



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 09:00 PM UTC

PDB ID : 3VUS / pdb_00003vus
Title : Escherichia coli PgaB N-terminal domain
Authors : Nishiyama, T.; Noguchi, H.; Yoshida, H.; Park, S.-Y.; Tame, J.R.H.
Deposited on : 2012-07-05
Resolution : 1.65 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

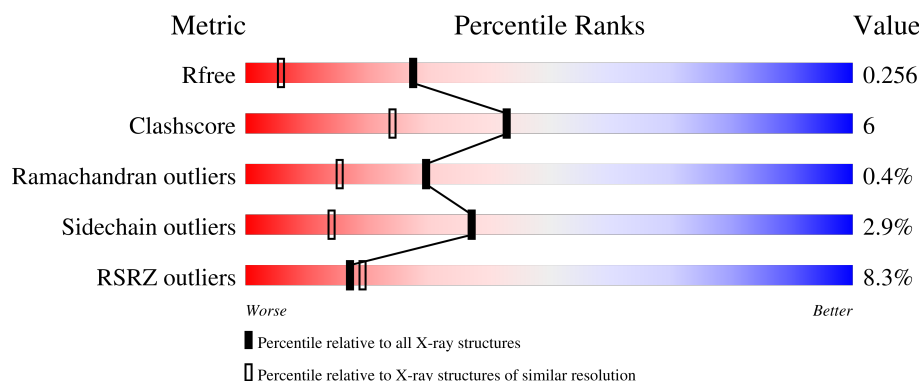
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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	2563 (1.66-1.66)
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)
RSRZ outliers	180081	2564 (1.66-1.66)

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	268	6% 78% 19% ..
1	B	268	10% 69% 22% ..

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
3	ACT	B	402	-	X	-	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4445 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.

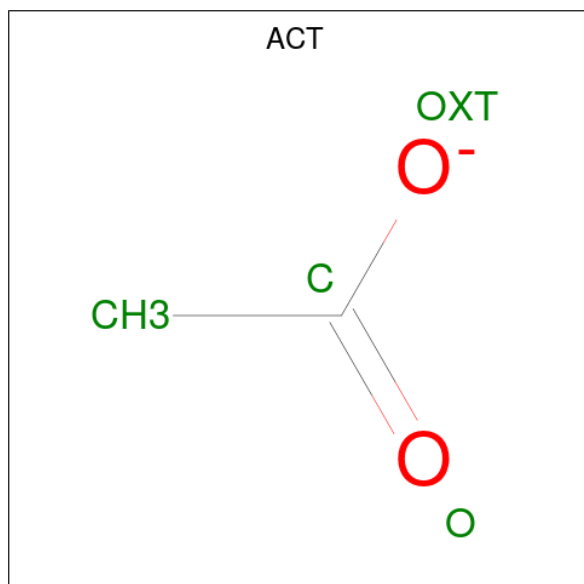
- Molecule 1 is a protein called Poly-beta-1,6-N-acetyl-D-glucosamine N-deacetylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	0	0	0
			2156	1386	376	391	3			
1	B	256	Total	C	N	O	S	0	0	0
			2103	1357	366	377	3			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

- Molecule 3 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



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

- Molecule 4 is MERCURY (II) ION (CCD ID: HG) (formula: Hg).

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

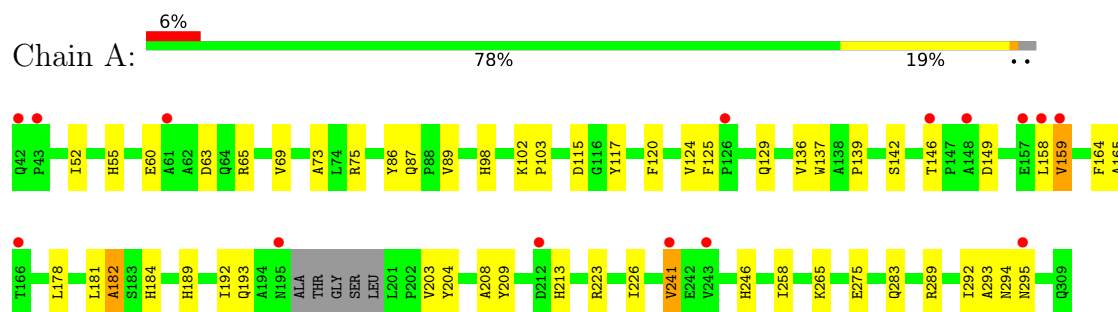
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	91	Total O 91 91	0	0
5	B	84	Total O 84 84	0	0

