



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 13, 2020 – 09:31 pm BST

PDB ID : 5QK5
Title : PanDDA analysis group deposition of models with modelled events (e.g. bound ligands) – Crystal Structure of NUDT5 in complex with Z1267773786
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 : 1.63 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 following versions of software and data (see [references ⓘ](#)) 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 : 2.11
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) : 2.11

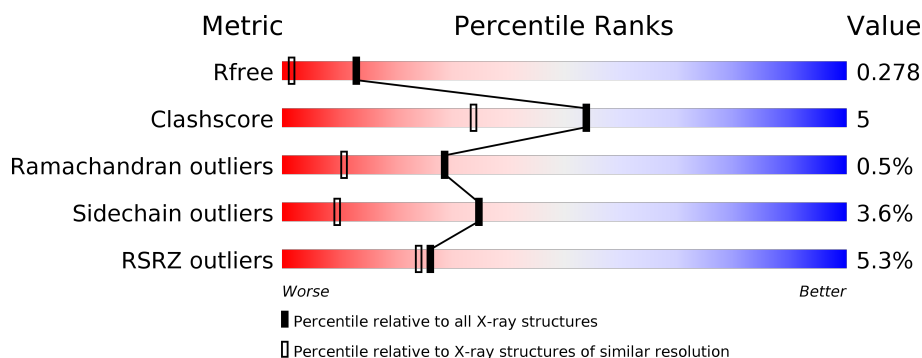
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.63 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	3122 (1.66-1.62)
Clashscore	141614	3268 (1.66-1.62)
Ramachandran outliers	138981	3215 (1.66-1.62)
Sidechain outliers	138945	3215 (1.66-1.62)
RSRZ outliers	127900	3079 (1.66-1.62)

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 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	<div> <div>4%</div> <div> <div></div> <div>81%</div> <div>11%</div> <div>7%</div> </div> </div>
1	B	209	<div> <div>%</div> <div> <div></div> <div>80%</div> <div>11%</div> <div>7%</div> </div> </div>
1	C	209	<div> <div>4%</div> <div> <div></div> <div>78%</div> <div>13%</div> <div>8%</div> </div> </div>
1	D	209	<div> <div>11%</div> <div> <div></div> <div>75%</div> <div>14%</div> <div>9%</div> </div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6086 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	923	244	283	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
			1406	887	235	277	7			
1	D	190	Total	C	N	O	S	0	1	0
			1419	899	240	272	8			

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

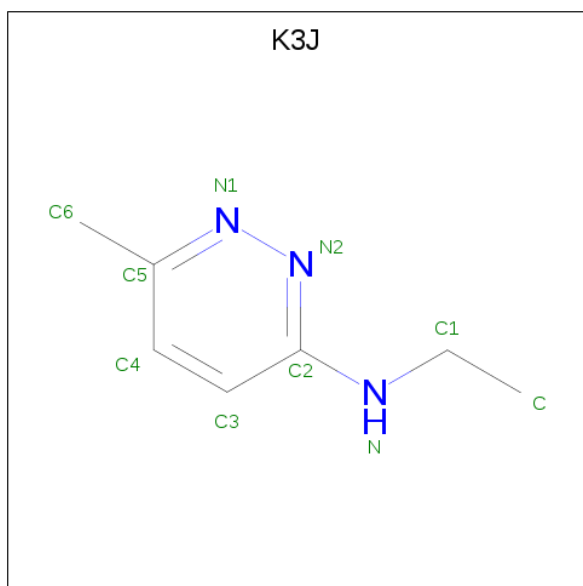
- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

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

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

- Molecule 4 is N-ethyl-6-methylpyridazin-3-amine (three-letter code: K3J) (formula: $C_7H_{11}N_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N 10 7 3	0	0
4	A	1	Total C N 10 7 3	0	0
4	C	1	Total C N 10 7 3	0	0
4	C	1	Total C N 10 7 3	0	0

