Full wwPDB X-ray Structure Validation Report (i)

Feb 28, 2024 – 01:21 AM EST

PDB ID	:	5QJ5
Title	:	PanDDA analysis group deposition – Crystal Structure of NUDT5 in complex with Z44592329
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Deposited on	:	2018-10-31
Resolution	:	1.72 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 (i) 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 (i)) were used in the production of this report:

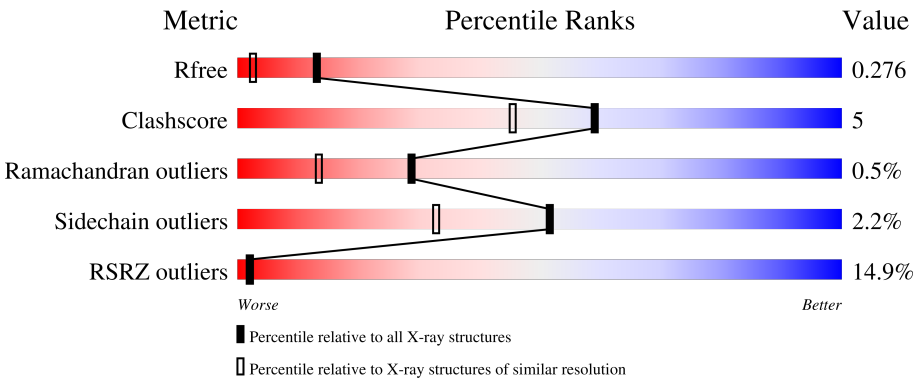
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 1.72 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}	130704	5722 (1.74-1.70)
Clashscore	141614	6152 (1.74-1.70)
Ramachandran outliers	138981	6051 (1.74-1.70)
Sidechain outliers	138945	6051 (1.74-1.70)
RSRZ outliers	127900	5629 (1.74-1.70)

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 <=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	209	13%; 84%; 8%; • 7%
1	B	209	14%; 83%; 10%; 7%
1	C	209	14%; 79%; 13%; 8%
1	D	209	14%; 84%; 6%; 9%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	EDO	B	301	-	-	X	-

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 6064 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 ADP-sugar pyrophosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	194	1458	922	243	285	8	0	0	0
1	B	194	1464	924	244	288	8	0	0	0
1	C	192	1402	884	234	277	7	0	0	0
1	D	190	1410	890	238	274	8	0	1	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP Q9UKK9
B	0	SER	-	expression tag	UNP Q9UKK9
C	0	SER	-	expression tag	UNP Q9UKK9
D	0	SER	-	expression tag	UNP Q9UKK9

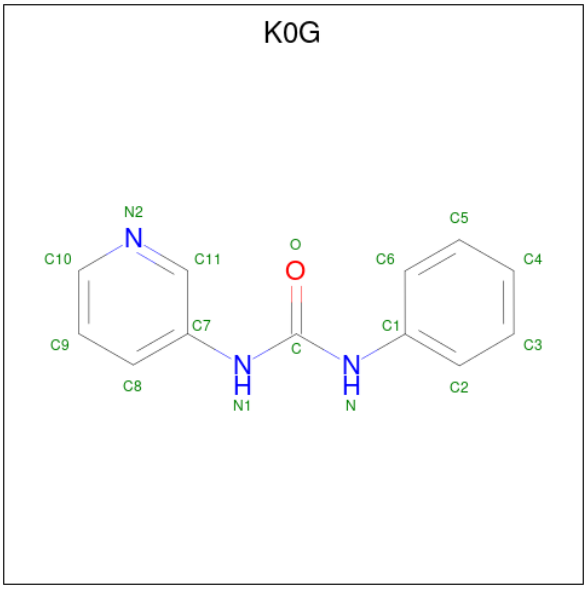
- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
2	A	2	2	2	0	0
2	B	2	2	2	0	0
2	C	2	2	2	0	0
2	D	2	2	2	0	0

