Full wwPDB X-ray Structure Validation Report ⓘ

Nov 16, 2023 – 03:17 AM JST

PDB ID	:	6KNM
Title	:	Apelin receptor in complex with single domain antibody
Authors	:	Ma, Y.B.; Ding, Y.; Song, X.; Ma, X.; Li, X.; Zhang, N.; Song, Y.; Sun, Y.; Shen, Y.; Zhong, W.; Hu, L.A.; Ma, Y.L.; Zhang, M.Y.
Deposited on	:	2019-08-06
Resolution	:	3.20 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:

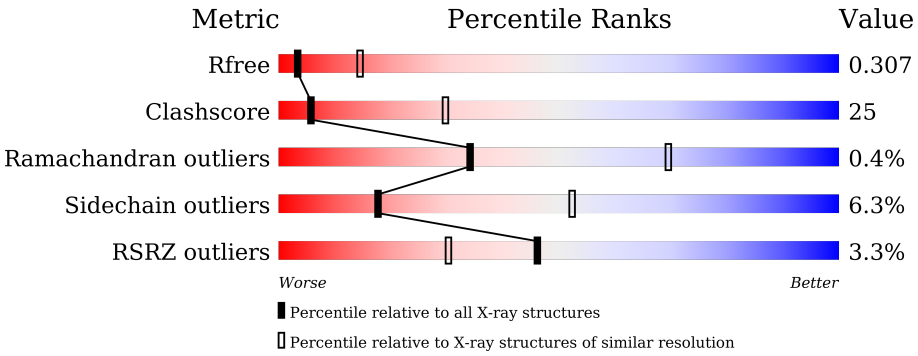
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson 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 3.20 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	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

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	B	407	3% 48% 35% . 14%
2	A	129	4% 55% 41% .

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 3747 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 Apelin receptor,Rubredoxin,Apelin receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	351	2787	1829	437	493	28	0	0	0

There are 47 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-17	MET	-	initiating methionine	UNP P35414
B	-16	LYS	-	expression tag	UNP P35414
B	-15	THR	-	expression tag	UNP P35414
B	-14	ILE	-	expression tag	UNP P35414
B	-13	ILE	-	expression tag	UNP P35414
B	-12	ALA	-	expression tag	UNP P35414
B	-11	LEU	-	expression tag	UNP P35414
B	-10	SER	-	expression tag	UNP P35414
B	-9	TYR	-	expression tag	UNP P35414
B	-8	ILE	-	expression tag	UNP P35414
B	-7	PHE	-	expression tag	UNP P35414
B	-6	CYS	-	expression tag	UNP P35414
B	-5	LEU	-	expression tag	UNP P35414
B	-4	VAL	-	expression tag	UNP P35414
B	-3	PHE	-	expression tag	UNP P35414
B	-2	ALA	-	expression tag	UNP P35414
B	-1	ASP	-	expression tag	UNP P35414
B	0	TYR	-	expression tag	UNP P35414
B	1	LYS	-	expression tag	UNP P35414
B	2	ASP	-	expression tag	UNP P35414
B	3	ASP	-	expression tag	UNP P35414
B	4	ASP	-	expression tag	UNP P35414
B	5	ASP	-	expression tag	UNP P35414
B	6	LYS	-	expression tag	UNP P35414
B	117	ALA	VAL	engineered mutation	UNP P35414
B	177	ASN	THR	engineered mutation	UNP P35414
B	261	LYS	TRP	engineered mutation	UNP P35414

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Chain	Residue	Modelled	Actual	Comment	Reference
B	325	LEU	CYS	engineered mutation	UNP P35414
B	326	MET	CYS	engineered mutation	UNP P35414
B	331	LEU	-	expression tag	UNP P35414
B	332	GLU	-	expression tag	UNP P35414
B	333	VAL	-	expression tag	UNP P35414
B	334	LEU	-	expression tag	UNP P35414
B	335	PHE	-	expression tag	UNP P35414
B	336	GLN	-	expression tag	UNP P35414
B	337	GLY	-	expression tag	UNP P35414
B	338	PRO	-	expression tag	UNP P35414
B	339	HIS	-	expression tag	UNP P35414
B	340	HIS	-	expression tag	UNP P35414
B	341	HIS	-	expression tag	UNP P35414
B	342	HIS	-	expression tag	UNP P35414
B	343	HIS	-	expression tag	UNP P35414
B	344	HIS	-	expression tag	UNP P35414
B	345	HIS	-	expression tag	UNP P35414
B	346	HIS	-	expression tag	UNP P35414
B	347	HIS	-	expression tag	UNP P35414
B	348	HIS	-	expression tag	UNP P35414

