



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 06:27 pm GMT

PDB ID : 4N8V
Title : Crystal structure of killer cell immunoglobulin-like receptor KIR2DS2 in complex with HLA-A
Authors : Liu, J.X.; Ren, E.C.
Deposited on : 2013-10-18
Resolution : 2.50 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
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

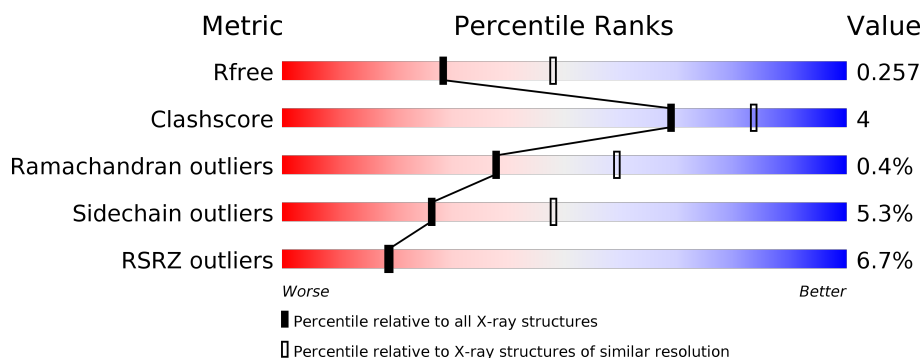
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	100719	3846 (2.50-2.50)
Clashscore	112137	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (2.50-2.50)
RSRZ outliers	101464	3876 (2.50-2.50)

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	G	201	<div> <div>6%</div> <div>80%</div> <div>12%</div> <div>5%</div> </div>
1	I	201	<div> <div>26%</div> <div>81%</div> <div>11%</div> <div>5%</div> </div>
2	A	274	<div> <div>%</div> <div>90%</div> <div>9%</div> <div>.</div> </div>
2	D	274	<div> <div>%</div> <div>92%</div> <div>7%</div> <div>.</div> </div>
3	B	100	<div> <div>3%</div> <div>89%</div> <div>10%</div> <div>.</div> </div>
3	E	100	<div> <div>%</div> <div>90%</div> <div>9%</div> <div>.</div> </div>

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Mol	Chain	Length	Quality of chain
4	C	9	 67%33%
4	F	9	 67%33%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9364 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 Killer cell immunoglobulin-like receptor 2DS2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	G	190	Total	C	N	O	S	0	0	0
			1473	930	252	284	7			
1	I	190	Total	C	N	O	S	0	0	0
			1473	930	252	284	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	0	MET	-	EXPRESSION TAG	UNP P43631
I	0	MET	-	EXPRESSION TAG	UNP P43631

- Molecule 2 is a protein called HLA class I histocompatibility antigen, A-11 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	274	Total	C	N	O	S	0	0	0
			2236	1389	407	431	9			
2	D	274	Total	C	N	O	S	0	0	0
			2236	1389	407	431	9			

- Molecule 3 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	B	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
3	E	99	Total	C	N	O	S	0	0	0
			829	528	140	158	3			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	EXPRESSION TAG	UNP P61769
E	0	MET	-	EXPRESSION TAG	UNP P61769

- Molecule 4 is a protein called peptide from Virion membrane protein A14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	C	9	Total	C	N	O	S	0	0	0
			78	53	11	12	2			
4	F	9	Total	C	N	O	S	0	0	0
			78	53	11	12	2			

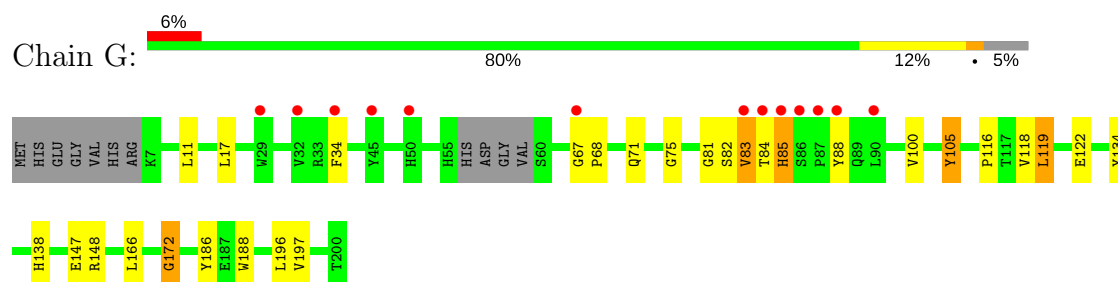
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	G	31	Total	O	0	0
			31	31		
5	I	29	Total	O	0	0
			29	29		
5	A	20	Total	O	0	0
			20	20		
5	B	7	Total	O	0	0
			7	7		
5	C	1	Total	O	0	0
			1	1		
5	D	26	Total	O	0	0
			26	26		
5	E	8	Total	O	0	0
			8	8		
5	F	2	Total	O	0	0
			2	2		

