Full wwPDB X-ray Structure Validation Report (i)

Aug 16, 2023 – 11:28 PM EDT

PDB ID	:	2B0O
Title	:	Crystal structure of UPLC1 GAP domain
Authors	:	Ismail, S.; Shen, L.; Arrowsmith, C.; Edwards, A.; Sundstrom, M.; Weigelt, J.; Bochkarev, A.; Park, H.; Structural Genomics Consortium (SGC)
Deposited on	:	2005-09-14
Resolution	:	2.06 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 (i) 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 (1)) were used in the production of this report:

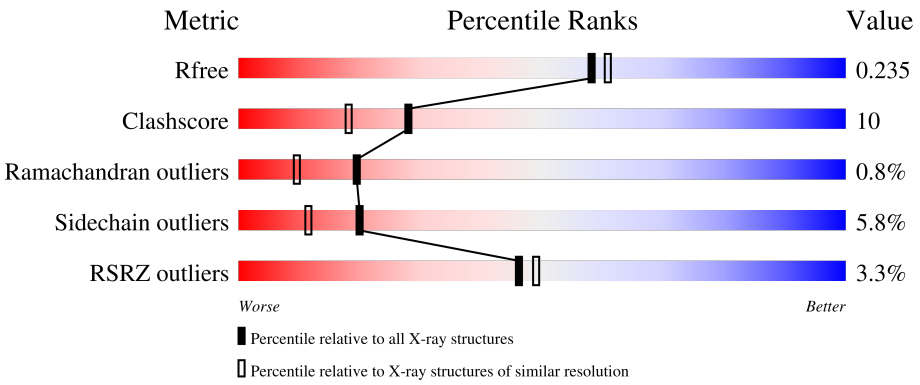
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 2.06 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	2684 (2.08-2.04)
Clashscore	141614	2801 (2.08-2.04)
Ramachandran outliers	138981	2768 (2.08-2.04)
Sidechain outliers	138945	2768 (2.08-2.04)
RSRZ outliers	127900	2646 (2.08-2.04)

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	E	301	5%; 71%; 11%; 5%; •; 12%
1	F	301	%; 74%; 13%; •; 12%
1	G	301	2%; 67%; 14%; • •; 16%
1	H	301	4%; 67%; 17%; • •; 13%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 8492 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 UPLC1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	E	264	2011	1257	359	381	14	0	0	0
1	F	266	2037	1278	361	384	14	0	0	0
1	G	253	1929	1203	347	366	13	0	0	0
1	H	262	2011	1264	357	377	13	0	0	0

There are 76 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	397	MET	-	cloning artifact	UNP Q6P9F4
E	398	GLY	-	cloning artifact	UNP Q6P9F4
E	399	SER	-	cloning artifact	UNP Q6P9F4
E	400	SER	-	cloning artifact	UNP Q6P9F4
E	401	HIS	-	expression tag	UNP Q6P9F4
E	402	HIS	-	expression tag	UNP Q6P9F4
E	403	HIS	-	expression tag	UNP Q6P9F4
E	404	HIS	-	expression tag	UNP Q6P9F4
E	405	HIS	-	expression tag	UNP Q6P9F4
E	406	HIS	-	expression tag	UNP Q6P9F4
E	407	SER	-	cloning artifact	UNP Q6P9F4
E	408	SER	-	cloning artifact	UNP Q6P9F4
E	409	GLY	-	cloning artifact	UNP Q6P9F4
E	410	LEU	-	cloning artifact	UNP Q6P9F4
E	411	VAL	-	cloning artifact	UNP Q6P9F4
E	412	PRO	-	cloning artifact	UNP Q6P9F4
E	413	ARG	-	cloning artifact	UNP Q6P9F4
E	414	GLY	-	cloning artifact	UNP Q6P9F4
E	415	SER	-	cloning artifact	UNP Q6P9F4
F	397	MET	-	cloning artifact	UNP Q6P9F4
F	398	GLY	-	cloning artifact	UNP Q6P9F4