- Molecule 2 is a protein called Single domain antibody JN241.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	129	Total	C	N	O	S	0	0	0
			959	587	170	193	9			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

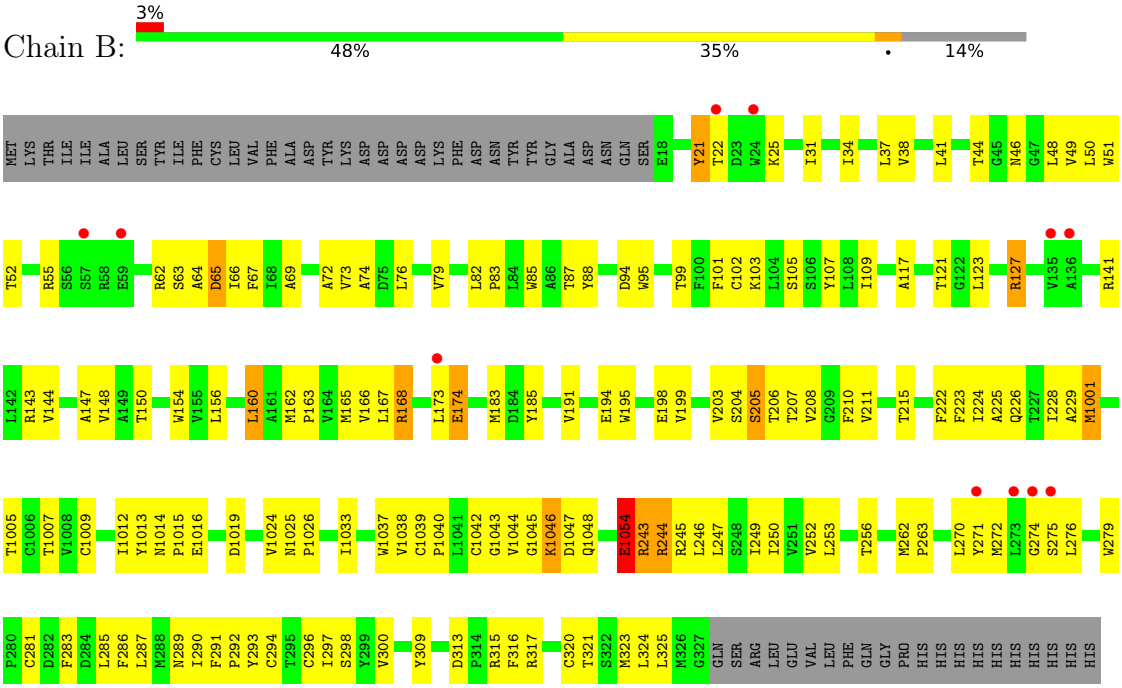
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		

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.

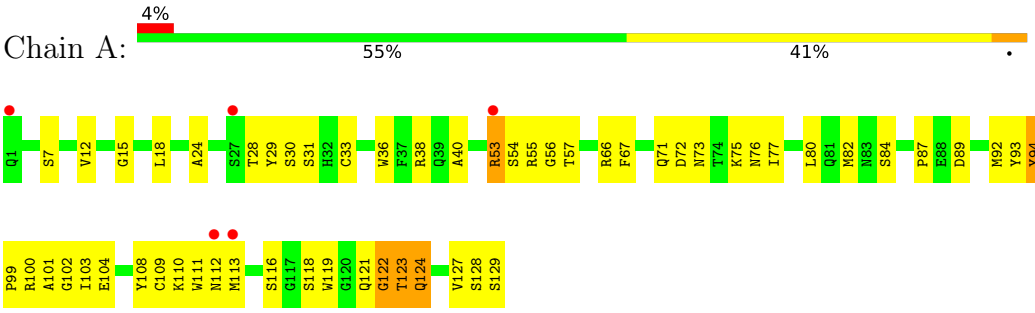
• Molecule 1: Apelin receptor,Rubredoxin,Apelin receptor