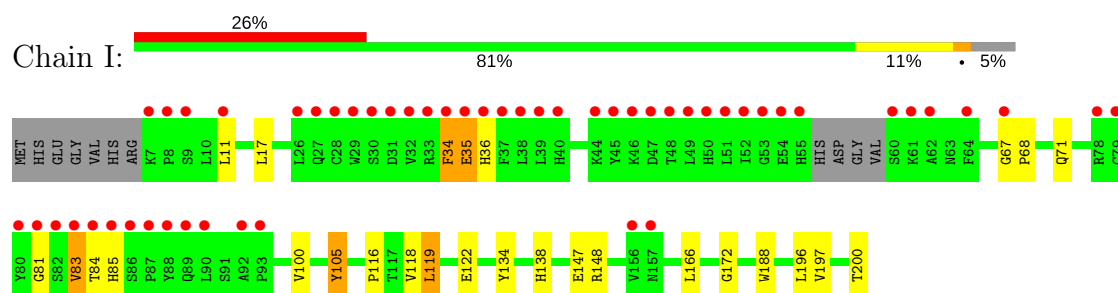
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

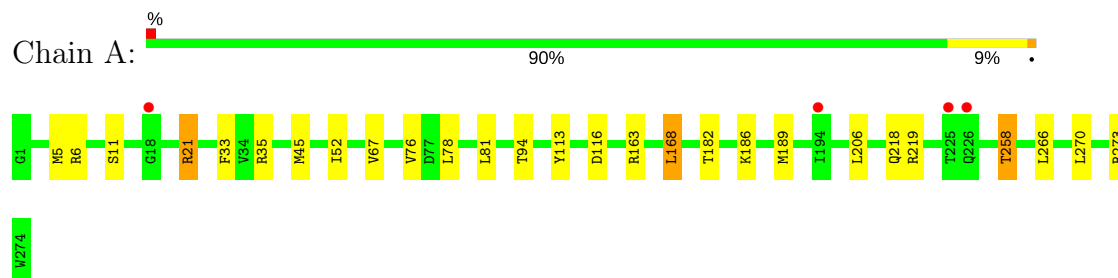
- Molecule 1: Killer cell immunoglobulin-like receptor 2DS2



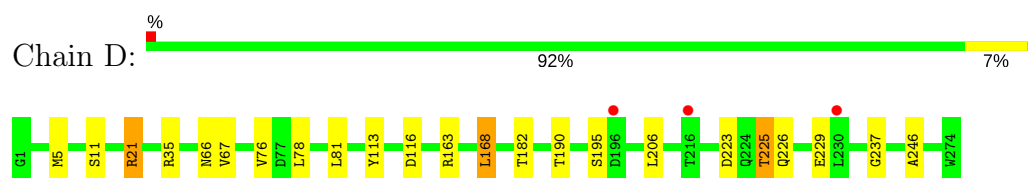
- Molecule 1: Killer cell immunoglobulin-like receptor 2DS2



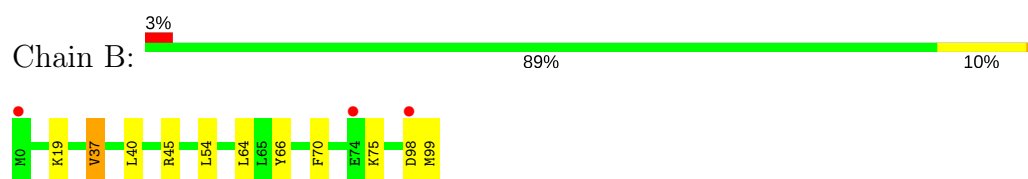
- Molecule 2: HLA class I histocompatibility antigen, A-11 alpha chain