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Chain	Residue	Modelled	Actual	Comment	Reference
F	399	SER	-	cloning artifact	UNP Q6P9F4
F	400	SER	-	cloning artifact	UNP Q6P9F4
F	401	HIS	-	expression tag	UNP Q6P9F4
F	402	HIS	-	expression tag	UNP Q6P9F4
F	403	HIS	-	expression tag	UNP Q6P9F4
F	404	HIS	-	expression tag	UNP Q6P9F4
F	405	HIS	-	expression tag	UNP Q6P9F4
F	406	HIS	-	expression tag	UNP Q6P9F4
F	407	SER	-	cloning artifact	UNP Q6P9F4
F	408	SER	-	cloning artifact	UNP Q6P9F4
F	409	GLY	-	cloning artifact	UNP Q6P9F4
F	410	LEU	-	cloning artifact	UNP Q6P9F4
F	411	VAL	-	cloning artifact	UNP Q6P9F4
F	412	PRO	-	cloning artifact	UNP Q6P9F4
F	413	ARG	-	cloning artifact	UNP Q6P9F4
F	414	GLY	-	cloning artifact	UNP Q6P9F4
F	415	SER	-	cloning artifact	UNP Q6P9F4
G	397	MET	-	cloning artifact	UNP Q6P9F4
G	398	GLY	-	cloning artifact	UNP Q6P9F4
G	399	SER	-	cloning artifact	UNP Q6P9F4
G	400	SER	-	cloning artifact	UNP Q6P9F4
G	401	HIS	-	expression tag	UNP Q6P9F4
G	402	HIS	-	expression tag	UNP Q6P9F4
G	403	HIS	-	expression tag	UNP Q6P9F4
G	404	HIS	-	expression tag	UNP Q6P9F4
G	405	HIS	-	expression tag	UNP Q6P9F4
G	406	HIS	-	expression tag	UNP Q6P9F4
G	407	SER	-	cloning artifact	UNP Q6P9F4
G	408	SER	-	cloning artifact	UNP Q6P9F4
G	409	GLY	-	cloning artifact	UNP Q6P9F4
G	410	LEU	-	cloning artifact	UNP Q6P9F4
G	411	VAL	-	cloning artifact	UNP Q6P9F4
G	412	PRO	-	cloning artifact	UNP Q6P9F4
G	413	ARG	-	cloning artifact	UNP Q6P9F4
G	414	GLY	-	cloning artifact	UNP Q6P9F4
G	415	SER	-	cloning artifact	UNP Q6P9F4
H	397	MET	-	cloning artifact	UNP Q6P9F4
H	398	GLY	-	cloning artifact	UNP Q6P9F4
H	399	SER	-	cloning artifact	UNP Q6P9F4
H	400	SER	-	cloning artifact	UNP Q6P9F4
H	401	HIS	-	expression tag	UNP Q6P9F4
H	402	HIS	-	expression tag	UNP Q6P9F4

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Chain	Residue	Modelled	Actual	Comment	Reference
H	403	HIS	-	expression tag	UNP Q6P9F4
H	404	HIS	-	expression tag	UNP Q6P9F4
H	405	HIS	-	expression tag	UNP Q6P9F4
H	406	HIS	-	expression tag	UNP Q6P9F4
H	407	SER	-	cloning artifact	UNP Q6P9F4
H	408	SER	-	cloning artifact	UNP Q6P9F4
H	409	GLY	-	cloning artifact	UNP Q6P9F4
H	410	LEU	-	cloning artifact	UNP Q6P9F4
H	411	VAL	-	cloning artifact	UNP Q6P9F4
H	412	PRO	-	cloning artifact	UNP Q6P9F4
H	413	ARG	-	cloning artifact	UNP Q6P9F4
H	414	GLY	-	cloning artifact	UNP Q6P9F4
H	415	SER	-	cloning artifact	UNP Q6P9F4

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
2	E	1	1	1	0	0
2	F	1	1	1	0	0
2	G	1	1	1	0	0
2	H	1	1	1	0	0

- Molecule 3 is water.

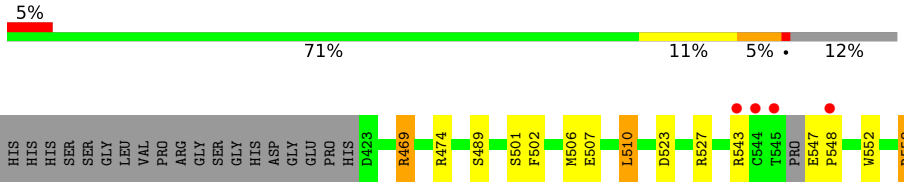
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	E	147	147	147	0	0
3	F	129	129	129	0	0
3	G	130	130	130	2	0
3	H	94	94	94	0	0

