



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:09 AM GMT

PDB ID : 1ZY0
Title : X-ray structure of peptide deformylase from Arabidopsis thaliana (AtPDF1A); crystals grown in PEG-6000
Authors : Fieulaine, S.; Juillan-Binard, C.; Serero, A.; Dardel, F.; Giglione, C.; Meinel, T.; Ferrer, J.-L.
Deposited on : 2005-06-09
Resolution : 2.90 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

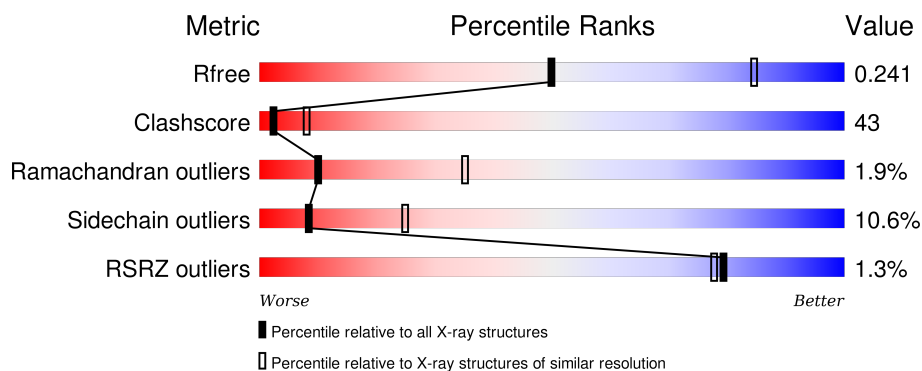
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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. 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	197	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 2%, orange 8%, yellow 50%, green 38%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 2% 38% 50% 8% • • </div> </div>
1	B	197	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 2%, orange 8%, yellow 49%, green 39%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 2% 39% 49% 8% • • </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3035 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 Peptide deformylase, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	191	Total	C	N	O	S	0	0	0
			1478	934	262	274	8			
1	B	191	Total	C	N	O	S	0	0	0
			1464	923	255	278	8			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	INITIATING METHIONINE	UNP Q9FV53
A	191	SER	-	EXPRESSION TAG	UNP Q9FV53
A	192	HIS	-	EXPRESSION TAG	UNP Q9FV53
A	193	HIS	-	EXPRESSION TAG	UNP Q9FV53
A	194	HIS	-	EXPRESSION TAG	UNP Q9FV53
A	195	HIS	-	EXPRESSION TAG	UNP Q9FV53
A	196	HIS	-	EXPRESSION TAG	UNP Q9FV53
A	197	HIS	-	EXPRESSION TAG	UNP Q9FV53
B	1	MET	-	INITIATING METHIONINE	UNP Q9FV53
B	191	SER	-	EXPRESSION TAG	UNP Q9FV53
B	192	HIS	-	EXPRESSION TAG	UNP Q9FV53
B	193	HIS	-	EXPRESSION TAG	UNP Q9FV53
B	194	HIS	-	EXPRESSION TAG	UNP Q9FV53
B	195	HIS	-	EXPRESSION TAG	UNP Q9FV53
B	196	HIS	-	EXPRESSION TAG	UNP Q9FV53
B	197	HIS	-	EXPRESSION TAG	UNP Q9FV53

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

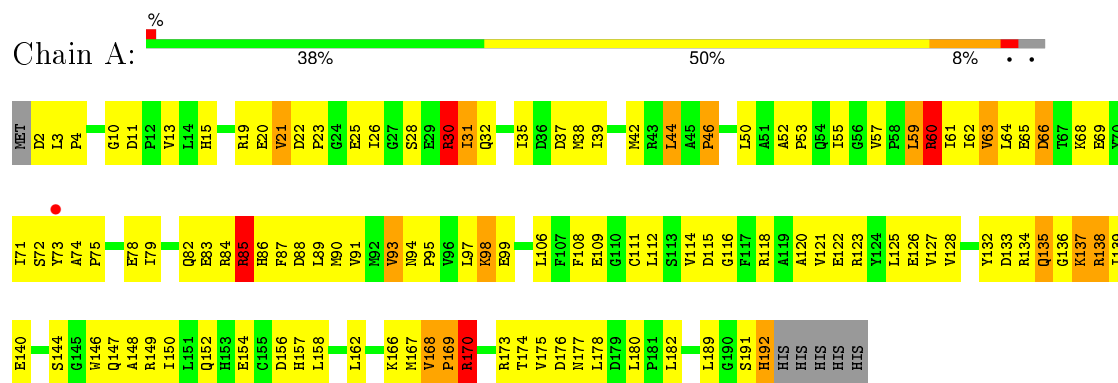
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	42	Total 42	O 42	0	0
3	B	49	Total 49	O 49	0	0

