



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:20 PM GMT

PDB ID : 4OIW
Title : Structural and kinetic bases for the metal preference of the M18 aminopeptidase from *Pseudomonas aeruginosa*
Authors : Nguyen, D.D.; Pandian, R.; Kim, D.D.; Ha, S.C.; Yoon, H.J.; Kim, K.S.; Yun, K.H.; Kim, J.H.; Kim, K.K.
Deposited on : 2014-01-20
Resolution : 2.44 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

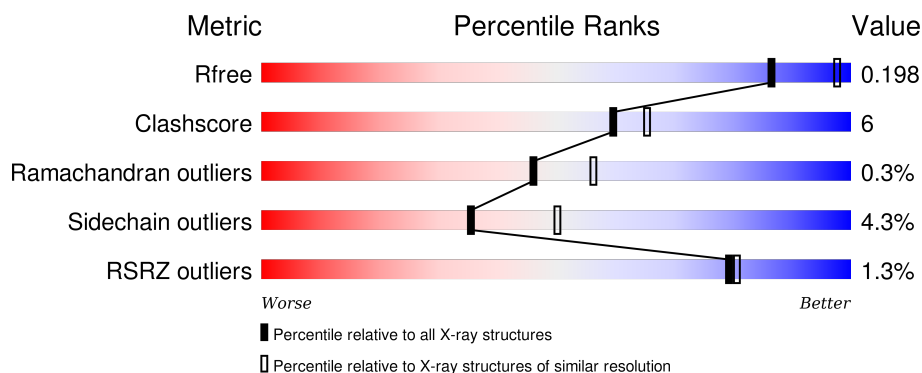
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.44 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	1003 (2.46-2.42)
Clashscore	102246	1071 (2.46-2.42)
Ramachandran outliers	100387	1065 (2.46-2.42)
Sidechain outliers	100360	1065 (2.46-2.42)
RSRZ outliers	91569	1005 (2.46-2.42)

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	431	<div> <div> <div>2%</div> <div>81%</div> <div>11%</div> <div>5%</div> </div> </div>
1	B	431	<div> <div> <div>2%</div> <div>83%</div> <div>9%</div> <div>5%</div> </div> </div>
1	C	431	<div> <div> <div>2%</div> <div>81%</div> <div>12%</div> <div>5%</div> </div> </div>
1	D	431	<div> <div> <div>2%</div> <div>84%</div> <div>10%</div> <div>5%</div> </div> </div>
1	E	431	<div> <div> <div>2%</div> <div>83%</div> <div>10%</div> <div>5%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	431	<div><div><div>%</div><div><div></div></div><div>81%</div><div>13%</div><div>• 5%</div></div></div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 19832 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 Probable M18 family aminopeptidase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	408	Total	C	N	O	S	0	0	0
			3156	1980	577	591	8			
1	B	408	Total	C	N	O	S	0	0	0
			3156	1980	577	591	8			
1	C	408	Total	C	N	O	S	0	0	0
			3156	1980	577	591	8			
1	D	408	Total	C	N	O	S	0	0	0
			3156	1980	577	591	8			
1	E	408	Total	C	N	O	S	0	0	0
			3156	1980	577	591	8			
1	F	408	Total	C	N	O	S	0	0	0
			3156	1980	577	591	8			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP Q9HYZ3
A	0	SER	-	EXPRESSION TAG	UNP Q9HYZ3
A	82	ALA	HIS	ENGINEERED MUTATION	UNP Q9HYZ3
B	-1	GLY	-	EXPRESSION TAG	UNP Q9HYZ3
B	0	SER	-	EXPRESSION TAG	UNP Q9HYZ3
B	82	ALA	HIS	ENGINEERED MUTATION	UNP Q9HYZ3
C	-1	GLY	-	EXPRESSION TAG	UNP Q9HYZ3
C	0	SER	-	EXPRESSION TAG	UNP Q9HYZ3
C	82	ALA	HIS	ENGINEERED MUTATION	UNP Q9HYZ3
D	-1	GLY	-	EXPRESSION TAG	UNP Q9HYZ3
D	0	SER	-	EXPRESSION TAG	UNP Q9HYZ3
D	82	ALA	HIS	ENGINEERED MUTATION	UNP Q9HYZ3
E	-1	GLY	-	EXPRESSION TAG	UNP Q9HYZ3
E	0	SER	-	EXPRESSION TAG	UNP Q9HYZ3
E	82	ALA	HIS	ENGINEERED MUTATION	UNP Q9HYZ3
F	-1	GLY	-	EXPRESSION TAG	UNP Q9HYZ3
F	0	SER	-	EXPRESSION TAG	UNP Q9HYZ3

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Chain	Residue	Modelled	Actual	Comment	Reference
F	82	ALA	HIS	ENGINEERED MUTATION	UNP Q9HYZ3

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	1	Total Zn 1 1	0	0
2	E	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	A	1	Total Zn 1 1	0	0
2	F	1	Total Zn 1 1	0	0