Chain B:



• Molecule 2: Single domain antibody JN241

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	44.78Å 48.39Å 350.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.64 – 3.20 46.64 – 3.20	Depositor EDS
% Data completeness (in resolution range)	92.2 (46.64-3.20) 92.5 (46.64-3.20)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.94 (at 3.19Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.258 , 0.305 0.258 , 0.307	Depositor DCC
R_{free} test set	621 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	81.1	Xtriage
Anisotropy	1.120	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 75.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.43$, $< L^2 > = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	3747	wwPDB-VP
Average B, all atoms (Å^2)	95.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.84% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	B	0.49	0/2865	0.69	1/3910 (0.0%)
2	A	0.65	1/976 (0.1%)	0.86	5/1315 (0.4%)
All	All	0.54	1/3841 (0.0%)	0.74	6/5225 (0.1%)

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
2	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	123	THR	C-O	5.56	1.33	1.23

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	116	SER	N-CA-C	-5.98	94.84	111.00
2	A	122	GLY	N-CA-C	-5.98	98.15	113.10
2	A	123	THR	N-CA-C	-5.42	96.36	111.00
1	B	325	LEU	CA-CB-CG	5.39	127.71	115.30
2	A	40	ALA	N-CA-C	5.33	125.39	111.00
2	A	122	GLY	CA-C-O	5.15	129.87	120.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	A	112	ASN	Mainchain

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	B	2787	0	2762	139	0
2	A	959	0	923	64	0
3	B	1	0	0	0	0
All	All	3747	0	3685	189	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 25.