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

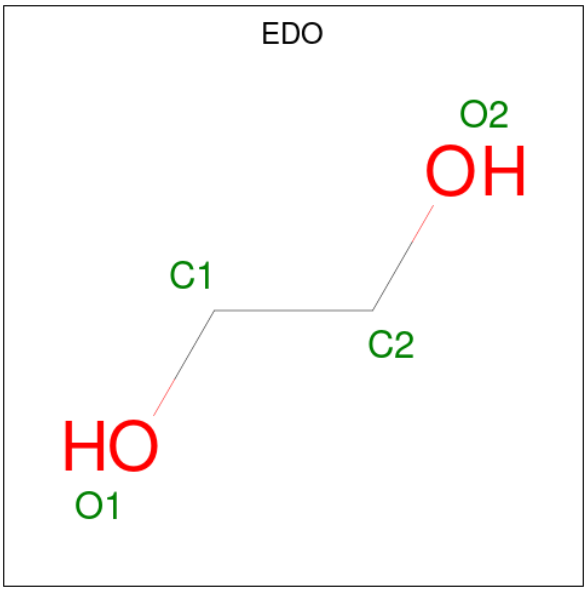
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Cl		
3	A	1	1	1	0	0

- Molecule 4 is N-phenyl-N'-pyridin-3-ylurea (three-letter code: K0G) (formula: $C_{12}H_{11}N_3O$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
4	A	1	16	12	3	1	0	0
			Total	C	N	O		
4	B	1	16	12	3	1	0	0
			Total	C	N	O		
4	C	1	16	12	3	1	0	0
			Total	C	N	O		
4	D	1	16	12	3	1	0	0

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
5	B	1	4	2	2	0	0
			Total	C	O		
5	C	1	4	2	2	0	0

- Molecule 6 is water.

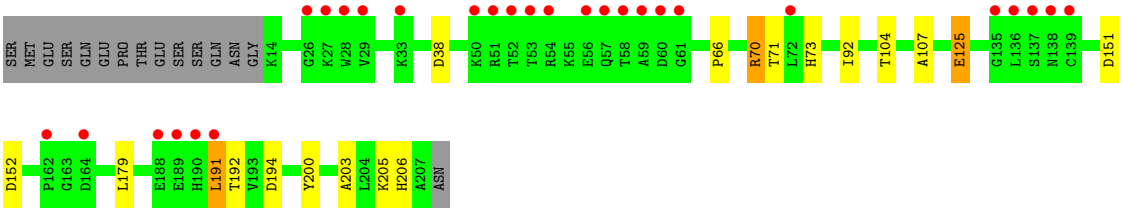
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	67	67	67	0	0
			Total	O		
6	B	73	73	73	0	0
			Total	O		
6	C	70	70	70	0	0
			Total	O		
6	D	39	39	39	0	0

3 Residue-property plots (i)

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.

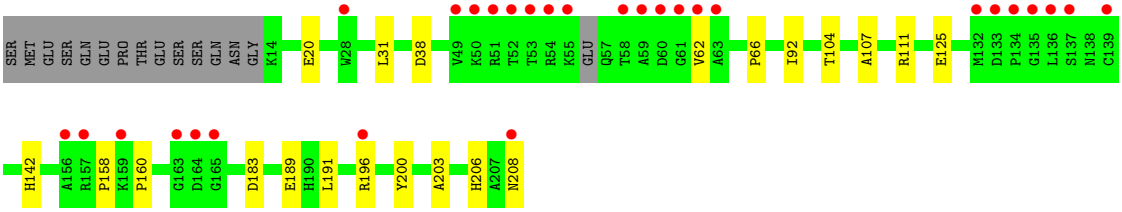
• Molecule 1: ADP-sugar pyrophosphatase

Chain A:



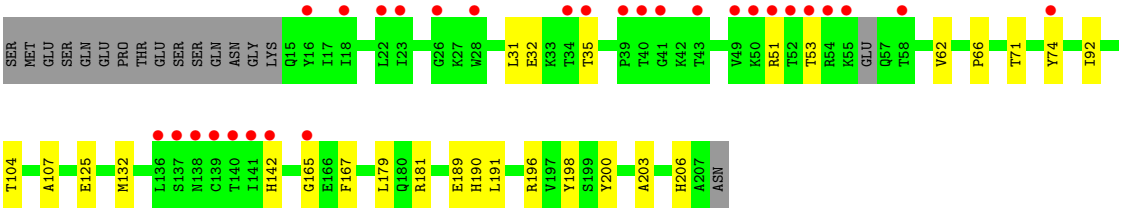
• Molecule 1: ADP-sugar pyrophosphatase

Chain B:



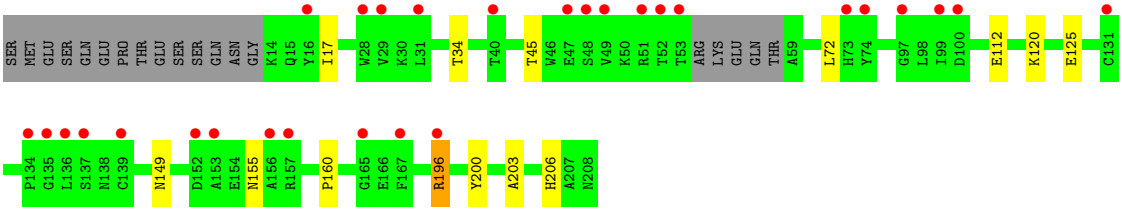
• Molecule 1: ADP-sugar pyrophosphatase

Chain C:



• Molecule 1: ADP-sugar pyrophosphatase

Chain D:


SER
MET
GLU
SER
GLN
GLU
PRO
THR
GLU
SER
SER
GLN
ASN
GLY
K14
Q15
Y16
I17
W28
V29
K30
L31
T34
T40
T45
W46
E47
S48
V49
K50
R51
T52
T53
ARG
LYS
GLU
GLN
THR
A59
L72
H73
Y74
G97
L98
I99
D100
E112
K120
E125
C131
P134
G135
L136
S137
N138
C139
N149
D152
A153
E154
N155
A156
R157
P160
G165
E166
F167
R196
Y200
A203
H206
A207
N208

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	49.15Å 59.77Å 80.09Å 79.47° 81.89° 75.78°	Depositor
Resolution (Å)	78.34 – 1.72 47.40 – 1.72	Depositor EDS
% Data completeness (in resolution range)	96.4 (78.34-1.72) 96.4 (47.40-1.72)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.44 (at 1.72Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.215 , 0.253 0.242 , 0.276	Depositor DCC
R_{free} test set	4450 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	30.0	Xtriage
Anisotropy	0.130	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 46.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	6064	wwPDB-VP
Average B, all atoms (Å^2)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.76% 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, 0.333 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: EDO, MG, K0G, CL

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	1.00	0/1486	1.01	5/2028 (0.2%)
1	B	0.90	0/1492	0.94	2/2035 (0.1%)
1	C	0.95	0/1429	0.98	3/1958 (0.2%)
1	D	0.90	0/1436	1.00	4/1961 (0.2%)
All	All	0.94	0/5843	0.98	14/7982 (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	D	0	1