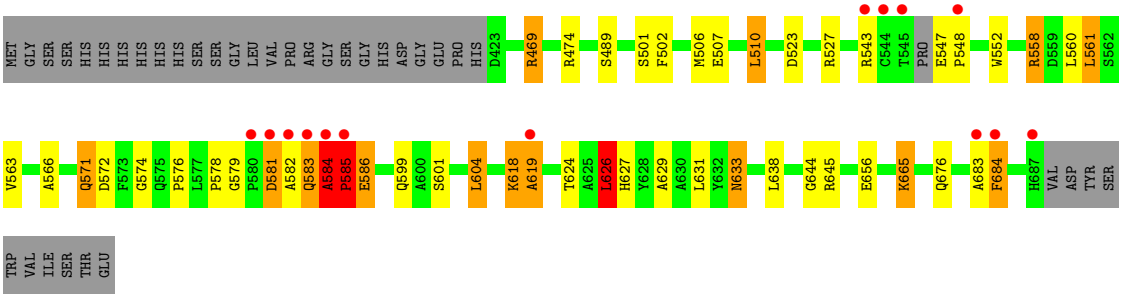
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: UPLC1

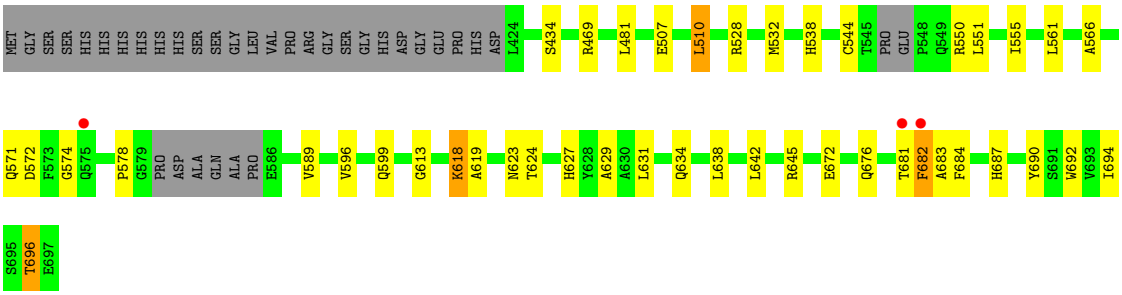
Chain E:





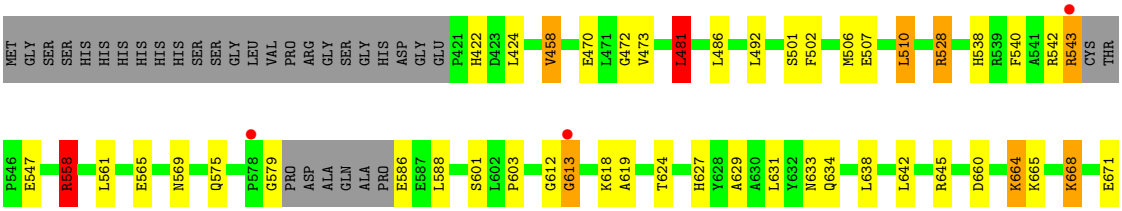
- Molecule 1: UPLC1

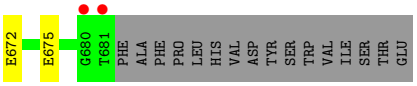
Chain F:



- Molecule 1: UPLC1

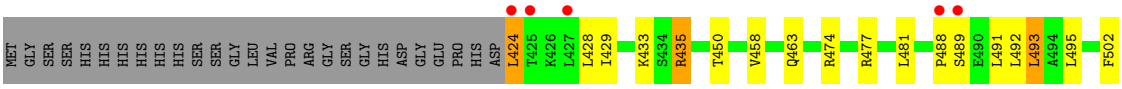
Chain G:

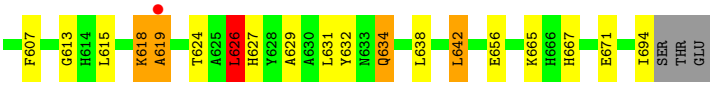




• Molecule 1: UPLC1

Chain H:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	54.05Å 106.72Å 100.36Å 90.00° 98.97° 90.00°	Depositor
Resolution (Å)	50.00 – 2.06 44.22 – 2.06	Depositor EDS
% Data completeness (in resolution range)	97.6 (50.00-2.06) 97.5 (44.22-2.06)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$< I/\sigma(I) >$ [1]	2.46 (at 2.06Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.183 , 0.233 0.183 , 0.235	Depositor DCC
R_{free} test set	3414 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	21.6	Xtriage
Anisotropy	0.307	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 58.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	8492	wwPDB-VP
Average B, all atoms (Å^2)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.41% 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 (i)

