



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 13, 2024 – 06:44 AM EST

PDB ID : 4KGH
Title : Crystal Structure of human splunc1 lacking the secretion signal sequence
Authors : Betts, L.; Walton, W.G.
Deposited on : 2013-04-29
Resolution : 2.81 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 : 4.02b-467
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 1.20.1
EDS : 3.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.003 (Gargrove)
Density-Fitness : 1.0.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

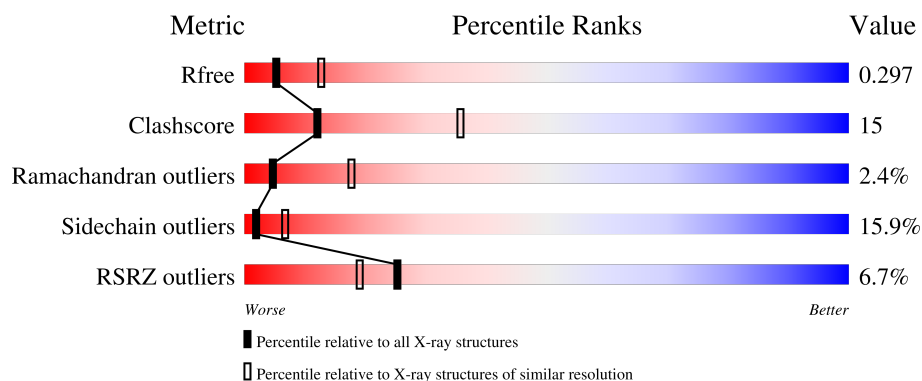
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.81 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}	164625	3657 (2.80-2.80)
Clashscore	180529	4123 (2.80-2.80)
Ramachandran outliers	177936	4071 (2.80-2.80)
Sidechain outliers	177891	4073 (2.80-2.80)
RSRZ outliers	164620	3659 (2.80-2.80)

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	240	<div> <div>5%</div> <div> <div></div> <div>50%</div> <div>24%</div> <div>6%</div> <div>20%</div> </div> </div>
1	B	240	<div> <div>6%</div> <div> <div></div> <div>54%</div> <div>25%</div> <div>•</div> <div>18%</div> </div> </div>

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 2781 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.

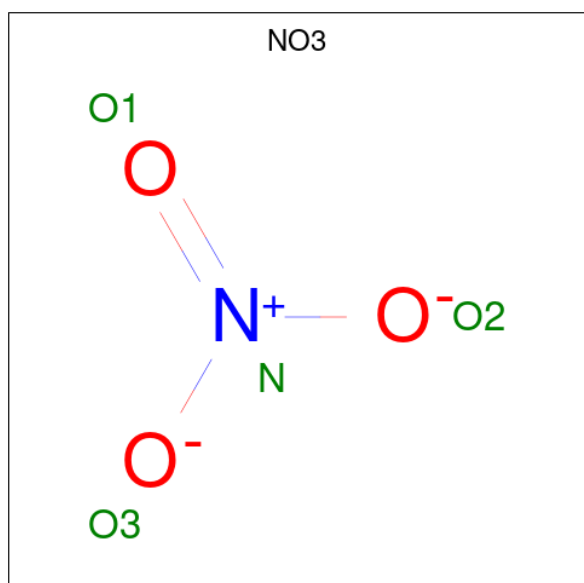
- Molecule 1 is a protein called BPI fold-containing family A member 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	197	Total	C	N	O	S	0	0	0
			1385	892	231	259	3			
1	A	193	Total	C	N	O	S	0	1	0
			1376	879	235	259	3			

There are 6 discrepancies between the modelled and reference sequences:

