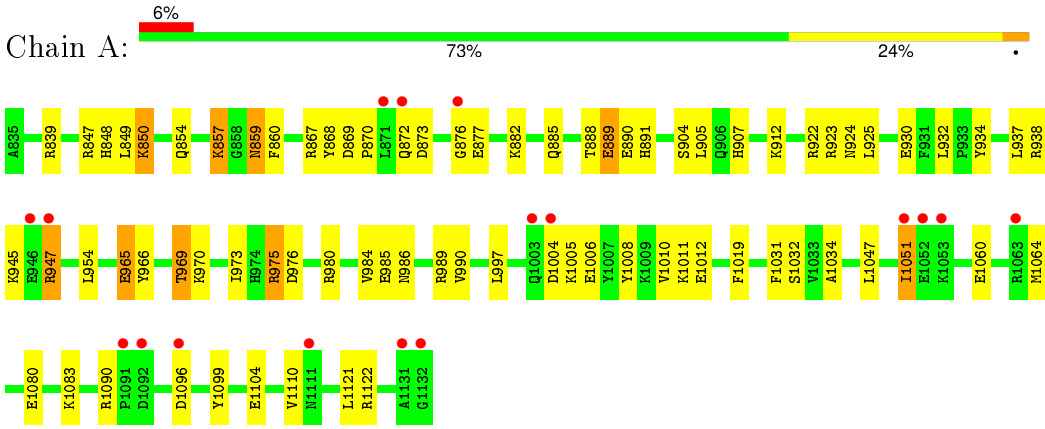
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	87	Total O 87 87	0	0
3	B	96	Total O 96 96	0	0

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.

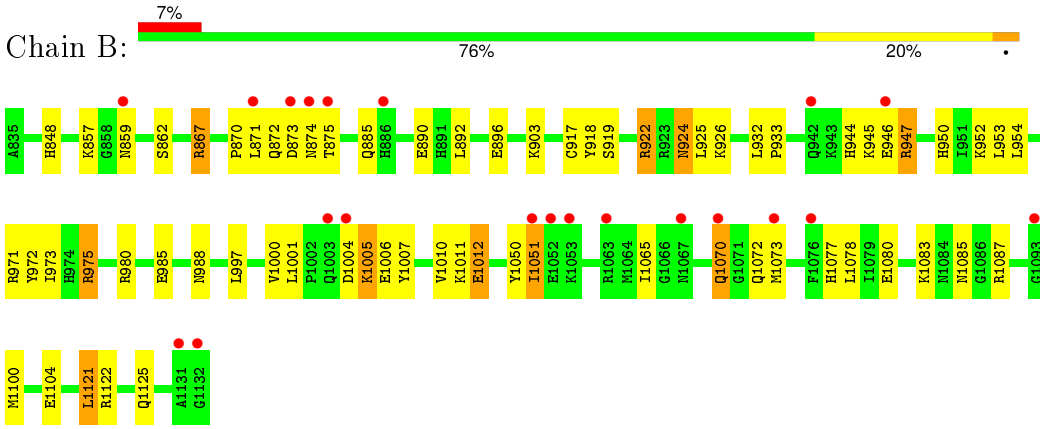
- Molecule 1: Tyrosine-protein kinase JAK2

Chain A:



- Molecule 1: Tyrosine-protein kinase JAK2

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	50.50Å 103.85Å 70.00Å 90.00° 90.02° 90.00°	Depositor
Resolution (Å)	50.50 – 1.95 50.50 – 1.95	Depositor EDS
% Data completeness (in resolution range)	98.4 (50.50-1.95) 98.4 (50.50-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	2.30 (at 1.95Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_991)	Depositor
R, R_{free}	0.186 , 0.234 0.189 , 0.214	Depositor DCC
R_{free} test set	2608 reflections (5.29%)	DCC
Wilson B-factor (Å^2)	18.8	Xtriage
Anisotropy	0.842	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.40 , 64.3	EDS
Estimated twinning fraction	0.220 for h,-k,-l 0.256 for h,-k,-l	Xtriage
Reported twinning fraction	0.220 for h,-k,-l	Depositor
L-test for twinning[2]	$< \|L\| > = 0.45$, $< L^2 > = 0.28$	Xtriage
Outliers	0 of 51908 reflections	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	5148	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.61% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: PTR, 0NV

