



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 03:12 PM UTC

PDB ID : 5QJM / pdb_00005qjm
Title : PanDDA analysis group deposition of models with modelled events (e.g. bound ligands) – Crystal Structure of NUDT5 in complex with Z328695024
Authors : Dubianok, Y.; Collins, P.; Krojer, T.; Wright, N.; Strain-Damerell, C.; Burgess-Brown, N.; Bountra, C.; Arrowsmith, C.H.; Edwards, A.; Huber, K.; von Delft, F.
Deposited on : 2018-10-31
Resolution : 1.75 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

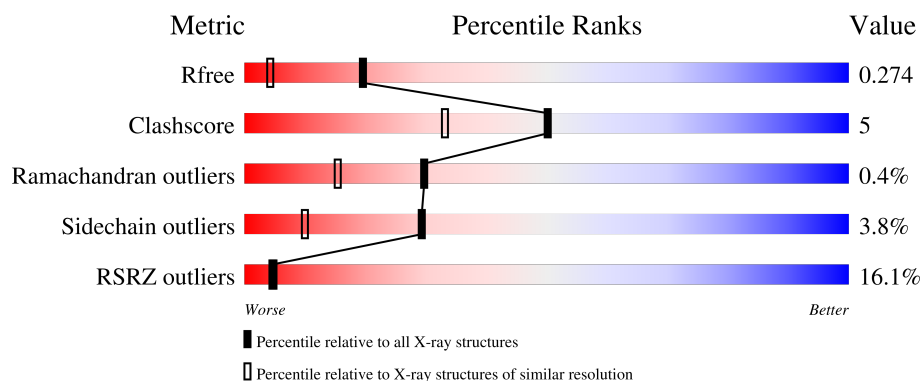
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3183 (1.76-1.76)
Clashscore	190562	3299 (1.76-1.76)
Ramachandran outliers	187476	3274 (1.76-1.76)
Sidechain outliers	187428	3274 (1.76-1.76)
RSRZ outliers	180081	3183 (1.76-1.76)

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

Mol	Chain	Length	Quality of chain
1	A	209	6% 83% 10% 7%
1	B	209	10% 82% 10% 7%
1	C	209	17% 81% 11% 8%
1	D	209	27% 74% 16% 9%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6090 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
1	A	194	Total	C	N	O	S	0	0	0
			1458	922	243	285	8			
1	B	194	Total	C	N	O	S	0	0	0
			1464	924	244	288	8			
1	C	192	Total	C	N	O	S	0	0	0
			1427	900	239	281	7			
1	D	190	Total	C	N	O	S	0	1	0
			1425	904	241	272	8			

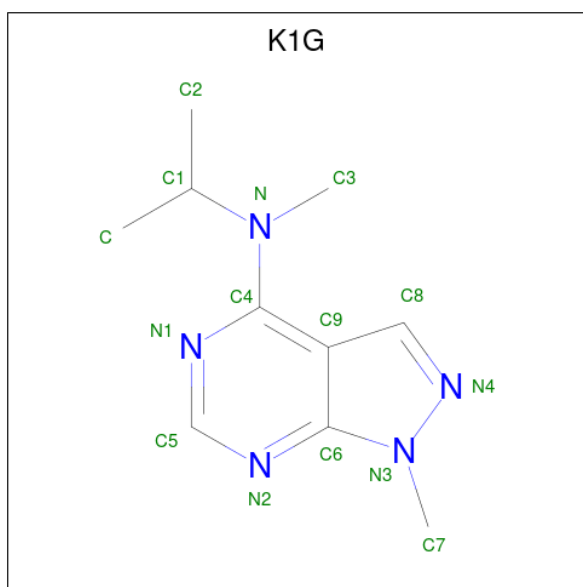
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP Q9UUKK9
B	0	SER	-	expression tag	UNP Q9UUKK9
C	0	SER	-	expression tag	UNP Q9UUKK9
D	0	SER	-	expression tag	UNP Q9UUKK9

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 3 is N,1-dimethyl-N-(propan-2-yl)-1H-pyrazolo[3,4-d]pyrimidin-4-amine (CCD ID: K1G) (formula: C₁₀H₁₅N₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			15	10	5		
3	B	1	Total	C	N	0	0
			15	10	5		
3	C	1	Total	C	N	0	0
			15	10	5		
3	D	1	Total	C	N	0	0
			15	10	5		