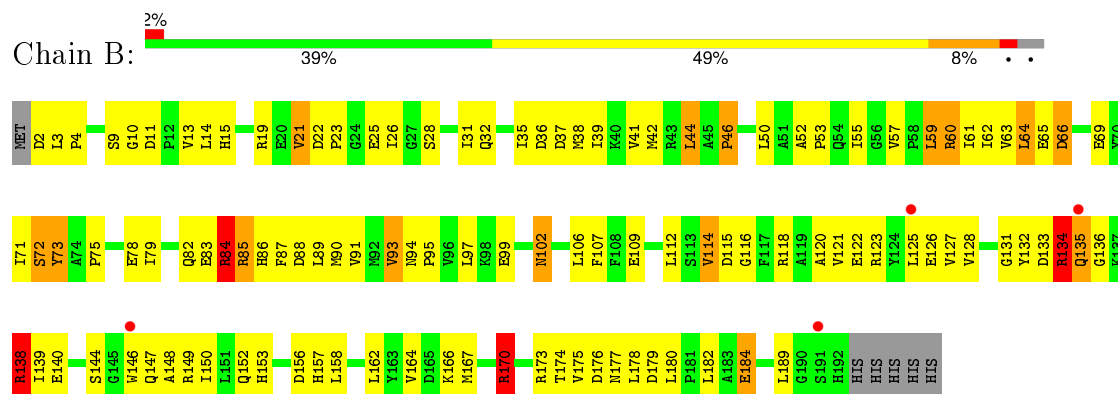
3 Residue-property plots

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



- Molecule 1: Peptide deformylase, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	51.50 Å 76.80 Å 109.30 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.90 46.59 – 2.89	Depositor EDS
% Data completeness (in resolution range)	87.2 (30.00-2.90) 99.3 (46.59-2.89)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.35 (at 2.91 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.243 , 0.283 0.261 , 0.241	Depositor DCC
R_{free} test set	503 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	59.0	Xtriage
Anisotropy	1.284	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 63.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 10135 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3035	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.74% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	A	2.07	28/1505 (1.9%)	1.72	33/2041 (1.6%)
1	B	1.69	19/1491 (1.3%)	1.65	24/2028 (1.2%)
All	All	1.89	47/2996 (1.6%)	1.69	57/4069 (1.4%)

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
1	A	0	4
1	B	0	4
All	All	0	8

All (47) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	138	ARG	CZ-NH2	-30.76	0.93	1.33
1	B	138	ARG	CZ-NH1	-28.22	0.96	1.33
1	A	140	GLU	CD-OE1	-26.82	0.96	1.25
1	A	140	GLU	CD-OE2	-25.51	0.97	1.25
1	B	140	GLU	CD-OE1	-24.14	0.99	1.25
1	B	140	GLU	CD-OE2	-23.59	0.99	1.25
1	A	138	ARG	CZ-NH1	-22.50	1.03	1.33
1	B	138	ARG	NE-CZ	-22.44	1.03	1.33
1	A	137	LYS	C-N	-20.53	0.86	1.34
1	A	138	ARG	NE-CZ	-19.03	1.08	1.33
1	A	60	ARG	CZ-NH1	-16.75	1.11	1.33
1	A	138	ARG	CD-NE	-14.01	1.22	1.46
1	A	138	ARG	CB-CG	-13.01	1.17	1.52
1	A	122	GLU	CD-OE1	-12.96	1.11	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	167	MET	C-N	12.64	1.63	1.34
1	B	122	GLU	CD-OE1	-12.59	1.11	1.25
1	B	138	ARG	CB-CG	-12.52	1.18	1.52
1	B	122	GLU	CD-OE2	-12.27	1.12	1.25
1	B	138	ARG	CZ-NH2	-11.40	1.18	1.33
1	A	167	MET	C-O	-11.13	1.02	1.23
1	A	122	GLU	CD-OE2	-10.98	1.13	1.25
1	B	60	ARG	CZ-NH1	-10.95	1.18	1.33
1	B	19	ARG	CZ-NH1	-10.48	1.19	1.33
1	A	19	ARG	CZ-NH1	-10.15	1.19	1.33
1	A	138	ARG	N-CA	9.49	1.65	1.46
1	A	137	LYS	CA-C	9.21	1.76	1.52
1	B	60	ARG	CZ-NH2	-8.48	1.22	1.33
1	A	85	ARG	CZ-NH1	-7.69	1.23	1.33
1	A	168	VAL	C-O	-7.66	1.08	1.23
1	A	170	ARG	CZ-NH1	-7.16	1.23	1.33
1	A	2	ASP	CG-OD2	-7.06	1.09	1.25
1	B	2	ASP	CG-OD1	-6.71	1.09	1.25
1	B	138	ARG	CD-NE	-6.58	1.35	1.46
1	A	64	LEU	C-N	-6.16	1.19	1.34
1	B	85	ARG	CZ-NH1	-6.08	1.25	1.33
1	B	170	ARG	NE-CZ	5.96	1.40	1.33
1	A	168	VAL	C-N	5.91	1.45	1.34
1	A	137	LYS	C-O	5.90	1.34	1.23
1	B	102	ASN	CB-CG	-5.68	1.38	1.51
1	B	140	GLU	CB-CG	-5.42	1.41	1.52
1	A	19	ARG	CZ-NH2	-5.41	1.26	1.33
1	B	85	ARG	NE-CZ	-5.28	1.26	1.33
1	A	63	VAL	C-O	5.25	1.33	1.23
1	A	63	VAL	C-N	-5.22	1.22	1.34
1	A	64	LEU	C-O	5.08	1.33	1.23
1	A	169	PRO	N-CA	-5.03	1.38	1.47
1	B	19	ARG	CG-CD	-5.02	1.39	1.51