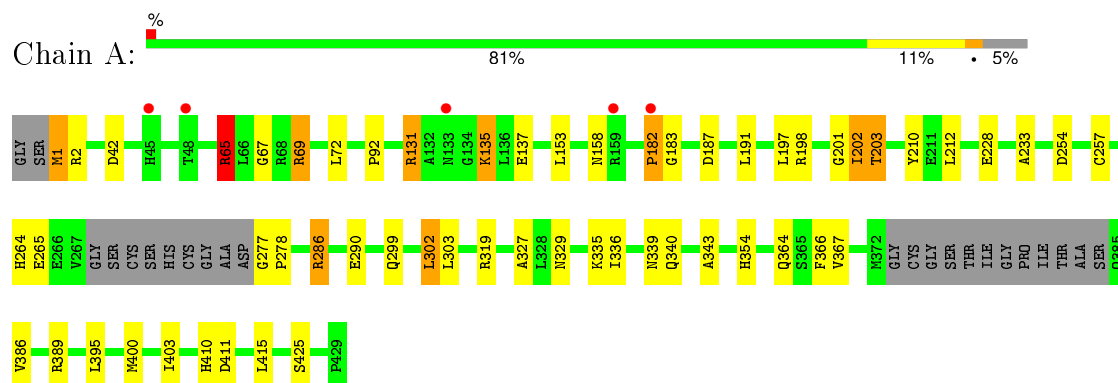
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	157	Total O 157 157	0	0
3	B	172	Total O 172 172	0	0
3	C	171	Total O 171 171	0	0
3	D	134	Total O 134 134	0	0
3	E	137	Total O 137 137	0	0
3	F	119	Total O 119 119	0	0

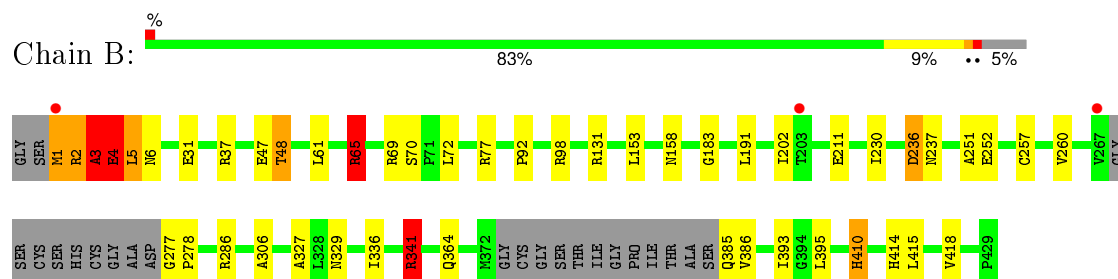
3 Residue-property plots

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.

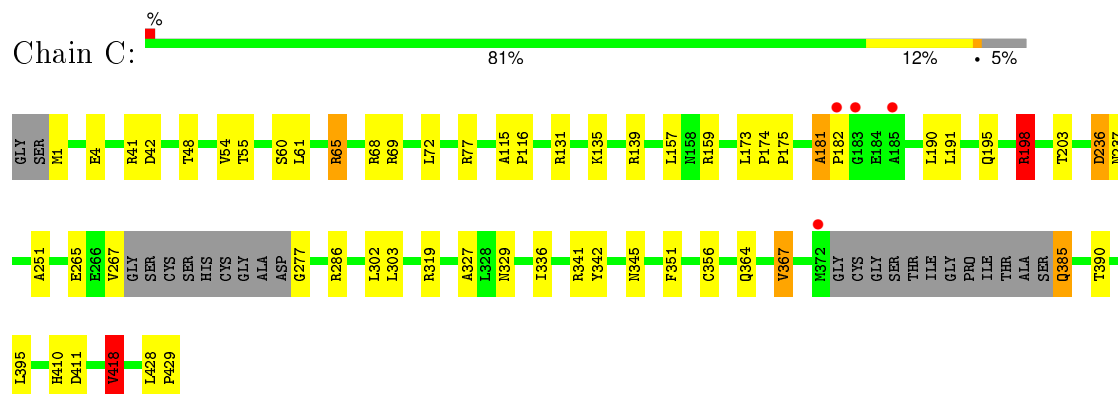
- Molecule 1: Probable M18 family aminopeptidase 2



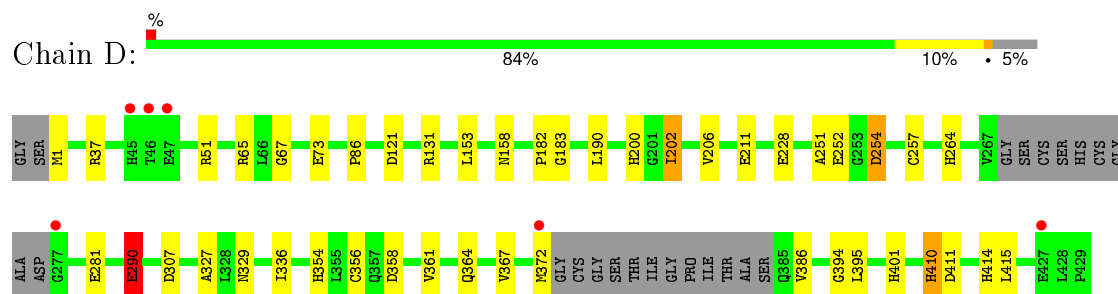
- Molecule 1: Probable M18 family aminopeptidase 2