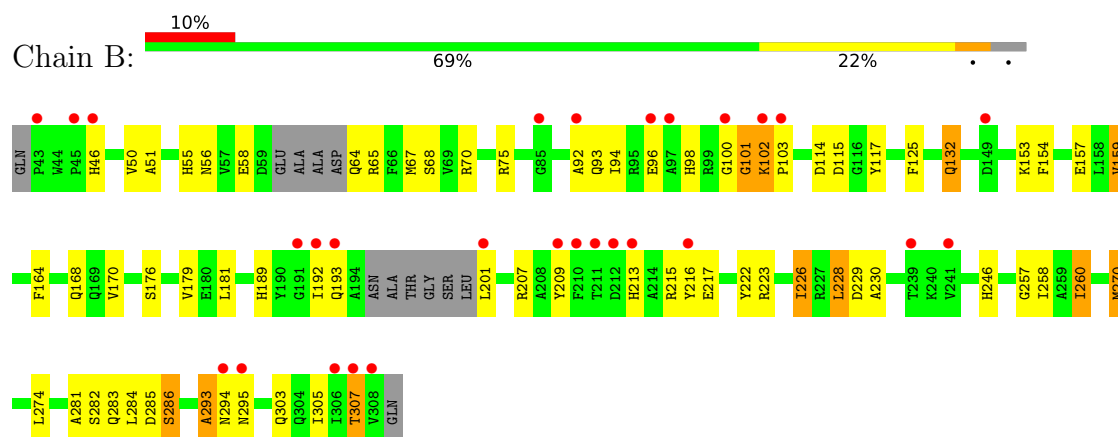
3 Residue-property plots [i](#)

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: Poly-beta-1,6-N-acetyl-D-glucosamine N-deacetylase



- Molecule 1: Poly-beta-1,6-N-acetyl-D-glucosamine N-deacetylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	39.56Å 53.11Å 144.17Å 90.00° 95.25° 90.00°	Depositor
Resolution (Å)	25.00 – 1.65 25.00 – 1.65	Depositor EDS
% Data completeness (in resolution range)	96.6 (25.00-1.65) 96.6 (25.00-1.65)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.75 (at 1.65Å)	Xtriage
Refinement program	REFMAC 5.6.0119	Depositor
R, R_{free}	0.205 , 0.256 0.204 , 0.256	Depositor DCC
R_{free} test set	996 reflections (1.39%)	wwPDB-VP
Wilson B-factor (Å ²)	25.9	Xtriage
Anisotropy	0.109	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 43.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.020 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4445	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.55% 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: HG, ZN, ACT

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	1.64	21/2219 (0.9%)	1.46	18/3023 (0.6%)
1	B	1.62	17/2165 (0.8%)	1.47	17/2947 (0.6%)
All	All	1.63	38/4384 (0.9%)	1.47	35/5970 (0.6%)