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	70	ARG	NE-CZ-NH2	-8.97	115.81	120.30
1	D	196[A]	ARG	NE-CZ-NH1	8.73	124.67	120.30
1	D	196[B]	ARG	NE-CZ-NH1	8.73	124.67	120.30
1	D	196[A]	ARG	NE-CZ-NH2	-8.14	116.23	120.30
1	D	196[B]	ARG	NE-CZ-NH2	-8.14	116.23	120.30
1	C	196	ARG	NE-CZ-NH1	7.71	124.16	120.30
1	A	70	ARG	NE-CZ-NH1	6.16	123.38	120.30
1	B	111	ARG	NE-CZ-NH1	5.92	123.26	120.30
1	B	183	ASP	CB-CG-OD1	5.76	123.48	118.30
1	A	194	ASP	CB-CG-OD1	-5.73	113.14	118.30
1	A	152	ASP	CB-CG-OD1	5.41	123.17	118.30
1	C	190	HIS	N-CA-CB	5.34	120.21	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	181	ARG	NE-CZ-NH1	5.23	122.91	120.30
1	A	152	ASP	CB-CG-OD2	-5.16	113.66	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	160	PRO	Peptide

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	1458	0	1420	18	0
1	B	1464	0	1414	16	0
1	C	1402	0	1298	18	0
1	D	1410	0	1349	11	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	1	0	0	0	0
4	A	16	0	0	0	0
4	B	16	0	0	0	0
4	C	16	0	0	0	0
4	D	16	0	0	0	0
5	B	4	0	6	4	0
5	C	4	0	6	0	0
6	A	67	0	0	1	0
6	B	73	0	0	5	0
6	C	70	0	0	4	0
6	D	39	0	0	1	0
All	All	6064	0	5493	53	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 5.

All (53) 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:C:125:GLU:OE1	6:C:401:HOH:O	1.90	0.89
1:B:92:ILE:HD11	1:B:191:LEU:HD13	1.53	0.89
1:D:120:LYS:H	1:D:155:ASN:HD21	1.27	0.82
1:A:104:THR:HG23	1:A:107:ALA:H	1.46	0.79
1:C:203:ALA:HB3	1:D:203:ALA:HB3	1.69	0.74
5:B:301:EDO:H22	6:B:409:HOH:O	1.88	0.73
1:B:196:ARG:NH2	6:B:401:HOH:O	2.23	0.70
1:A:92:ILE:HD11	1:A:191:LEU:HD11	1.72	0.70
5:B:301:EDO:C2	6:B:409:HOH:O	2.41	0.68
1:A:203:ALA:HB3	1:B:203:ALA:HB3	1.80	0.63
1:C:179:LEU:HD11	1:C:198:TYR:CZ	2.33	0.62
1:A:179:LEU:HD23	1:A:205:LYS:HD2	1.83	0.60
1:C:104:THR:HG23	1:C:107:ALA:H	1.67	0.60
6:C:401:HOH:O	1:D:206:HIS:HE1	1.86	0.58
1:A:191:LEU:C	1:A:191:LEU:HD23	2.25	0.57
1:C:200:TYR:CD1	1:D:203:ALA:HB2	2.40	0.56