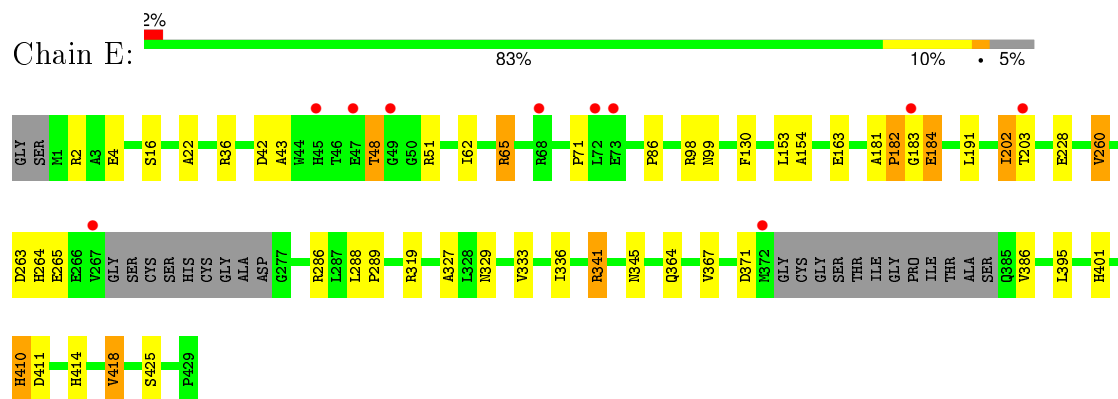
- Molecule 1: Probable M18 family aminopeptidase 2



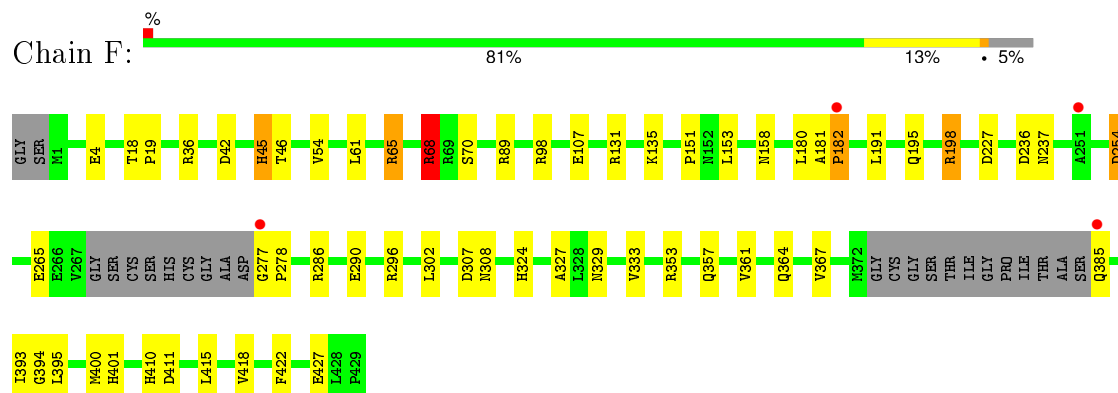
- Molecule 1: Probable M18 family aminopeptidase 2



- Molecule 1: Probable M18 family aminopeptidase 2



- Molecule 1: Probable M18 family aminopeptidase 2



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	150.02Å 218.27Å 172.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.79 – 2.44 48.74 – 2.44	Depositor EDS
% Data completeness (in resolution range)	96.0 (48.79-2.44) 96.0 (48.74-2.44)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.70 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.148 , 0.197 0.156 , 0.198	Depositor DCC
R_{free} test set	5026 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	39.8	Xtriage
Anisotropy	0.117	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 36.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 100620 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	19832	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.18% 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, 0.375 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	0.78	1/3220 (0.0%)	0.93	8/4372 (0.2%)
1	B	0.82	0/3220	0.97	10/4372 (0.2%)
1	C	0.77	0/3220	0.91	6/4372 (0.1%)
1	D	0.75	1/3220 (0.0%)	0.90	3/4372 (0.1%)
1	E	0.74	1/3220 (0.0%)	0.89	5/4372 (0.1%)
1	F	0.74	0/3220	0.90	6/4372 (0.1%)
All	All	0.77	3/19320 (0.0%)	0.92	38/26232 (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
1	A	0	3
1	D	0	1
1	E	0	1
All	All	0	5

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	228	GLU	CD-OE1	7.37	1.33	1.25
1	D	228	GLU	CD-OE1	5.81	1.32	1.25
1	A	228	GLU	CD-OE1	5.33	1.31	1.25