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	170	ARG	NE-CZ-NH2	-32.99	103.81	120.30
1	B	60	ARG	NE-CZ-NH1	26.85	133.73	120.30
1	A	138	ARG	NE-CZ-NH2	-26.69	106.95	120.30
1	A	30	ARG	NE-CZ-NH2	-25.55	107.52	120.30
1	A	85	ARG	NE-CZ-NH2	-24.16	108.22	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	84	ARG	NE-CZ-NH2	-22.63	108.98	120.30
1	A	170	ARG	NE-CZ-NH2	-21.53	109.54	120.30
1	B	140	GLU	OE1-CD-OE2	-18.85	100.68	123.30
1	A	140	GLU	OE1-CD-OE2	-17.29	102.55	123.30
1	A	170	ARG	NE-CZ-NH1	15.62	128.11	120.30
1	B	138	ARG	NE-CZ-NH2	-15.17	112.71	120.30
1	A	85	ARG	NE-CZ-NH1	12.92	126.76	120.30
1	B	138	ARG	CD-NE-CZ	-12.55	106.03	123.60
1	B	60	ARG	NH1-CZ-NH2	-12.53	105.61	119.40
1	A	138	ARG	NE-CZ-NH1	12.09	126.35	120.30
1	B	85	ARG	NE-CZ-NH1	-11.94	114.33	120.30
1	A	137	LYS	CA-C-O	-10.62	97.79	120.10
1	A	137	LYS	O-C-N	9.92	138.57	122.70
1	B	170	ARG	NH1-CZ-NH2	-9.87	108.54	119.40
1	A	60	ARG	NE-CZ-NH2	8.89	124.75	120.30
1	B	122	GLU	OE1-CD-OE2	-8.67	112.90	123.30
1	A	170	ARG	NH1-CZ-NH2	-8.28	110.30	119.40
1	B	84	ARG	NH1-CZ-NH2	-8.03	110.57	119.40
1	A	122	GLU	OE1-CD-OE2	-8.00	113.70	123.30
1	B	138	ARG	CG-CD-NE	-7.87	95.28	111.80
1	A	60	ARG	NH1-CZ-NH2	-7.86	110.75	119.40
1	B	2	ASP	CB-CG-OD2	7.69	125.22	118.30
1	A	30	ARG	NH1-CZ-NH2	-7.65	110.99	119.40
1	A	170	ARG	CD-NE-CZ	-7.53	113.06	123.60
1	A	2	ASP	CB-CG-OD1	7.41	124.97	118.30
1	B	118	ARG	NE-CZ-NH2	-7.07	116.77	120.30
1	B	19	ARG	CG-CD-NE	6.84	126.17	111.80
1	A	168	VAL	N-CA-CB	-6.82	96.49	111.50
1	B	102	ASN	CB-CA-C	-6.78	96.84	110.40
1	A	138	ARG	NH1-CZ-NH2	6.63	126.69	119.40
1	A	167	MET	CA-C-N	-6.61	102.66	117.20
1	A	60	ARG	NE-CZ-NH1	6.52	123.56	120.30
1	A	168	VAL	CA-C-O	6.50	133.74	120.10
1	A	169	PRO	N-CD-CG	-6.45	93.53	103.20
1	B	60	ARG	NE-CZ-NH2	-6.31	117.14	120.30
1	B	118	ARG	CG-CD-NE	-6.10	98.99	111.80
1	A	138	ARG	N-CA-CB	6.04	121.47	110.60
1	A	85	ARG	CD-NE-CZ	-5.81	115.47	123.60
1	B	140	GLU	CG-CD-OE1	5.74	129.77	118.30
1	A	169	PRO	CA-N-CD	5.73	119.72	111.70
1	A	19	ARG	NE-CZ-NH2	5.69	123.14	120.30
1	B	140	GLU	CG-CD-OE2	5.62	129.53	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	19	ARG	NE-CZ-NH1	-5.50	117.55	120.30
1	A	2	ASP	OD1-CG-OD2	-5.35	113.13	123.30
1	A	192	HIS	N-CA-C	-5.34	96.59	111.00
1	A	140	GLU	CG-CD-OE2	5.33	128.97	118.30
1	B	138	ARG	NH1-CZ-NH2	-5.18	113.70	119.40
1	A	168	VAL	N-CA-C	-5.17	97.04	111.00
1	B	118	ARG	NE-CZ-NH1	5.16	122.88	120.30
1	B	2	ASP	C-N-CA	-5.11	108.93	121.70
1	A	140	GLU	CG-CD-OE1	5.09	128.48	118.30
1	B	2	ASP	OD1-CG-OD2	-5.05	113.71	123.30