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Cl	0	0
			1	1		

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

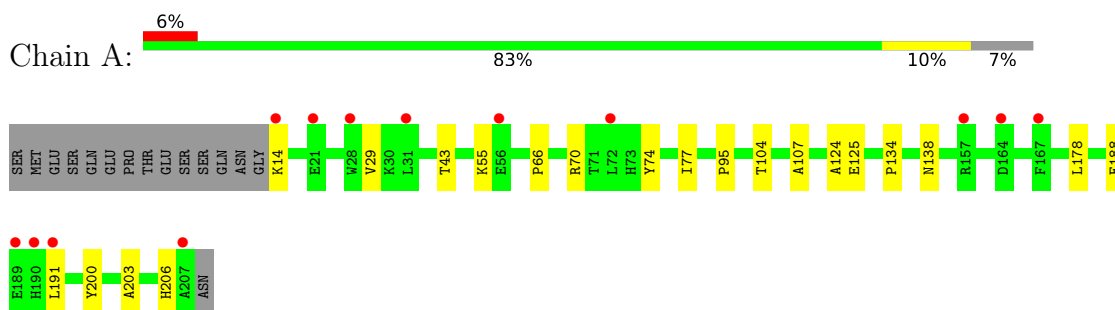
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	68	Total	O	0	0
			68	68		
6	B	71	Total	O	0	0
			71	71		
6	C	67	Total	O	0	0
			67	67		
6	D	29	Total	O	0	0
			29	29		

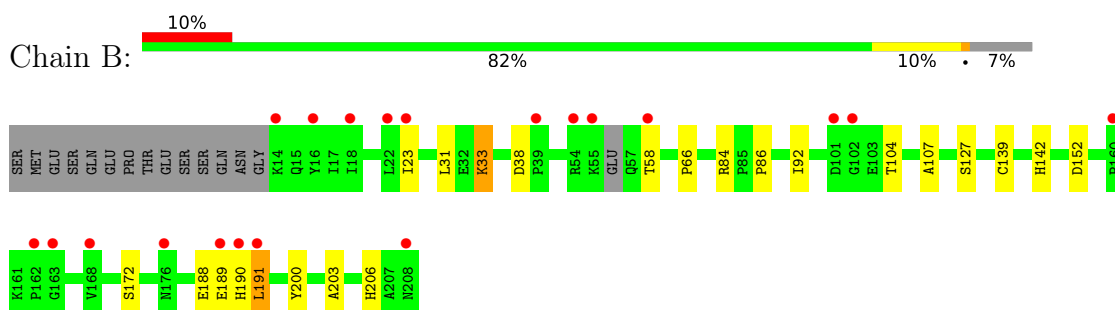
3 Residue-property plots [i](#)

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.