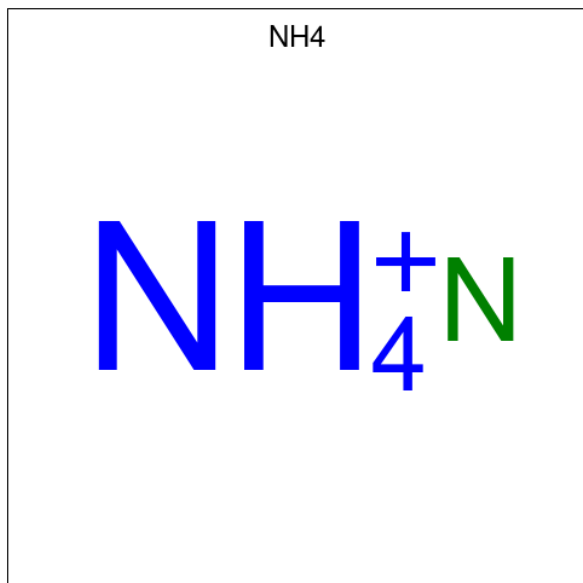
Chain	Residue	Modelled	Actual	Comment	Reference
B	17	SER	-	expression tag	UNP Q9NP55
B	18	ASN	-	expression tag	UNP Q9NP55
B	105	VAL	ILE	SEE REMARK 999	UNP Q9NP55
A	17	SER	-	expression tag	UNP Q9NP55
A	18	ASN	-	expression tag	UNP Q9NP55
A	105	VAL	ILE	SEE REMARK 999	UNP Q9NP55

- Molecule 2 is NITRATE ION (three-letter code: NO3) (formula: NO₃).



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

- Molecule 3 is AMMONIUM ION (three-letter code: NH4) (formula: H_4N).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	N	0	0
			1	1		

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

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

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $\text{C}_3\text{H}_8\text{O}_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			6	3	3		

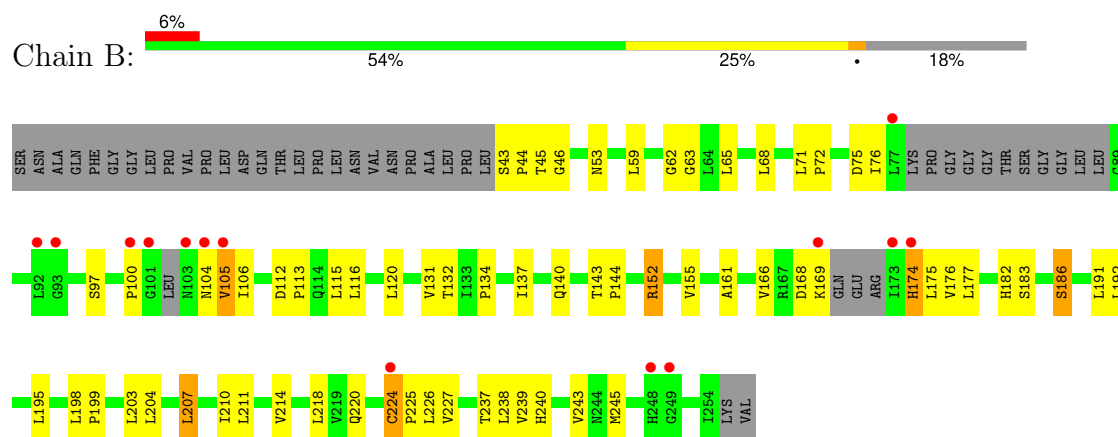
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	5	Total	O	0	0
			5	5		
6	A	3	Total	O	0	0
			3	3		

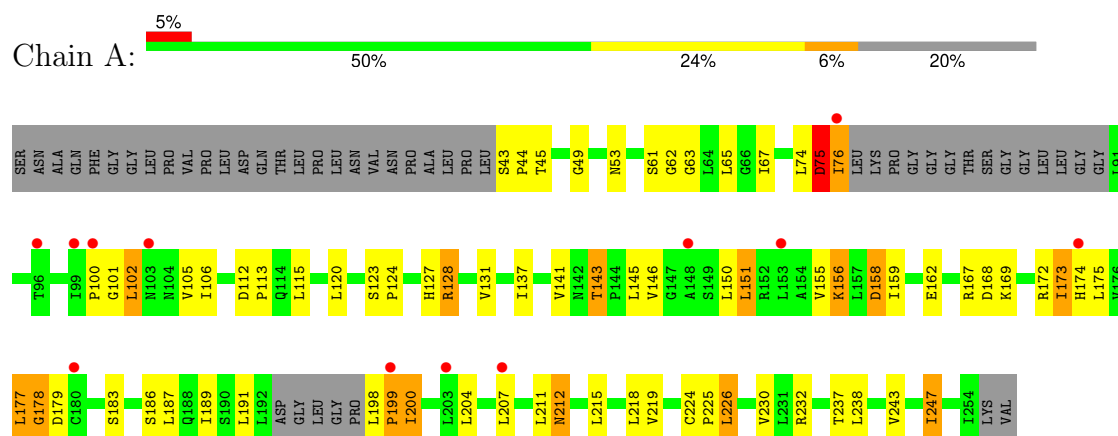
3 Residue-property plots [i](#)