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	0.91	2/2478 (0.1%)	0.93	3/3333 (0.1%)
1	B	0.92	1/2475 (0.0%)	0.95	5/3330 (0.2%)
All	All	0.91	3/4953 (0.1%)	0.94	8/6663 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	965	GLU	CG-CD	5.96	1.60	1.51
1	A	1019	PHE	CB-CG	5.20	1.60	1.51
1	B	917	CYS	CB-SG	5.03	1.90	1.82

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	975	ARG	NE-CZ-NH2	-7.56	116.52	120.30
1	B	975	ARG	NE-CZ-NH1	7.12	123.86	120.30
1	A	975	ARG	NE-CZ-NH2	-6.75	116.92	120.30
1	A	975	ARG	NE-CZ-NH1	6.56	123.58	120.30
1	B	1087	ARG	NE-CZ-NH1	5.98	123.29	120.30
1	A	976	ASP	CB-CG-OD1	5.45	123.20	118.30
1	B	922	ARG	NE-CZ-NH1	-5.38	117.61	120.30
1	B	867	ARG	NE-CZ-NH1	5.00	122.80	120.30

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	2459	0	2415	47	3
1	B	2456	0	2408	51	3
2	A	25	0	21	1	0
2	B	25	0	21	0	0
3	A	87	0	0	0	0
3	B	96	0	0	5	0
All	All	5148	0	4865	96	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 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:1073:MET:CE	1:B:1077:HIS:HE1	1.87	0.87
1:B:1073:MET:HE3	1:B:1077:HIS:HE1	1.51	0.75
1:B:867:ARG:NH2	1:B:874:ASN:HA	2.03	0.73
1:B:1073:MET:HE1	1:B:1077:HIS:HE1	1.57	0.69
1:B:1073:MET:CE	1:B:1077:HIS:CE1	2.74	0.68
1:A:937:LEU:HD21	1:A:1047:LEU:HD21	1.75	0.68
1:B:975:ARG:HD3	1:B:997:LEU:O	1.92	0.68
1:A:867:ARG:NH1	1:B:933:PRO:O	2.28	0.67
1:A:975:ARG:HD3	1:A:997:LEU:O	1.97	0.64
1:A:1005:LYS:HD3	1:A:1006:GLU:H	1.62	0.64
1:A:868:TYR:O	1:A:876:GLY:HA3	1.98	0.63
1:B:952:LYS:NZ	1:B:988:ASN:OD1	2.30	0.63
1:A:984:VAL:HG22	1:A:990:VAL:HG12	1.81	0.62
1:B:975:ARG:HH22	1:B:1012:GLU:CD	2.01	0.62
1:B:1073:MET:HE3	1:B:1077:HIS:CE1	2.33	0.62
1:B:947:ARG:O	1:B:947:ARG:HG2	2.01	0.61
1:B:975:ARG:NH2	1:B:1012:GLU:OE1	2.33	0.60
1:B:1080:GLU:OE2	1:B:1083:LYS:HE2	2.01	0.60
1:B:953:LEU:HD11	1:B:1050:TYR:CZ	2.37	0.60
1:A:985:GLU:OE1	1:A:989:ARG:HD2	2.03	0.59
1:B:1104:GLU:CD	1:B:1122:ARG:HH22	2.07	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:972:TYR:CD1	1:B:1000:VAL:HG22	2.39	0.58
1:B:1073:MET:HE1	1:B:1077:HIS:CE1	2.38	0.57
1:B:1010:VAL:HG12	1:B:1012:GLU:HG2	1.87	0.57
1:A:1104:GLU:OE1	1:A:1122:ARG:NH2	2.38	0.57
1:B:1104:GLU:OE1	1:B:1122:ARG:NH2	2.38	0.55
1:B:1122:ARG:HD2	3:B:1370:HOH:O	2.07	0.55
1:B:971:ARG:NH2	1:B:1005:LYS:O	2.40	0.54
1:B:972:TYR:CE1	1:B:1000:VAL:HG22	2.42	0.54
1:A:947:ARG:HG2	1:A:947:ARG:O	2.08	0.53
1:B:1121:LEU:O	1:B:1125:GLN:HG3	2.09	0.52
1:A:1005:LYS:CD	1:A:1006:GLU:H	2.23	0.52
1:A:1010:VAL:HG12	1:A:1012:GLU:HG2	1.90	0.52
1:A:1104:GLU:CD	1:A:1122:ARG:HH22	2.12	0.52
1:A:888:THR:HB	1:A:891:HIS:H	1.75	0.51
1:B:919:SER:OG	1:B:926:LYS:HE3	2.10	0.51
1:B:973:ILE:HD11	1:B:1001:LEU:HD11	1.92	0.51
1:A:854:GLN:HE22	1:A:857:LYS:HE2	1.75	0.51
1:A:934:TYR:OH	1:B:988:ASN:ND2	2.44	0.50