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



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

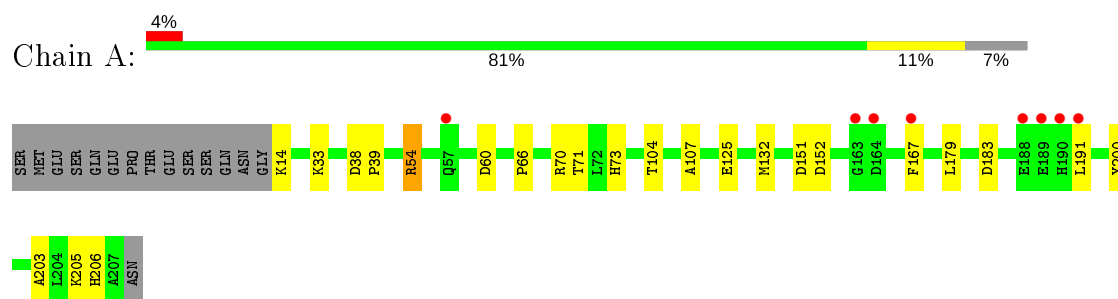
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	80	Total	O	0	0
			80	80		
6	B	82	Total	O	0	0
			82	82		
6	C	75	Total	O	0	0
			75	75		
6	D	33	Total	O	0	0
			33	33		

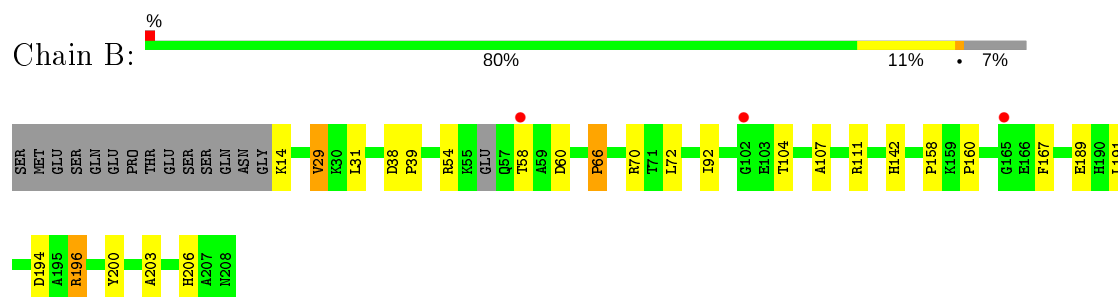
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

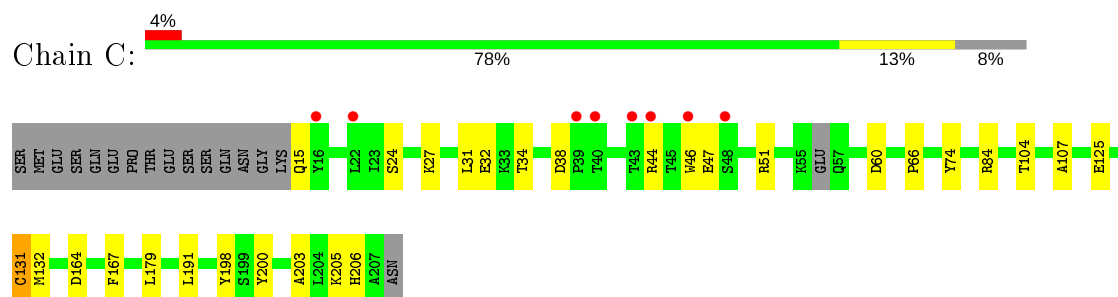
- Molecule 1: ADP-sugar pyrophosphatase



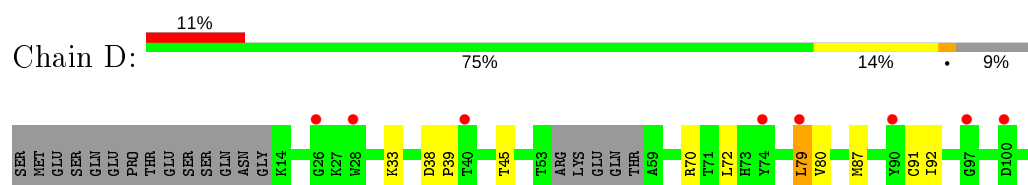
- Molecule 1: ADP-sugar pyrophosphatase

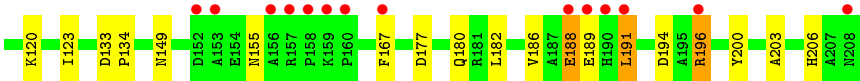


- Molecule 1: ADP-sugar pyrophosphatase



- 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	49.44Å 59.96Å 79.98Å 79.36° 81.22° 75.51°	Depositor
Resolution (Å)	78.11 – 1.63 50.04 – 1.63	Depositor EDS
% Data completeness (in resolution range)	96.3 (78.11-1.63) 96.3 (50.04-1.63)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.60 (at 1.63Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.231 , 0.266 0.245 , 0.278	Depositor DCC
R_{free} test set	5259 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	26.4	Xtriage
Anisotropy	0.217	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 41.1	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	6086	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.87% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, CL, EDO, K3J