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	65	ARG	NE-CZ-NH1	8.39	124.49	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	68	ARG	NE-CZ-NH1	8.04	124.32	120.30
1	A	65	ARG	NE-CZ-NH2	-7.88	116.36	120.30
1	C	139	ARG	NE-CZ-NH2	-7.81	116.39	120.30
1	D	254	ASP	CB-CG-OD1	7.62	125.15	118.30
1	B	341	ARG	NE-CZ-NH1	6.99	123.80	120.30
1	F	68	ARG	NE-CZ-NH2	-6.59	117.01	120.30
1	A	198	ARG	NE-CZ-NH1	6.42	123.51	120.30
1	B	183	GLY	N-CA-C	-6.30	97.34	113.10
1	B	3	ALA	C-N-CA	6.08	136.91	121.70
1	A	69	ARG	NE-CZ-NH1	6.06	123.33	120.30
1	F	254	ASP	CB-CG-OD1	6.04	123.74	118.30
1	B	69	ARG	NE-CZ-NH1	5.89	123.25	120.30
1	D	183	GLY	N-CA-C	-5.83	98.53	113.10
1	C	418	VAL	CB-CA-C	-5.81	100.36	111.40
1	E	418	VAL	CB-CA-C	-5.67	100.62	111.40
1	C	69	ARG	NE-CZ-NH1	5.63	123.11	120.30
1	C	236	ASP	CB-CG-OD2	-5.46	113.38	118.30
1	F	198	ARG	NE-CZ-NH1	5.43	123.02	120.30
1	B	286	ARG	NE-CZ-NH2	-5.42	117.59	120.30
1	B	260	VAL	CB-CA-C	-5.35	101.24	111.40
1	B	65	ARG	NE-CZ-NH2	-5.34	117.63	120.30
1	D	121	ASP	CB-CG-OD1	5.30	123.07	118.30
1	A	187	ASP	CB-CG-OD1	5.29	123.06	118.30
1	B	236	ASP	CB-CG-OD1	5.27	123.04	118.30
1	E	65	ARG	CG-CD-NE	5.19	122.70	111.80
1	F	68	ARG	CG-CD-NE	5.19	122.69	111.80
1	A	131	ARG	NE-CZ-NH1	5.18	122.89	120.30
1	A	286	ARG	NE-CZ-NH2	-5.11	117.75	120.30
1	F	227	ASP	CB-CG-OD1	5.10	122.89	118.30
1	E	260	VAL	CB-CA-C	-5.10	101.72	111.40
1	E	371	ASP	CB-CG-OD1	5.08	122.87	118.30
1	C	198	ARG	NE-CZ-NH1	5.06	122.83	120.30
1	B	48	THR	C-N-CA	-5.06	111.68	122.30
1	A	65	ARG	CG-CD-NE	5.06	122.42	111.80
1	B	37	ARG	NE-CZ-NH1	5.04	122.82	120.30
1	E	51	ARG	NE-CZ-NH1	5.03	122.81	120.30
1	C	367	VAL	CB-CA-C	-5.00	101.90	111.40

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1	MET	Peptide
1	A	182	PRO	Peptide
1	A	366	PHE	Peptide
1	D	182	PRO	Peptide
1	E	182	PRO	Peptide

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	3156	0	3099	36	0
1	B	3156	0	3099	61	0
1	C	3156	0	3099	40	0
1	D	3156	0	3099	27	0
1	E	3156	0	3099	26	0
1	F	3156	0	3099	38	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	157	0	0	5	0
3	B	172	0	0	9	0
3	C	171	0	0	9	0
3	D	134	0	0	6	0
3	E	137	0	0	4	0
3	F	119	0	0	9	0
All	All	19832	0	18594	214	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 (214) 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:2:ARG:HB3	1:B:3:ALA:CB	1.58	1.31