1:B:872:GLN:NE2	3:B:1358:HOH:O	2.45	0.50
1:B:950:HIS:O	1:B:954:LEU:HG	2.11	0.50
1:A:941:LEU:O	1:A:945:LYS:HB3	2.12	0.49
1:A:1080:GLU:OE2	1:A:1083:LYS:HE2	2.12	0.49
1:B:1051:ILE:O	1:B:1051:ILE:HG22	2.13	0.49
1:A:1034:ALA:CB	1:A:1110:VAL:HG13	2.44	0.48
1:B:922:ARG:O	1:B:922:ARG:HG3	2.14	0.48
1:B:1005:LYS:HD3	1:B:1006:GLU:H	1.78	0.48
1:B:932:LEU:HD13	1:B:985:GLU:HA	1.96	0.47
1:A:849:LEU:C	1:A:850:LYS:HD2	2.35	0.47
1:A:888:THR:HG1	1:A:891:HIS:CE1	2.33	0.47
1:A:888:THR:HG22	1:A:890:GLU:H	1.80	0.47
1:B:918:TYR:HA	1:B:924:ASN:O	2.16	0.46
1:A:907:HIS:HB2	1:A:966:TYR:CD2	2.50	0.46
1:B:1070:GLN:HG3	1:B:1070:GLN:H	1.60	0.46
1:B:1005:LYS:HE3	1:B:1007:PTR:CE1	2.46	0.46
1:A:1090:ARG:HG3	1:A:1099:TYR:HB2	1.97	0.46
1:A:966:TYR:O	1:A:969:THR:HG22	2.16	0.46
1:A:973:ILE:HD12	1:A:1032:SER:HA	1.98	0.45
1:B:848:HIS:CD2	1:B:870:PRO:HA	2.51	0.45
1:A:847:ARG:HG3	1:A:848:HIS:CD2	2.51	0.45
1:A:904:SER:O	1:A:970:LYS:HE3	2.16	0.45
1:A:869:ASP:HB3	1:A:872:GLN:HA	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:903:LYS:NZ	3:B:1321:HOH:O	2.39	0.44
1:A:912:LYS:HB2	1:A:930:GLU:CD	2.37	0.44
1:A:954:LEU:HD23	1:A:954:LEU:HA	1.73	0.44
1:A:934:TYR:CE2	1:A:986:ASN:HA	2.52	0.44
1:B:947:ARG:HD3	3:B:1338:HOH:O	2.17	0.44
1:A:1008:PTR:N	1:A:1031:PHE:O	2.50	0.44
1:A:934:TYR:CD2	1:A:986:ASN:HA	2.52	0.44
1:A:839:ARG:HA	1:A:839:ARG:HD2	1.83	0.44
1:B:857:LYS:HE3	1:B:885:GLN:NE2	2.32	0.44
1:B:932:LEU:HA	1:B:933:PRO:HD3	1.87	0.44
1:B:862:SER:HB2	1:B:885:GLN:OE1	2.18	0.44
1:A:938:ARG:HG3	1:A:1051:ILE:HG21	2.00	0.44
1:B:1065:ILE:HD11	1:B:1078:LEU:HB2	2.00	0.43
1:A:860:PHE:HA	1:A:885:GLN:O	2.18	0.43
1:B:1085:ASN:HB3	3:B:1366:HOH:O	2.18	0.43
1:B:1097:GLU:O	1:B:1100:MET:HB3	2.17	0.43
1:B:848:HIS:CG	1:B:870:PRO:HA	2.54	0.43
1:A:868:TYR:CE2	1:A:870:PRO:HG3	2.54	0.43
1:B:971:ARG:HB3	1:B:1001:LEU:HB2	2.02	0.42
1:B:871:LEU:HB3	1:B:873:ASP:OD1	2.20	0.42
1:B:944:HIS:O	1:B:946:GLU:N	2.52	0.42
1:B:892:LEU:O	1:B:896:GLU:HG3	2.20	0.42
1:A:1011:LYS:HE3	1:A:1011:LYS:HB2	1.71	0.41
1:A:923:ARG:HD2	1:A:923:ARG:HH11	1.71	0.41
1:A:905:LEU:HD23	1:A:970:LYS:HD2	2.03	0.41
1:B:871:LEU:HD12	1:B:875:THR:OG1	2.21	0.41
1:A:877:GLU:HG2	1:A:877:GLU:H	1.39	0.41
1:B:1005:LYS:CD	1:B:1006:GLU:H	2.33	0.41
1:A:854:GLN:OE1	1:A:857:LYS:HG3	2.20	0.41
1:A:932:LEU:HD22	1:A:985:GLU:HG3	2.02	0.41
1:A:882:LYS:HE3	2:A:1201:0NV:H19	2.03	0.41
1:A:937:LEU:CD2	1:A:1047:LEU:HD21	2.48	0.40
1:A:859:ASN:HB2	1:A:860:PHE:CD2	2.56	0.40
1:A:1060:GLU:O	1:A:1064:MET:HG3	2.21	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:889:GLU:OE2	1:B:1073:MET:CE[2_546]	1.59	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:839:ARG:NH2	1:B:890:GLU:OE1[2_546]	2.12	0.08
1:A:889:GLU:OE2	1:B:1073:MET:CG[2_546]	2.13	0.07