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	170	ARG	Sidechain
1	A	30	ARG	Sidechain
1	A	60	ARG	Sidechain
1	A	85	ARG	Sidechain
1	B	138	ARG	Sidechain
1	B	170	ARG	Sidechain
1	B	60	ARG	Sidechain
1	B	84	ARG	Sidechain

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	1478	0	1465	127	0
1	B	1464	0	1421	126	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	42	0	0	2	0
3	B	49	0	0	4	0
All	All	3035	0	2886	249	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 43.

All (249) 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:137:LYS:CA	1:A:137:LYS:C	1.76	1.52
1:A:21:VAL:HG23	1:A:59:LEU:HD12	1.31	1.11
1:A:30:ARG:HH11	1:A:30:ARG:HG3	1.17	1.07
1:A:137:LYS:C	1:A:138:ARG:CA	2.23	1.06
1:B:78:GLU:HG3	3:B:241:HOH:O	1.55	1.03
1:A:177:ASN:HB3	1:A:180:LEU:HD12	1.42	1.01
1:B:177:ASN:HB3	1:B:180:LEU:HD12	1.43	0.95
1:B:75:PRO:HB2	1:B:78:GLU:HB2	1.50	0.93
1:B:73:TYR:H	1:B:73:TYR:HD1	1.10	0.93
1:A:137:LYS:C	1:A:138:ARG:N	0.86	0.91
1:B:21:VAL:HG23	1:B:59:LEU:HD12	1.51	0.91
1:A:147:GLN:NE2	3:A:226:HOH:O	2.04	0.89
1:A:133:ASP:HB3	1:A:139:ILE:HD13	1.55	0.88
1:A:137:LYS:CA	1:A:138:ARG:N	2.36	0.88
1:A:137:LYS:O	1:A:138:ARG:N	2.07	0.87
1:A:21:VAL:HG23	1:A:59:LEU:CD1	2.04	0.87
1:B:28:SER:O	1:B:32:GLN:HG2	1.76	0.83
1:B:112:LEU:HD12	1:B:112:LEU:N	1.94	0.83
1:A:123:ARG:HG3	1:A:149:ARG:HG3	1.60	0.83
1:B:138:ARG:NH1	1:B:138:ARG:CG	2.27	0.82
1:B:10:GLY:H	1:B:15:HIS:HE1	1.27	0.81
1:A:75:PRO:HB2	1:A:78:GLU:HB2	1.61	0.81
1:A:21:VAL:CG2	1:A:31:ILE:HD11	2.12	0.80
1:A:30:ARG:NH1	1:A:30:ARG:HG3	1.98	0.79
1:B:10:GLY:H	1:B:15:HIS:CE1	2.01	0.79
1:A:30:ARG:CG	1:A:30:ARG:HH11	1.96	0.77
1:B:123:ARG:HG3	1:B:149:ARG:HG3	1.66	0.77
1:A:10:GLY:H	1:A:15:HIS:CE1	2.03	0.77
1:B:75:PRO:HB2	1:B:78:GLU:CB	2.15	0.77
1:A:10:GLY:H	1:A:15:HIS:HE1	1.31	0.77
1:A:21:VAL:CG2	1:A:59:LEU:HD12	2.14	0.76
1:A:93:VAL:HG22	1:A:132:TYR:HB2	1.67	0.76
1:B:21:VAL:HG23	1:B:59:LEU:CD1	2.16	0.75
1:A:134:ARG:HH11	1:A:134:ARG:HG2	1.52	0.74
1:A:109:GLU:OE2	1:A:123:ARG:NE	2.21	0.74
1:A:75:PRO:HB2	1:A:78:GLU:CB	2.18	0.73
1:A:137:LYS:CA	1:A:137:LYS:O	2.36	0.73
1:B:179:ASP:OD1	3:B:236:HOH:O	2.07	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:133:ASP:HB3	1:B:139:ILE:HD13	1.71	0.72
1:B:107:PHE:HA	1:B:189:LEU:HD22	1.72	0.71
1:A:25:GLU:O	1:A:28:SER:HB3	1.90	0.71
1:B:73:TYR:HD1	1:B:73:TYR:N	1.88	0.70
1:B:25:GLU:O	1:B:28:SER:HB3	1.92	0.70
1:B:69:GLU:HA	1:B:72:SER:HB3	1.74	0.69
1:B:174:THR:HB	1:B:177:ASN:ND2	2.09	0.68
1:B:93:VAL:HG22	1:B:132:TYR:HB2	1.75	0.68
1:A:174:THR:HB	1:A:177:ASN:ND2	2.09	0.68
1:A:71:ILE:HG22	1:A:79:ILE:HD13	1.75	0.67
1:B:3:LEU:HD12	1:B:4:PRO:HD2	1.76	0.67
1:A:21:VAL:HG22	1:A:31:ILE:HD11	1.75	0.67
1:B:65:GLU:HB2	3:B:227:HOH:O	1.95	0.67
1:B:23:PRO:HA	1:B:26:ILE:HG13	1.79	0.65
1:A:11:ASP:OD2	1:A:13:VAL:HG12	1.97	0.65
1:A:108:PHE:HE1	1:A:189:LEU:HD11	1.62	0.64