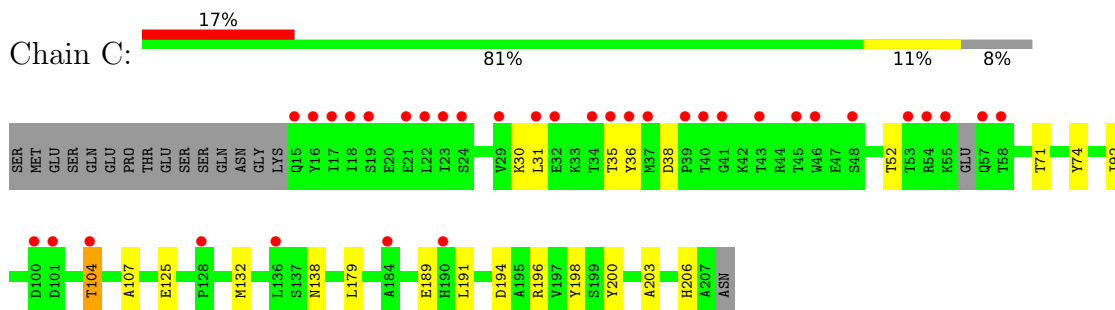
- Molecule 1: ADP-sugar pyrophosphatase



- Molecule 1: ADP-sugar pyrophosphatase

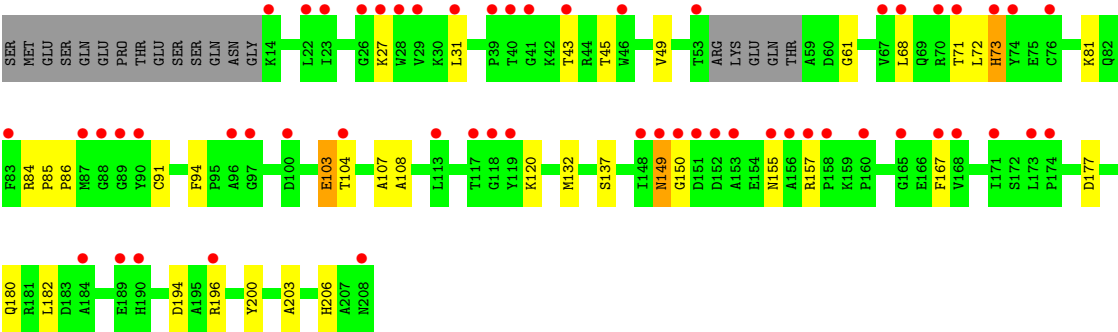


- Molecule 1: ADP-sugar pyrophosphatase



- Molecule 1: ADP-sugar pyrophosphatase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	49.64Å 59.95Å 80.27Å 79.09° 81.28° 75.44°	Depositor
Resolution (Å)	78.34 – 1.75 78.34 – 1.75	Depositor EDS
% Data completeness (in resolution range)	87.3 (78.34-1.75) 87.3 (78.34-1.75)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.25 (at 1.75Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.223 , 0.271 0.235 , 0.274	Depositor DCC
R_{free} test set	4452 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	29.0	Xtriage
Anisotropy	0.151	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 28.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6090	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.57% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: K1G, MG, CL, EDO

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	1.29	4/1486 (0.3%)	1.09	4/2028 (0.2%)
1	B	1.15	2/1492 (0.1%)	1.08	1/2035 (0.0%)
1	C	1.18	1/1454 (0.1%)	1.09	0/1987
1	D	1.20	2/1453 (0.1%)	1.13	4/1983 (0.2%)
All	All	1.21	9/5885 (0.2%)	1.10	9/8033 (0.1%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	55	LYS	CA-C	9.99	1.57	1.52
1	D	84	ARG	C-O	-9.01	1.17	1.25
1	D	61	GLY	N-CA	5.72	1.50	1.45
1	B	139	CYS	C-O	5.62	1.30	1.23
1	B	127	SER	C-O	-5.46	1.16	1.23
1	C	104	THR	CA-C	5.45	1.59	1.53
1	A	178	LEU	CA-C	5.23	1.59	1.52
1	A	95	PRO	C-O	5.07	1.29	1.23
1	A	124	ALA	CA-C	-5.05	1.47	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	94	PHE	CA-C-N	-7.38	112.15	119.76
1	D	94	PHE	C-N-CA	-7.38	112.15	119.76
1	A	188	GLU	N-CA-C	6.70	121.19	112.89
1	D	85	PRO	N-CA-C	6.17	118.22	110.70
1	A	77	ILE	N-CA-C	-5.69	100.21	108.17
1	D	182	LEU	N-CA-C	5.53	116.99	111.07
1	A	29	VAL	CB-CA-C	-5.40	101.37	110.71
1	B	188	GLU	N-CA-C	5.10	119.58	112.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	74	TYR	N-CA-C	-5.02	103.06	110.48

There are no chirality outliers.

There are no planarity outliers.