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	0.98	0/1486	0.99	5/2027 (0.2%)
1	B	0.90	0/1492	1.01	6/2035 (0.3%)
1	C	0.99	1/1433 (0.1%)	1.01	4/1962 (0.2%)
1	D	0.92	0/1447	1.07	7/1976 (0.4%)
All	All	0.94	1/5858 (0.0%)	1.02	22/8000 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	131	CYS	CB-SG	-5.08	1.73	1.81

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	60	ASP	CB-CG-OD2	7.91	125.41	118.30
1	A	183	ASP	CB-CG-OD1	7.69	125.22	118.30
1	A	70	ARG	NE-CZ-NH2	-7.50	116.55	120.30
1	B	196	ARG	NE-CZ-NH1	7.38	123.99	120.30
1	D	196[A]	ARG	NE-CZ-NH1	7.33	123.97	120.30
1	D	196[B]	ARG	NE-CZ-NH1	7.33	123.97	120.30
1	C	60	ASP	CB-CG-OD1	-7.13	111.89	118.30
1	D	133	ASP	CB-CG-OD1	6.46	124.11	118.30
1	A	54	ARG	NE-CZ-NH2	-6.46	117.07	120.30
1	B	194	ASP	CB-CG-OD2	-6.29	112.64	118.30
1	D	196[A]	ARG	NE-CZ-NH2	-6.28	117.16	120.30
1	D	196[B]	ARG	NE-CZ-NH2	-6.28	117.16	120.30
1	D	91	CYS	CA-CB-SG	-6.08	103.06	114.00
1	A	54	ARG	NE-CZ-NH1	6.07	123.34	120.30
1	B	29	VAL	CB-CA-C	-6.02	99.96	111.40
1	C	38	ASP	CB-CG-OD1	5.86	123.57	118.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	152	ASP	CB-CG-OD2	-5.75	113.13	118.30
1	C	51	ARG	NE-CZ-NH1	5.68	123.14	120.30
1	B	72	LEU	CB-CG-CD1	-5.63	101.42	111.00
1	B	111	ARG	NE-CZ-NH1	5.42	123.01	120.30
1	D	87	MET	CG-SD-CE	5.29	108.67	100.20
1	B	70	ARG	NE-CZ-NH2	-5.24	117.68	120.30