1:A:200:TYR:CD1	1:B:203:ALA:HB2	2.40	0.56
1:B:104:THR:HG23	1:B:107:ALA:H	1.70	0.56
1:A:125:GLU:OE1	1:B:206:HIS:HE1	1.90	0.55
1:C:132:MET:O	1:D:196[B]:ARG:NH1	2.41	0.54
5:B:301:EDO:H21	6:B:452:HOH:O	2.09	0.53
1:A:71:THR:HG23	1:A:151:ASP:OD2	2.08	0.53
1:A:206:HIS:HE1	1:B:125:GLU:OE1	1.92	0.53
1:D:112:GLU:OE2	6:D:401:HOH:O	2.19	0.51
1:C:200:TYR:OH	1:D:206:HIS:HD2	1.94	0.50
1:C:62:VAL:HG12	1:C:142:HIS:HB2	1.93	0.49
1:B:62:VAL:HG12	1:B:142:HIS:HB2	1.94	0.49
1:A:206:HIS:HD2	1:B:200:TYR:OH	1.96	0.49
1:C:206:HIS:HE1	1:D:125:GLU:OE1	1.96	0.49
1:C:125:GLU:HG2	6:C:405:HOH:O	2.12	0.48
1:B:208:ASN:OD1	1:B:208:ASN:C	2.53	0.47
5:B:301:EDO:C2	6:B:452:HOH:O	2.61	0.47
1:B:62:VAL:HA	1:B:142:HIS:O	2.16	0.46
1:C:203:ALA:HB2	1:D:200:TYR:CD1	2.51	0.45
1:C:92:ILE:HD12	1:C:191:LEU:HD11	1.99	0.45
1:A:203:ALA:HB2	1:B:200:TYR:CD1	2.52	0.45
1:C:191:LEU:HD12	1:C:191:LEU:C	2.37	0.45
1:D:17:ILE:HD12	1:D:34:THR:CG2	2.48	0.44
1:A:92:ILE:HD12	1:A:191:LEU:HD21	1.99	0.44
1:C:206:HIS:HD2	1:D:200:TYR:OH	2.01	0.43
1:A:92:ILE:CD1	1:A:191:LEU:HD11	2.46	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:70:ARG:HD3	6:A:433:HOH:O	2.18	0.42
1:B:189:GLU:OE2	1:B:191:LEU:HD11	2.18	0.42
1:A:200:TYR:OH	1:B:206:HIS:HD2	2.00	0.42
1:A:191:LEU:HD23	1:A:192:THR:N	2.34	0.42
1:C:31:LEU:HD23	1:C:32:GLU:H	1.85	0.41
1:C:165:GLY:HA2	1:C:167:PHE:CE1	2.55	0.41
1:A:191:LEU:C	1:A:191:LEU:CD2	2.88	0.41
1:C:92:ILE:HD12	1:C:191:LEU:CD1	2.51	0.41
1:B:158:PRO:O	1:B:160:PRO:HD3	2.21	0.41
1:A:71:THR:O	1:A:73:HIS:HD2	2.04	0.40
1:B:92:ILE:CD1	1:B:191:LEU:HD13	2.37	0.40
1:C:74:TYR:HB3	6:C:424:HOH:O	2.21	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	192/209 (92%)	186 (97%)	5 (3%)	1 (0%)	29 13
1	B	190/209 (91%)	185 (97%)	4 (2%)	1 (0%)	29 13
1	C	188/209 (90%)	181 (96%)	5 (3%)	2 (1%)	14 3
1	D	187/209 (90%)	180 (96%)	7 (4%)	0	100 100
All	All	757/836 (91%)	732 (97%)	21 (3%)	4 (0%)	29 13