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	1458	0	1420	10	0
1	B	1464	0	1414	20	0
1	C	1427	0	1348	16	0
1	D	1425	0	1382	28	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	15	0	0	0	0
3	B	15	0	0	0	0
3	C	15	0	0	0	0
3	D	15	0	0	0	0
4	B	1	0	0	0	0
5	B	8	0	12	1	0
5	C	4	0	6	0	0
6	A	68	0	0	0	0
6	B	71	0	0	1	0
6	C	67	0	0	4	0
6	D	29	0	0	1	0
All	All	6090	0	5582	60	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 (60) 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:D:196[B]:ARG:HH21	1:D:196[B]:ARG:CG	1.53	1.17
1:D:196[B]:ARG:HG3	1:D:196[B]:ARG:NH2	1.56	1.09
1:B:189:GLU:HB3	1:B:191:LEU:HD12	1.48	0.92
1:D:196[B]:ARG:HH21	1:D:196[B]:ARG:HG3	0.74	0.87
1:C:125:GLU:OE1	6:C:401:HOH:O	2.00	0.79
1:D:196[B]:ARG:NH2	1:D:196[B]:ARG:N	2.36	0.72
1:C:132:MET:HE1	1:C:196:ARG:CZ	2.20	0.72
1:D:196[B]:ARG:CG	1:D:196[B]:ARG:NH2	2.25	0.68
1:C:104:THR:HG23	1:C:107:ALA:H	1.60	0.67
1:C:203:ALA:HB3	1:D:203:ALA:HB3	1.77	0.66
1:A:203:ALA:HB3	1:B:203:ALA:HB3	1.78	0.66
1:B:189:GLU:OE2	1:B:191:LEU:HD11	1.95	0.65
1:D:120:LYS:H	1:D:155:ASN:HD21	1.43	0.64
1:A:104:THR:HG23	1:A:107:ALA:H	1.65	0.62
1:A:206:HIS:HD2	1:B:200:TYR:OH	1.84	0.59
1:B:104:THR:HG23	1:B:107:ALA:H	1.67	0.58
1:D:196[B]:ARG:NH2	1:D:196[B]:ARG:H	2.00	0.57
1:C:206:HIS:HD2	1:D:200:TYR:OH	1.87	0.57
1:B:189:GLU:CB	1:B:191:LEU:HD12	2.30	0.56
1:C:194:ASP:OD2	1:C:196:ARG:NH2	2.40	0.54
1:D:196[B]:ARG:NH2	1:D:196[B]:ARG:CB	2.71	0.52
1:C:200:TYR:OH	1:D:206:HIS:HD2	1.92	0.52
1:D:149:ASN:C	1:D:149:ASN:HD22	2.17	0.52
1:C:200:TYR:CD1	1:D:203:ALA:HB2	2.46	0.51
1:D:132:MET:HE1	1:D:196[A]:ARG:CZ	2.41	0.51
1:D:150:GLY:O	6:D:401:HOH:O	2.19	0.50
1:D:196[B]:ARG:HH21	1:D:196[B]:ARG:CB	2.20	0.49
1:C:179:LEU:HD11	1:C:198:TYR:CZ	2.47	0.49
1:A:200:TYR:OH	1:B:206:HIS:HD2	1.96	0.49
1:B:190:HIS:O	1:B:190:HIS:CG	2.65	0.48
1:D:71:THR:O	1:D:73:HIS:ND1	2.45	0.48
1:A:200:TYR:CD1	1:B:203:ALA:HB2	2.49	0.46
1:B:23:ILE:HD11	1:B:33:LYS:HB2	1.98	0.46
1:B:92:ILE:HD11	1:B:191:LEU:HD13	1.97	0.46
1:B:189:GLU:HG2	1:B:191:LEU:HD11	1.98	0.46
1:B:189:GLU:OE2	1:B:191:LEU:CD1	2.61	0.46
1:A:125:GLU:OE1	1:B:206:HIS:HE1	1.98	0.46
1:D:177:ASP:OD1	1:D:180:GLN:CB	2.64	0.46
1:C:138:ASN:HB3	1:D:86:PRO:O	2.16	0.45
1:A:104:THR:CG2	1:A:107:ALA:H	2.27	0.45
1:D:196[B]:ARG:NH2	1:D:196[B]:ARG:CA	2.79	0.45
1:A:134:PRO:HB2	1:B:84:ARG:HG3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:132:MET:O	1:D:196[B]:ARG:NH1	2.50	0.43
1:A:138:ASN:HB3	1:B:86:PRO:O	2.18	0.43
1:D:104:THR:HG23	1:D:107:ALA:H	1.84	0.43
1:B:58:THR:CB	1:B:142:HIS:NE2	2.82	0.43
1:D:103:GLU:CD	1:D:108:ALA:HB2	2.43	0.43
1:C:30:LYS:HG3	1:C:52:THR:HG22	2.01	0.42
1:C:38:ASP:HB2	1:D:167:PHE:CE2	2.54	0.42
1:D:194:ASP:OD1	1:D:196[A]:ARG:HD3	2.19	0.42
1:C:92:ILE:HG13	1:C:191:LEU:HD13	2.02	0.42
6:C:401:HOH:O	1:D:206:HIS:HE1	2.02	0.42
1:C:36:TYR:OH	6:C:402:HOH:O	2.21	0.42
1:C:74:TYR:HB3	6:C:439:HOH:O	2.19	0.42
5:B:302:EDO:C2	6:B:405:HOH:O	2.69	0.41
1:B:189:GLU:CG	1:B:191:LEU:HD11	2.51	0.41
1:D:49:VAL:O	1:D:137:SER:HA	2.21	0.41
1:D:81:LYS:HA	1:D:91:CYS:O	2.20	0.41
1:B:189:GLU:CG	1:B:191:LEU:CD1	3.00	0.40
1:A:203:ALA:HB2	1:B:200:TYR:CD1	2.56	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%)	24	11
1	B	190/209 (91%)	185 (97%)	4 (2%)	1 (0%)	24	11
1	C	188/209 (90%)	185 (98%)	3 (2%)	0	100	100
1	D	187/209 (90%)	183 (98%)	3 (2%)	1 (0%)	24	11
All	All	757/836 (91%)	739 (98%)	15 (2%)	3 (0%)	30	15