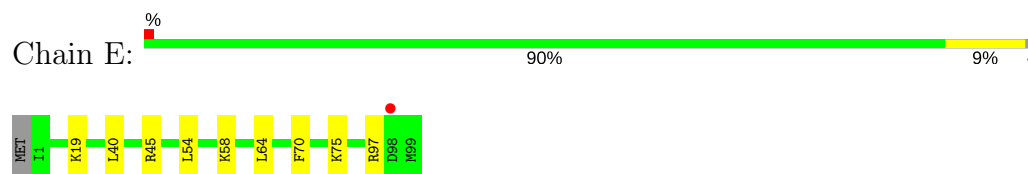
- Molecule 2: HLA class I histocompatibility antigen, A-11 alpha chain



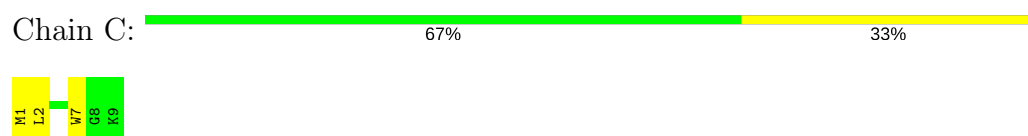
- Molecule 3: Beta-2-microglobulin



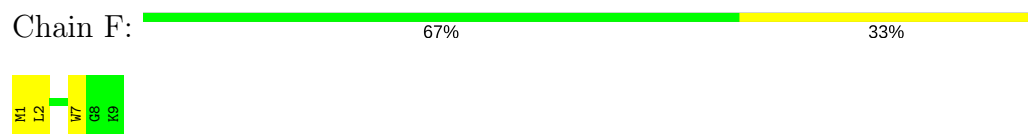
- Molecule 3: Beta-2-microglobulin



- Molecule 4: peptide from Virion membrane protein A14



- Molecule 4: peptide from Virion membrane protein A14



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.55Å 94.67Å 228.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	66.53 – 2.50 66.53 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.3 (66.53-2.50) 99.3 (66.53-2.50)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	9.75 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.222 , 0.257 0.225 , 0.257	Depositor DCC
R_{free} test set	2679 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	34.8	Xtriage
Anisotropy	0.222	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 40.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9364	wwPDB-VP
Average B, all atoms (Å ²)	36.0	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 28.69 % 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 1.7650e-03. 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 [i](#)

5.1 Standard geometry [i](#)

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	G	0.50	0/1516	0.71	3/2058 (0.1%)
1	I	0.47	0/1516	0.69	2/2058 (0.1%)
2	A	0.46	0/2297	0.71	0/3117
2	D	0.47	0/2297	0.71	0/3117
3	B	0.43	0/860	0.61	0/1162
3	E	0.45	0/852	0.63	0/1152
4	C	0.58	0/80	0.67	0/104
4	F	0.53	0/80	0.64	0/104
All	All	0.47	0/9498	0.69	5/12872 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	35	GLU	CB-CA-C	-5.96	98.49	110.40
1	G	85	HIS	N-CA-C	5.80	126.65	111.00
1	G	85	HIS	CB-CA-C	-5.45	99.49	110.40
1	I	148	ARG	NE-CZ-NH2	-5.28	117.66	120.30
1	G	148	ARG	NE-CZ-NH2	-5.22	117.69	120.30

There are no chirality outliers.

There are no planarity outliers.

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	G	1473	0	1404	19	0
1	I	1473	0	1404	23	0
2	A	2236	0	2085	22	0
2	D	2236	0	2085	8	0
3	B	837	0	803	3	0
3	E	829	0	794	3	0
4	C	78	0	82	5	0
4	F	78	0	82	3	0
5	A	20	0	0	1	0
5	B	7	0	0	0	0
5	C	1	0	0	0	0
5	D	26	0	0	1	0
5	E	8	0	0	0	0
5	F	2	0	0	0	0
5	G	31	0	0	1	0
5	I	29	0	0	1	0
All	All	9364	0	8739	78	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 4.