5.1 Standard geometry (i)

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	E	0.92	1/2048 (0.0%)	0.95	8/2776 (0.3%)
1	F	0.92	1/2074 (0.0%)	0.90	3/2809 (0.1%)
1	G	0.86	1/1962 (0.1%)	0.92	8/2654 (0.3%)
1	H	0.78	0/2048	0.82	3/2774 (0.1%)
All	All	0.87	3/8132 (0.0%)	0.90	22/11013 (0.2%)

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	E	0	3
1	F	0	2
1	G	0	1
1	H	0	2
All	All	0	8

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	613	GLY	N-CA	6.31	1.55	1.46
1	F	469	ARG	CZ-NH1	5.89	1.40	1.33
1	E	619	ALA	N-CA	5.41	1.57	1.46

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	E	469	ARG	NE-CZ-NH2	-13.76	113.42	120.30
1	G	528	ARG	NE-CZ-NH2	-11.34	114.63	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	E	469	ARG	NE-CZ-NH1	9.51	125.05	120.30
1	F	469	ARG	NE-CZ-NH2	-8.91	115.84	120.30
1	G	458	VAL	CG1-CB-CG2	7.40	122.74	110.90
1	G	528	ARG	NE-CZ-NH1	7.28	123.94	120.30
1	G	558	ARG	NE-CZ-NH2	-7.16	116.72	120.30
1	G	612	GLY	C-N-CA	-6.63	108.38	122.30
1	H	618	LYS	C-N-CA	-6.47	105.53	121.70
1	G	558	ARG	NE-CZ-NH1	6.22	123.41	120.30
1	F	618	LYS	C-N-CA	-6.21	106.18	121.70
1	E	618	LYS	C-N-CA	-6.11	106.42	121.70
1	E	527	ARG	NE-CZ-NH2	-5.86	117.37	120.30
1	G	481	LEU	CA-CB-CG	5.76	128.56	115.30
1	E	626	LEU	CB-CG-CD2	5.66	120.63	111.00
1	H	626	LEU	CA-CB-CG	5.48	127.91	115.30
1	H	527	ARG	NE-CZ-NH2	-5.37	117.61	120.30
1	E	584	ALA	C-N-CD	5.25	139.44	128.40
1	E	626	LEU	CA-CB-CG	5.22	127.30	115.30
1	E	604	LEU	CA-CB-CG	5.12	127.08	115.30
1	G	458	VAL	N-CA-CB	-5.05	100.39	111.50
1	F	510	LEU	CA-CB-CG	5.02	126.85	115.30

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	584	ALA	Peptide
1	E	585	PRO	Peptide
1	E	683	ALA	Peptide
1	F	613	GLY	Peptide
1	F	681	THR	Peptide
1	G	613	GLY	Peptide
1	H	613	GLY	Peptide
1	H	619	ALA	Peptide

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	E	2011	0	1992	49	0
1	F	2037	0	2021	37	0
1	G	1929	0	1918	45	0
1	H	2011	0	1996	36	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
3	E	147	0	0	7	0
3	F	129	0	0	5	0
3	G	130	0	0	5	0
3	H	94	0	0	5	0
All	All	8492	0	7927	158	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 10.