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2:ARG:CG	1:B:3:ALA:HA	1.79	1.11
1:B:2:ARG:CB	1:B:3:ALA:HB2	1.82	1.07
1:B:4:GLU:H	1:B:5:LEU:CB	1.81	0.94
1:B:4:GLU:H	1:B:5:LEU:HB3	1.32	0.93
1:B:327:ALA:H	1:B:364:GLN:HE22	1.13	0.91
1:A:327:ALA:H	1:A:364:GLN:HE22	1.11	0.90
1:B:2:ARG:HG2	1:B:3:ALA:HA	1.52	0.89
1:A:265:GLU:OE1	3:A:621:HOH:O	1.90	0.87
1:B:2:ARG:HB3	1:B:3:ALA:HB2	0.89	0.87
1:B:2:ARG:CD	1:B:3:ALA:HA	2.05	0.84
1:C:327:ALA:H	1:C:364:GLN:HE22	1.22	0.84
1:D:327:ALA:H	1:D:364:GLN:HE22	1.23	0.83
1:B:2:ARG:CB	1:B:3:ALA:CB	2.50	0.81
1:F:265:GLU:OE1	3:F:610:HOH:O	1.97	0.81
1:B:2:ARG:CB	1:B:3:ALA:CA	2.58	0.81
1:B:2:ARG:HB3	1:B:3:ALA:CA	2.09	0.80
1:E:265:GLU:OE1	3:E:633:HOH:O	1.98	0.80
1:A:131:ARG:CD	3:B:653:HOH:O	2.32	0.77
1:A:131:ARG:HD3	3:B:653:HOH:O	1.83	0.77
1:C:265:GLU:OE2	3:C:625:HOH:O	2.03	0.77
1:B:6:ASN:OD1	1:B:230:ILE:HD11	1.85	0.76
1:E:36:ARG:NH2	3:E:690:HOH:O	2.19	0.75
1:B:3:ALA:HB3	1:B:4:GLU:OE1	1.86	0.75
1:B:4:GLU:N	1:B:5:LEU:CB	2.50	0.75
1:F:327:ALA:H	1:F:364:GLN:HE22	1.33	0.73
1:B:1:MET:O	1:B:3:ALA:HB2	1.89	0.72
1:F:70:SER:HB2	3:F:661:HOH:O	1.88	0.71
1:C:181:ALA:HB1	1:C:182:PRO:HD2	1.70	0.71
1:B:2:ARG:CB	1:B:3:ALA:HA	2.20	0.71
1:C:181:ALA:HB1	1:C:182:PRO:CD	2.20	0.70
1:C:329:ASN:ND2	1:C:410:HIS:H	1.89	0.70
1:E:327:ALA:H	1:E:364:GLN:HE22	1.36	0.70
1:F:195:GLN:OE1	1:F:198:ARG:NH1	2.23	0.70
1:B:4:GLU:N	1:B:5:LEU:HB2	2.06	0.70
1:B:6:ASN:CG	1:B:230:ILE:HD11	2.12	0.69
1:B:2:ARG:CG	1:B:3:ALA:CA	2.64	0.69
1:B:65:ARG:NH1	3:B:632:HOH:O	2.27	0.67
1:A:299:GLN:O	1:A:389:ARG:NH2	2.29	0.66
1:C:65:ARG:NH2	1:C:251:ALA:O	2.26	0.65
1:B:277:GLY:N	1:B:278:PRO:CD	2.60	0.64
1:B:4:GLU:N	1:B:4:GLU:CD	2.52	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:395:LEU:HD22	1:E:411:ASP:HB3	1.80	0.63
1:B:3:ALA:CB	1:B:4:GLU:OE1	2.46	0.63
1:B:2:ARG:HD3	1:B:3:ALA:HA	1.81	0.62
1:A:339:ASN:HA	1:C:267:VAL:HG11	1.81	0.62
1:B:341:ARG:HH11	1:B:341:ARG:HG3	1.65	0.61
1:F:153:LEU:HD23	1:F:158:ASN:HB2	1.83	0.61
1:A:354:HIS:CE1	1:C:131:ARG:HD2	2.36	0.60
1:F:395:LEU:HD22	1:F:411:ASP:HB3	1.83	0.60
1:B:1:MET:O	1:B:3:ALA:CB	2.49	0.60
1:C:181:ALA:CB	1:C:182:PRO:HD2	2.31	0.60
1:F:98:ARG:HD3	3:F:710:HOH:O	2.02	0.60
1:A:277:GLY:N	1:A:278:PRO:HD2	2.16	0.59
1:A:135:LYS:HE3	1:A:137:GLU:OE2	2.01	0.59
1:D:65:ARG:NH2	1:D:251:ALA:O	2.35	0.59
1:A:131:ARG:HD2	3:B:653:HOH:O	2.00	0.59
1:A:395:LEU:HD22	1:A:411:ASP:HB3	1.85	0.58
1:D:65:ARG:HB3	1:D:257:CYS:HB2	1.84	0.58
1:D:67:GLY:HA3	1:D:254:ASP:O	2.03	0.58
1:D:395:LEU:HD22	1:D:411:ASP:HB3	1.85	0.58
1:D:329:ASN:ND2	1:D:410:HIS:H	2.01	0.58
1:F:70:SER:CB	3:F:661:HOH:O	2.47	0.57
1:A:69:ARG:HD3	3:A:717:HOH:O	2.05	0.57
1:F:361:VAL:HG21	1:F:418:VAL:HG13	1.86	0.57
1:F:308:ASN:HB2	1:F:400:MET:CE	2.35	0.57
1:B:277:GLY:N	1:B:278:PRO:HD2	2.19	0.56
1:D:254:ASP:HB3	3:D:653:HOH:O	2.05	0.56
1:E:329:ASN:ND2	1:E:410:HIS:H	2.02	0.56
1:B:131:ARG:HG2	3:B:757:HOH:O	2.05	0.56
1:F:308:ASN:HB2	1:F:400:MET:HE3	1.86	0.56
1:F:65:ARG:NH1	3:F:717:HOH:O	2.37	0.56
1:C:265:GLU:OE2	3:C:616:HOH:O	2.18	0.56
1:A:290:GLU:O	3:A:625:HOH:O	2.17	0.56
1:C:41:ARG:HD2	3:C:672:HOH:O	2.06	0.55
1:A:329:ASN:ND2	1:A:410:HIS:H	2.05	0.55