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	295	ASN	CA-C	8.69	1.60	1.53
1	A	295	ASN	CA-CB	8.14	1.58	1.52
1	B	55	HIS	N-CA	7.89	1.51	1.46
1	B	189	HIS	ND1-CE1	7.52	1.40	1.32
1	A	52	ILE	C-O	7.28	1.31	1.24
1	A	295	ASN	C-O	-7.07	1.20	1.23
1	A	184	HIS	CG-CD2	6.93	1.43	1.35
1	B	226	ILE	N-CA	6.84	1.54	1.46
1	A	246	HIS	CG-CD2	6.66	1.43	1.35
1	A	209	TYR	CA-C	-6.44	1.45	1.52
1	A	182	ALA	CA-C	-6.24	1.44	1.52
1	A	117	TYR	N-CA	6.24	1.53	1.46
1	B	189	HIS	CE1-NE2	6.16	1.38	1.32
1	B	229	ASP	C-O	6.07	1.31	1.24
1	B	68	SER	CA-C	-5.99	1.45	1.52
1	B	179	VAL	N-CA	5.93	1.53	1.46
1	A	136	VAL	CA-CB	5.92	1.62	1.54
1	A	75	ARG	C-O	5.83	1.30	1.24
1	B	230	ALA	N-CA	5.80	1.53	1.46
1	A	289	ARG	N-CA	5.63	1.53	1.46
1	A	178	LEU	CA-C	5.62	1.60	1.52
1	A	265	LYS	N-CA	5.51	1.53	1.46
1	B	270	MET	SD-CE	-5.41	1.66	1.79
1	B	293	ALA	CA-C	-5.36	1.46	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	125	PHE	CA-CB	5.35	1.61	1.53
1	A	120	PHE	C-O	-5.29	1.18	1.24
1	B	246	HIS	CG-ND1	-5.27	1.32	1.38
1	A	86	TYR	C-N	5.26	1.40	1.33
1	B	114	ASP	CG-OD2	5.21	1.35	1.25
1	A	89	VAL	N-CA	5.18	1.51	1.46
1	B	164	PHE	C-O	5.17	1.30	1.23
1	A	86	TYR	CA-CB	5.15	1.61	1.53
1	A	137	TRP	CD2-CE2	5.12	1.50	1.41
1	B	46	HIS	CG-CD2	5.11	1.41	1.35
1	B	246	HIS	C-N	5.08	1.39	1.33
1	B	216	TYR	CA-C	-5.03	1.46	1.52
1	A	125	PHE	N-CA	5.03	1.51	1.46
1	A	189	HIS	ND1-CE1	5.01	1.37	1.32

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	125	PHE	CA-C-N	-8.06	110.64	119.19
1	A	125	PHE	C-N-CA	-8.06	110.64	119.19
1	A	124	VAL	N-CA-C	-7.45	104.52	111.45
1	A	295	ASN	CB-CA-C	6.80	116.92	111.00
1	A	73	ALA	O-C-N	6.67	129.76	122.15
1	A	73	ALA	CA-C-O	-6.64	113.38	120.42
1	A	55	HIS	N-CA-CB	-6.53	105.52	111.59
1	B	50	VAL	CB-CA-C	-6.51	101.80	110.98
1	A	129	GLN	N-CA-C	-6.37	104.26	111.07
1	A	75	ARG	NE-CZ-NH1	6.28	127.78	121.50
1	B	94	ILE	CA-C-O	-6.24	114.12	121.05
1	A	98	HIS	O-C-N	5.97	128.22	122.07
1	B	201	LEU	CA-C-N	-5.95	113.40	119.83
1	B	201	LEU	C-N-CA	-5.95	113.40	119.83
1	B	257	GLY	N-CA-C	5.94	119.86	112.73
1	B	125	PHE	O-C-N	5.87	126.51	120.70
1	B	93	GLN	N-CA-C	5.81	117.61	111.28
1	B	260	ILE	CB-CA-C	5.80	119.40	111.97
1	B	228	LEU	N-CA-C	-5.74	105.11	111.36
1	A	241	VAL	N-CA-C	-5.52	107.89	113.47
1	A	295	ASN	CA-C-N	5.45	125.39	119.78
1	A	295	ASN	C-N-CA	5.45	125.39	119.78
1	A	87	GLN	CA-C-O	-5.43	114.52	119.62
1	B	286	SER	CA-C-N	-5.35	117.45	123.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	286	SER	C-N-CA	-5.35	117.45	123.02
1	A	165	ALA	N-CA-C	5.32	117.20	110.33
1	B	94	ILE	O-C-N	5.32	127.43	121.90
1	A	292	ILE	CA-C-N	-5.31	114.34	122.62
1	A	292	ILE	C-N-CA	-5.31	114.34	122.62
1	B	222	TYR	N-CA-C	-5.26	105.46	111.14
1	B	176	SER	N-CA-C	5.18	117.67	111.71
1	B	51	ALA	N-CA-C	-5.10	100.86	109.07
1	B	305	ILE	N-CA-C	5.10	115.87	110.72
1	B	132	GLN	N-CA-C	-5.08	104.38	111.39
1	A	69	VAL	N-CA-C	-5.03	100.66	108.86

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	2156	0	2075	18	0
1	B	2103	0	2033	35	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	4	0	3	0	0
3	B	4	0	3	1	0
4	B	1	0	0	0	0
5	A	91	0	0	5	0
5	B	84	0	0	4	0
All	All	4445	0	4114	52	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 6.