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: BPI fold-containing family A member 1



- Molecule 1: BPI fold-containing family A member 1



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	47.48Å 203.73Å 118.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.64 – 2.81 38.64 – 2.81	Depositor EDS
% Data completeness (in resolution range)	94.4 (38.64-2.81) 90.8 (38.64-2.81)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.65 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.8.1_1168	Depositor
R, R_{free}	0.246 , 0.295 0.251 , 0.297	Depositor DCC
R_{free} test set	1380 reflections (9.95%)	wwPDB-VP
Wilson B-factor (Å ²)	66.7	Xtriage
Anisotropy	0.691	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 72.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	2781	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.75% 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: NO3, NH4, NA, GOL

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	0.22	0/1394	0.45	0/1907
1	B	0.29	1/1399 (0.1%)	0.48	1/1913 (0.1%)
All	All	0.26	1/2793 (0.0%)	0.47	1/3820 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	225	PRO	N-CD	5.22	1.55	1.47

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	224	CYS	C-N-CD	5.71	140.38	128.40

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1376	0	1391	50	1
1	B	1385	0	1421	36	0
2	B	4	0	0	0	0
3	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	1	0	0	0	0
5	B	6	0	8	2	0
6	A	3	0	0	1	0
6	B	5	0	0	0	0
All	All	2781	0	2820	86	1

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 15.