1:A:197:LEU:O	1:A:201:GLY:HA2	2.07	0.55
1:B:4:GLU:N	1:B:5:LEU:HB3	2.11	0.55
1:B:4:GLU:HA	1:B:6:ASN:H	1.72	0.55
1:B:70:SER:HB2	3:B:662:HOH:O	2.06	0.55
1:C:329:ASN:HD21	1:C:410:HIS:H	1.55	0.54
1:D:414:HIS:HE1	3:D:660:HOH:O	1.90	0.54
1:B:2:ARG:HD3	1:B:3:ALA:CA	2.36	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:329:ASN:ND2	1:B:410:HIS:H	2.05	0.54
1:A:264:HIS:HE1	3:A:651:HOH:O	1.90	0.53
1:F:329:ASN:ND2	1:F:410:HIS:H	2.06	0.53
1:B:3:ALA:H	1:B:4:GLU:HB2	1.74	0.52
1:B:327:ALA:H	1:B:364:GLN:NE2	1.95	0.52
1:B:6:ASN:OD1	1:B:230:ILE:CD1	2.56	0.52
1:C:181:ALA:CB	1:C:182:PRO:CD	2.86	0.52
1:B:65:ARG:HB3	1:B:257:CYS:HB2	1.90	0.52
1:F:254:ASP:HB3	3:F:647:HOH:O	2.10	0.51
1:E:414:HIS:HD2	3:E:644:HOH:O	1.93	0.50
1:D:37:ARG:NH2	3:D:699:HOH:O	2.45	0.50
1:B:3:ALA:H	1:B:4:GLU:CB	2.25	0.50
1:B:3:ALA:HB3	1:B:4:GLU:HG2	1.94	0.50
1:B:4:GLU:H	1:B:5:LEU:HB2	1.61	0.50
1:C:159:ARG:NH2	3:C:767:HOH:O	2.45	0.49
1:F:400:MET:O	1:F:401:HIS:HB2	2.13	0.49
1:C:173:LEU:N	1:C:174:PRO:CD	2.76	0.48
1:D:354:HIS:CE1	1:F:131:ARG:HD3	2.48	0.48
1:A:277:GLY:N	1:A:278:PRO:CD	2.76	0.48
1:F:277:GLY:N	1:F:278:PRO:CD	2.76	0.48
1:F:36:ARG:HH12	1:F:45:HIS:CE1	2.31	0.48
1:C:68:ARG:HD2	3:C:692:HOH:O	2.14	0.48
1:A:72:LEU:O	1:A:72:LEU:HD23	2.13	0.48
1:C:157:LEU:O	1:E:319:ARG:NH1	2.47	0.48
1:B:131:ARG:CG	3:B:757:HOH:O	2.62	0.48
1:C:319:ARG:HG3	1:E:153:LEU:HD21	1.96	0.47
1:A:65:ARG:HB3	1:A:257:CYS:HB2	1.96	0.47
1:A:65:ARG:HG2	1:A:65:ARG:HH11	1.79	0.47
1:B:153:LEU:HD23	1:B:158:ASN:HB2	1.95	0.47
1:F:307:ASP:O	1:F:394:GLY:HA3	2.14	0.47
1:B:395:LEU:HD21	1:B:415:LEU:HB2	1.95	0.47
1:C:115:ALA:N	1:C:116:PRO:CD	2.78	0.47
1:A:354:HIS:ND1	1:C:131:ARG:HD2	2.30	0.47
1:F:181:ALA:HB1	1:F:182:PRO:HD2	1.97	0.47
1:B:211:GLU:OE2	1:C:345:ASN:HB2	2.13	0.47
1:C:356:CYS:SG	1:C:418:VAL:HG11	2.55	0.46
1:D:65:ARG:HD2	1:D:257:CYS:HB2	1.98	0.46
1:F:254:ASP:HA	3:F:717:HOH:O	2.14	0.46
1:D:153:LEU:HD23	1:D:158:ASN:HB2	1.97	0.46
1:C:385:GLN:HG3	3:C:650:HOH:O	2.15	0.46
1:F:18:THR:HB	1:F:19:PRO:HD2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:395:LEU:HD22	1:C:411:ASP:HB3	1.98	0.46
1:F:45:HIS:ND1	1:F:45:HIS:N	2.63	0.46
1:C:195:GLN:OE1	1:C:198:ARG:HD2	2.15	0.46
1:F:395:LEU:HD21	1:F:415:LEU:HB2	1.98	0.46
1:D:131:ARG:HG2	3:D:724:HOH:O	2.16	0.46
1:A:42:ASP:O	1:A:286:ARG:NH2	2.48	0.46
1:C:351:PHE:CD1	1:C:429:PRO:HD3	2.51	0.46
1:C:159:ARG:HD2	3:C:702:HOH:O	2.15	0.46
1:B:65:ARG:CG	1:B:65:ARG:HH11	2.29	0.45
1:D:264:HIS:CD2	1:E:341:ARG:HD3	2.51	0.45
1:F:89:ARG:HD2	3:F:618:HOH:O	2.16	0.45
3:D:650:HOH:O	1:F:131:ARG:CD	2.63	0.45
1:A:153:LEU:HD23	1:A:158:ASN:HB2	1.98	0.45
1:D:202:ILE:HD12	3:D:640:HOH:O	2.16	0.45
1:F:329:ASN:HD21	1:F:410:HIS:H	1.65	0.45
1:F:302:LEU:HD23	1:F:422:PHE:CE1	2.52	0.45
1:C:342:TYR:O	1:C:390:THR:HB	2.16	0.45
1:E:202:ILE:HD12	1:E:203:THR:N	2.31	0.45
1:D:51:ARG:CZ	1:D:65:ARG:HG2	2.46	0.45
1:A:65:ARG:CG	1:A:65:ARG:HH11	2.30	0.44
1:E:181:ALA:HB3	1:E:184:GLU:HG2	1.99	0.44
1:C:61:LEU:C	1:C:61:LEU:HD23	2.38	0.44
1:B:414:HIS:HD2	3:B:652:HOH:O	1.98	0.44
1:A:400:MET:HE1	3:A:630:HOH:O	2.16	0.44
1:A:233:ALA:HB2	1:A:403:ILE:O	2.18	0.44
1:E:43:ALA:HB2	1:F:296:ARG:NH1	2.33	0.44