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	1427	18	0
1	B	1464	0	1414	16	0
1	C	1406	0	1309	18	0
1	D	1419	0	1364	25	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	20	0	0	0	0
4	C	20	0	0	0	0
5	A	4	0	6	0	0
5	B	8	0	12	1	0
5	C	8	0	12	0	0
6	A	80	0	0	0	0
6	B	82	0	0	1	0
6	C	75	0	0	4	0
6	D	33	0	0	0	0
All	All	6086	0	5544	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:79:LEU:HD21	1:D:92:ILE:HG23	1.31	1.12
1:B:92:ILE:HD11	1:B:191:LEU:HD13	1.47	0.97
1:C:125:GLU:OE1	6:C:401:HOH:O	1.98	0.81
1:D:191:LEU:HD12	1:D:191:LEU:O	1.81	0.80
1:D:79:LEU:HD23	1:D:80:VAL:N	1.97	0.78
1:D:120:LYS:H	1:D:155:ASN:HD21	1.30	0.77
1:D:79:LEU:C	1:D:79:LEU:HD23	2.03	0.77
1:D:191:LEU:C	1:D:191:LEU:HD12	2.05	0.76
1:C:104:THR:HG23	1:C:107:ALA:H	1.50	0.74
1:D:79:LEU:CD2	1:D:92:ILE:HG23	2.15	0.71
1:A:191:LEU:C	1:A:191:LEU:HD12	2.14	0.68
1:A:203:ALA:HB3	1:B:203:ALA:HB3	1.77	0.66
1:A:206:HIS:HD2	1:B:200:TYR:OH	1.79	0.64
1:C:15:GLN:N	6:C:403:HOH:O	2.31	0.62
1:A:104:THR:HG23	1:A:107:ALA:H	1.65	0.61
1:C:206:HIS:HD2	1:D:200:TYR:OH	1.82	0.61
1:B:104:THR:HG23	1:B:107:ALA:H	1.65	0.60
1:D:79:LEU:HD11	1:D:92:ILE:CG2	2.32	0.59
1:C:200:TYR:CD1	1:D:203:ALA:HB2	2.41	0.55
1:D:191:LEU:C	1:D:191:LEU:CD1	2.73	0.55
1:A:71:THR:HG23	1:A:151:ASP:OD2	2.09	0.53
1:C:203:ALA:HB2	1:D:200:TYR:CD1	2.44	0.52
1:D:79:LEU:CD2	1:D:79:LEU:C	2.74	0.52
1:B:189:GLU:HB3	1:B:191:LEU:HD12	1.92	0.52
1:C:203:ALA:HB3	1:D:203:ALA:HB3	1.93	0.51
1:A:125:GLU:OE1	1:B:206:HIS:HE1	1.94	0.51
1:D:110:LEU:HG	1:D:123:ILE:HD11	1.93	0.50
1:A:104:THR:CG2	1:A:107:ALA:H	2.25	0.48
1:B:158:PRO:O	1:B:160:PRO:HD3	2.13	0.48
1:C:200:TYR:OH	1:D:206:HIS:HD2	1.95	0.48
6:C:401:HOH:O	1:D:206:HIS:HE1	1.96	0.48
1:A:200:TYR:CD1	1:B:203:ALA:HB2	2.49	0.47
1:B:54:ARG:NH2	1:B:60:ASP:OD2	2.39	0.47
1:C:32:GLU:O	1:C:47:GLU:HA	2.15	0.47
1:C:131:CYS:SG	1:D:194:ASP:HA	2.55	0.46
1:A:167:PHE:CD2	1:B:39:PRO:HG2	2.50	0.46
1:A:54:ARG:HD3	1:A:60:ASP:OD1	2.15	0.46
1:D:104:THR:HG23	1:D:107:ALA:H	1.81	0.46
1:C:179:LEU:HD23	1:C:205:LYS:HD2	1.97	0.45
1:B:14:LYS:HE3	1:B:39:PRO:HB2	1.97	0.45
1:A:200:TYR:OH	1:B:206:HIS:HD2	2.00	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:182:LEU:O	1:D:186:VAL:HG23	2.17	0.44
1:C:132:MET:O	1:D:196[B]:ARG:NH1	2.50	0.44
1:A:39:PRO:HG2	1:B:167:PHE:CD2	2.53	0.44
1:C:34:THR:N	1:C:46:TRP:O	2.51	0.44
5:B:301:EDO:C2	6:B:402:HOH:O	2.66	0.44
1:C:74:TYR:HB3	6:C:431:HOH:O	2.18	0.43
1:A:179:LEU:HD23	1:A:205:LYS:HD2	2.02	0.42
1:C:179:LEU:HD11	1:C:198:TYR:CZ	2.55	0.42
1:D:188:GLU:HG3	1:D:189:GLU:HG3	2.02	0.42
1:C:84:ARG:HG3	1:D:134:PRO:HB2	2.02	0.42
1:C:167:PHE:CD2	1:D:39:PRO:HG2	2.55	0.41
1:A:203:ALA:HB2	1:B:200:TYR:CD1	2.55	0.41
1:C:191:LEU:HD12	1:C:191:LEU:C	2.40	0.41
1:A:71:THR:O	1:A:73:HIS:HD2	2.04	0.41
1:A:132:MET:O	1:B:196:ARG:NH2	2.55	0.40
1:B:58:THR:CB	1:B:142:HIS:NE2	2.85	0.40
1:A:191:LEU:HD12	1:A:191:LEU:O	2.21	0.40
1:D:177:ASP:OD1	1:D:180:GLN:CB	2.68	0.40
1:A:191:LEU:CD1	1:A:191:LEU:C	2.85	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%)	188 (98%)	3 (2%)	1 (0%)	29	11
1	B	190/209 (91%)	184 (97%)	5 (3%)	1 (0%)	29	11
1	C	188/209 (90%)	182 (97%)	4 (2%)	2 (1%)	14	2
1	D	187/209 (90%)	180 (96%)	7 (4%)	0	100	100
All	All	757/836 (91%)	734 (97%)	19 (2%)	4 (0%)	29	11