All (189) 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:1037:TRP:O	1:B:1046:LYS:CE	1.66	1.42
2:A:33:CYS:SG	2:A:109:CYS:SG	1.43	1.31
1:B:1037:TRP:O	1:B:1046:LYS:HE3	1.38	1.10
1:B:1037:TRP:O	1:B:1046:LYS:HE2	1.28	1.04
2:A:29:TYR:O	2:A:100:ARG:NH1	1.92	1.03
2:A:93:TYR:N	2:A:123:THR:O	1.96	0.98
2:A:113:MET:HA	2:A:113:MET:CE	2.02	0.89
1:B:244:ARG:HH21	1:B:244:ARG:HG3	1.38	0.89
1:B:22:THR:CG2	2:A:56:GLY:HA2	2.03	0.88
1:B:22:THR:HG21	2:A:56:GLY:HA2	1.56	0.88
2:A:94:TYR:CD1	2:A:122:GLY:HA3	2.08	0.87
1:B:263:PRO:HG3	1:B:297:ILE:HD11	1.57	0.87
2:A:33:CYS:CB	2:A:109:CYS:SG	2.62	0.87
1:B:274:GLY:HA2	1:B:279:TRP:HD1	1.42	0.85
1:B:207:THR:HA	1:B:211:VAL:HG12	1.58	0.84
1:B:271:TYR:CD2	2:A:103:ILE:HD11	2.13	0.84
2:A:113:MET:HA	2:A:113:MET:HE3	1.59	0.84
2:A:99:PRO:O	2:A:110:LYS:HG3	1.79	0.83
1:B:1046:LYS:CE	1:B:1046:LYS:H	1.94	0.79
1:B:203:VAL:O	1:B:207:THR:OG1	2.01	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:67:PHE:HB2	1:B:123:LEU:HD13	1.65	0.79
1:B:174:GLU:OE2	2:A:53:ARG:N	2.18	0.76
2:A:53:ARG:HG2	2:A:54:SER:N	1.99	0.76
1:B:243:ARG:HB2	1:B:245:ARG:HH12	1.51	0.75
1:B:1045:GLY:CA	1:B:1046:LYS:NZ	2.52	0.73
1:B:1046:LYS:H	1:B:1046:LYS:HZ2	1.37	0.72
1:B:1045:GLY:HA3	1:B:1046:LYS:NZ	2.04	0.72
2:A:92:MET:HA	2:A:124:GLN:HA	1.72	0.72
1:B:1046:LYS:H	1:B:1046:LYS:NZ	1.90	0.70
2:A:33:CYS:SG	2:A:109:CYS:CB	2.76	0.70
1:B:244:ARG:HG3	1:B:244:ARG:NH2	2.06	0.70
2:A:87:PRO:HA	2:A:127:VAL:HB	1.75	0.69
2:A:94:TYR:CD1	2:A:122:GLY:CA	2.77	0.68
2:A:96:ALA:HB2	2:A:119:TRP:CZ3	2.29	0.68
1:B:185:TYR:HB2	1:B:194:GLU:HG3	1.74	0.67
1:B:1046:LYS:HZ2	1:B:1046:LYS:N	1.93	0.67
2:A:93:TYR:O	2:A:123:THR:N	2.28	0.66
1:B:274:GLY:HA2	1:B:279:TRP:CD1	2.29	0.65
1:B:127:ARG:HD2	1:B:250:ILE:HD12	1.77	0.65
1:B:22:THR:HG21	2:A:56:GLY:CA	2.25	0.63
1:B:1045:GLY:HA3	1:B:1046:LYS:HZ1	1.61	0.63
2:A:94:TYR:HD1	2:A:122:GLY:CA	2.12	0.63
2:A:121:GLN:HA	2:A:121:GLN:OE1	1.99	0.62
1:B:207:THR:HA	1:B:211:VAL:CG1	2.28	0.61
1:B:1045:GLY:CA	1:B:1046:LYS:HZ1	2.10	0.61
1:B:243:ARG:HB2	1:B:245:ARG:NH1	2.14	0.61
2:A:92:MET:SD	2:A:124:GLN:HB2	2.41	0.61
1:B:224:ILE:O	1:B:228:ILE:HG13	2.01	0.59
2:A:93:TYR:O	2:A:122:GLY:HA2	2.02	0.59
2:A:111:TRP:HA	2:A:111:TRP:CE3	2.38	0.59
1:B:107:TYR:CD1	1:B:165:MET:HB2	2.38	0.59
1:B:46:ASN:HB2	1:B:76:LEU:HD13	1.85	0.58
1:B:321:THR:O	1:B:324:LEU:HB3	2.04	0.58
1:B:51:TRP:O	1:B:55:ARG:N	2.29	0.58
1:B:174:GLU:CD	2:A:53:ARG:HD3	2.25	0.57
1:B:69:ALA:O	1:B:73:VAL:HG23	2.05	0.57
1:B:205:SER:OG	1:B:206:THR:N	2.38	0.56
1:B:313:ASP:O	1:B:317:ARG:HG3	2.05	0.56
1:B:163:PRO:O	1:B:167:LEU:HB2	2.05	0.56
1:B:1007:THR:HG22	1:B:1048:GLN:O	2.06	0.56
2:A:94:TYR:HA	2:A:122:GLY:HA2	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:96:ALA:HB2	2:A:119:TRP:CE3	2.41	0.56
2:A:75:LYS:HZ3	2:A:77:ILE:HD12	1.71	0.56