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	53	THR
1	A	66	PRO
1	C	66	PRO

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Mol	Chain	Res	Type
1	B	66	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	153/179 (86%)	150 (98%)	3 (2%)	55	37
1	B	154/179 (86%)	151 (98%)	3 (2%)	57	39
1	C	139/179 (78%)	135 (97%)	4 (3%)	42	22
1	D	144/179 (80%)	141 (98%)	3 (2%)	53	35
All	All	590/716 (82%)	577 (98%)	13 (2%)	52	33

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	ASP
1	A	125	GLU
1	A	191	LEU
1	B	20	GLU
1	B	31	LEU
1	B	38	ASP
1	C	35	THR
1	C	51	ARG
1	C	71	THR
1	C	189	GLU
1	D	45	THR
1	D	72	LEU
1	D	149	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	HIS
1	A	206	HIS

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Mol	Chain	Res	Type
1	B	15	GLN
1	B	73	HIS
1	B	206	HIS
1	C	206	HIS
1	D	15	GLN
1	D	149	ASN
1	D	155	ASN
1	D	206	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)

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 15 ligands modelled in this entry, 9 are monoatomic - leaving 6 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
4	K0G	C	304	-	17,17,17	1.13	1 (5%)	21,21,21	1.08	1 (4%)
5	EDO	B	301	-	3,3,3	0.61	0	2,2,2	1.09	0
4	K0G	D	303	-	17,17,17	1.42	2 (11%)	21,21,21	0.98	0
4	K0G	A	304	-	17,17,17	1.14	2 (11%)	21,21,21	0.94	1 (4%)
5	EDO	C	301	-	3,3,3	0.80	0	2,2,2	0.10	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
4	K0G	B	304	-	17,17,17	1.26	2 (11%)	21,21,21	0.59	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
4	K0G	C	304	-	-	2/8/8/8	0/2/2/2
5	EDO	B	301	-	-	1/1/1/1	-
4	K0G	D	303	-	-	1/8/8/8	0/2/2/2
4	K0G	A	304	-	-	1/8/8/8	0/2/2/2
5	EDO	C	301	-	-	1/1/1/1	-
4	K0G	B	304	-	-	2/8/8/8	0/2/2/2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	303	K0G	C2-C1	2.74	1.43	1.39
4	B	304	K0G	C2-C1	2.71	1.43	1.39
4	D	303	K0G	C-N1	2.56	1.42	1.37
4	C	304	K0G	C8-C7	2.38	1.43	1.39
4	A	304	K0G	C8-C7	2.26	1.43	1.39
4	A	304	K0G	C6-C1	2.23	1.43	1.39
4	B	304	K0G	C-N1	2.06	1.41	1.37

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	C	304	K0G	C7-N1-C	-3.35	119.76	126.61
4	A	304	K0G	C7-N1-C	-2.16	122.19	126.61

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	304	K0G	C11-C7-N1-C
4	B	304	K0G	C11-C7-N1-C
4	D	303	K0G	C11-C7-N1-C
5	B	301	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
4	A	304	K0G	C11-C7-N1-C
4	C	304	K0G	C8-C7-N1-C
4	B	304	K0G	C8-C7-N1-C
5	C	301	EDO	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	301	EDO	4	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	194/209 (92%)	0.50	28 (14%)	2	2	21, 32, 58, 87	26 (13%)
1	B	194/209 (92%)	0.62	29 (14%)	2	2	21, 37, 64, 109	21 (10%)
1	C	192/209 (91%)	0.45	29 (15%)	2	2	22, 39, 71, 93	21 (10%)
1	D	190/209 (90%)	0.65	29 (15%)	2	2	25, 41, 70, 78	26 (13%)
All	All	770/836 (92%)	0.56	115 (14%)	2	2	21, 37, 69, 109	94 (12%)

All (115) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	53	THR	9.3
1	B	52	THR	9.1
1	A	191	LEU	8.8
1	D	74	TYR	8.0
1	A	139	CYS	7.5
1	B	63	ALA	7.0
1	A	136	LEU	6.9
1	A	28	TRP	6.7
1	A	53	THR	6.7
1	B	135	GLY	6.6
1	B	54	ARG	6.6
1	B	208	ASN	6.5
1	B	139	CYS	6.4
1	B	134	PRO	6.4
1	B	49	VAL	6.3
1	D	31	LEU	6.3
1	B	136	LEU	6.2
1	B	132	MET	5.9
1	A	59	ALA	5.9
1	C	52	THR	5.7
1	A	29	VAL	5.5