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	66	PRO
1	B	66	PRO
1	D	157	ARG

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%)	149 (97%)	4 (3%)	40	20
1	B	154/179 (86%)	148 (96%)	6 (4%)	28	9
1	C	145/179 (81%)	141 (97%)	4 (3%)	38	18
1	D	147/179 (82%)	138 (94%)	9 (6%)	17	3
All	All	599/716 (84%)	576 (96%)	23 (4%)	29	10

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

Mol	Chain	Res	Type
1	A	14	LYS
1	A	43	THR
1	A	70	ARG
1	A	191	LEU
1	B	31	LEU
1	B	33	LYS
1	B	38	ASP
1	B	152	ASP
1	B	172	SER
1	B	191	LEU
1	C	31	LEU
1	C	35	THR
1	C	71	THR
1	C	189	GLU
1	D	27	LYS
1	D	31	LEU
1	D	43	THR
1	D	45	THR
1	D	68	LEU

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Mol	Chain	Res	Type
1	D	72	LEU
1	D	73	HIS
1	D	103	GLU
1	D	149	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	HIS
1	A	206	HIS
1	B	15	GLN
1	B	69	GLN
1	B	190	HIS
1	B	206	HIS
1	C	206	HIS
1	D	82	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 9 are monoatomic - leaving 7 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
3	K1G	B	305	-	15,16,16	2.22	6 (40%)	16,23,23	1.89	4 (25%)
3	K1G	C	304	-	15,16,16	1.85	3 (20%)	16,23,23	2.37	5 (31%)
3	K1G	A	303	-	15,16,16	2.20	7 (46%)	16,23,23	1.93	5 (31%)
5	EDO	C	301	-	3,3,3	0.56	0	2,2,2	0.51	0
5	EDO	B	306	-	3,3,3	0.46	0	2,2,2	0.38	0
5	EDO	B	302	-	3,3,3	0.74	0	2,2,2	0.16	0
3	K1G	D	303	-	15,16,16	1.79	5 (33%)	16,23,23	2.02	5 (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
3	K1G	B	305	-	-	2/8/8/8	0/2/2/2
3	K1G	C	304	-	-	2/8/8/8	0/2/2/2
3	K1G	A	303	-	-	1/8/8/8	0/2/2/2
5	EDO	C	301	-	-	1/1/1/1	-
5	EDO	B	306	-	-	0/1/1/1	-
5	EDO	B	302	-	-	0/1/1/1	-
3	K1G	D	303	-	-	5/8/8/8	0/2/2/2

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	303	K1G	C4-N	4.31	1.48	1.36
3	B	305	K1G	C4-N	4.23	1.48	1.36
3	C	304	K1G	C4-N	3.55	1.46	1.36
3	A	303	K1G	C5-N1	3.43	1.40	1.33
3	B	305	K1G	C9-C8	3.42	1.46	1.42
3	B	305	K1G	C4-N1	3.41	1.40	1.34
3	D	303	K1G	C4-N	3.32	1.45	1.36
3	C	304	K1G	C9-C8	2.86	1.45	1.42
3	A	303	K1G	C9-C6	2.84	1.44	1.40
3	B	305	K1G	C5-N1	2.63	1.38	1.33
3	B	305	K1G	C6-N2	2.58	1.39	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	305	K1G	C9-C6	2.52	1.44	1.40
3	D	303	K1G	C8-N4	2.50	1.34	1.32
3	A	303	K1G	C9-C4	2.48	1.46	1.42
3	A	303	K1G	C9-C8	2.45	1.44	1.42
3	D	303	K1G	C3-N	2.44	1.51	1.45
3	D	303	K1G	C9-C6	2.43	1.44	1.40
3	C	304	K1G	C6-N3	2.11	1.38	1.35
3	D	303	K1G	C-C1	2.10	1.57	1.52
3	A	303	K1G	C6-N3	2.04	1.38	1.35
3	A	303	K1G	C6-N2	2.01	1.38	1.34