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	294/298 (99%)	288 (98%)	6 (2%)	0	100 100
1	B	294/298 (99%)	282 (96%)	11 (4%)	1 (0%)	46 35
All	All	588/596 (99%)	570 (97%)	17 (3%)	1 (0%)	52 42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1051	ILE

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	266/269 (99%)	248 (93%)	18 (7%)	20 6
1	B	266/269 (99%)	250 (94%)	16 (6%)	24 9
All	All	532/538 (99%)	498 (94%)	34 (6%)	22 8

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	850	LYS
1	A	857	LYS
1	A	859	ASN
1	A	873	ASP
1	A	889	GLU
1	A	922	ARG
1	A	924	ASN
1	A	925	LEU
1	A	947	ARG
1	A	965	GLU
1	A	969	THR
1	A	980	ARG
1	A	1004	ASP
1	A	1051	ILE
1	A	1070	GLN
1	A	1072	GLN
1	A	1096	ASP
1	A	1121	LEU
1	B	859	ASN
1	B	924	ASN
1	B	925	LEU
1	B	945	LYS
1	B	947	ARG
1	B	958	SER
1	B	969	THR
1	B	980	ARG
1	B	1004	ASP
1	B	1005	LYS
1	B	1011	LYS
1	B	1012	GLU
1	B	1070	GLN
1	B	1072	GLN
1	B	1096	ASP
1	B	1121	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	924	ASN
1	B	924	ASN
1	B	955	GLN
1	B	1077	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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	1.23	1 (7%)	18,22,24	1.01	2 (11%)
1	PTR	A	1008	1	14,16,17	1.15	1 (7%)	18,22,24	1.13	2 (11%)
1	PTR	B	1007	1	14,16,17	1.14	1 (7%)	18,22,24	1.12	1 (5%)
1	PTR	B	1008	1	14,16,17	1.19	1 (7%)	18,22,24	1.00	2 (11%)

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

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1007	PTR	OH-CZ	-4.18	1.30	1.40
1	B	1008	PTR	OH-CZ	-3.88	1.31	1.40
1	B	1007	PTR	OH-CZ	-3.73	1.31	1.40
1	A	1008	PTR	OH-CZ	-3.50	1.32	1.40

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	1008	PTR	O-C-CA	-2.64	118.62	125.49
1	B	1008	PTR	O-C-CA	-2.52	118.93	125.49
1	A	1007	PTR	O-C-CA	-2.01	120.25	125.49
1	B	1007	PTR	O3P-P-OH	2.51	114.23	105.22
1	A	1007	PTR	P-OH-CZ	2.58	131.19	123.76
1	B	1008	PTR	O3P-P-OH	3.00	115.97	105.22
1	A	1008	PTR	O3P-P-OH	3.06	116.18	105.22

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	1008	PTR	1	0
1	B	1007	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	0NV	A	1201	-	22,29,29	3.38	7 (31%)	27,41,41	2.32	9 (33%)
2	0NV	B	1201	-	22,29,29	3.15	7 (31%)	27,41,41	2.58	10 (37%)

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	0NV	A	1201	-	-	0/8/18/18	0/5/5/5
2	0NV	B	1201	-	-	0/8/18/18	0/5/5/5