1:B:38:VAL:HG21	1:B:300:VAL:HG12	1.87	0.56
1:B:166:VAL:HG12	1:B:167:LEU:HD12	1.87	0.55
1:B:1014:ASN:OD1	1:B:1016:GLU:HB2	2.07	0.55
1:B:1042:CYS:SG	1:B:1044:VAL:HG22	2.47	0.55
1:B:286:PHE:CE1	1:B:290:ILE:HD13	2.43	0.54
1:B:66:ILE:HD11	1:B:150:THR:HG21	1.90	0.54
1:B:1025:ASN:N	1:B:1025:ASN:OD1	2.40	0.54
1:B:99:THR:HG22	1:B:103:LYS:HD2	1.89	0.54
1:B:79:VAL:HA	1:B:82:LEU:HD12	1.90	0.53
1:B:21:TYR:O	1:B:22:THR:HG23	2.09	0.53
1:B:48:LEU:O	1:B:52:THR:HG23	2.08	0.53
1:B:62:ARG:NH1	1:B:315:ARG:HB3	2.23	0.53
1:B:225:ALA:O	1:B:229:ALA:N	2.40	0.53
1:B:270:LEU:O	1:B:274:GLY:N	2.42	0.53
1:B:51:TRP:CH2	1:B:323:MET:SD	3.03	0.52
1:B:62:ARG:O	1:B:65:ASP:HB2	2.09	0.52
1:B:101:PHE:O	1:B:105:SER:N	2.33	0.52
2:A:53:ARG:NH2	2:A:54:SER:OG	2.43	0.52
1:B:67:PHE:HB2	1:B:123:LEU:CD1	2.38	0.52
1:B:1045:GLY:CA	1:B:1046:LYS:HZ2	2.23	0.52
1:B:62:ARG:HH21	1:B:246:LEU:HG	1.73	0.52
2:A:101:ALA:O	2:A:102:GLY:C	2.47	0.52
2:A:53:ARG:NE	2:A:108:TYR:CZ	2.78	0.52
1:B:44:THR:O	1:B:48:LEU:HD13	2.09	0.52
1:B:275:SER:O	1:B:276:LEU:HD13	2.10	0.52
1:B:144:VAL:O	1:B:148:VAL:N	2.37	0.52
1:B:211:VAL:O	1:B:215:THR:HG23	2.09	0.52
1:B:117:ALA:O	1:B:121:THR:HG23	2.10	0.51
1:B:275:SER:HB3	1:B:276:LEU:HD22	1.92	0.51
1:B:83:PRO:O	1:B:87:THR:HG23	2.11	0.51
1:B:62:ARG:HG2	1:B:65:ASP:CG	2.31	0.51
1:B:1046:LYS:H	1:B:1046:LYS:CD	2.23	0.51
1:B:21:TYR:CD1	2:A:29:TYR:CE1	2.98	0.51
1:B:67:PHE:CB	1:B:123:LEU:HD13	2.40	0.51
2:A:12:VAL:HG21	2:A:18:LEU:HG	1.93	0.50
1:B:21:TYR:CD1	2:A:29:TYR:HE1	2.29	0.50
1:B:51:TRP:HH2	1:B:323:MET:SD	2.34	0.50
2:A:93:TYR:O	2:A:122:GLY:CA	2.58	0.50
1:B:1019:ASP:HB3	1:B:1024:VAL:HG13	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:94:TYR:CD1	2:A:121:GLN:O	2.65	0.49
1:B:199:VAL:O	1:B:203:VAL:HG23	2.12	0.49
1:B:49:VAL:HG12	1:B:316:PHE:CE1	2.47	0.49
2:A:66:ARG:NH1	2:A:89:ASP:OD2	2.36	0.49
1:B:195:TRP:CH2	1:B:276:LEU:HD11	2.48	0.49
1:B:252:VAL:O	1:B:256:THR:HG23	2.13	0.49
2:A:72:ASP:O	2:A:76:ASN:N	2.45	0.49
2:A:15:GLY:HA2	2:A:84:SER:HA	1.94	0.48
1:B:168:ARG:HH11	2:A:104:GLU:HB3	1.78	0.48
1:B:225:ALA:O	1:B:229:ALA:HB2	2.13	0.48
1:B:51:TRP:CH2	1:B:55:ARG:HD2	2.47	0.48
1:B:297:ILE:O	1:B:300:VAL:HG22	2.12	0.48
2:A:128:SER:OG	2:A:129:SER:N	2.46	0.48
2:A:53:ARG:HD2	2:A:108:TYR:CG	2.49	0.48
2:A:94:TYR:CE1	2:A:121:GLN:O	2.66	0.48
1:B:1045:GLY:HA3	1:B:1046:LYS:HZ2	1.78	0.47
1:B:243:ARG:HG3	1:B:243:ARG:O	2.15	0.47
1:B:253:LEU:HD11	1:B:309:TYR:CE1	2.50	0.47
1:B:191:VAL:HG13	2:A:99:PRO:HG3	1.97	0.47
1:B:37:LEU:O	1:B:41:LEU:HG	2.14	0.47
1:B:204:SER:O	1:B:208:VAL:HG12	2.15	0.47
1:B:1016:GLU:O	1:B:1026:PRO:HB3	2.15	0.47
1:B:1037:TRP:CZ2	1:B:1040:PRO:HD3	2.50	0.46
1:B:31:ILE:HD11	1:B:293:TYR:CE1	2.51	0.46
1:B:63:SER:OG	1:B:64:ALA:N	2.48	0.46
2:A:100:ARG:HA	2:A:108:TYR:O	2.16	0.46
1:B:291:PHE:N	1:B:292:PRO:HD2	2.31	0.46