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	304	K1G	C2-C1-N	6.59	121.04	111.45
3	B	305	K1G	C7-N3-C6	-4.81	124.92	128.30
3	A	303	K1G	C2-C1-N	4.64	118.20	111.45
3	D	303	K1G	C7-N3-C6	-4.61	125.05	128.30
3	C	304	K1G	C-C1-N	-3.46	106.40	111.45
3	D	303	K1G	C7-N3-N4	3.34	124.19	120.05
3	D	303	K1G	C5-N1-C4	3.23	119.73	111.83
3	D	303	K1G	C-C1-N	3.09	115.94	111.45
3	B	305	K1G	C5-N1-C4	3.08	119.35	111.83
3	A	303	K1G	C5-N1-C4	3.00	119.15	111.83
3	C	304	K1G	C5-N1-C4	2.89	118.90	111.83
3	A	303	K1G	C9-C4-N1	-2.88	117.40	122.53
3	C	304	K1G	C9-C4-N1	-2.79	117.55	122.53
3	C	304	K1G	N1-C4-N	2.64	119.88	116.82
3	A	303	K1G	C-C1-N	-2.60	107.67	111.45
3	B	305	K1G	C9-C4-N1	-2.48	118.11	122.53
3	D	303	K1G	C9-C4-N1	-2.42	118.22	122.53
3	B	305	K1G	C7-N3-N4	2.32	122.92	120.05
3	A	303	K1G	N1-C4-N	2.21	119.39	116.82

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	303	K1G	C2-C1-N-C3
3	C	304	K1G	C2-C1-N-C3
3	D	303	K1G	C-C1-N-C3
3	D	303	K1G	C2-C1-N-C3

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Mol	Chain	Res	Type	Atoms
5	C	301	EDO	O1-C1-C2-O2
3	C	304	K1G	C2-C1-N-C4
3	D	303	K1G	C-C1-N-C4
3	B	305	K1G	C-C1-N-C3
3	D	303	K1G	C9-C4-N-C1
3	D	303	K1G	N1-C4-N-C1
3	B	305	K1G	C2-C1-N-C3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	302	EDO	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, 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%)	0.69	13 (6%) 24 28	18, 30, 52, 64	8 (4%)
1	B	194/209 (92%)	0.74	20 (10%) 12 13	21, 34, 58, 68	3 (1%)
1	C	192/209 (91%)	1.06	35 (18%) 3 4	22, 36, 60, 67	1 (0%)
1	D	190/209 (90%)	1.62	56 (29%) 1 1	8, 42, 63, 70	3 (1%)
All	All	770/836 (92%)	1.02	124 (16%) 4 5	8, 36, 59, 70	15 (1%)

All (124) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	196[A]	ARG	11.5
1	A	191	LEU	11.4
1	A	190	HIS	10.2
1	A	189	GLU	8.8
1	D	190	HIS	8.6
1	C	54	ARG	7.7
1	D	189	GLU	6.8
1	C	22	LEU	5.9
1	A	21	GLU	5.7
1	B	191	LEU	5.6
1	C	23	ILE	5.3
1	D	40	THR	4.8
1	C	18	ILE	4.7
1	C	16	TYR	4.5
1	C	40	THR	4.3
1	C	43	THR	4.3
1	D	74	TYR	4.2
1	D	53	THR	4.2
1	C	21	GLU	4.0
1	D	167	PHE	3.9
1	C	58	THR	3.9

