



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 29, 2020 – 03:42 PM BST

PDB ID : 6DWH
Title : Crystal structure of complex of BBKI and Bovine Trypsin
Authors : Li, M.; Wlodawer, A.
Deposited on : 2018-06-26
Resolution : 2.00 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 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	:	2.13
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)	:	2.13

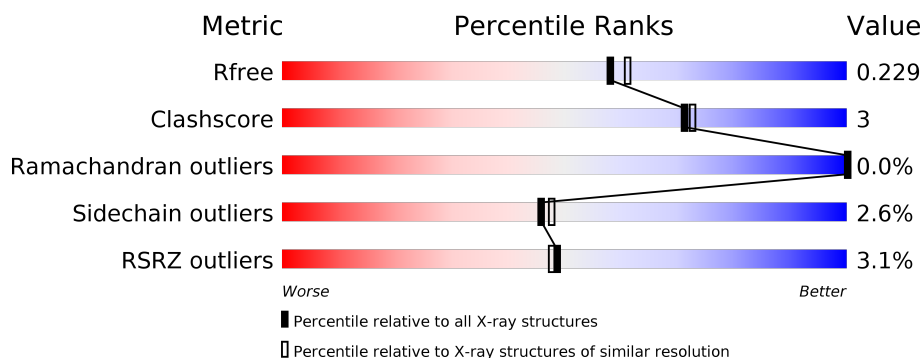
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	223	<div> <div>2%</div> <div> <div></div> <div>94%</div> <div>5%</div> </div> </div>
1	B	223	<div> <div>4%</div> <div> <div></div> <div>91%</div> <div>9%</div> </div> </div>
1	C	223	<div> <div>2%</div> <div> <div></div> <div>91%</div> <div>8%</div> </div> </div>
1	D	223	<div> <div>3%</div> <div> <div></div> <div>88%</div> <div>12%</div> </div> </div>
1	E	223	<div> <div>3%</div> <div> <div></div> <div>89%</div> <div>10%</div> </div> </div>
1	F	223	<div> <div>4%</div> <div> <div></div> <div>86%</div> <div>14%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
2	G	175	
2	H	175	
2	I	175	
2	J	175	
2	K	175	
2	L	175	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 18208 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 Cationic trypsin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	223	Total	C	N	O	S	0	1	0
			1631	1013	279	325	14			
1	B	223	Total	C	N	O	S	0	1	0
			1631	1013	279	325	14			
1	C	223	Total	C	N	O	S	0	1	0
			1631	1013	279	325	14			
1	D	223	Total	C	N	O	S	0	1	0
			1631	1013	279	325	14			
1	E	223	Total	C	N	O	S	0	1	0
			1631	1013	279	325	14			
1	F	223	Total	C	N	O	S	0	1	0
			1631	1013	279	325	14			

- Molecule 2 is a protein called Kunitz-type inhibitor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	163	Total	C	N	O	S	0	3	0
			1267	810	217	239	1			
2	H	163	Total	C	N	O	S	0	3	0
			1267	810	217	239	1			
2	I	163	Total	C	N	O	S	0	3	0
			1267	810	217	239	1			
2	J	163	Total	C	N	O	S	0	3	0
			1267	810	217	239	1			
2	K	163	Total	C	N	O	S	0	3	0
			1267	810	217	239	1			
2	L	163	Total	C	N	O	S	0	3	0
			1267	810	217	239	1			

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	J	1	Total Na 1 1	0	0
3	I	2	Total Na 2 2	0	0
3	L	2	Total Na 2 2	0	0
3	C	1	Total Na 1 1	0	0
3	E	2	Total Na 2 2	0	0

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	D	1	Total Cl 1 1	0	0
4	F	1	Total Cl 1 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	61	Total O 61 61	0	0
5	G	57	Total O 57 57	0	0
5	B	59	Total O 59 59	0	0
5	C	65	Total O 65 65	0	0
5	D	72	Total O 72 72	0	0
5	E	90	Total O 90 90	0	0
5	F	89	Total O 89 89	0	0
5	H	54	Total O 54 54	0	0
5	I	45	Total O 45 45	0	0
5	J	71	Total O 71 71	0	0
5	K	81	Total O 81 81	0	0