All (158) 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:H:618:LYS:O	1:H:619:ALA:CB	1.90	1.14
1:F:618:LYS:O	1:F:619:ALA:HB3	1.40	1.11
1:E:618:LYS:O	1:E:619:ALA:HB3	1.37	1.10
1:E:618:LYS:O	1:E:619:ALA:CB	1.93	1.08
1:G:618:LYS:O	1:G:619:ALA:HB3	1.48	1.08
1:F:618:LYS:O	1:F:619:ALA:CB	1.98	1.03
1:G:543:ARG:HA	1:G:569:ASN:OD1	1.61	1.00
1:H:618:LYS:O	1:H:619:ALA:HB3	1.22	0.99
1:G:618:LYS:O	1:G:619:ALA:CB	2.13	0.95
1:G:575:GLN:HG2	3:H:727:HOH:O	1.68	0.94
1:G:422:HIS:HD2	1:G:424:LEU:H	1.15	0.90
1:F:544:CYS:HB3	3:F:812:HOH:O	1.71	0.88
1:E:581:ASP:O	1:E:583:GLN:N	2.05	0.88
1:E:583:GLN:O	1:E:585:PRO:HD2	1.78	0.83
1:G:668:LYS:NZ	1:G:672:GLU:HG3	1.96	0.81
1:H:542:ARG:O	1:H:543:ARG:HB2	1.85	0.77
1:F:599:GLN:HG2	3:F:758:HOH:O	1.84	0.76
1:E:583:GLN:HB2	3:E:837:HOH:O	1.86	0.76
1:E:629:ALA:HB3	1:E:638:LEU:HD13	1.69	0.74
1:G:660:ASP:O	1:G:664:LYS:HD2	1.86	0.74
1:F:551:LEU:HD22	1:F:571:GLN:HE22	1.54	0.70
1:G:542:ARG:O	1:G:543:ARG:CG	2.39	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:629:ALA:HB3	1:F:638:LEU:HD13	1.74	0.70
1:E:644:GLY:O	1:E:645:ARG:HG2	1.92	0.69
1:G:668:LYS:HZ3	1:G:672:GLU:HG3	1.58	0.68
1:E:507:GLU:HB3	1:E:510:LEU:HD22	1.76	0.67
1:F:692:TRP:O	1:F:696:THR:HB	1.93	0.67
1:G:668:LYS:HD3	1:G:668:LYS:C	2.16	0.66
1:F:578:PRO:HA	3:F:824:HOH:O	1.96	0.66
1:E:626:LEU:HD22	1:E:638:LEU:HD12	1.78	0.65
1:H:528:ARG:O	1:H:532:MET:HG3	1.96	0.65
1:H:555:ILE:HD11	1:H:589:VAL:HG13	1.78	0.65
1:E:676:GLN:HB3	1:F:682:PHE:HE2	1.59	0.65
1:H:543:ARG:HA	1:H:569:ASN:OD1	1.97	0.65
1:G:542:ARG:O	1:G:543:ARG:HG2	1.96	0.64
1:E:676:GLN:HB3	1:F:682:PHE:CE2	2.32	0.64
1:F:538:HIS:HE1	3:F:746:HOH:O	1.81	0.63
1:F:672:GLU:O	1:F:676:GLN:HG3	1.99	0.62
1:G:507:GLU:HB3	1:G:510:LEU:HD22	1.82	0.62
1:F:566:ALA:O	1:F:571:GLN:HB2	2.00	0.61
1:G:422:HIS:CD2	1:G:424:LEU:H	2.07	0.61
1:E:684:PHE:H	1:E:684:PHE:HD1	1.43	0.61
1:H:542:ARG:HG2	1:H:565:GLU:OE1	2.00	0.61
1:E:560:LEU:HA	1:E:604:LEU:CD2	2.31	0.61
1:E:558:ARG:HG2	1:E:601:SER:HB3	1.82	0.61
1:G:634:GLN:NE2	3:G:706:HOH:O	2.34	0.61
1:E:684:PHE:CD1	1:E:684:PHE:N	2.54	0.60
1:F:551:LEU:HD22	1:F:571:GLN:NE2	2.16	0.60
1:G:624:THR:OG1	1:G:627:HIS:HD2	1.84	0.60
1:H:667:HIS:O	1:H:671:GLU:HG3	2.01	0.60
1:G:558:ARG:HG2	1:G:601:SER:HB3	1.85	0.59
1:H:629:ALA:HB3	1:H:638:LEU:HD13	1.85	0.59
1:G:472:GLY:HA3	3:G:811:HOH:O	2.03	0.58
1:G:558:ARG:CG	1:G:558:ARG:HH11	2.16	0.58
1:H:538:HIS:HE1	3:H:715:HOH:O	1.87	0.57