1:B:320:CYS:O	1:B:321:THR:C	2.53	0.46
1:B:287:LEU:O	1:B:290:ILE:HG22	2.16	0.46
1:B:226:GLN:HA	1:B:229:ALA:HB2	1.98	0.45
1:B:143:ARG:O	1:B:147:ALA:N	2.50	0.45
1:B:173:LEU:HD12	2:A:108:TYR:HB3	1.98	0.45
2:A:33:CYS:O	2:A:97:ALA:HA	2.17	0.45
1:B:1005:THR:HG22	1:B:1012:ILE:HG12	1.99	0.44
1:B:275:SER:HB2	2:A:30:SER:OG	2.17	0.44
2:A:67:PHE:CE1	2:A:82:MET:HB3	2.52	0.44
1:B:50:LEU:HD23	1:B:50:LEU:HA	1.67	0.44
1:B:1007:THR:CG2	1:B:1048:GLN:O	2.65	0.44
1:B:247:LEU:HD12	1:B:247:LEU:HA	1.71	0.44
1:B:297:ILE:HG13	1:B:298:SER:N	2.31	0.44
2:A:24:ALA:HB3	2:A:76:ASN:OD1	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:53:ARG:HG3	2:A:108:TYR:CE1	2.53	0.44
2:A:28:THR:OG1	2:A:31:SER:N	2.43	0.43
1:B:173:LEU:HD12	2:A:108:TYR:CB	2.48	0.43
1:B:272:MET:O	1:B:276:LEU:HD23	2.18	0.43
2:A:36:TRP:CG	2:A:80:LEU:HD22	2.53	0.43
2:A:97:ALA:HB3	2:A:118:SER:HB2	2.01	0.43
1:B:262:MET:N	1:B:263:PRO:HD2	2.34	0.43
2:A:92:MET:CG	2:A:124:GLN:HB2	2.49	0.43
1:B:88:TYR:CD2	1:B:95:TRP:HB2	2.52	0.43
1:B:253:LEU:HD11	1:B:309:TYR:HE1	1.83	0.43
1:B:51:TRP:CZ3	1:B:55:ARG:HD2	2.54	0.43
2:A:18:LEU:HD23	2:A:18:LEU:HA	1.85	0.43
2:A:12:VAL:O	2:A:127:VAL:HA	2.19	0.43
1:B:85:TRP:NE1	1:B:109:ILE:HD11	2.33	0.42
1:B:1038:VAL:HG12	1:B:1045:GLY:HA2	2.01	0.42
1:B:1039:CYS:HB3	1:B:1043:GLY:H	1.84	0.42
1:B:31:ILE:O	1:B:34:ILE:N	2.53	0.42
1:B:48:LEU:O	1:B:51:TRP:N	2.52	0.42
1:B:195:TRP:CZ3	1:B:276:LEU:HD21	2.54	0.42
1:B:147:ALA:O	1:B:150:THR:HG22	2.20	0.42
1:B:1001:MET:HG2	1:B:1015:PRO:CB	2.49	0.42
1:B:185:TYR:CE2	1:B:198:GLU:HB2	2.55	0.42
1:B:168:ARG:HG2	1:B:183:MET:HG2	2.01	0.42
1:B:156:LEU:O	1:B:160:LEU:HD23	2.20	0.42
1:B:162:MET:O	1:B:165:MET:N	2.51	0.41
1:B:1013:TYR:CZ	1:B:1033:ILE:HD11	2.55	0.41
1:B:285:LEU:O	1:B:285:LEU:HG	2.20	0.41
2:A:96:ALA:HB2	2:A:119:TRP:CH2	2.55	0.41
1:B:22:THR:HG23	2:A:56:GLY:HA2	1.96	0.41
2:A:103:ILE:H	2:A:103:ILE:HD12	1.86	0.41
1:B:1045:GLY:HA2	1:B:1046:LYS:NZ	2.33	0.41
1:B:76:LEU:HD12	1:B:76:LEU:HA	1.75	0.41
1:B:72:ALA:O	1:B:76:LEU:HB2	2.21	0.41
2:A:38:ARG:HA	2:A:92:MET:O	2.21	0.41
1:B:249:ILE:O	1:B:252:VAL:HG12	2.20	0.41
1:B:263:PRO:HB2	1:B:294:CYS:SG	2.61	0.41
1:B:74:ALA:HB2	1:B:154:TRP:CH2	2.56	0.40
1:B:206:THR:O	1:B:210:PHE:N	2.40	0.40
1:B:222:PHE:HD2	1:B:223:PHE:CD1	2.38	0.40
1:B:225:ALA:HA	1:B:228:ILE:HB	2.02	0.40
1:B:1045:GLY:O	1:B:1047:ASP:N	2.55	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:144:VAL:O	1:B:147:ALA:N	2.55	0.40
1:B:199:VAL:HG13	1:B:272:MET:HB3	2.03	0.40
1:B:1054:GLU:H	1:B:1054:GLU:HG2	1.46	0.40
2:A:29:TYR:HB3	2:A:73:ASN:ND2	2.37	0.40
1:B:296:CYS:O	1:B:300:VAL:HG13	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	B	349/407 (86%)	307 (88%)	40 (12%)	2 (1%)	25 64
2	A	127/129 (98%)	109 (86%)	18 (14%)	0	100 100
All	All	476/536 (89%)	416 (87%)	58 (12%)	2 (0%)	34 69