1:B:112:LEU:CD1	1:B:112:LEU:N	2.61	0.64
1:A:23:PRO:HA	1:A:26:ILE:HG13	1.80	0.64
1:B:116:GLY:C	1:B:175:VAL:HG13	2.19	0.62
1:B:11:ASP:OD2	1:B:13:VAL:HG12	1.98	0.62
1:B:84:ARG:HG3	1:B:146:TRP:CB	2.29	0.62
1:B:84:ARG:HG3	1:B:146:TRP:CG	2.34	0.62
1:A:38:MET:HE1	1:A:61:ILE:HB	1.81	0.62
1:B:35:ILE:O	1:B:39:ILE:HG13	2.00	0.62
1:B:65:GLU:HB3	1:B:89:LEU:HA	1.81	0.62
1:B:84:ARG:HG3	1:B:146:TRP:CD1	2.35	0.62
1:B:35:ILE:HD13	1:B:91:VAL:HG11	1.82	0.61
1:A:3:LEU:HD12	1:A:4:PRO:HD2	1.82	0.61
1:A:97:LEU:HG	1:A:127:VAL:HG21	1.81	0.61
1:A:174:THR:HG22	1:A:176:ASP:H	1.64	0.61
1:B:10:GLY:N	1:B:15:HIS:HE1	1.98	0.61
1:A:134:ARG:HH11	1:A:134:ARG:CG	2.14	0.60
1:B:138:ARG:HG2	1:B:138:ARG:NH1	2.14	0.60
1:A:55:ILE:HG13	1:A:57:VAL:HG23	1.81	0.60
1:A:10:GLY:N	1:A:15:HIS:HE1	1.98	0.60
1:A:32:GLN:HA	1:A:32:GLN:NE2	2.15	0.60
1:B:55:ILE:HG13	1:B:57:VAL:HG23	1.83	0.60
1:B:174:THR:HB	1:B:177:ASN:HD22	1.66	0.60
1:B:32:GLN:HE21	1:B:32:GLN:HA	1.66	0.60
1:B:174:THR:HG22	1:B:176:ASP:H	1.67	0.60
1:A:52:ALA:HB3	1:A:53:PRO:HD3	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:LEU:O	1:A:46:PRO:HD3	2.01	0.59
1:A:174:THR:HB	1:A:177:ASN:HD22	1.66	0.59
1:A:98:LYS:HG3	1:A:99:GLU:OE1	2.02	0.59
1:B:21:VAL:CG2	1:B:59:LEU:HD12	2.30	0.59
1:A:169:PRO:C	1:A:170:ARG:HG2	2.22	0.59
1:B:3:LEU:CD1	1:B:4:PRO:HD2	2.32	0.59
1:A:112:LEU:HD12	1:A:112:LEU:H	1.67	0.59
1:A:35:ILE:O	1:A:39:ILE:HG13	2.03	0.58
1:A:32:GLN:OE1	1:A:134:ARG:NH1	2.36	0.58
1:A:156:ASP:OD2	1:A:166:LYS:NZ	2.37	0.58
1:B:134:ARG:HH11	1:B:134:ARG:CG	2.17	0.57
1:B:44:LEU:O	1:B:46:PRO:HD3	2.04	0.57
1:B:66:ASP:OD2	1:B:84:ARG:HD3	2.04	0.57
1:A:116:GLY:C	1:A:175:VAL:HG13	2.25	0.57
1:A:127:VAL:HG22	1:A:128:VAL:N	2.19	0.57
1:A:118:ARG:NH1	1:A:175:VAL:HG12	2.20	0.56
1:A:137:LYS:C	1:A:138:ARG:HA	2.18	0.56
1:B:127:VAL:HG22	1:B:128:VAL:N	2.21	0.56
1:B:52:ALA:HB3	1:B:53:PRO:HD3	1.86	0.56
1:A:175:VAL:HG23	1:A:176:ASP:N	2.20	0.56
1:A:71:ILE:HG12	1:A:84:ARG:HH11	1.71	0.56
1:B:97:LEU:HG	1:B:127:VAL:HG21	1.88	0.56
1:A:65:GLU:HB3	1:A:89:LEU:HA	1.88	0.55
1:A:3:LEU:CD1	1:A:4:PRO:HD2	2.36	0.55
1:B:99:GLU:OE1	1:B:99:GLU:N	2.40	0.55
1:B:175:VAL:HG23	1:B:176:ASP:N	2.22	0.55
1:B:177:ASN:OD1	3:B:211:HOH:O	2.18	0.55
1:A:177:ASN:HB3	1:A:180:LEU:CD1	2.28	0.55
1:B:106:LEU:CD2	1:B:120:ALA:HB1	2.37	0.54
1:A:106:LEU:HD22	1:A:189:LEU:HD23	1.89	0.54
1:B:32:GLN:NE2	1:B:32:GLN:HA	2.22	0.54
1:A:66:ASP:HB2	1:A:87:PHE:CE1	2.43	0.53
1:B:114:VAL:O	1:B:114:VAL:HG12	2.08	0.53
1:B:148:ALA:O	1:B:152:GLN:HG3	2.08	0.53
1:B:72:SER:OG	1:B:73:TYR:N	2.41	0.53
1:B:135:GLN:HE21	1:B:135:GLN:CA	2.20	0.53
1:A:69:GLU:HA	1:A:72:SER:HB2	1.90	0.53
1:B:38:MET:HE1	1:B:61:ILE:O	2.08	0.53
1:B:177:ASN:HB3	1:B:180:LEU:CD1	2.29	0.52
1:B:66:ASP:HB2	1:B:87:PHE:CE1	2.44	0.52
1:A:21:VAL:HG21	1:A:31:ILE:HD11	1.89	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:PHE:CE1	1:A:189:LEU:HD11	2.43	0.52