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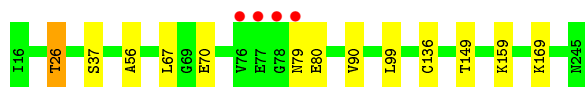
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	L	66	Total	O	0	0
			66	66		

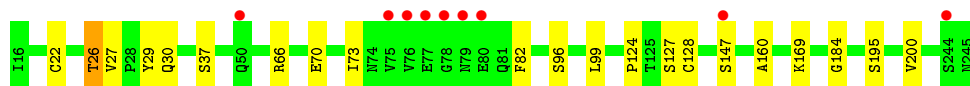
3 Residue-property plots [i](#)

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: Cationic trypsin



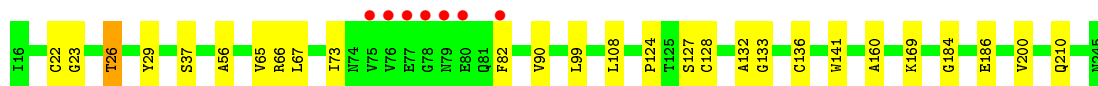
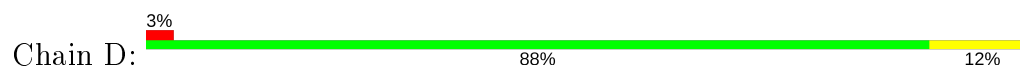
- Molecule 1: Cationic trypsin