1:B:65:ARG:HH21	1:B:251:ALA:HB3	1.82	0.44
1:B:277:GLY:HA2	3:B:664:HOH:O	2.17	0.44
1:C:236:ASP:HA	1:C:237:ASN:HA	1.73	0.44
1:E:86:PRO:HG2	1:E:401:HIS:CG	2.52	0.44
1:C:277:GLY:N	3:C:682:HOH:O	2.51	0.44
1:B:92:PRO:HG3	1:C:336:ILE:HG13	1.99	0.44
1:C:115:ALA:HB3	1:C:116:PRO:HD3	2.00	0.44
1:E:182:PRO:HA	1:E:183:GLY:HA2	1.87	0.43
1:F:324:HIS:CD2	1:F:367:VAL:HG22	2.53	0.43
1:A:202:ILE:HD12	1:A:203:THR:N	2.34	0.43
1:A:340:GLN:HE21	1:A:343:ALA:CB	2.30	0.43
1:A:182:PRO:HA	1:A:183:GLY:HA2	1.88	0.43
1:C:55:THR:HA	1:C:60:SER:O	2.19	0.43
1:E:99:ASN:OD1	3:E:725:HOH:O	2.21	0.43
1:F:333:VAL:O	1:F:393:ILE:HA	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:327:ALA:H	1:D:364:GLN:NE2	2.03	0.42
1:D:395:LEU:HD21	1:D:415:LEU:HB2	2.01	0.42
1:F:61:LEU:C	1:F:61:LEU:HD23	2.40	0.42
1:F:68:ARG:CG	1:F:68:ARG:HH11	2.32	0.42
1:C:303:LEU:O	1:C:390:THR:HA	2.19	0.42
1:E:16:SER:HB3	1:E:22:ALA:HA	2.00	0.42
1:A:339:ASN:CA	1:C:267:VAL:HG11	2.49	0.42
1:A:210:TYR:HB2	1:A:212:LEU:HG	2.00	0.42
1:B:65:ARG:NH2	1:B:251:ALA:O	2.53	0.42
1:A:395:LEU:HD21	1:A:415:LEU:HB2	2.00	0.42
1:F:353:ARG:O	1:F:357:GLN:HG3	2.20	0.42
1:B:4:GLU:N	1:B:4:GLU:OE2	2.52	0.42
1:C:174:PRO:HA	1:C:175:PRO:HD3	1.89	0.42
1:D:356:CYS:HB3	1:D:361:VAL:O	2.19	0.42
1:B:236:ASP:HA	1:B:237:ASN:HA	1.72	0.42
1:D:86:PRO:HG2	1:D:401:HIS:CG	2.55	0.42
1:D:329:ASN:HD21	1:D:410:HIS:H	1.66	0.41
1:D:200:HIS:HB2	1:D:202:ILE:HG22	2.02	0.41
1:E:329:ASN:HD21	1:E:410:HIS:H	1.67	0.41
1:E:130:PHE:CE2	1:E:202:ILE:HD11	2.55	0.41
1:E:62:ILE:HG12	1:E:260:VAL:HG13	2.01	0.41
1:C:42:ASP:O	1:C:286:ARG:NH2	2.53	0.41
1:E:333:VAL:HG22	1:E:364:GLN:HB2	2.02	0.41
1:E:42:ASP:O	1:E:286:ARG:NH2	2.53	0.41
1:D:281:GLU:HB2	1:D:386:VAL:HG12	2.02	0.41
1:B:3:ALA:C	1:B:4:GLU:CD	2.78	0.41
3:C:610:HOH:O	1:E:154:ALA:HB2	2.21	0.41
1:F:42:ASP:O	1:F:286:ARG:NH2	2.54	0.41
1:B:4:GLU:CA	1:B:5:LEU:CB	2.98	0.41
1:E:327:ALA:H	1:E:364:GLN:NE2	2.10	0.41
1:D:131:ARG:HB3	1:D:206:VAL:HG12	2.02	0.41
1:E:263:ASP:OD1	1:E:264:HIS:N	2.53	0.41
1:D:290:GLU:HG3	1:F:290:GLU:OE1	2.20	0.41
1:B:341:ARG:HH11	1:B:341:ARG:CG	2.33	0.41
1:C:428:LEU:HB3	1:C:429:PRO:HD2	2.02	0.41
1:F:236:ASP:HA	1:F:237:ASN:HA	1.75	0.41
1:D:211:GLU:OE2	1:E:345:ASN:HB2	2.20	0.41
1:B:306:ALA:HA	1:B:393:ILE:O	2.21	0.41
1:B:61:LEU:HD23	1:B:61:LEU:C	2.41	0.41
1:A:67:GLY:HA3	1:A:254:ASP:O	2.21	0.41
1:C:302:LEU:HG	1:C:303:LEU:N	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:PRO:HG3	1:B:336:ILE:HG13	2.03	0.41
1:B:3:ALA:HB3	1:B:4:GLU:CD	2.40	0.40
1:A:302:LEU:HG	1:A:303:LEU:N	2.36	0.40
1:E:71:PRO:HB2	1:E:288:LEU:HD21	2.03	0.40
1:B:72:LEU:C	1:B:72:LEU:HD23	2.41	0.40
1:F:89:ARG:CD	3:F:618:HOH:O	2.69	0.40
1:A:336:ILE:O	1:A:367:VAL:HG12	2.21	0.40
1:D:307:ASP:O	1:D:394:GLY:HA3	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	402/431 (93%)	387 (96%)	15 (4%)	0	100	100
1	B	402/431 (93%)	379 (94%)	19 (5%)	4 (1%)	19	22
1	C	402/431 (93%)	386 (96%)	15 (4%)	1 (0%)	52	64
1	D	402/431 (93%)	387 (96%)	14 (4%)	1 (0%)	52	64
1	E	402/431 (93%)	379 (94%)	22 (6%)	1 (0%)	52	64
1	F	402/431 (93%)	378 (94%)	23 (6%)	1 (0%)	52	64
All	All	2412/2586 (93%)	2296 (95%)	108 (4%)	8 (0%)	46	56