1:A:106:LEU:CD2	1:A:120:ALA:HB1	2.39	0.52
1:B:134:ARG:HH11	1:B:134:ARG:HG2	1.73	0.52
1:A:87:PHE:CE1	1:A:147:GLN:HG3	2.45	0.52
1:B:146:TRP:O	1:B:150:ILE:HG13	2.10	0.52
1:A:25:GLU:HA	1:A:25:GLU:OE1	2.10	0.51
1:A:106:LEU:HD22	1:A:189:LEU:CD2	2.40	0.51
1:B:125:LEU:O	1:B:144:SER:HA	2.09	0.51
1:B:21:VAL:CG2	1:B:31:ILE:CD1	2.88	0.51
1:A:114:VAL:O	1:A:114:VAL:HG12	2.10	0.51
1:B:112:LEU:HD12	1:B:112:LEU:H	1.74	0.51
1:A:32:GLN:HE22	1:A:35:ILE:HD12	1.76	0.51
1:B:87:PHE:CE1	1:B:147:GLN:HG3	2.46	0.51
1:B:39:ILE:HD13	1:B:89:LEU:HD21	1.92	0.51
1:A:146:TRP:O	1:A:150:ILE:HG13	2.11	0.50
1:A:125:LEU:O	1:A:144:SER:HA	2.12	0.50
1:B:66:ASP:HB3	1:B:86:HIS:HA	1.92	0.50
1:B:84:ARG:HG3	1:B:146:TRP:HB3	1.92	0.50
1:B:63:VAL:HG12	1:B:91:VAL:HG22	1.93	0.50
1:A:97:LEU:HG	1:A:127:VAL:CG2	2.40	0.50
1:A:65:GLU:HA	1:A:87:PHE:CE2	2.46	0.50
1:B:112:LEU:H	1:B:112:LEU:CD1	2.24	0.50
1:B:21:VAL:HG21	1:B:31:ILE:HD13	1.94	0.49
1:A:148:ALA:O	1:A:152:GLN:HG3	2.11	0.49
1:A:175:VAL:HG23	1:A:176:ASP:H	1.76	0.49
1:B:175:VAL:HG23	1:B:176:ASP:H	1.77	0.49
1:A:52:ALA:HB1	1:A:57:VAL:HB	1.94	0.49
1:B:3:LEU:HD21	1:B:41:VAL:HG22	1.95	0.49
1:A:30:ARG:NH1	1:A:30:ARG:CG	2.64	0.48
1:B:131:GLY:O	1:B:139:ILE:HG12	2.13	0.48
1:B:73:TYR:CD1	1:B:73:TYR:N	2.61	0.48
1:B:157:HIS:CE1	1:B:162:LEU:HD21	2.48	0.48
1:A:63:VAL:HG12	1:A:91:VAL:HG22	1.95	0.48
1:A:62:ILE:HG23	1:A:158:LEU:HD11	1.96	0.48
1:B:109:GLU:HG3	1:B:149:ARG:HH11	1.78	0.48
1:A:68:LYS:HG2	1:A:86:HIS:CG	2.48	0.48
1:B:42:MET:HE2	1:B:63:VAL:CG2	2.44	0.48
1:B:64:LEU:HD12	1:B:87:PHE:HE2	1.78	0.48
1:B:153:HIS:C	1:B:153:HIS:CD2	2.87	0.48
1:B:52:ALA:HB1	1:B:57:VAL:HB	1.96	0.47
1:A:75:PRO:HB2	1:A:78:GLU:HB3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:128:VAL:HG13	1:B:128:VAL:O	2.14	0.47
1:A:15:HIS:HB3	1:B:170:ARG:HB3	1.95	0.47
1:B:3:LEU:CG	1:B:4:PRO:HD2	2.44	0.47
1:A:21:VAL:HG21	1:A:31:ILE:CD1	2.45	0.47
1:B:21:VAL:HG22	1:B:31:ILE:CD1	2.44	0.47
1:A:66:ASP:HB3	1:A:86:HIS:HA	1.95	0.47
1:A:66:ASP:O	1:A:86:HIS:HB2	2.15	0.47
1:B:106:LEU:HD22	1:B:189:LEU:HD23	1.97	0.47
1:A:22:ASP:OD2	1:A:25:GLU:HG2	2.14	0.47
1:B:21:VAL:HG22	1:B:31:ILE:HD11	1.96	0.46
1:A:3:LEU:HG	1:A:4:PRO:HD2	1.97	0.46
1:A:99:GLU:O	1:A:99:GLU:HG2	2.15	0.46
1:A:134:ARG:CG	1:A:134:ARG:NH1	2.73	0.46
1:A:128:VAL:O	1:A:128:VAL:HG13	2.14	0.46
1:A:82:GLN:O	1:A:83:GLU:HB2	2.15	0.46
1:B:21:VAL:CG2	1:B:31:ILE:HD11	2.45	0.46
1:A:21:VAL:CG2	1:A:31:ILE:CD1	2.89	0.46
1:A:154:GLU:OE2	3:A:199:HOH:O	2.21	0.46
1:B:62:ILE:HG23	1:B:158:LEU:HD11	1.98	0.46
1:B:21:VAL:CG2	1:B:31:ILE:HD13	2.46	0.46
1:A:156:ASP:CG	1:A:166:LYS:NZ	2.70	0.46
1:A:111:CYS:SG	1:A:157:HIS:HE1	2.38	0.46
1:A:112:LEU:N	1:A:112:LEU:HD12	2.31	0.45
1:B:64:LEU:HD12	1:B:87:PHE:CE2	2.51	0.45
1:B:82:GLN:O	1:B:83:GLU:HB2	2.17	0.45
1:A:168:VAL:O	1:A:169:PRO:C	2.53	0.45