- Molecule 1: Cationic trypsin



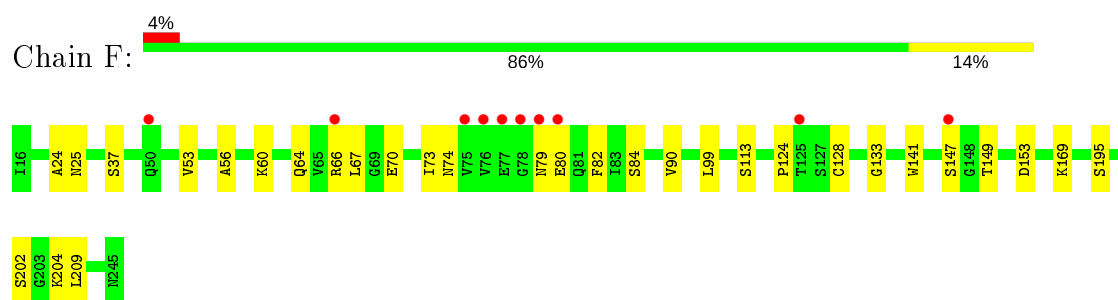
- Molecule 1: Cationic trypsin



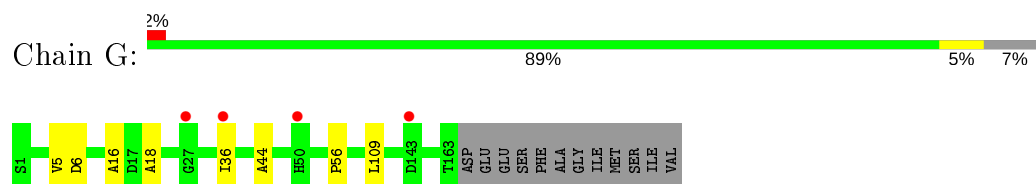
- Molecule 1: Cationic trypsin



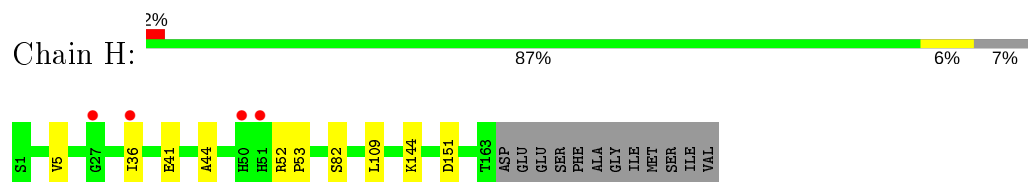
- Molecule 1: Cationic trypsin



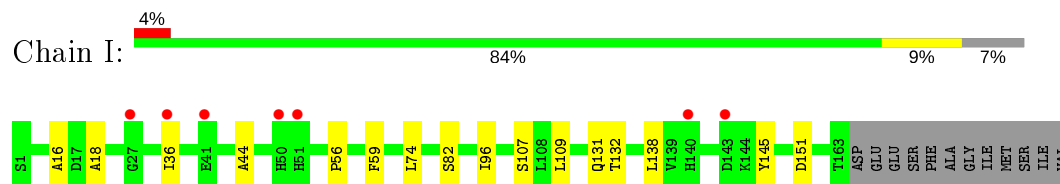
- Molecule 2: Kunitz-type inhibitor



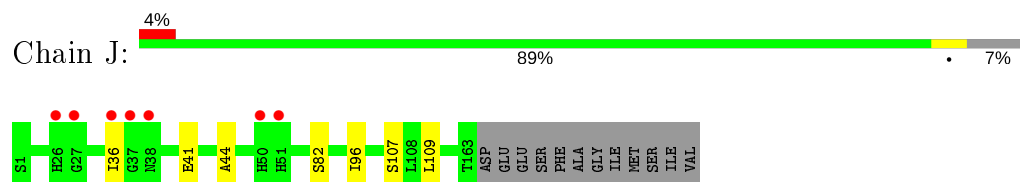
- Molecule 2: Kunitz-type inhibitor



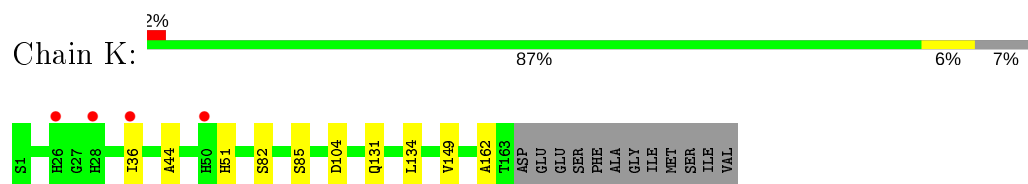
- Molecule 2: Kunitz-type inhibitor



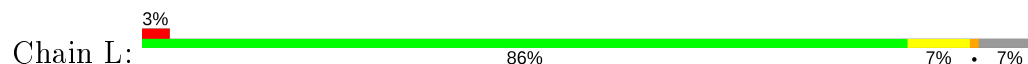
- Molecule 2: Kunitz-type inhibitor



- Molecule 2: Kunitz-type inhibitor



- Molecule 2: Kunitz-type inhibitor





4 Data and refinement statistics

Property	Value	Source
Space group	P 64	Depositor
Cell constants a, b, c, α , β , γ	207.81Å 207.81Å 107.17Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.91 – 2.00 49.91 – 2.01	Depositor EDS
% Data completeness (in resolution range)	80.4 (49.91-2.00) 80.4 (49.91-2.01)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.190 , 0.223 0.197 , 0.229	Depositor DCC
R_{free} test set	1397 reflections (0.99%)	wwPDB-VP
Wilson B-factor (Å ²)	24.9	Xtriage
Anisotropy	0.010	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 29.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.488 for h,-h-k,-l	Xtriage
Reported twinning fraction	0.479 for H, K, L 0.521 for -K, -H, -L	Depositor
Outliers	0 of 141523 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	18208	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 52.31 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.9517e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	0.59	0/1666	0.67	0/2258
1	B	0.66	2/1666 (0.1%)	0.72	1/2258 (0.0%)
1	C	0.73	2/1666 (0.1%)	0.70	1/2258 (0.0%)
1	D	0.72	3/1666 (0.2%)	0.75	3/2258 (0.1%)
1	E	0.70	1/1666 (0.1%)	0.73	1/2258 (0.0%)
1	F	0.80	3/1666 (0.2%)	0.79	3/2258 (0.1%)
2	G	0.44	0/1311	0.73	0/1781
2	H	0.43	0/1311	0.72	0/1781
2	I	0.43	0/1311	0.70	0/1781
2	J	0.43	0/1311	0.70	0/1781
2	K	0.46	0/1311	0.72	0/1781
2	L	0.47	0/1311	0.74	0/1781
All	All	0.60	11/17862 (0.1%)	0.72	9/24234 (0.0%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	67	LEU	C-N	14.14	1.58	1.33
1	F	128	CYS	C-N	11.13	1.59	1.34
1	F	133	GLY	C-N	11.03	1.59	1.34
1	E	127	SER	C-N	10.37	1.57	1.34
1	D	133	GLY	C-N	10.10	1.57	1.34
1	F	202	SER	C-N	10.09	1.51	1.33
1	B	127	SER	C-N	6.51	1.49	1.34
1	C	127	SER	C-N	6.33	1.48	1.34
1	D	128	CYS	C-N	6.29	1.48	1.34
1	D	127	SER	C-N	5.49	1.46	1.34
1	B	128	CYS	C-N	5.21	1.46	1.34