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	2	ARG
1	B	3	ALA
1	B	4	GLU
1	B	5	LEU

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Mol	Chain	Res	Type
1	E	48	THR
1	D	290	GLU
1	C	181	ALA
1	F	182	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	329/344 (96%)	317 (96%)	12 (4%)	42	58
1	B	329/344 (96%)	313 (95%)	16 (5%)	31	43
1	C	329/344 (96%)	313 (95%)	16 (5%)	31	43
1	D	329/344 (96%)	318 (97%)	11 (3%)	45	61
1	E	329/344 (96%)	312 (95%)	17 (5%)	29	40
1	F	329/344 (96%)	316 (96%)	13 (4%)	38	53
All	All	1974/2064 (96%)	1889 (96%)	85 (4%)	35	50

All (85) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	2	ARG
1	A	65	ARG
1	A	135	LYS
1	A	191	LEU
1	A	202	ILE
1	A	203	THR
1	A	302	LEU
1	A	319	ARG
1	A	335	LYS
1	A	386	VAL
1	A	425	SER
1	B	1	MET
1	B	4	GLU

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Mol	Chain	Res	Type
1	B	31	GLU
1	B	47	GLU
1	B	48	THR
1	B	65	ARG
1	B	77	ARG
1	B	98	ARG
1	B	191	LEU
1	B	202	ILE
1	B	252	GLU
1	B	341	ARG
1	B	385	GLN
1	B	386	VAL
1	B	410	HIS
1	B	418	VAL
1	C	1	MET
1	C	4	GLU
1	C	48	THR
1	C	54	VAL
1	C	65	ARG
1	C	72	LEU
1	C	77	ARG
1	C	135	LYS
1	C	190	LEU
1	C	191	LEU
1	C	198	ARG
1	C	203	THR
1	C	341	ARG
1	C	367	VAL
1	C	385	GLN
1	C	418	VAL
1	D	1	MET
1	D	73	GLU
1	D	190	LEU
1	D	202	ILE
1	D	252	GLU
1	D	290	GLU
1	D	336	ILE
1	D	358	ASP
1	D	367	VAL
1	D	372	MET
1	D	410	HIS
1	E	2	ARG

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Mol	Chain	Res	Type
1	E	4	GLU
1	E	48	THR
1	E	65	ARG
1	E	98	ARG
1	E	163	GLU
1	E	184	GLU
1	E	191	LEU
1	E	202	ILE
1	E	289	PRO
1	E	336	ILE
1	E	341	ARG
1	E	367	VAL
1	E	386	VAL
1	E	410	HIS
1	E	418	VAL
1	E	425	SER
1	F	4	GLU
1	F	45	HIS
1	F	46	THR
1	F	54	VAL
1	F	65	ARG
1	F	68	ARG
1	F	107	GLU
1	F	135	LYS
1	F	151	PRO
1	F	180	LEU
1	F	191	LEU
1	F	385	GLN
1	F	427	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	93	ASN
1	A	250	ASN
1	A	264	HIS
1	A	329	ASN
1	A	340	GLN
1	A	364	GLN
1	B	250	ASN
1	B	329	ASN
1	B	354	HIS

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Mol	Chain	Res	Type
1	B	364	GLN
1	B	414	HIS
1	C	45	HIS
1	C	329	ASN
1	C	340	GLN
1	C	364	GLN
1	D	133	ASN
1	D	264	HIS
1	D	329	ASN
1	D	337	ASN
1	D	364	GLN
1	D	414	HIS
1	E	133	ASN
1	E	329	ASN
1	E	340	GLN
1	E	364	GLN
1	F	250	ASN
1	F	329	ASN
1	F	340	GLN
1	F	364	GLN
1	F	385	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 6 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 ⓘ

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	408/431 (94%)	-0.33	5 (1%) 81 82	23, 33, 58, 87	0
1	B	408/431 (94%)	-0.36	3 (0%) 89 90	25, 33, 61, 109	0
1	C	408/431 (94%)	-0.39	4 (0%) 84 85	25, 34, 56, 94	0
1	D	408/431 (94%)	-0.39	6 (1%) 76 77	25, 36, 60, 87	0
1	E	408/431 (94%)	-0.18	10 (2%) 61 59	26, 37, 65, 86	0
1	F	408/431 (94%)	-0.30	4 (0%) 84 85	30, 39, 63, 88	0
All	All	2448/2586 (94%)	-0.32	32 (1%) 79 80	23, 36, 61, 109	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	277	GLY	4.1
1	E	267	VAL	3.8
1	C	182	PRO	3.8
1	F	182	PRO	3.7
1	E	372	MET	3.7
1	B	1	MET	3.6
1	B	267	VAL	3.2
1	E	68	ARG	3.1
1	D	372	MET	3.0
1	E	183	GLY	2.9
1	A	182	PRO	2.8
1	F	251	ALA	2.8
1	A	133	ASN	2.7
1	E	49	GLY	2.7
1	E	73	GLU	2.5
1	C	185	ALA	2.5
1	D	47	GLU	2.5
1	C	372	MET	2.4
1	A	45	HIS	2.4

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Mol	Chain	Res	Type	RSRZ
1	E	47	GLU	2.4
1	C	183	GLY	2.3
1	E	203	THR	2.2
1	F	385	GLN	2.2
1	A	48	THR	2.2
1	A	159	ARG	2.1
1	D	46	THR	2.1
1	B	203	THR	2.1
1	D	277	GLY	2.1
1	D	45	HIS	2.0
1	E	45	HIS	2.0
1	E	72	LEU	2.0
1	D	427	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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(Å ²)	Q<0.9
2	ZN	F	501	1/1	0.97	0.11	-1.42	65,65,65,65	0
2	ZN	B	501	1/1	0.99	0.05	-2.24	55,55,55,55	0
2	ZN	D	501	1/1	0.99	0.04	-2.28	60,60,60,60	0
2	ZN	E	501	1/1	0.98	0.06	-2.44	59,59,59,59	0
2	ZN	C	501	1/1	0.99	0.07	-3.17	56,56,56,56	0
2	ZN	A	501	1/1	0.99	0.05	-3.98	59,59,59,59	0

6.5 Other polymers [i](#)

There are no such residues in this entry.