1:B:156:ASP:OD2	1:B:166:LYS:NZ	2.49	0.45
1:B:22:ASP:OD2	1:B:25:GLU:HG2	2.17	0.45
1:A:173:ARG:HA	1:B:9:SER:HB2	1.98	0.45
1:A:3:LEU:CG	1:A:4:PRO:HD2	2.47	0.45
1:A:162:LEU:HD23	1:A:162:LEU:HA	1.77	0.45
1:A:26:ILE:HD13	1:A:93:VAL:HG21	1.99	0.45
1:B:134:ARG:CG	1:B:134:ARG:NH1	2.75	0.44
1:A:168:VAL:HA	1:A:169:PRO:HD3	1.78	0.44
1:B:132:TYR:N	1:B:132:TYR:CD2	2.85	0.44
1:A:42:MET:HE2	1:A:63:VAL:CG2	2.48	0.44
1:B:97:LEU:C	1:B:97:LEU:HD23	2.38	0.44
1:A:154:GLU:OE1	1:A:154:GLU:HA	2.18	0.44
1:B:65:GLU:HA	1:B:87:PHE:CE2	2.53	0.44
1:A:26:ILE:HD12	1:A:136:GLY:HA2	2.00	0.44
1:A:71:ILE:O	1:A:72:SER:C	2.56	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:ASP:O	1:B:86:HIS:HB2	2.19	0.43
1:B:3:LEU:HG	1:B:4:PRO:HD2	1.98	0.43
1:B:97:LEU:HG	1:B:127:VAL:CG2	2.47	0.43
1:B:106:LEU:HD21	1:B:120:ALA:HB1	2.01	0.43
1:A:20:GLU:HA	1:A:60:ARG:HG3	2.00	0.43
1:B:88:ASP:O	1:B:89:LEU:C	2.56	0.43
1:B:184:GLU:HG3	1:B:184:GLU:O	2.14	0.43
1:B:32:GLN:HE21	1:B:32:GLN:CA	2.29	0.43
1:A:97:LEU:C	1:A:97:LEU:HD23	2.39	0.43
1:B:133:ASP:C	1:B:133:ASP:OD1	2.57	0.43
1:B:89:LEU:HD12	1:B:90:MET:H	1.84	0.43
1:A:137:LYS:C	1:A:137:LYS:CB	2.75	0.43
1:A:30:ARG:HG2	1:A:31:ILE:N	2.34	0.42
1:B:23:PRO:HA	1:B:26:ILE:CG1	2.48	0.42
1:A:174:THR:O	1:A:175:VAL:C	2.58	0.42
1:A:88:ASP:O	1:A:89:LEU:C	2.58	0.42
1:A:84:ARG:HG3	1:A:146:TRP:CD1	2.55	0.42
1:B:94:ASN:N	1:B:95:PRO:CD	2.82	0.42
1:A:65:GLU:CB	1:A:89:LEU:HA	2.50	0.41
1:B:26:ILE:HD12	1:B:136:GLY:HA2	2.01	0.41
1:B:36:ASP:OD1	1:B:134:ARG:NH2	2.53	0.41
1:A:23:PRO:HA	1:A:26:ILE:CG1	2.49	0.41
1:B:26:ILE:HD13	1:B:93:VAL:HG21	2.02	0.41
1:A:173:ARG:CD	1:A:178:LEU:HA	2.50	0.41
1:B:127:VAL:HG22	1:B:128:VAL:H	1.86	0.41
1:A:73:TYR:O	1:A:74:ALA:HB2	2.19	0.41
1:A:94:ASN:N	1:A:95:PRO:CD	2.83	0.41
1:A:135:GLN:H	1:A:135:GLN:NE2	2.18	0.41
1:B:71:ILE:O	1:B:72:SER:C	2.59	0.41
1:B:71:ILE:O	1:B:79:ILE:HD11	2.21	0.41
1:A:89:LEU:HD12	1:A:90:MET:H	1.84	0.41
1:B:65:GLU:CB	1:B:89:LEU:HA	2.48	0.41
1:A:173:ARG:HA	1:B:9:SER:CB	2.50	0.41
1:A:191:SER:HB2	1:A:192:HIS:H	1.55	0.41
1:A:74:ALA:HB3	1:A:79:ILE:HD11	2.03	0.41
1:B:157:HIS:CE1	1:B:162:LEU:CD2	3.03	0.41
1:B:173:ARG:CD	1:B:178:LEU:HA	2.51	0.41
1:B:97:LEU:CG	1:B:127:VAL:HG21	2.51	0.41
1:B:174:THR:O	1:B:175:VAL:C	2.60	0.40
1:B:162:LEU:HA	1:B:162:LEU:HD23	1.75	0.40
1:A:174:THR:HG22	1:A:176:ASP:N	2.33	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:170:ARG:HB2	1:B:15:HIS:HB3	2.02	0.40
1:A:127:VAL:CG2	1:A:128:VAL:N	2.84	0.40
1:B:13:VAL:HG13	1:B:14:LEU:N	2.36	0.40
1:A:173:ARG:HD3	1:A:178:LEU:HA	2.04	0.40
1:A:106:LEU:HD21	1:A:120:ALA:HB1	2.03	0.40
1:B:164:VAL:O	1:B:167:MET:HG2	2.22	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	189/197 (96%)	159 (84%)	28 (15%)	2 (1%)	17	51
1	B	189/197 (96%)	159 (84%)	25 (13%)	5 (3%)	7	26
All	All	378/394 (96%)	318 (84%)	53 (14%)	7 (2%)	10	35