1:E:576:PRO:HB2	1:E:586:GLU:HG2	1.86	0.57
1:H:493:LEU:HD11	1:H:607:PHE:HB2	1.85	0.57
1:G:558:ARG:HG2	1:G:558:ARG:HH11	1.70	0.57
1:G:668:LYS:HZ2	1:G:672:GLU:HG3	1.68	0.57
1:G:579:GLY:HA2	3:G:766:HOH:O	2.05	0.56
1:F:682:PHE:O	1:F:683:ALA:HB3	2.05	0.56
1:G:542:ARG:HD2	1:G:565:GLU:OE1	2.06	0.56
1:F:618:LYS:O	1:F:623:ASN:O	2.24	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:422:HIS:CD2	1:G:424:LEU:HB2	2.40	0.56
1:H:424:LEU:O	1:H:428:LEU:HG	2.07	0.55
3:E:842:HOH:O	1:F:687:HIS:HD2	1.90	0.55
1:G:671:GLU:O	1:G:675:GLU:HG3	2.07	0.55
1:E:558:ARG:CG	1:E:558:ARG:HH11	2.21	0.54
1:H:435:ARG:HH11	1:H:435:ARG:HA	1.71	0.54
1:H:595:LYS:HE2	1:H:632:TYR:CZ	2.42	0.54
1:E:676:GLN:OE1	1:F:682:PHE:HZ	1.91	0.54
1:G:543:ARG:HA	1:G:569:ASN:CG	2.27	0.54
1:E:558:ARG:HG2	1:E:558:ARG:HH11	1.73	0.54
1:H:450:THR:O	1:H:481:LEU:HB2	2.09	0.53
1:H:624:THR:H	1:H:627:HIS:CD2	2.26	0.53
1:E:629:ALA:CB	1:E:638:LEU:HD13	2.36	0.53
1:F:507:GLU:HB3	1:F:510:LEU:HD22	1.90	0.53
1:E:633:ASN:ND2	3:E:729:HOH:O	2.32	0.53
1:G:668:LYS:HD3	1:G:668:LYS:O	2.09	0.53
1:G:624:THR:H	1:G:627:HIS:CD2	2.27	0.52
1:H:538:HIS:HD2	3:H:753:HOH:O	1.91	0.52
1:G:668:LYS:HD2	1:G:672:GLU:OE2	2.09	0.52
1:G:629:ALA:HB3	1:G:638:LEU:HD13	1.92	0.52
1:H:624:THR:OG1	1:H:627:HIS:HD2	1.92	0.52
1:G:470:GLU:OE1	1:G:528:ARG:HD3	2.09	0.51
1:E:583:GLN:O	1:E:585:PRO:CD	2.55	0.51
1:G:538:HIS:HE1	3:G:716:HOH:O	1.92	0.51
1:G:510:LEU:HD13	1:G:540:PHE:HB3	1.91	0.51
1:H:474:ARG:HD2	1:H:474:ARG:H	1.75	0.51
1:E:563:VAL:HG21	1:E:604:LEU:HD13	1.91	0.51
1:E:502:PHE:CE2	1:E:506:MET:HG3	2.46	0.51
1:E:578:PRO:HB3	1:E:586:GLU:OE1	2.10	0.50
1:H:491:LEU:O	1:H:495:LEU:HD12	2.12	0.50
1:H:627:HIS:HE1	1:H:656:GLU:O	1.95	0.50
1:F:624:THR:OG1	1:F:627:HIS:HD2	1.95	0.50
1:G:502:PHE:CE2	1:G:506:MET:HG3	2.47	0.50
1:F:528:ARG:O	1:F:532:MET:HG3	2.12	0.49
1:E:624:THR:OG1	1:E:627:HIS:HD2	1.95	0.49
1:E:560:LEU:HA	1:E:604:LEU:HD22	1.93	0.49
1:E:469:ARG:HD3	3:E:716:HOH:O	2.12	0.48
1:F:624:THR:H	1:F:627:HIS:CD2	2.31	0.48
1:F:538:HIS:HD2	3:F:740:HOH:O	1.95	0.47
1:F:634:GLN:HE21	1:F:634:GLN:HA	1.78	0.47
1:H:578:PRO:HG3	1:H:586:GLU:OE2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:572:ASP:C	1:F:574:GLY:H	2.18	0.47
1:H:429:ILE:O	1:H:433:LYS:HG3	2.16	0.46
1:E:558:ARG:CG	1:E:558:ARG:NH1	2.78	0.46
1:E:645:ARG:HE	1:F:645:ARG:HE	1.64	0.46
1:E:560:LEU:CA	1:E:604:LEU:HD21	2.46	0.46