All (86) 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:A:159:ILE:HD11	1:A:219:VAL:HG12	1.62	0.81
1:A:49:GLY:O	1:A:53:ASN:ND2	2.14	0.80
1:B:183:SER:O	1:B:186:SER:OG	2.01	0.78
1:B:239:VAL:O	1:B:243:VAL:HG23	1.88	0.73
1:A:183:SER:O	1:A:186:SER:OG	2.02	0.72
1:A:62:GLY:N	1:A:63:GLY:HA2	2.05	0.71
1:B:220:GLN:HG3	1:B:224:CYS:SG	2.32	0.69
1:A:101:GLY:O	1:A:102:LEU:CB	2.43	0.67
1:B:220:GLN:O	1:B:224:CYS:HB2	1.96	0.65
1:A:106:ILE:HG22	1:A:143:THR:HG23	1.78	0.64
1:B:62:GLY:N	1:B:63:GLY:HA2	2.12	0.64
1:A:74:LEU:O	1:A:76:ILE:N	2.31	0.64
1:A:173:ILE:HG21	1:A:243:VAL:CG1	2.28	0.64
1:B:182:HIS:ND1	5:B:304:GOL:O1	2.31	0.62
1:A:158:ASP:HB2	1:A:186:SER:HB3	1.82	0.61
1:B:220:GLN:O	1:B:224:CYS:SG	2.59	0.60
1:A:168:ASP:OD1	1:A:169:LYS:N	2.35	0.60
1:B:220:GLN:O	1:B:224:CYS:CB	2.51	0.58
1:A:173:ILE:CG2	1:A:243:VAL:CG1	2.82	0.57
1:A:75:ASP:O	1:A:76:ILE:HG13	2.05	0.57
1:B:43:SER:N	1:B:44:PRO:HD2	2.20	0.57
1:A:128:ARG:N	1:A:128:ARG:HD2	2.21	0.56
1:B:43:SER:N	1:B:44:PRO:CD	2.71	0.54
1:B:68:LEU:HA	1:B:71:LEU:HD11	1.89	0.54
1:A:173:ILE:HG22	1:A:243:VAL:HG11	1.89	0.53
1:B:105:VAL:HG12	1:B:106:ILE:H	1.74	0.52
1:A:143:THR:O	1:A:146:VAL:O	2.28	0.52
1:A:173:ILE:CG2	1:A:243:VAL:HG11	2.41	0.51
1:B:143:THR:O	1:B:143:THR:HG23	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:105:VAL:O	1:B:144:PRO:HD2	2.11	0.51
1:B:131:VAL:HG13	1:B:161:ALA:HB3	1.93	0.50
1:B:166:VAL:O	1:B:174:HIS:ND1	2.44	0.49
1:B:116:LEU:HB2	1:B:134:PRO:HB2	1.94	0.49
1:A:167:ARG:HG3	1:A:173:ILE:CD1	2.42	0.49
1:B:59:LEU:HD21	1:B:227:VAL:HG12	1.95	0.49
1:A:123:SER:CB	1:A:128:ARG:HG2	2.42	0.48
1:B:175:LEU:HB2	1:B:239:VAL:HG13	1.96	0.48
1:B:105:VAL:HG12	1:B:106:ILE:N	2.29	0.48
1:A:224:CYS:N	1:A:225:PRO:CD	2.76	0.48
1:A:43:SER:N	1:A:44:PRO:CD	2.76	0.48
1:A:177:LEU:O	1:A:178:GLY:C	2.51	0.47
1:B:198:LEU:HB3	1:B:199:PRO:HD3	1.97	0.47
1:A:167:ARG:HG3	1:A:173:ILE:HD11	1.96	0.47
1:B:182:HIS:ND1	1:B:182:HIS:O	2.48	0.47
1:B:176:VAL:HG12	1:B:177:LEU:N	2.30	0.46
1:A:123:SER:HB3	1:A:128:ARG:HG2	1.97	0.46
1:A:173:ILE:CD1	1:A:247:ILE:HG13	2.46	0.46
1:B:112:ASP:N	1:B:113:PRO:CD	2.79	0.46
1:B:140:GLN:OE1	1:B:152:ARG:HG3	2.15	0.45
1:A:105:VAL:O	1:A:143:THR:HG22	2.16	0.45
1:A:156:LYS:NZ	6:A:301:HOH:O	2.49	0.45
1:A:112:ASP:N	1:A:113:PRO:CD	2.79	0.45
1:A:178:GLY:HA2	1:A:179:ASP:HA	1.68	0.45
1:B:143:THR:O	1:B:143:THR:CG2	2.65	0.45
1:B:72:PRO:HB2	1:B:75:ASP:HB2	1.99	0.45
1:A:183:SER:HB3	1:A:186:SER:OG	2.16	0.45
1:A:198:LEU:N	1:A:199:PRO:HD2	2.32	0.45
1:A:226:LEU:HD22	1:A:230:VAL:HG23	1.99	0.45
1:A:123:SER:HB2	1:A:124:PRO:HD2	2.00	0.44
1:A:162:GLU:O	1:A:177:LEU:O	2.35	0.44
1:A:187:LEU:CD2	1:A:212:ASN:HA	2.47	0.44
1:B:182:HIS:CE1	5:B:304:GOL:HO1	2.34	0.44
1:B:45:THR:OG1	1:B:46:GLY:N	2.47	0.44
1:A:75:ASP:OD1	1:A:75:ASP:N	2.49	0.44
1:B:97:SER:O	1:B:100:PRO:CB	2.67	0.43
1:A:191:LEU:C	1:A:191:LEU:HD13	2.39	0.43
1:A:151:LEU:HD22	1:A:204:LEU:CD1	2.48	0.43
1:A:175:LEU:HB2	1:A:243:VAL:CG2	2.49	0.42
1:A:151:LEU:CD2	1:A:204:LEU:CD1	2.96	0.42
1:A:75:ASP:O	1:A:76:ILE:CG1	2.65	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:VAL:HG23	1:A:106:ILE:H	1.83	0.42
1:B:168:ASP:O	1:B:169:LYS:CB	2.67	0.41
1:A:137:ILE:HG13	1:A:155:VAL:HB	2.02	0.41
1:B:76:ILE:HG23	1:B:214:VAL:HG11	2.02	0.41
1:B:210:ILE:O	1:B:214:VAL:HG23	2.21	0.41
1:A:141:VAL:HG21	1:A:207:LEU:HD11	2.01	0.41
1:B:137:ILE:HG13	1:B:155:VAL:HB	2.03	0.41
1:A:168:ASP:HB3	1:A:172:ARG:O	2.20	0.41
1:B:204:LEU:O	1:B:207:LEU:N	2.54	0.41
1:A:141:VAL:HG21	1:A:151:LEU:HD11	2.03	0.41
1:B:140:GLN:OE1	1:B:152:ARG:CG	2.69	0.41
1:A:199:PRO:O	1:A:200:ILE:CB	2.69	0.40
1:A:127:HIS:C	1:A:128:ARG:HD2	2.42	0.40
1:A:198:LEU:N	1:A:199:PRO:CD	2.85	0.40
1:A:43:SER:N	1:A:44:PRO:HD2	2.36	0.40
1:A:63:GLY:O	1:A:67:ILE:HD13	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:ARG:NH2	1:A:168:ASP:OD1[4_555]	2.09	0.11

