



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 08:05 AM GMT

PDB ID : 4E5X
Title : Crystal structure of a complex between the human adenovirus type 2 E3-19K protein and MHC class I molecule HLA-A2/Tax
Authors : Li, L.; Bouvier, M.
Deposited on : 2012-03-14
Resolution : 1.95 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

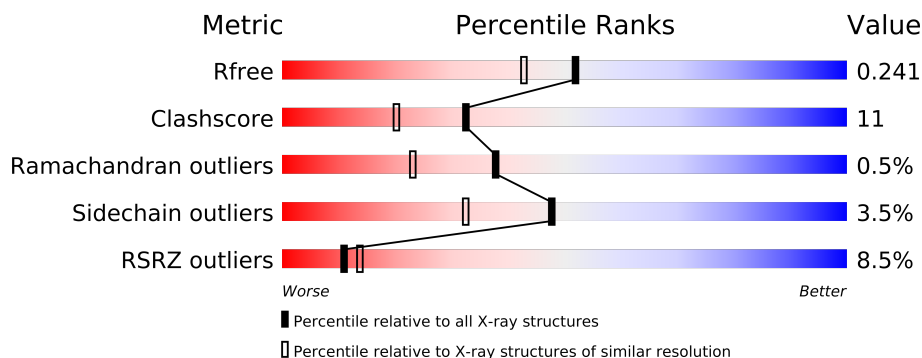
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.95 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}	66092	1321 (1.96-1.96)
Clashscore	79885	1488 (1.96-1.96)
Ramachandran outliers	78287	1475 (1.96-1.96)
Sidechain outliers	78261	1475 (1.96-1.96)
RSRZ outliers	66119	1321 (1.96-1.96)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	275	
1	D	275	
2	B	100	
2	E	100	
3	C	9	
3	F	9	
4	G	100	
4	H	100	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8439 atoms, of which 0 are hydrogen and 0 are deuterium.

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 HLA class I histocompatibility antigen, A-2 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	S	0	0	0
			2247	1403	409	426	9			
1	D	275	Total	C	N	O	S	0	0	0
			2247	1403	409	426	9			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	E	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	INITIATING METHIONINE	UNP P61769
E	0	MET	-	INITIATING METHIONINE	UNP P61769

- Molecule 3 is a protein called Protein Tax-1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	0	0	0
			77	56	9	12			
3	F	9	Total	C	N	O	0	0	0
			77	56	9	12			

- Molecule 4 is a protein called Early E3 18.5 kDa glycoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	G	99	Total	C	N	O	S	0	0	0
			832	539	134	151	8			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	97	Total 813	C 528	N 130	O 147	S 8	0	0	0

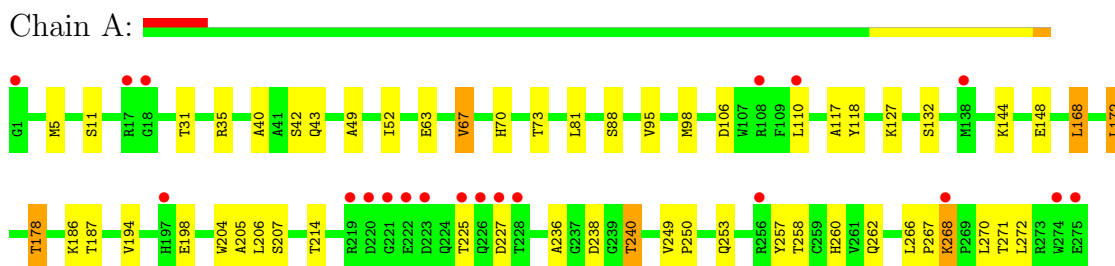
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	107	Total 107	O 107	0	0
5	B	52	Total 52	O 52	0	0
5	C	10	Total 10	O 10	0	0
5	D	101	Total 101	O 101	0	0
5	E	35	Total 35	O 35	0	0
5	F	4	Total 4	O 4	0	0
5	G	82	Total 82	O 82	0	0
5	H	81	Total 81	O 81	0	0