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	25	ASN	N-CA-CB	-8.73	94.88	110.60
1	D	124	PRO	O-C-N	8.68	136.58	122.70
1	C	67	LEU	C-N-CA	-7.84	105.84	122.30
1	D	124	PRO	CA-C-N	-6.43	103.05	117.20
1	B	124	PRO	O-C-N	6.19	132.60	122.70
1	E	128	CYS	O-C-N	6.00	132.31	122.70
1	F	24	ALA	N-CA-C	5.95	127.05	111.00
1	F	124	PRO	O-C-N	5.16	130.96	122.70
1	D	124	PRO	C-N-CA	-5.10	108.96	121.70

There are no chirality outliers.

There are no planarity outliers.

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	1631	0	1589	9	0
1	B	1631	0	1589	15	0
1	C	1631	0	1589	13	0
1	D	1631	0	1589	18	0
1	E	1631	0	1589	19	0
1	F	1631	0	1589	15	0
2	G	1267	0	1271	10	0
2	H	1267	0	1271	6	1
2	I	1267	0	1271	12	2
2	J	1267	0	1271	4	0
2	K	1267	0	1271	8	0
2	L	1267	0	1271	11	0
3	C	1	0	0	0	0
3	E	2	0	0	0	0
3	I	2	0	0	0	0
3	J	1	0	0	0	0
3	L	2	0	0	0	0
4	D	1	0	0	0	0
4	F	1	0	0	0	0
5	A	61	0	0	0	0
5	B	59	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	C	65	0	0	1	0
5	D	72	0	0	5	0
5	E	90	0	0	6	0
5	F	89	0	0	2	0
5	G	57	0	0	2	0
5	H	54	0	0	1	0
5	I	45	0	0	2	0
5	J	71	0	0	0	0
5	K	81	0	0	3	0
5	L	66	0	0	3	0
All	All	18208	0	17160	120	3

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 3.