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Mol	Chain	Res	Type	RSRZ
1	D	153	ALA	3.8
1	D	113	LEU	3.6
1	A	56	GLU	3.6
1	A	207	ALA	3.5
1	B	101	ASP	3.5
1	B	190	HIS	3.4
1	D	14	LYS	3.4
1	D	68	LEU	3.4
1	D	23	ILE	3.3
1	D	168	VAL	3.3
1	B	208	ASN	3.3
1	D	26	GLY	3.2
1	C	57	GLN	3.2
1	C	17	ILE	3.2
1	D	97	GLY	3.2
1	D	71	THR	3.2
1	D	73	HIS	3.2
1	B	14	LYS	3.2
1	D	29	VAL	3.1
1	A	28	TRP	3.1
1	D	88	GLY	3.1
1	C	35	THR	3.0
1	B	58	THR	3.0
1	D	100	ASP	3.0
1	B	102	GLY	3.0
1	B	163	GLY	3.0
1	B	189	GLU	2.9
1	D	158	PRO	2.9
1	C	55	LYS	2.9
1	D	90	TYR	2.9
1	C	136	LEU	2.9
1	B	55	LYS	2.9
1	D	27	LYS	2.9
1	C	24	SER	2.9
1	D	119	TYR	2.8
1	C	53	THR	2.8
1	C	48	SER	2.8
1	B	54	ARG	2.8
1	D	149	ASN	2.8
1	C	45	THR	2.8
1	B	176	ASN	2.7
1	C	36	TYR	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	165	GLY	2.7
1	D	151	ASP	2.7
1	D	28	TRP	2.7
1	D	148	ILE	2.7
1	D	174	PRO	2.6
1	D	43	THR	2.6
1	A	31	LEU	2.6
1	C	184	ALA	2.6
1	D	155	ASN	2.6
1	D	152	ASP	2.5
1	C	37	MET	2.5
1	C	46	TRP	2.5
1	D	46	TRP	2.5
1	D	160	PRO	2.5
1	C	100	ASP	2.5
1	D	184	ALA	2.5
1	D	150	GLY	2.5
1	B	23	ILE	2.5
1	C	34	THR	2.5
1	D	39	PRO	2.4
1	B	168	VAL	2.4
1	D	208	ASN	2.4
1	A	157	ARG	2.4
1	B	22	LEU	2.4
1	D	89	GLY	2.4
1	D	156	ALA	2.4
1	D	22	LEU	2.4
1	C	15	GLN	2.3
1	D	41	GLY	2.3
1	B	160	PRO	2.3
1	D	83	PHE	2.3
1	B	16	TYR	2.3
1	B	18	ILE	2.3
1	D	171	ILE	2.3
1	B	162	PRO	2.3
1	D	70	ARG	2.3
1	D	104	THR	2.3
1	D	117	THR	2.3
1	D	76	CYS	2.2
1	B	39	PRO	2.2
1	C	19	SER	2.2
1	A	164	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	173	LEU	2.2
1	C	41	GLY	2.2
1	D	118	GLY	2.2
1	C	39	PRO	2.2
1	D	87	MET	2.2
1	D	96	ALA	2.2
1	C	32	GLU	2.2
1	A	72	LEU	2.2
1	D	31	LEU	2.2
1	C	190	HIS	2.2
1	C	101	ASP	2.1
1	D	157	ARG	2.1
1	D	67	VAL	2.1
1	A	14	LYS	2.1
1	C	29	VAL	2.0
1	C	31	LEU	2.0
1	C	104	THR	2.0
1	C	128	PRO	2.0
1	A	167	PHE	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	EDO	B	302	4/4	0.81	0.12	39,42,48,48	0
5	EDO	B	306	4/4	0.86	0.15	27,30,31,33	4
5	EDO	C	301	4/4	0.86	0.13	36,51,52,60	0
3	K1G	C	304	15/15	0.87	0.17	42,46,52,52	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	CL	B	301	1/1	0.89	0.13	62,62,62,62	0
3	K1G	D	303	15/15	0.89	0.14	42,46,50,51	0
2	MG	D	301	1/1	0.90	0.08	39,39,39,39	0
2	MG	A	302	1/1	0.93	0.08	53,53,53,53	0
3	K1G	B	305	15/15	0.93	0.14	30,34,40,40	15
2	MG	B	304	1/1	0.94	0.07	54,54,54,54	0
3	K1G	A	303	15/15	0.94	0.12	36,42,50,51	0
2	MG	C	302	1/1	0.94	0.06	60,60,60,60	0
2	MG	D	302	1/1	0.97	0.08	41,41,41,41	0
2	MG	B	303	1/1	0.98	0.06	35,35,35,35	0
2	MG	C	303	1/1	0.98	0.04	35,35,35,35	0
2	MG	A	301	1/1	0.99	0.03	34,34,34,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.