All (52) 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:132:GLN:HG3	5:B:545:HOH:O	1.59	1.01
1:B:192:ILE:HG21	5:B:544:HOH:O	1.72	0.89
1:B:283:GLN:HE22	1:B:286:SER:H	1.24	0.81
1:A:213:HIS:HE1	5:A:540:HOH:O	1.68	0.76
1:B:303:GLN:O	1:B:307:THR:HB	1.86	0.76
1:A:181:LEU:HD13	1:A:241:VAL:HG11	1.67	0.75
1:A:275:GLU:HG3	5:A:535:HOH:O	1.86	0.74
1:A:181:LEU:CD1	1:A:241:VAL:HG11	2.17	0.74
1:B:283:GLN:HE22	1:B:286:SER:N	1.89	0.70
1:A:146:THR:HG21	5:A:527:HOH:O	1.94	0.66
1:B:75:ARG:HD3	5:B:570:HOH:O	1.95	0.65
1:A:60:GLU:CD	1:B:168:GLN:HG3	2.26	0.60
1:B:100:GLY:O	1:B:101:GLY:O	2.19	0.60
1:B:56:ASN:ND2	1:B:58:GLU:CG	2.65	0.60
1:B:270:MET:HG2	1:B:284:LEU:HD13	1.85	0.58
1:A:146:THR:CG2	5:A:527:HOH:O	2.52	0.56
1:B:56:ASN:ND2	1:B:58:GLU:HG3	2.22	0.54
1:B:64:GLN:HA	1:B:64:GLN:OE1	2.04	0.54
1:A:213:HIS:CE1	5:A:540:HOH:O	2.50	0.54
1:B:65:ARG:NH1	1:B:154:PHE:O	2.41	0.53
1:B:58:GLU:O	1:B:70:ARG:HA	2.09	0.53
1:B:98:HIS:ND1	1:B:281:ALA:O	2.41	0.52
1:B:213:HIS:CE1	5:B:582:HOH:O	2.63	0.52
1:A:192:ILE:HD13	1:A:203:VAL:HG23	1.92	0.52
1:B:70:ARG:HG3	1:B:295:ASN:HD21	1.74	0.51
1:B:170:VAL:HG13	1:B:181:LEU:HD11	1.94	0.49
1:B:223:ARG:NH2	1:B:258:ILE:HD12	2.27	0.49
1:A:204:TYR:CE1	1:A:226:ILE:HG13	2.48	0.49
1:B:274:LEU:HD11	3:B:402:ACT:H3	1.95	0.48
1:B:283:GLN:NE2	1:B:285:ASP:H	2.12	0.48
1:B:92:ALA:O	1:B:96:GLU:HG3	2.14	0.47
1:B:65:ARG:HB3	1:B:154:PHE:CE2	2.49	0.47
1:B:56:ASN:HD21	1:B:58:GLU:CD	2.23	0.47
1:A:293:ALA:O	1:A:294:ASN:HB2	2.14	0.46
1:A:142:SER:O	1:A:146:THR:HG23	2.16	0.45
1:B:228:LEU:HD23	1:B:228:LEU:HA	1.81	0.45
1:A:63:ASP:OD2	1:A:65:ARG:NH2	2.50	0.44
1:A:192:ILE:HD12	1:A:208:ALA:HA	2.00	0.44
1:B:102:LYS:HA	1:B:103:PRO:HD2	1.85	0.44
1:B:65:ARG:HB3	1:B:154:PHE:HE2	1.83	0.43
1:B:193:GLN:HG2	1:B:209:TYR:O	2.17	0.43
1:A:159:VAL:HG11	1:A:164:PHE:HE2	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:157:GLU:HG2	1:B:159:VAL:HG23	2.00	0.42
1:B:115:ASP:HB2	1:B:117:TYR:CE2	2.54	0.42
1:B:283:GLN:NE2	1:B:285:ASP:N	2.68	0.42
1:B:293:ALA:O	1:B:294:ASN:HB2	2.20	0.42
1:A:223:ARG:NH2	1:A:258:ILE:HD12	2.34	0.41
1:B:207:ARG:HD2	1:B:217:GLU:O	2.20	0.41
1:B:283:GLN:HE21	1:B:285:ASP:H	1.66	0.41
1:A:102:LYS:HA	1:A:103:PRO:HD3	1.97	0.41
1:B:98:HIS:HE1	1:B:282:SER:O	2.03	0.41
1:A:139:PRO:HD2	1:A:182:ALA: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	259/268 (97%)	249 (96%)	9 (4%)	1 (0%)	30	15
1	B	250/268 (93%)	236 (94%)	13 (5%)	1 (0%)	30	15
All	All	509/536 (95%)	485 (95%)	22 (4%)	2 (0%)	30	15