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	24	SER
1	A	66	PRO
1	B	66	PRO
1	C	66	PRO

5.3.2 Protein sidechains ⓘ

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	29
1	B	154/179 (86%)	150 (97%)	4 (3%)	46	19
1	C	140/179 (78%)	136 (97%)	4 (3%)	42	15
1	D	145/179 (81%)	135 (93%)	10 (7%)	15	2
All	All	592/716 (83%)	571 (96%)	21 (4%)	35	10

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	LYS
1	A	33	LYS
1	A	38	ASP
1	B	29	VAL
1	B	31	LEU
1	B	38	ASP
1	B	66	PRO
1	C	27	LYS
1	C	31	LEU
1	C	44	ARG
1	C	164	ASP
1	D	33	LYS
1	D	38	ASP
1	D	45	THR
1	D	70	ARG
1	D	72	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	79	LEU
1	D	149	ASN
1	D	167	PHE
1	D	188	GLU
1	D	191	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	HIS
1	A	180	GLN
1	A	206	HIS
1	B	73	HIS
1	B	206	HIS
1	C	206	HIS
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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 9 are monoatomic - leaving 9 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$
5	EDO	C	306	-	3,3,3	0.48	0	2,2,2	0.62	0
5	EDO	B	301	-	3,3,3	0.58	0	2,2,2	0.12	0
5	EDO	A	306	-	3,3,3	0.64	0	2,2,2	0.36	0
4	K3J	A	305	-	10,10,10	2.04	2 (20%)	11,12,12	2.63	2 (18%)
4	K3J	C	305	-	10,10,10	2.58	4 (40%)	11,12,12	2.11	4 (36%)
5	EDO	C	301	-	3,3,3	0.55	0	2,2,2	0.19	0
4	K3J	A	304	-	10,10,10	1.06	1 (10%)	11,12,12	2.05	5 (45%)
5	EDO	B	304	-	3,3,3	0.54	0	2,2,2	0.49	0
4	K3J	C	304	-	10,10,10	1.58	3 (30%)	11,12,12	2.33	3 (27%)

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
5	EDO	C	306	-	-	0/1/1/1	-
5	EDO	B	301	-	-	1/1/1/1	-
5	EDO	A	306	-	-	0/1/1/1	-
4	K3J	A	305	-	-	0/3/3/3	0/1/1/1
4	K3J	C	305	-	-	0/3/3/3	0/1/1/1
5	EDO	C	301	-	-	1/1/1/1	-
4	K3J	A	304	-	-	0/3/3/3	0/1/1/1
5	EDO	B	304	-	-	0/1/1/1	-
4	K3J	C	304	-	-	0/3/3/3	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	305	K3J	C5-N1	6.29	1.40	1.33
4	A	305	K3J	C5-N1	4.76	1.38	1.33
4	C	304	K3J	C5-N1	3.27	1.37	1.33
4	A	305	K3J	C2-N2	3.24	1.43	1.33
4	C	305	K3J	C2-N2	2.62	1.41	1.33
4	C	305	K3J	N2-N1	-2.60	1.27	1.34
4	C	304	K3J	C2-N2	2.38	1.40	1.33
4	C	305	K3J	C3-C4	2.34	1.43	1.38
4	A	304	K3J	C2-N2	2.07	1.39	1.33

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	304	K3J	C3-C2	2.06	1.44	1.39