All (78) 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:I:34:PHE:CE1	1:I:81:GLY:HA3	1.73	1.22
1:I:84:THR:HG22	1:I:85:HIS:H	1.15	1.06
1:I:34:PHE:CD1	1:I:81:GLY:HA3	1.95	1.00
1:I:34:PHE:HE1	1:I:81:GLY:HA3	1.24	0.95
2:A:45:MET:HE2	2:A:67:VAL:HG21	1.54	0.90
1:I:34:PHE:CE1	1:I:81:GLY:CA	2.61	0.84
2:A:45:MET:CE	2:A:67:VAL:HG21	2.06	0.83
1:G:82:SER:O	1:G:83:VAL:HG22	1.79	0.82
2:A:6:ARG:HG2	2:A:113:TYR:OH	1.80	0.82
1:I:34:PHE:HE1	1:I:81:GLY:CA	1.97	0.77
1:G:172:GLY:HA3	5:G:321:HOH:O	1.90	0.71
1:I:84:THR:HG22	1:I:85:HIS:N	1.99	0.71
1:I:34:PHE:O	1:I:35:GLU:HB2	1.90	0.71
1:G:82:SER:O	1:G:83:VAL:CG2	2.41	0.68
1:I:84:THR:CG2	1:I:85:HIS:H	1.93	0.68
1:I:34:PHE:O	1:I:35:GLU:CB	2.43	0.66
2:A:163:ARG:NH2	4:C:1:MET:SD	2.69	0.65
2:A:5:MET:CE	2:A:33:PHE:HZ	2.10	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:118:VAL:CG1	1:I:122:GLU:HB2	2.26	0.65
1:G:118:VAL:CG1	1:G:122:GLU:HB2	2.26	0.65
2:A:218:GLN:HB2	2:A:258:THR:HG23	1.80	0.64
1:I:71:GLN:HG2	2:D:76:VAL:HG11	1.80	0.64
2:A:5:MET:CE	2:A:33:PHE:CZ	2.81	0.63
1:G:82:SER:C	1:G:83:VAL:CG2	2.68	0.62
1:G:85:HIS:CD2	1:G:85:HIS:C	2.71	0.62
2:A:266:LEU:HD13	2:A:270:LEU:HD13	1.82	0.61
1:I:83:VAL:O	1:I:83:VAL:HG13	2.02	0.59
1:I:34:PHE:CD1	1:I:81:GLY:CA	2.78	0.58
3:B:37:VAL:HG13	3:B:66:TYR:CE1	2.40	0.57
1:G:85:HIS:O	1:G:85:HIS:HD2	1.89	0.54
2:D:5:MET:HB2	2:D:168:LEU:HG	1.89	0.54
2:A:67:VAL:HG23	4:C:2:LEU:HD21	1.89	0.53
1:G:82:SER:C	1:G:83:VAL:HG23	2.27	0.53
2:A:5:MET:HB2	2:A:168:LEU:HG	1.91	0.52
2:A:5:MET:HE3	2:A:33:PHE:CZ	2.46	0.51
2:A:45:MET:HE1	2:A:67:VAL:HG21	1.89	0.49
2:A:186:LYS:NZ	5:A:318:HOH:O	2.46	0.49
2:A:5:MET:HE2	2:A:33:PHE:HZ	1.78	0.48
2:A:33:PHE:HD2	2:A:52:ILE:HD12	1.78	0.48
1:I:118:VAL:HG12	1:I:119:LEU:O	2.13	0.48
2:D:229:GLU:HG3	2:D:246:ALA:HB3	1.96	0.48
1:G:118:VAL:HG12	1:G:119:LEU:O	2.15	0.47
2:A:266:LEU:HD22	2:A:270:LEU:HD11	1.97	0.47
3:B:54:LEU:HA	3:B:64:LEU:HD13	1.96	0.46
1:I:67:GLY:O	1:I:68:PRO:C	2.54	0.46
1:I:34:PHE:HD1	1:I:81:GLY:HA3	1.68	0.46
1:G:67:GLY:O	1:G:68:PRO:C	2.53	0.46
3:E:54:LEU:HA	3:E:64:LEU:HD13	1.99	0.45
2:A:189:MET:HE1	2:A:273:ARG:C	2.37	0.45
1:G:83:VAL:HA	1:G:84:THR:HA	1.58	0.45
2:A:5:MET:HE1	4:C:1:MET:H2	1.83	0.44
1:I:100:VAL:HG11	1:I:188:TRP:CD2	2.52	0.44
1:G:34:PHE:CE1	1:G:81:GLY:HA3	2.53	0.44
2:A:67:VAL:CG2	4:C:2:LEU:HD21	2.47	0.44
1:G:138:HIS:HB3	1:G:147:GLU:HG3	2.00	0.44
1:G:116:PRO:O	1:G:197:VAL:HA	2.18	0.43
4:F:2:LEU:HD12	4:F:2:LEU:C	2.39	0.43
2:D:163:ARG:NH2	4:F:1:MET:SD	2.91	0.43
1:G:100:VAL:HG11	1:G:188:TRP:CD2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:11:SER:HA	2:A:21:ARG:O	2.19	0.42
1:I:138:HIS:HB3	1:I:147:GLU:HG3	2.01	0.42
2:A:67:VAL:CG2	4:C:2:LEU:CD2	2.98	0.42
2:D:11:SER:HA	2:D:21:ARG:O	2.20	0.42
3:E:58:LYS:HD2	3:E:58:LYS:HA	1.84	0.42
1:I:116:PRO:O	1:I:197:VAL:HA	2.19	0.42
1:I:105:TYR:HB2	1:I:134:TYR:OH	2.20	0.42
2:D:225:THR:O	2:D:226:GLN:HB2	2.20	0.41
2:A:33:PHE:CD2	2:A:52:ILE:HD12	2.56	0.41
1:G:105:TYR:HB2	1:G:134:TYR:OH	2.21	0.41
1:G:85:HIS:CD2	1:G:88:TYR:CE1	3.09	0.41
3:B:40:LEU:HD23	3:B:45:ARG:HA	2.03	0.41
1:G:85:HIS:CD2	1:G:85:HIS:O	2.69	0.41
1:I:200:THR:OXT	5:I:325:HOH:O	2.22	0.41
3:E:40:LEU:HD23	3:E:45:ARG:HA	2.02	0.41
1:G:75:GLY:HA2	1:G:186:TYR:CD2	2.55	0.41
1:I:83:VAL:HA	1:I:84:THR:HA	1.73	0.41
2:D:237:GLY:N	5:D:315:HOH:O	2.47	0.40
2:D:66:ASN:CB	4:F:2:LEU:HD11	2.52	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	G	186/201 (92%)	175 (94%)	9 (5%)	2 (1%)	17	29
1	I	186/201 (92%)	171 (92%)	13 (7%)	2 (1%)	17	29
2	A	272/274 (99%)	267 (98%)	5 (2%)	0	100	100
2	D	272/274 (99%)	269 (99%)	3 (1%)	0	100	100
3	B	98/100 (98%)	96 (98%)	2 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	E	97/100 (97%)	96 (99%)	1 (1%)	0	100	100
4	C	7/9 (78%)	7 (100%)	0	0	100	100
4	F	7/9 (78%)	7 (100%)	0	0	100	100
All	All	1125/1168 (96%)	1088 (97%)	33 (3%)	4 (0%)	38	59