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	101	GLY
1	A	115	ASP

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	224/227 (99%)	219 (98%)	5 (2%)	45	23
1	B	219/227 (96%)	211 (96%)	8 (4%)	30	8
All	All	443/454 (98%)	430 (97%)	13 (3%)	37	14

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

Mol	Chain	Res	Type
1	A	149	ASP
1	A	158	LEU
1	A	159	VAL
1	A	193	GLN
1	A	283	GLN
1	B	67	MET
1	B	102	LYS
1	B	153	LYS
1	B	159	VAL
1	B	215	ARG
1	B	226	ILE
1	B	260	ILE
1	B	307	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	151	GLN
1	B	56	ASN
1	B	132	GLN
1	B	151	GLN
1	B	213	HIS
1	B	283	GLN
1	B	295	ASN

5.3.3 RNA ⓘ

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 5 ligands modelled in this entry, 3 are monoatomic - leaving 2 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	ACT	B	402	2	3,3,3	1.99	1 (33%)	3,3,3	3.86	2 (66%)
3	ACT	A	402	2	3,3,3	1.73	1 (33%)	3,3,3	0.18	0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	ACT	CH3-C	3.03	1.61	1.49
3	A	402	ACT	CH3-C	2.02	1.57	1.49

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	ACT	OXT-C-CH3	4.96	135.87	115.05
3	B	402	ACT	OXT-C-O	-4.23	106.33	122.03

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
3	B	402	ACT	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	263/268 (98%)	0.49	15 (5%)	29	33	17, 29, 57, 101	0
1	B	256/268 (95%)	0.62	28 (10%)	10	11	17, 31, 57, 79	0
All	All	519/536 (96%)	0.56	43 (8%)	17	19	17, 30, 58, 101	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	61	ALA	5.5
1	B	43	PRO	4.4
1	B	210	PHE	3.5
1	B	192	ILE	3.3
1	B	45	PRO	3.3
1	A	148	ALA	3.3
1	B	209	TYR	3.2
1	B	191	GLY	3.1
1	B	46	HIS	2.9
1	B	201	LEU	2.8
1	A	146	THR	2.7
1	B	295	ASN	2.7
1	B	307	THR	2.7
1	B	308	VAL	2.7
1	A	195	ASN	2.6
1	A	212	ASP	2.6
1	A	157	GLU	2.6
1	B	212	ASP	2.6
1	B	97	ALA	2.6
1	B	213	HIS	2.4
1	A	43	PRO	2.4
1	A	158	LEU	2.4
1	B	92	ALA	2.4
1	B	96	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	102	LYS	2.4
1	B	193	GLN	2.3
1	A	42	GLN	2.3
1	B	216	TYR	2.3
1	B	241	VAL	2.3
1	A	159	VAL	2.2
1	B	211	THR	2.2
1	B	100	GLY	2.2
1	B	103	PRO	2.2
1	B	294	ASN	2.2
1	A	126	PRO	2.2
1	A	295	ASN	2.1
1	B	149	ASP	2.1
1	B	306	ILE	2.1
1	B	85	GLY	2.1
1	A	243	VAL	2.1
1	A	166	THR	2.0
1	B	239	THR	2.0
1	A	241	VAL	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
3	ACT	A	402	4/4	0.86	0.17	21,30,30,38	0
3	ACT	B	402	4/4	0.86	0.18	14,16,24,25	4
4	HG	B	403	1/1	0.93	0.07	33,33,33,33	1
2	ZN	B	401	1/1	0.98	0.05	24,24,24,24	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	A	401	1/1	1.00	0.06	26,26,26,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.