1:G:538:HIS:HD2	3:G:738:HOH:O	1.99	0.46
1:H:634:GLN:HE21	1:H:634:GLN:HA	1.80	0.45
1:E:560:LEU:HA	1:E:604:LEU:HD21	1.96	0.45
1:E:558:ARG:HG2	1:E:601:SER:CB	2.45	0.45
1:G:502:PHE:CZ	1:G:506:MET:HG3	2.52	0.45
1:E:583:GLN:HG3	3:E:781:HOH:O	2.17	0.44
1:F:551:LEU:CD2	1:F:571:GLN:HE22	2.26	0.44
1:H:561:LEU:HD23	3:H:742:HOH:O	2.17	0.44
1:H:572:ASP:C	1:H:574:GLY:H	2.21	0.44
1:E:501:SER:HB3	1:E:561:LEU:HD11	1.99	0.44
1:E:624:THR:H	1:E:627:HIS:CD2	2.35	0.44
1:E:684:PHE:CD1	1:E:684:PHE:O	2.71	0.44
1:G:473:VAL:HG11	1:H:615:LEU:HB3	1.99	0.44
1:E:627:HIS:HE1	1:E:656:GLU:O	2.01	0.43
1:F:690:TYR:CE1	1:F:694:ILE:CD1	3.01	0.43
1:F:550:ARG:HE	1:F:550:ARG:HB2	1.61	0.43
1:E:584:ALA:O	1:E:586:GLU:N	2.51	0.43
1:E:645:ARG:HE	1:F:645:ARG:NE	2.17	0.43
1:F:555:ILE:HD11	1:F:589:VAL:HG13	2.00	0.43
1:G:501:SER:HB3	1:G:561:LEU:HD11	2.00	0.43
1:H:543:ARG:H	1:H:569:ASN:HD21	1.66	0.43
1:G:588:LEU:HD12	1:G:619:ALA:HB2	2.00	0.43
1:H:588:LEU:HD12	1:H:619:ALA:HB2	2.01	0.43
3:E:842:HOH:O	1:F:687:HIS:CD2	2.69	0.42
1:G:481:LEU:HA	1:G:486:LEU:HD11	2.01	0.42
1:G:492:LEU:HG	1:G:603:PRO:HG3	2.01	0.42
1:G:645:ARG:HD3	1:H:477:ARG:NH1	2.34	0.42
1:G:634:GLN:HE21	1:G:634:GLN:HA	1.84	0.42
1:H:626:LEU:HD22	1:H:638:LEU:CD1	2.48	0.42
1:E:552:TRP:CD2	1:E:579:GLY:HA2	2.54	0.42
1:E:581:ASP:HA	3:E:781:HOH:O	2.19	0.42
1:H:502:PHE:CE2	1:H:506:MET:HG3	2.54	0.42
1:F:682:PHE:HD1	1:F:684:PHE:H	1.67	0.42
1:E:572:ASP:C	1:E:574:GLY:H	2.23	0.41
1:H:626:LEU:HD21	1:H:642:LEU:HD13	2.01	0.41
1:G:645:ARG:HD3	1:H:477:ARG:HH12	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:463:GLN:HG3	3:H:782:HOH:O	2.20	0.41
1:H:492:LEU:CD1	1:H:603:PRO:HG3	2.51	0.41
1:G:424:LEU:HD23	1:G:424:LEU:HA	1.87	0.41
1:G:542:ARG:O	1:G:543:ARG:CB	2.64	0.41
1:E:581:ASP:O	1:E:583:GLN:CA	2.68	0.41
1:E:599:GLN:HG2	1:F:692:TRP:CD2	2.55	0.41
1:E:665:LYS:N	1:E:665:LYS:HD2	2.36	0.41
1:E:548:PRO:HB3	1:E:578:PRO:HD2	2.03	0.41
1:F:690:TYR:CE1	1:F:694:ILE:HD11	2.55	0.41
1:E:684:PHE:HD1	1:E:684:PHE:N	2.08	0.40
1:E:566:ALA:O	1:E:571:GLN:HB2	2.20	0.40
1:F:682:PHE:O	1:F:683:ALA:CB	2.70	0.40
1:F:561:LEU:HD23	1:F:561:LEU:HA	1.87	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	E	260/301 (86%)	250 (96%)	5 (2%)	5 (2%)	8 2
1	F	260/301 (86%)	253 (97%)	7 (3%)	0	100 100
1	G	247/301 (82%)	242 (98%)	5 (2%)	0	100 100
1	H	256/301 (85%)	240 (94%)	13 (5%)	3 (1%)	13 4
All	All	1023/1204 (85%)	985 (96%)	30 (3%)	8 (1%)	19 9