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	243	ARG
1	B	1054	GLU

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	B	307/358 (86%)	288 (94%)	19 (6%)	18 53
2	A	103/103 (100%)	96 (93%)	7 (7%)	16 49
All	All	410/461 (89%)	384 (94%)	26 (6%)	18 52

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	21	TYR
1	B	25	LYS
1	B	65	ASP
1	B	94	ASP
1	B	102	CYS
1	B	127	ARG
1	B	141	ARG
1	B	160	LEU
1	B	168	ARG
1	B	174	GLU
1	B	205	SER
1	B	1001	MET
1	B	1009	CYS
1	B	1046	LYS
1	B	1054	GLU
1	B	244	ARG
1	B	281	CYS
1	B	283	PHE
1	B	289	ASN
2	A	7	SER
2	A	53	ARG
2	A	55	ARG
2	A	57	THR
2	A	71	GLN
2	A	94	TYR
2	A	124	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	112	ASN
1	B	301	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

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

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 95^{th} 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	B	351/407 (86%)	0.03	11 (3%) 49 32	73, 91, 122, 154	0
2	A	129/129 (100%)	0.19	5 (3%) 39 25	79, 97, 123, 132	0
All	All	480/536 (89%)	0.07	16 (3%) 46 30	73, 93, 123, 154	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	135	VAL	4.1
2	A	1	GLN	4.1
1	B	275	SER	3.7
1	B	274	GLY	3.6
1	B	22	THR	3.0
1	B	271	TYR	2.8
2	A	27	SER	2.8
2	A	112	ASN	2.8
1	B	57	SER	2.4
1	B	173	LEU	2.4
1	B	136	ALA	2.3
1	B	59	GLU	2.3
1	B	273	LEU	2.3
1	B	24	TRP	2.2
2	A	53	ARG	2.1
2	A	113	MET	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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)