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	44	LEU
1	B	44	LEU
1	A	46	PRO
1	B	46	PRO
1	B	72	SER
1	B	134	ARG
1	B	114	VAL

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	157/170 (92%)	143 (91%)	14 (9%)	12	35
1	B	154/170 (91%)	135 (88%)	19 (12%)	6	17
All	All	311/340 (92%)	278 (89%)	33 (11%)	8	25

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	VAL
1	A	31	ILE
1	A	37	ASP
1	A	50	LEU
1	A	59	LEU
1	A	66	ASP
1	A	85	ARG
1	A	93	VAL
1	A	98	LYS
1	A	115	ASP
1	A	121	VAL
1	A	126	GLU
1	A	135	GLN
1	A	182	LEU
1	B	21	VAL
1	B	37	ASP
1	B	50	LEU
1	B	59	LEU
1	B	64	LEU
1	B	66	ASP
1	B	73	TYR
1	B	85	ARG
1	B	93	VAL
1	B	102	ASN
1	B	115	ASP
1	B	121	VAL
1	B	126	GLU

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Mol	Chain	Res	Type
1	B	134	ARG
1	B	135	GLN
1	B	138	ARG
1	B	170	ARG
1	B	182	LEU
1	B	184	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	HIS
1	A	135	GLN
1	A	147	GLN
1	A	177	ASN
1	A	192	HIS
1	B	15	HIS
1	B	32	GLN
1	B	135	GLN
1	B	147	GLN
1	B	177	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 are 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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	191/197 (96%)	0.22	1 (0%) 91 90	47, 64, 81, 95	0
1	B	191/197 (96%)	0.31	4 (2%) 67 62	47, 68, 85, 98	0
All	All	382/394 (96%)	0.27	5 (1%) 79 78	47, 66, 83, 98	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	135	GLN	2.6
1	A	73	TYR	2.3
1	B	191	SER	2.3
1	B	125	LEU	2.1
1	B	146	TRP	2.1

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	ZN	A	198	1/1	0.98	0.14	-1.52	60,60,60,60	0
2	ZN	B	198	1/1	0.96	0.12	-2.18	64,64,64,64	0

6.5 Other polymers [i](#)

There are no such residues in this entry.