All (120) 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:E:66:ARG:HG2	1:E:82:PHE:CD1	1.92	1.03
1:E:66:ARG:HD3	1:E:82:PHE:CE1	2.09	0.87
1:D:66:ARG:HD3	1:D:82:PHE:CE1	2.15	0.80
1:E:210:GLN:HG2	5:E:435:HOH:O	1.85	0.76
1:A:26:THR:HG22	1:D:26:THR:HG22	1.68	0.75
1:E:33:LEU:CD2	5:E:450:HOH:O	2.37	0.73
1:B:26:THR:HG22	1:C:26:THR:HG22	1.72	0.71
1:D:136:CYS:HB3	5:D:402:HOH:O	1.89	0.71
1:B:27:VAL:HG12	5:B:301:HOH:O	1.88	0.71
1:D:66:ARG:HD3	1:D:82:PHE:CZ	2.26	0.70
1:E:33:LEU:HD21	5:E:450:HOH:O	1.90	0.69
2:G:36:ILE:HG12	2:G:44:ALA:HB2	1.77	0.67
1:B:66:ARG:HD3	1:B:82:PHE:CE1	2.29	0.66
2:L:36:ILE:HG12	2:L:44:ALA:HB2	1.78	0.66
1:E:70:GLU:HG3	1:E:80:GLU:CG	2.26	0.65
1:D:200:VAL:N	5:D:402:HOH:O	2.28	0.64
1:B:70:GLU:OE2	1:B:73:ILE:HG22	1.99	0.62
1:E:70:GLU:HG3	1:E:80:GLU:HG2	1.80	0.61
2:H:144:LYS:HB3	5:H:206:HOH:O	1.99	0.61
1:D:99:LEU:HD11	2:J:109:LEU:HD11	1.81	0.61
1:E:66:ARG:CG	1:E:82:PHE:CD1	2.79	0.60
1:B:29:TYR:N	5:B:301:HOH:O	2.34	0.60
5:K:224:HOH:O	2:L:53:PRO:HD2	2.00	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:74:LEU:HD22	5:I:311:HOH:O	2.01	0.59
1:C:66:ARG:CZ	1:C:82:PHE:CE2	2.87	0.58
1:E:58:CYS:O	5:E:401:HOH:O	2.17	0.57
1:E:44:GLY:HA3	5:E:450:HOH:O	2.05	0.57
2:L:92:GLN:NE2	5:L:301:HOH:O	2.39	0.56
1:C:137:LEU:HA	5:C:444:HOH:O	2.06	0.56
1:B:66:ARG:HD3	1:B:82:PHE:CZ	2.41	0.56
1:E:66:ARG:CD	1:E:82:PHE:CE1	2.85	0.55
1:F:149:THR:HG21	2:L:16:ALA:HB1	1.88	0.55
2:K:36:ILE:HG12	2:K:44:ALA:HB2	1.88	0.55
2:I:36:ILE:HG12	2:I:44:ALA:HB2	1.88	0.55
1:B:22:CYS:O	1:B:26:THR:HG21	2.08	0.54
2:K:162:ALA:HB2	5:K:220:HOH:O	2.07	0.54
2:L:36:ILE:CD1	2:L:44:ALA:HB2	2.39	0.52
1:D:99:LEU:CD1	2:J:109:LEU:HD11	2.40	0.52
2:J:36:ILE:HG12	2:J:44:ALA:HB2	1.91	0.52
1:C:149:THR:HG21	2:I:16:ALA:CB	2.40	0.51
1:E:22:CYS:O	1:E:26:THR:HG21	2.12	0.50
1:C:72:ASN:HA	1:C:153:ASP:O	2.11	0.50
1:B:26:THR:HB	1:C:26:THR:HA	1.92	0.50
1:F:70:GLU:HG3	1:F:80:GLU:HG2	1.94	0.50
1:C:99:LEU:CD1	2:I:109:LEU:HD11	2.43	0.49
1:D:29:TYR:CZ	1:D:200:VAL:HG21	2.46	0.49
1:F:70:GLU:HG3	1:F:80:GLU:CG	2.42	0.49
2:L:10:GLN:NE2	5:L:303:HOH:O	2.46	0.49
1:E:53:VAL:HG11	1:E:209:LEU:HD21	1.94	0.49
1:F:73:ILE:HG23	1:F:141:TRP:CE2	2.49	0.48
2:L:36:ILE:CG1	2:L:44:ALA:HB2	2.43	0.48
1:F:149:THR:HG21	2:L:16:ALA:CB	2.42	0.48
1:D:132:ALA:C	5:D:401:HOH:O	2.52	0.48
1:F:66:ARG:HD3	1:F:82:PHE:CZ	2.48	0.48
2:G:36:ILE:CG1	2:G:44:ALA:HB2	2.44	0.48
1:D:22:CYS:O	1:D:26:THR:HG21	2.12	0.48
1:A:149:THR:HG21	2:G:16:ALA:CB	2.44	0.48
1:C:160:ALA:HB1	1:C:184:GLY:HA2	1.96	0.48
1:D:132:ALA:O	5:D:401:HOH:O	2.20	0.48
2:G:5:VAL:HB	5:G:201:HOH:O	2.13	0.47
2:K:36:ILE:CG1	2:K:44:ALA:HB2	2.43	0.47
1:B:27:VAL:C	5:B:301:HOH:O	2.51	0.47