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	582	ALA
1	E	584	ALA

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Mol	Chain	Res	Type
1	H	542	ARG
1	E	581	ASP
1	E	583	GLN
1	H	489	SER
1	E	585	PRO
1	H	488	PRO

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	E	215/247 (87%)	200 (93%)	15 (7%)	15 7
1	F	219/247 (89%)	212 (97%)	7 (3%)	39 32
1	G	207/247 (84%)	194 (94%)	13 (6%)	18 10
1	H	215/247 (87%)	200 (93%)	15 (7%)	15 7
All	All	856/988 (87%)	806 (94%)	50 (6%)	20 11

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	474	ARG
1	E	489	SER
1	E	510	LEU
1	E	523	ASP
1	E	543	ARG
1	E	547	GLU
1	E	558	ARG
1	E	561	LEU
1	E	571	GLN
1	E	586	GLU
1	E	626	LEU
1	E	631	LEU
1	E	633	ASN
1	E	665	LYS
1	E	684	PHE

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Mol	Chain	Res	Type
1	F	434	SER
1	F	481	LEU
1	F	596	VAL
1	F	631	LEU
1	F	642	LEU
1	F	682	PHE
1	F	696	THR
1	G	458	VAL
1	G	481	LEU
1	G	510	LEU
1	G	543	ARG
1	G	547	GLU
1	G	558	ARG
1	G	586	GLU
1	G	631	LEU
1	G	633	ASN
1	G	642	LEU
1	G	664	LYS
1	G	665	LYS
1	G	668	LYS
1	H	424	LEU
1	H	435	ARG
1	H	458	VAL
1	H	493	LEU
1	H	524	MET
1	H	542	ARG
1	H	558	ARG
1	H	571	GLN
1	H	599	GLN
1	H	626	LEU
1	H	631	LEU
1	H	634	GLN
1	H	642	LEU
1	H	665	LYS
1	H	694	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	627	HIS
1	E	634	GLN
1	E	687	HIS

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Mol	Chain	Res	Type
1	F	513	HIS
1	F	538	HIS
1	F	571	GLN
1	F	627	HIS
1	F	633	ASN
1	F	634	GLN
1	G	422	HIS
1	G	538	HIS
1	G	627	HIS
1	G	633	ASN
1	G	634	GLN
1	H	538	HIS
1	H	571	GLN
1	H	627	HIS
1	H	633	ASN
1	H	634	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 4 ligands modelled in this entry, 4 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 (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 (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	E	264/301 (87%)	0.01	14 (5%)	26	27	10, 19, 43, 57	0
1	F	266/301 (88%)	-0.17	3 (1%)	80	81	11, 19, 39, 50	0
1	G	253/301 (84%)	-0.04	5 (1%)	65	67	11, 20, 40, 56	0
1	H	262/301 (87%)	0.07	13 (4%)	28	29	14, 25, 49, 63	0
All	All	1045/1204 (86%)	-0.03	35 (3%)	46	49	10, 21, 44, 63	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	512	SER	8.4
1	H	513	HIS	6.2
1	E	582	ALA	4.8
1	G	681	THR	4.7
1	H	514	GLY	4.6
1	E	684	PHE	4.5
1	H	427	LEU	4.2
1	F	682	PHE	4.2
1	E	581	ASP	3.9
1	E	580	PRO	3.8
1	E	584	ALA	3.5
1	H	543	ARG	3.4
1	E	545	THR	3.4
1	E	687	HIS	3.1
1	G	543	ARG	3.0
1	H	488	PRO	2.9
1	H	489	SER	2.8
1	E	583	GLN	2.7
1	E	585	PRO	2.6
1	H	424	LEU	2.6
1	E	683	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	F	681	THR	2.5
1	H	548	PRO	2.5
1	E	543	ARG	2.4
1	H	511	PRO	2.4
1	H	516	PRO	2.4
1	E	544	CYS	2.4
1	H	425	THR	2.2
1	G	578	PRO	2.2
1	G	613	GLY	2.2
1	H	619	ALA	2.1
1	F	575	GLN	2.1
1	E	548	PRO	2.1
1	E	619	ALA	2.0
1	G	680	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(Å^2)	Q<0.9
2	ZN	F	698	1/1	0.99	0.08	16,16,16,16	0
2	ZN	E	698	1/1	1.00	0.11	17,17,17,17	0
2	ZN	G	698	1/1	1.00	0.08	16,16,16,16	0
2	ZN	H	698	1/1	1.00	0.06	22,22,22,22	0

6.5 Other polymers (i)

There are no such residues in this entry.