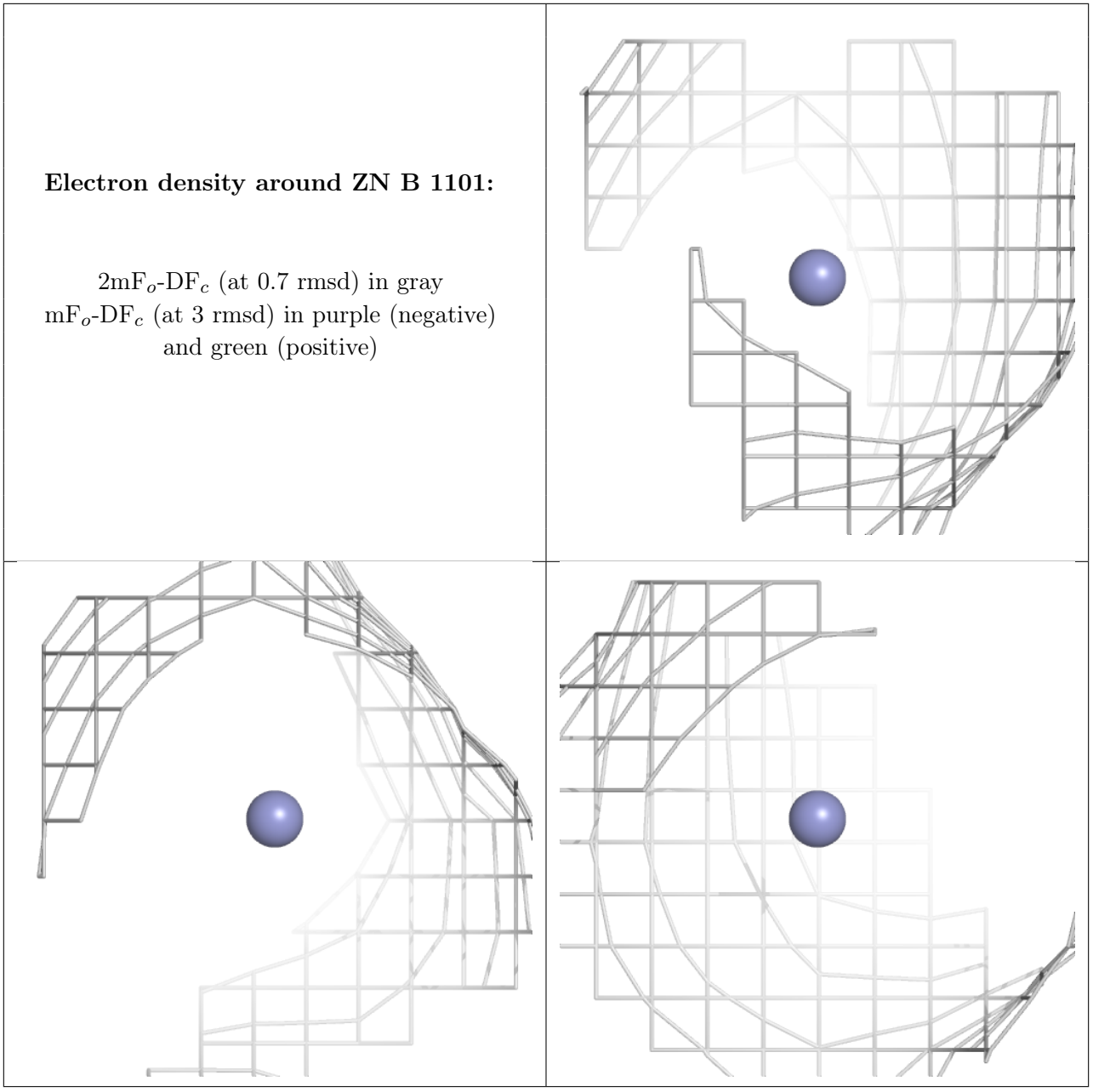
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' 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
3	ZN	B	1101	1/1	0.95	0.21	90,90,90,90	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around ZN B 1101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



6.5 Other polymers (i)

There are no such residues in this entry.