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Mol	Chain	Res	Type	RSRZ
1	B	62	VAL	5.5
1	A	52	THR	5.4
1	C	51	ARG	5.4
1	A	58	THR	5.4
1	D	73	HIS	5.2
1	B	59	ALA	5.1
1	A	137	SER	5.0
1	A	190	HIS	5.0
1	A	54	ARG	5.0
1	A	51	ARG	4.9
1	A	135	GLY	4.9
1	C	141	ILE	4.8
1	B	61	GLY	4.7
1	B	133	ASP	4.6
1	B	137	SER	4.6
1	C	139	CYS	4.6
1	A	61	GLY	4.5
1	B	50	LYS	4.4
1	A	189	GLU	4.3
1	A	50	LYS	4.3
1	C	41	GLY	4.3
1	D	53	THR	4.0
1	D	29	VAL	3.9
1	C	18	ILE	3.9
1	B	196	ARG	3.9
1	C	23	ILE	3.8
1	C	28	TRP	3.8
1	A	138	ASN	3.8
1	D	136	LEU	3.7
1	B	157	ARG	3.6
1	D	131	CYS	3.6
1	D	196[A]	ARG	3.6
1	D	157	ARG	3.6
1	C	49	VAL	3.6
1	B	51	ARG	3.5
1	D	97	GLY	3.4
1	B	60	ASP	3.4
1	C	54	ARG	3.4
1	B	28	TRP	3.3
1	C	16	TYR	3.2
1	C	58	THR	3.2
1	D	134	PRO	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	99	ILE	3.1
1	D	167	PHE	3.1
1	A	26	GLY	3.1
1	C	50	LYS	3.0
1	B	58	THR	3.0
1	C	26	GLY	3.0
1	C	140	THR	3.0
1	C	142	HIS	3.0
1	C	165	GLY	2.9
1	C	136	LEU	2.9
1	D	28	TRP	2.9
1	D	153	ALA	2.9
1	D	152	ASP	2.8
1	A	57	GLN	2.8
1	C	43	THR	2.8
1	D	139	CYS	2.8
1	D	48	SER	2.7
1	B	163	GLY	2.7
1	C	74	TYR	2.7
1	A	164	ASP	2.7
1	B	164	ASP	2.6
1	C	137	SER	2.6
1	B	159	LYS	2.5
1	C	34	THR	2.5
1	A	60	ASP	2.5
1	A	27	LYS	2.5
1	D	156	ALA	2.5
1	B	165	GLY	2.5
1	C	53	THR	2.4
1	C	22	LEU	2.4
1	B	55	LYS	2.4
1	D	16	TYR	2.4
1	A	56	GLU	2.4
1	D	49	VAL	2.3
1	A	188	GLU	2.3
1	C	35	THR	2.3
1	C	39	PRO	2.3
1	D	47	GLU	2.3
1	A	72	LEU	2.3
1	C	55	LYS	2.3
1	D	165	GLY	2.3
1	D	40	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	52	THR	2.2
1	C	138	ASN	2.2
1	D	135	GLY	2.2
1	D	100	ASP	2.2
1	A	33	LYS	2.2
1	D	51	ARG	2.2
1	D	137	SER	2.2
1	A	162	PRO	2.1
1	B	156	ALA	2.1
1	C	40	THR	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides 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(Å^2)	Q<0.9
4	K0G	B	304	16/16	0.64	0.26	33,36,40,41	16
4	K0G	A	304	16/16	0.65	0.34	32,39,44,44	16
4	K0G	D	303	16/16	0.72	0.23	39,48,52,53	16
4	K0G	C	304	16/16	0.82	0.18	35,42,49,50	16
5	EDO	C	301	4/4	0.86	0.11	39,52,52,58	0
2	MG	D	301	1/1	0.89	0.05	43,43,43,43	0
2	MG	B	303	1/1	0.89	0.14	58,58,58,58	0
5	EDO	B	301	4/4	0.90	0.18	38,40,41,42	0
3	CL	A	303	1/1	0.91	0.08	70,70,70,70	0
2	MG	A	302	1/1	0.93	0.04	52,52,52,52	0
2	MG	D	302	1/1	0.96	0.11	41,41,41,41	0
2	MG	C	303	1/1	0.96	0.06	42,42,42,42	0
2	MG	B	302	1/1	0.98	0.03	35,35,35,35	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	MG	A	301	1/1	0.98	0.06	35,35,35,35	0
2	MG	C	302	1/1	0.99	0.06	62,62,62,62	0

6.5 Other polymers (i)

There are no such residues in this entry.