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	188/240 (78%)	170 (90%)	11 (6%)	7 (4%)	2	9
1	B	189/240 (79%)	177 (94%)	10 (5%)	2 (1%)	12	37
All	All	377/480 (78%)	347 (92%)	21 (6%)	9 (2%)	5	18

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	104	ASN
1	B	105	VAL
1	A	75	ASP
1	A	100	PRO
1	A	102	LEU
1	A	178	GLY
1	A	199	PRO
1	A	200	ILE
1	A	145	LEU

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	148/202 (73%)	120 (81%)	28 (19%)	1	4
1	B	149/202 (74%)	129 (87%)	20 (13%)	3	10
All	All	297/404 (74%)	249 (84%)	48 (16%)	2	6

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	53	ASN
1	B	65	LEU
1	B	115	LEU
1	B	120	LEU
1	B	132	THR
1	B	152	ARG
1	B	174	HIS
1	B	186	SER
1	B	191	LEU
1	B	192	LEU
1	B	195	LEU
1	B	203	LEU
1	B	207	LEU
1	B	211	LEU

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Mol	Chain	Res	Type
1	B	218	LEU
1	B	226	LEU
1	B	237	THR
1	B	238	LEU
1	B	240	HIS
1	B	245	MET
1	A	45	THR
1	A	61	SER
1	A	65	LEU
1	A	75	ASP
1	A	76	ILE
1	A	115	LEU
1	A	120	LEU
1	A	128	ARG
1	A	131	VAL
1	A	143	THR
1	A	150	LEU
1	A	151	LEU
1	A	156	LYS
1	A	158	ASP
1	A	173	ILE
1	A	174[A]	HIS
1	A	174[B]	HIS
1	A	177	LEU
1	A	189	ILE
1	A	211	LEU
1	A	212	ASN
1	A	215	LEU
1	A	218	LEU
1	A	226	LEU
1	A	232	ARG
1	A	237	THR
1	A	238	LEU
1	A	247	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	127	HIS
1	A	127	HIS
1	A	182	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 4 ligands modelled in this entry, 1 is modelled with single atom and 1 is 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	$\# Z > 2$
2	NO3	B	301	-	1,3,3	3.26	1 (100%)	0,3,3	-	-
5	GOL	B	304	-	5,5,5	0.45	0	5,5,5	0.35	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
5	GOL	B	304	-	-	4/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	NO3	O1-N	3.26	1.40	1.24

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	304	GOL	O1-C1-C2-C3
5	B	304	GOL	C1-C2-C3-O3
5	B	304	GOL	O1-C1-C2-O2
5	B	304	GOL	O2-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	304	GOL	2	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	193/240 (80%)	0.49	12 (6%)	28 21	40, 77, 123, 134	1 (0%)
1	B	197/240 (82%)	0.22	14 (7%)	23 18	43, 65, 112, 139	0
All	All	390/480 (81%)	0.36	26 (6%)	25 19	40, 69, 118, 139	1 (0%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	105	VAL	4.9
1	B	169	LYS	4.7
1	B	103	ASN	4.6
1	B	173	ILE	4.5
1	B	101	GLY	4.0
1	B	77	LEU	3.6
1	A	180	CYS	3.2
1	B	100	PRO	3.2
1	A	174[A]	HIS	3.1
1	B	174	HIS	2.8
1	A	207	LEU	2.8
1	B	104	ASN	2.8
1	A	100	PRO	2.7
1	A	103	ASN	2.7
1	A	153	LEU	2.6
1	A	76	ILE	2.5
1	B	92	LEU	2.5
1	A	199	PRO	2.5
1	A	203	LEU	2.4
1	A	99	ILE	2.4
1	B	248	HIS	2.3
1	A	148	ALA	2.3
1	B	249	GLY	2.2
1	B	224	CYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	96	THR	2.0
1	B	93	GLY	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(\AA^2)	Q<0.9
5	GOL	B	304	6/6	0.74	0.18	53,57,58,59	6
3	NH4	B	302	1/1	0.86	0.12	41,41,41,41	0
2	NO3	B	301	4/4	0.87	0.12	73,75,77,80	0
4	NA	B	303	1/1	0.92	0.20	56,56,56,56	0

6.5 Other polymers [i](#)

There are no such residues in this entry.