2:H:36:ILE:CG1	2:H:44:ALA:HB2	2.44	0.47
1:F:74:ASN:ND2	1:F:153:ASP:OD1	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:53:VAL:CG1	1:E:209:LEU:HD21	2.45	0.47
1:B:160:ALA:HB1	1:B:184:GLY:HA2	1.97	0.47
1:B:30:GLN:N	5:B:301:HOH:O	2.47	0.47
1:B:99:LEU:HD11	2:H:109:LEU:HD11	1.98	0.46
2:K:51:HIS:N	5:K:204:HOH:O	2.49	0.46
1:A:99:LEU:HD11	2:G:109:LEU:HD11	1.97	0.46
2:H:36:ILE:HD11	2:H:44:ALA:HB2	1.98	0.45
1:A:26:THR:HA	1:D:26:THR:HB	1.98	0.45
1:C:73:ILE:HG23	1:C:141:TRP:CE2	2.52	0.45
2:H:36:ILE:HG12	2:H:44:ALA:HB2	1.98	0.44
2:H:52:ARG:HG3	2:H:53:PRO:HA	1.98	0.44
1:B:29:TYR:CD1	5:B:301:HOH:O	2.70	0.44
2:J:96:ILE:O	2:K:131:GLN:NE2	2.49	0.44
1:A:136:CYS:O	1:A:159:LYS:HA	2.17	0.44
1:C:149:THR:HG21	2:I:16:ALA:HB2	1.99	0.44
2:K:134:LEU:HB3	2:K:149:VAL:HG13	2.00	0.44
1:D:160:ALA:HB1	1:D:184:GLY:HA2	2.00	0.43
2:I:36:ILE:CD1	2:I:44:ALA:HB2	2.48	0.43
2:K:36:ILE:HD11	2:K:44:ALA:HB2	2.00	0.43
1:D:73:ILE:HG23	1:D:141:TRP:CE2	2.53	0.43
1:F:64:GLN:HA	1:F:84:SER:HA	2.00	0.43
2:I:59:PHE:C	5:I:311:HOH:O	2.57	0.43
1:A:70:GLU:HG3	1:A:80:GLU:HG2	2.01	0.43
1:E:210:GLN:CG	5:E:435:HOH:O	2.56	0.43
1:B:29:TYR:CZ	1:B:200:VAL:HG21	2.54	0.43
1:C:99:LEU:HD11	2:I:109:LEU:HD11	2.01	0.43
1:A:149:THR:HG21	2:G:16:ALA:HB1	2.01	0.43
1:A:56:ALA:HB1	1:A:90:VAL:HG13	2.01	0.43
1:F:195:SER:OG	2:L:64:ARG:C	2.58	0.43
1:E:160:ALA:HB1	1:E:184:GLY:HA2	2.01	0.42
2:I:138:LEU:HB3	2:I:145:TYR:HB3	2.01	0.42
1:D:65:VAL:HG21	1:D:108:LEU:HD21	2.01	0.42
1:F:60:LYS:HE2	2:L:3:VAL:HG21	2.00	0.42
1:F:99:LEU:HD13	5:F:459:HOH:O	2.20	0.42
1:E:66:ARG:HD3	1:E:82:PHE:CZ	2.53	0.42
1:E:73:ILE:HG23	1:E:141:TRP:CE2	2.54	0.42
2:G:36:ILE:CD1	2:G:44:ALA:HB2	2.49	0.42
2:K:36:ILE:CD1	2:K:44:ALA:HB2	2.50	0.42
1:F:53:VAL:HG11	1:F:209:LEU:HD21	2.02	0.42
1:F:53:VAL:CG1	1:F:209:LEU:HD21	2.49	0.41
1:E:75:VAL:HG23	1:E:77:GLU:HG3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:77:LYS:HD2	5:L:312:HOH:O	2.20	0.41
2:G:18:ALA:HB1	2:G:56:PRO:HB2	2.02	0.41
2:I:36:ILE:CG1	2:I:44:ALA:HB2	2.49	0.41
1:C:53:VAL:CG1	1:C:209:LEU:HD21	2.51	0.41
1:C:72:ASN:HB3	1:C:75:VAL:HG22	2.02	0.41
1:F:56:ALA:HB1	1:F:90:VAL:HG13	2.03	0.41
1:D:56:ALA:HB1	1:D:90:VAL:HG13	2.03	0.41
2:G:6:ASP:N	5:G:201:HOH:O	2.53	0.41
2:I:18:ALA:HB1	2:I:56:PRO:HB2	2.02	0.41
1:A:99:LEU:CD1	2:G:109:LEU:HD11	2.51	0.41
1:B:70:GLU:CD	1:B:73:ILE:HG22	2.41	0.41
1:D:23:GLY:O	1:D:26:THR:HG23	2.21	0.40
1:F:204:LYS:HD3	5:F:453:HOH:O	2.20	0.40
2:I:36:ILE:HD11	2:I:44:ALA:HB2	2.04	0.40
1:D:210:GLN:HG2	5:D:418:HOH:O	2.22	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:151:ASP:OD2	2:I:151:ASP:OD2[4_655]	1.77	0.43
2:I:96:ILE:O	2:I:131:GLN:NE2[4_655]	2.03	0.17
2:H:151:ASP:OD2	2:H:151:ASP:OD2[4_765]	2.14	0.06