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	I	172	GLY
1	G	83	VAL
1	G	172	GLY
1	I	83	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	G	164/173 (95%)	157 (96%)	7 (4%)	33	58
1	I	164/173 (95%)	156 (95%)	8 (5%)	29	52
2	A	231/231 (100%)	219 (95%)	12 (5%)	27	49
2	D	231/231 (100%)	217 (94%)	14 (6%)	22	40
3	B	95/95 (100%)	89 (94%)	6 (6%)	21	38
3	E	94/95 (99%)	90 (96%)	4 (4%)	33	58
4	C	8/8 (100%)	7 (88%)	1 (12%)	5	10
4	F	8/8 (100%)	7 (88%)	1 (12%)	5	10
All	All	995/1014 (98%)	942 (95%)	53 (5%)	26	48

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

Mol	Chain	Res	Type
1	G	11	LEU
1	G	17	LEU

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Mol	Chain	Res	Type
1	G	71	GLN
1	G	105	TYR
1	G	119	LEU
1	G	166	LEU
1	G	196	LEU
1	I	11	LEU
1	I	17	LEU
1	I	34	PHE
1	I	36	HIS
1	I	105	TYR
1	I	119	LEU
1	I	166	LEU
1	I	196	LEU
2	A	21	ARG
2	A	35	ARG
2	A	76	VAL
2	A	78	LEU
2	A	81	LEU
2	A	94	THR
2	A	116	ASP
2	A	168	LEU
2	A	182	THR
2	A	206	LEU
2	A	219	ARG
2	A	258	THR
3	B	19	LYS
3	B	37	VAL
3	B	70	PHE
3	B	75	LYS
3	B	98	ASP
3	B	99	MET
4	C	7	TRP
2	D	21	ARG
2	D	35	ARG
2	D	67	VAL
2	D	78	LEU
2	D	81	LEU
2	D	113	TYR
2	D	116	ASP
2	D	168	LEU
2	D	182	THR
2	D	190	THR