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	305	K3J	C6-C5-N1	6.61	119.36	116.24
4	C	304	K3J	C5-N1-N2	5.99	123.70	120.03
4	A	305	K3J	C5-N1-N2	4.81	122.97	120.03
4	C	305	K3J	C5-N1-N2	4.34	122.69	120.03
4	A	304	K3J	C2-N2-N1	3.38	121.77	119.37
4	A	304	K3J	C5-N1-N2	3.26	122.03	120.03
4	C	304	K3J	C6-C5-N1	3.14	117.72	116.24
4	A	304	K3J	C6-C5-N1	3.03	117.67	116.24
4	C	305	K3J	C2-N2-N1	3.00	121.50	119.37
4	C	304	K3J	C4-C5-N1	-2.53	119.65	121.94
4	C	305	K3J	C3-C2-N	2.43	126.26	121.04
4	A	304	K3J	C4-C5-N1	-2.32	119.84	121.94
4	A	304	K3J	C3-C2-N2	-2.23	118.28	122.56
4	C	305	K3J	C4-C5-N1	-2.10	120.04	121.94

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	301	EDO	O1-C1-C2-O2
5	C	301	EDO	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	301	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%)	-0.18	8 (4%) 37 35	19, 29, 51, 60	9 (4%)
1	B	194/209 (92%)	-0.28	3 (1%) 73 74	19, 33, 59, 80	2 (1%)
1	C	192/209 (91%)	-0.03	8 (4%) 36 33	19, 35, 63, 71	5 (2%)
1	D	190/209 (90%)	0.70	22 (11%) 4 3	23, 41, 65, 83	16 (8%)
All	All	770/836 (92%)	0.05	41 (5%) 26 24	19, 34, 61, 83	32 (4%)

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	191	LEU	8.1
1	D	79	LEU	8.1
1	D	160	PRO	7.4
1	A	191	LEU	7.2
1	D	156	ALA	6.6
1	D	158	PRO	6.4
1	D	153	ALA	6.3
1	D	157	ARG	4.9
1	D	190	HIS	4.9
1	D	159	LYS	4.7
1	B	58	THR	4.3
1	C	40	THR	4.2
1	A	190	HIS	3.9
1	C	46	TRP	3.8
1	C	44	ARG	3.8
1	D	189	GLU	3.7
1	A	189	GLU	3.6
1	D	208	ASN	3.4
1	D	188	GLU	3.3
1	A	188	GLU	3.3
1	D	90	TYR	3.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	43	THR	3.2
1	D	26	GLY	2.9
1	A	167	PHE	2.9
1	A	57	GLN	2.9
1	D	40	THR	2.8
1	C	39	PRO	2.8
1	D	100	ASP	2.7
1	D	97	GLY	2.6
1	C	48	SER	2.6
1	D	196[A]	ARG	2.5
1	C	16	TYR	2.5
1	A	163	GLY	2.5
1	B	102	GLY	2.4
1	D	167	PHE	2.3
1	A	164	ASP	2.2
1	B	165	GLY	2.2
1	D	28	TRP	2.2
1	D	152	ASP	2.1
1	C	22	LEU	2.1
1	D	74	TYR	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 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. 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	K3J	C	305	10/10	0.75	0.17	37,40,45,45	10
2	MG	A	302	1/1	0.82	0.12	64,64,64,64	0
2	MG	C	302	1/1	0.87	0.12	69,69,69,69	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	CL	A	303	1/1	0.88	0.08	65,65,65,65	0
5	EDO	C	306	4/4	0.89	0.08	41,43,46,47	0
5	EDO	C	301	4/4	0.89	0.12	35,49,56,62	0
2	MG	D	301	1/1	0.90	0.08	39,39,39,39	0
5	EDO	B	304	4/4	0.90	0.08	30,30,30,31	4
4	K3J	A	305	10/10	0.91	0.11	30,35,36,37	10
5	EDO	B	301	4/4	0.93	0.13	40,45,48,49	0
2	MG	B	303	1/1	0.93	0.14	75,75,75,75	0
4	K3J	C	304	10/10	0.94	0.10	39,41,44,48	0
5	EDO	A	306	4/4	0.94	0.06	25,26,26,26	0
4	K3J	A	304	10/10	0.96	0.09	27,28,30,30	10
2	MG	D	302	1/1	0.97	0.07	39,39,39,39	0
2	MG	C	303	1/1	0.97	0.05	33,33,33,33	0
2	MG	B	302	1/1	0.99	0.03	27,27,27,27	0
2	MG	A	301	1/1	1.00	0.06	29,29,29,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.