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	222/223 (100%)	216 (97%)	6 (3%)	0	100	100
1	B	222/223 (100%)	215 (97%)	7 (3%)	0	100	100
1	C	222/223 (100%)	215 (97%)	7 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	222/223 (100%)	217 (98%)	5 (2%)	0	100	100
1	E	222/223 (100%)	218 (98%)	4 (2%)	0	100	100
1	F	222/223 (100%)	214 (96%)	8 (4%)	0	100	100
2	G	164/175 (94%)	155 (94%)	9 (6%)	0	100	100
2	H	164/175 (94%)	157 (96%)	7 (4%)	0	100	100
2	I	164/175 (94%)	154 (94%)	10 (6%)	0	100	100
2	J	164/175 (94%)	155 (94%)	9 (6%)	0	100	100
2	K	164/175 (94%)	156 (95%)	8 (5%)	0	100	100
2	L	164/175 (94%)	159 (97%)	4 (2%)	1 (1%)	25	19
All	All	2316/2388 (97%)	2231 (96%)	84 (4%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	L	27	GLY

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	185/184 (100%)	180 (97%)	5 (3%)	44	46
1	B	185/184 (100%)	179 (97%)	6 (3%)	39	38
1	C	185/184 (100%)	180 (97%)	5 (3%)	44	46
1	D	185/184 (100%)	180 (97%)	5 (3%)	44	46
1	E	185/184 (100%)	178 (96%)	7 (4%)	33	31
1	F	185/184 (100%)	179 (97%)	6 (3%)	39	38
2	G	140/147 (95%)	140 (100%)	0	100	100
2	H	140/147 (95%)	137 (98%)	3 (2%)	53	57
2	I	140/147 (95%)	136 (97%)	4 (3%)	42	43

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	J	140/147 (95%)	136 (97%)	4 (3%)	42	43
2	K	140/147 (95%)	137 (98%)	3 (2%)	53	57
2	L	140/147 (95%)	135 (96%)	5 (4%)	35	34
All	All	1950/1986 (98%)	1897 (97%)	53 (3%)	46	46

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	THR
1	A	37	SER
1	A	67	LEU
1	A	79	ASN
1	A	169	LYS
1	B	26	THR
1	B	37	SER
1	B	96	SER
1	B	147	SER
1	B	169	LYS
1	B	195	SER
1	C	26	THR
1	C	37	SER
1	C	67	LEU
1	C	79	ASN
1	C	169	LYS
1	D	26	THR
1	D	37	SER
1	D	67	LEU
1	D	169	LYS
1	D	186	GLU
1	E	26	THR
1	E	37	SER
1	E	67	LEU
1	E	79	ASN
1	E	80	GLU
1	E	147	SER
1	E	169	LYS
1	F	37	SER
1	F	67	LEU
1	F	79	ASN
1	F	113	SER
1	F	147	SER

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Mol	Chain	Res	Type
1	F	169	LYS
2	H	5	VAL
2	H	41	GLU
2	H	82	SER
2	I	82	SER
2	I	107[A]	SER
2	I	107[B]	SER
2	I	132	THR
2	J	41	GLU
2	J	82	SER
2	J	107[A]	SER
2	J	107[B]	SER
2	K	82	SER
2	K	85	SER
2	K	104	ASP
2	L	36	ILE
2	L	81	SER
2	L	82	SER
2	L	107[A]	SER
2	L	107[B]	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	ASN
1	A	95	ASN
1	B	50	GLN
1	C	34	ASN
1	C	240	GLN
1	E	34	ASN
1	E	50	GLN
1	E	175	GLN
2	H	66	ASN
2	K	66	ASN
2	L	10	GLN
2	L	92	GLN

5.3.3 RNA ⓘ

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 10 ligands modelled in this entry, 10 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 ⓘ