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Mol	Chain	Res	Type
2	D	195	SER
2	D	206	LEU
2	D	223	ASP
2	D	225	THR
3	E	19	LYS
3	E	70	PHE
3	E	75	LYS
3	E	97	ARG
4	F	7	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	G	85	HIS
1	G	157	ASN
2	D	174	ASN

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](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

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	G	190/201 (94%)	0.54	13 (6%) 18 18	9, 32, 73, 126	0
1	I	190/201 (94%)	1.79	53 (27%) 1 0	10, 34, 141, 181	0
2	A	274/274 (100%)	0.21	4 (1%) 74 75	15, 30, 54, 76	0
2	D	274/274 (100%)	0.21	3 (1%) 80 81	15, 28, 55, 78	0
3	B	100/100 (100%)	0.27	3 (3%) 51 53	19, 31, 60, 85	0
3	E	99/100 (99%)	0.30	1 (1%) 82 83	20, 30, 57, 93	0
4	C	9/9 (100%)	0.47	0 100 100	20, 30, 38, 46	0
4	F	9/9 (100%)	0.17	0 100 100	12, 20, 33, 35	0
All	All	1145/1168 (98%)	0.54	77 (6%) 19 19	9, 30, 85, 181	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	86	SER	24.1
1	I	87	PRO	19.4
1	I	90	LEU	11.5
1	I	89	GLN	11.0
1	I	81	GLY	10.1
1	I	38	LEU	10.1
1	I	34	PHE	9.6
1	I	51	LEU	9.3
1	I	83	VAL	8.4
1	G	87	PRO	8.1
1	I	37	PHE	8.1
3	E	98	ASP	7.5
1	I	88	TYR	7.3
1	I	60	SER	6.7
1	I	35	GLU	6.6
1	I	36	HIS	6.4

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Mol	Chain	Res	Type	RSRZ
1	I	93	PRO	6.3
1	G	83	VAL	6.2
1	I	62	ALA	6.0
1	I	54	GLU	5.9
1	G	86	SER	5.7
1	I	50	HIS	5.6
1	I	92	ALA	5.6
1	I	32	VAL	5.6
1	I	30	SER	5.4
1	I	84	THR	5.4
1	I	29	TRP	4.9
1	G	84	THR	4.8
3	B	98	ASP	4.6
1	I	48	THR	4.5
1	I	31	ASP	4.5
1	I	82	SER	4.2
1	I	53	GLY	4.1
1	I	45	TYR	4.1
3	B	0	MET	4.1
1	I	28	CYS	4.0
1	I	61	LYS	3.9
1	I	49	LEU	3.9
1	G	45	TYR	3.7
1	I	8	PRO	3.7
1	I	85	HIS	3.6
1	I	64	PHE	3.6
1	I	79	CYS	3.6
1	G	85	HIS	3.5
1	G	67	GLY	3.5
1	I	27	GLN	3.3
1	I	55	HIS	3.2
2	D	196	ASP	3.1
1	I	33	ARG	3.1
1	I	52	ILE	3.0
2	A	194	ILE	3.0
1	I	78	ARG	2.9
1	I	67	GLY	2.9
1	I	156	VAL	2.9
1	G	88	TYR	2.7
1	G	90	LEU	2.7
1	I	157	ASN	2.7
1	I	40	HIS	2.7

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Mol	Chain	Res	Type	RSRZ
1	I	44	LYS	2.7
1	I	11	LEU	2.6
1	I	46	LYS	2.6
2	A	226	GLN	2.6
1	I	39	LEU	2.5
1	I	80	TYR	2.4
3	B	74	GLU	2.4
1	I	47	ASP	2.4
1	I	7	LYS	2.4
1	G	50	HIS	2.4
1	I	26	LEU	2.3
2	A	18	GLY	2.2
2	D	230	LEU	2.2
2	D	216	THR	2.2
2	A	225	THR	2.1
1	I	9	SER	2.1
1	G	32	VAL	2.0
1	G	34	PHE	2.0
1	G	29	TRP	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.