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1201	0NV	C11-N12	-9.91	1.31	1.47
2	A	1201	0NV	C17-C18	-8.67	1.41	1.54
2	B	1201	0NV	C11-N12	-8.60	1.33	1.47
2	B	1201	0NV	C17-C18	-8.47	1.42	1.54
2	B	1201	0NV	C18-N7	-5.41	1.38	1.49
2	A	1201	0NV	C18-N7	-5.17	1.38	1.49
2	A	1201	0NV	C10-N9	-4.48	1.36	1.48
2	B	1201	0NV	C10-N9	-4.02	1.38	1.48
2	A	1201	0NV	C8-N7	-3.12	1.37	1.46
2	B	1201	0NV	C15-N14	-3.03	1.38	1.48
2	A	1201	0NV	C15-N14	-2.87	1.38	1.48
2	B	1201	0NV	C8-N7	-2.62	1.39	1.46
2	B	1201	0NV	C16-C17	-2.19	1.47	1.54
2	A	1201	0NV	C16-C17	-2.15	1.48	1.54

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1201	0NV	C4-C5-C6	-4.55	103.89	110.80
2	B	1201	0NV	C2-C1-C6	-4.51	103.95	110.80
2	A	1201	0NV	C2-C1-C6	-3.47	105.52	110.80
2	B	1201	0NV	C24-C25-C20	-2.44	116.77	120.65
2	B	1201	0NV	C23-C22-C21	-2.39	116.69	120.19
2	A	1201	0NV	C23-C22-C21	-2.23	116.92	120.19
2	B	1201	0NV	C19-N3-C4	-2.15	106.29	111.08
2	B	1201	0NV	C16-C15-N14	2.17	110.89	105.48
2	A	1201	0NV	C16-C15-N14	2.53	111.80	105.48
2	A	1201	0NV	C15-N14-C13	2.88	108.68	105.40
2	A	1201	0NV	C8-N7-C18	3.06	108.07	104.75
2	A	1201	0NV	C11-C10-N9	3.31	134.55	117.69
2	B	1201	0NV	C8-N7-C18	3.32	108.36	104.75
2	B	1201	0NV	C11-C10-N9	3.43	135.17	117.69
2	B	1201	0NV	C11-C10-C18	3.88	118.11	111.15
2	A	1201	0NV	C11-C10-C18	3.90	118.14	111.15
2	B	1201	0NV	C15-N14-C13	4.70	110.76	105.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1201	0NV	C8-N7-C6	6.45	124.99	114.28
2	B	1201	0NV	C8-N7-C6	7.75	127.15	114.28

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	1201	0NV	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

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, 95^{th} 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	293/298 (98%)	0.44	18 (6%)	25	32	14, 24, 55, 102	0
1	B	296/298 (99%)	0.39	21 (7%)	19	27	11, 21, 53, 95	0
All	All	589/596 (98%)	0.41	39 (6%)	22	29	11, 23, 54, 102	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1132	GLY	13.2
1	A	871	LEU	10.8
1	A	872	GLN	9.2
1	B	875	THR	6.9
1	A	1132	GLY	6.9
1	B	873	ASP	6.8
1	B	1131	ALA	5.1
1	B	874	ASN	4.0
1	A	946	GLU	3.8
1	B	1004	ASP	3.6
1	A	1004	ASP	3.5
1	B	946	GLU	3.1
1	B	886	HIS	3.1
1	B	1063	ARG	3.0
1	B	871	LEU	3.0
1	A	1003	GLN	2.9
1	B	1053	LYS	2.8
1	A	1053	LYS	2.8
1	A	1131	ALA	2.6
1	A	1063	ARG	2.6
1	B	1070	GLN	2.6
1	A	1092	ASP	2.5
1	B	1051	ILE	2.5
1	B	1093	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	1003	GLN	2.4
1	B	1052	GLU	2.3
1	A	1091	PRO	2.3
1	A	1096	ASP	2.3
1	A	1111	ASN	2.3
1	B	859	ASN	2.3
1	A	942	GLN	2.2
1	B	1076	PHE	2.2
1	B	1073	MET	2.1
1	A	876	GLY	2.1
1	B	1067	ASN	2.1
1	A	1052	GLU	2.0
1	A	947	ARG	2.0
1	A	1051	ILE	2.0
1	B	942	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, 95^{th} 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	0.95	0.11	-	15,26,50,51	0
1	PTR	B	1008	16/17	0.91	0.15	-	15,22,83,86	0
1	PTR	A	1008	16/17	0.90	0.15	-	17,24,49,67	0
1	PTR	A	1007	16/17	0.94	0.11	-	17,36,97,98	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, 95^{th} 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
2	0NV	A	1201	25/25	0.95	0.10	-0.31	6,16,31,31	0
2	0NV	B	1201	25/25	0.95	0.10	-0.37	9,17,30,34	0

6.5 Other polymers (i)

There are no such residues in this entry.