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	223/223 (100%)	-0.04	4 (1%) 68 66	18, 29, 46, 82	0
1	B	223/223 (100%)	0.13	9 (4%) 38 37	17, 34, 52, 92	0
1	C	223/223 (100%)	0.02	5 (2%) 62 60	20, 31, 48, 82	0
1	D	223/223 (100%)	0.05	7 (3%) 49 48	17, 31, 49, 86	0
1	E	223/223 (100%)	0.16	6 (2%) 54 53	17, 30, 48, 93	0
1	F	223/223 (100%)	0.23	10 (4%) 33 32	16, 31, 48, 89	0
2	G	163/175 (93%)	-0.02	4 (2%) 57 56	16, 28, 57, 71	0
2	H	163/175 (93%)	0.02	4 (2%) 57 56	16, 30, 54, 72	0
2	I	163/175 (93%)	0.04	7 (4%) 35 34	18, 31, 62, 81	0
2	J	163/175 (93%)	0.03	7 (4%) 35 34	16, 29, 56, 68	0
2	K	163/175 (93%)	0.06	4 (2%) 57 56	16, 29, 57, 73	0
2	L	163/175 (93%)	0.09	5 (3%) 49 48	15, 27, 53, 70	0
All	All	2316/2388 (96%)	0.07	72 (3%) 49 48	15, 30, 55, 93	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	76	VAL	6.6
1	F	76	VAL	6.2
1	D	75	VAL	6.1
1	A	77	GLU	5.7
2	H	27	GLY	5.6
1	F	77	GLU	5.5
1	B	77	GLU	5.4
1	E	79	ASN	5.4
1	B	75	VAL	5.3
1	B	79	ASN	5.1
1	E	77	GLU	5.1

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Mol	Chain	Res	Type	RSRZ
2	I	50	HIS	5.0
2	G	50	HIS	4.8
1	D	76	VAL	4.7
1	C	76	VAL	4.6
2	H	50	HIS	4.5
1	F	78	GLY	4.5
2	K	50	HIS	4.2
1	E	76	VAL	4.1
2	J	37	GLY	4.0
1	F	75	VAL	3.9
1	F	79	ASN	3.9
1	C	79	ASN	3.8
2	K	36	ILE	3.8
2	G	27	GLY	3.8
2	J	50	HIS	3.7
1	D	77	GLU	3.7
1	A	78	GLY	3.5
2	L	36	ILE	3.4
2	L	27	GLY	3.4
2	J	36	ILE	3.3
1	B	244	SER	3.3
1	C	77	GLU	3.3
1	F	147	SER	3.1
2	J	26	HIS	3.1
2	L	26	HIS	3.1
2	L	50	HIS	3.0
1	E	78	GLY	3.0
2	H	36	ILE	2.9
2	L	143	ASP	2.9
1	D	78	GLY	2.8
1	D	79	ASN	2.6
2	I	27	GLY	2.6
2	J	51	HIS	2.6
1	E	74	ASN	2.6
1	B	147	SER	2.5
2	K	28	HIS	2.5
1	F	125	THR	2.5
1	B	78	GLY	2.5
2	I	41	GLU	2.4
2	I	143	ASP	2.4
1	A	79	ASN	2.4
1	A	76	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	E	75	VAL	2.3
2	I	51	HIS	2.3
1	F	66	ARG	2.3
2	I	36	ILE	2.3
2	G	36	ILE	2.3
2	H	51	HIS	2.2
1	D	80	GLU	2.2
1	B	50	GLN	2.2
2	J	38	ASN	2.2
2	K	26	HIS	2.2
2	I	140	HIS	2.1
2	J	27	GLY	2.1
1	F	80	GLU	2.1
1	D	82	PHE	2.1
1	F	50	GLN	2.0
1	B	80	GLU	2.0
1	C	117	ARG	2.0
1	C	80	GLU	2.0
2	G	143	ASP	2.0

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	NA	J	201	1/1	0.92	0.32	38,38,38,38	0
3	NA	L	202	1/1	0.92	0.28	42,42,42,42	0
4	CL	D	301	1/1	0.93	0.15	55,55,55,55	0
4	CL	F	301	1/1	0.94	0.13	54,54,54,54	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	NA	I	201	1/1	0.94	0.21	49,49,49,49	0
3	NA	I	202	1/1	0.95	0.14	49,49,49,49	0
3	NA	E	302	1/1	0.96	0.09	32,32,32,32	0
3	NA	E	301	1/1	0.99	0.34	21,21,21,21	0
3	NA	C	301	1/1	0.99	0.10	28,28,28,28	0
3	NA	L	201	1/1	0.99	0.07	22,22,22,22	0

6.5 Other polymers [i](#)

There are no such residues in this entry.