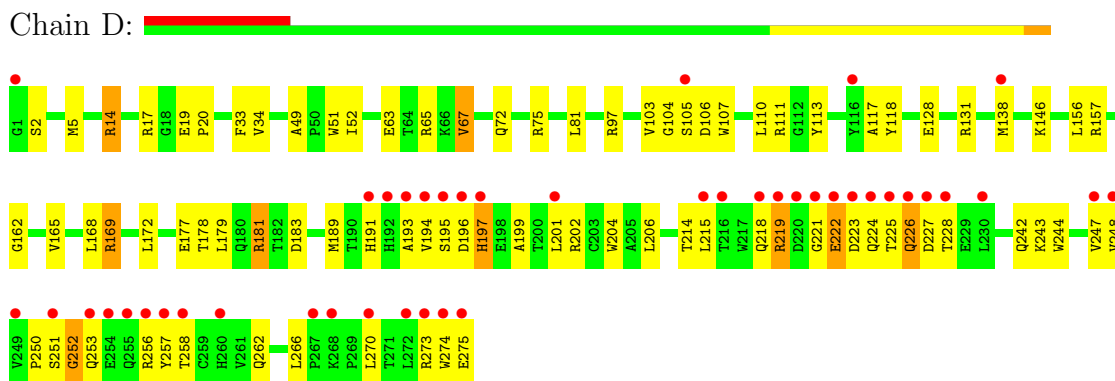
3 Residue-property plots

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.

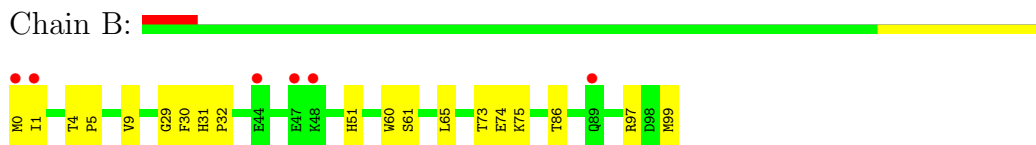
- Molecule 1: HLA class I histocompatibility antigen, A-2 alpha chain



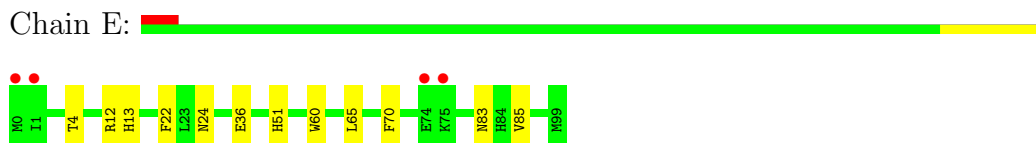
- Molecule 1: HLA class I histocompatibility antigen, A-2 alpha chain



- Molecule 2: Beta-2-microglobulin

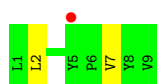


- Molecule 2: Beta-2-microglobulin



- Molecule 3: Protein Tax-1





- Molecule 3: Protein Tax-1

Chain F:



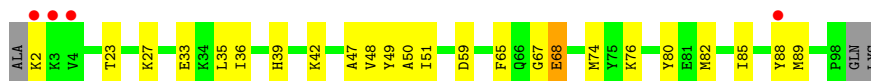
- Molecule 4: Early E3 18.5 kDa glycoprotein

Chain G:



- Molecule 4: Early E3 18.5 kDa glycoprotein

Chain H:



4 Data and refinement statistics

Property	Value	Source
Space group	I 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	113.19Å 133.67Å 196.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.06 – 1.95 49.18 – 1.95	Depositor EDS
% Data completeness (in resolution range)	94.9 (37.06-1.95) 94.6 (49.18-1.95)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.38 (at 1.95Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.210 , 0.241 0.210 , 0.241	Depositor DCC
R_{free} test set	10300 reflections (10.01%)	DCC
Wilson B-factor (Å ²)	25.4	Xtriage
Anisotropy	0.497	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 43.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 108334 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8439	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.26% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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	0.32	0/2312	0.59	0/3137
1	D	0.33	0/2312	0.61	0/3137
2	B	0.34	0/860	0.62	0/1162
2	E	0.32	0/860	0.59	0/1162
3	C	0.51	0/80	0.81	0/108
3	F	0.44	0/80	0.85	0/108
4	G	0.39	0/855	0.61	0/1150
4	H	0.37	0/836	0.57	0/1127
All	All	0.34	0/8195	0.61	0/11091

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
4	H	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	H	80	TYR	Sidechain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit,

and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2247	0	2096	47	0
1	D	2247	0	2096	82	0
2	B	837	0	803	14	0
2	E	837	0	803	13	0
3	C	77	0	79	1	0
3	F	77	0	79	3	0
4	G	832	0	817	7	0
4	H	813	0	796	30	1
5	A	107	0	0	6	0
5	B	52	0	0	1	0
5	C	10	0	0	0	0
5	D	101	0	0	3	0
5	E	35	0	0	3	0
5	F	4	0	0	0	0
5	G	82	0	0	0	0
5	H	81	0	0	2	0
All	All	8439	0	7569	176	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 11.

All (176) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:219:ARG:HE	1:D:224:GLN:NE2	1.59	1.00
1:D:63:GLU:O	1:D:67:VAL:HG23	1.61	0.99
1:A:63:GLU:O	1:A:67:VAL:HG23	1.72	0.89
1:D:14:ARG:HD2	1:D:19:GLU:O	1.76	0.84
1:A:266:LEU:HD13	1:A:270:LEU:HD13	1.59	0.82
2:B:73:THR:HG22	2:B:75:LYS:H	1.46	0.81
2:B:0:MET:HG2	2:B:1:ILE:H	1.47	0.79
1:D:228:THR:HG22	1:D:247:VAL:CG1	2.14	0.77
1:D:191:HIS:CE1	1:D:199:ALA:HB1	2.23	0.74
1:D:258:THR:HG22	1:D:273:ARG:HG2	1.68	0.73
1:A:168:LEU:HD22	1:A:172:LEU:HD22	1.72	0.72
2:E:85:VAL:HG13	5:E:101:HOH:O	1.88	0.72
1:A:187:THR:HG22	5:A:392:HOH:O	1.90	0.72
4:H:68:GLU:H	4:H:68:GLU:CD	1.94	0.71
1:D:228:THR:HG22	1:D:247:VAL:HG12	1.72	0.71
1:D:219:ARG:NE	1:D:224:GLN:NE2	2.36	0.71

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:63:GLU:O	1:A:67:VAL:CG2	2.38	0.70
1:A:127:LYS:HD2	1:A:132:SER:OG	1.91	0.70
1:D:20:PRO:HD2	1:D:75:ARG:HD3	1.75	0.69
1:D:65:ARG:HD2	5:D:401:HOH:O	1.93	0.69
1:D:63:GLU:O	1:D:67:VAL:CG2	2.41	0.68
1:D:226:GLN:O	1:D:227:ASP:HB2	1.94	0.67
1:D:67:VAL:HG22	3:F:2:LEU:HD21	1.75	0.67
1:D:67:VAL:HG22	3:F:2:LEU:CD2	2.24	0.67
1:A:194:VAL:CG2	1:A:198:GLU:HG3	2.26	0.66
1:A:238:ASP:OD1	1:A:240:THR:HG22	1.97	0.65
1:D:146:LYS:HE3	3:F:8:TYR:O	1.97	0.65
1:A:249:VAL:HG13	1:A:250:PRO:HD2	1.79	0.65
1:A:205:ALA:C	1:A:206:LEU:HD12	2.17	0.64
4:H:23:THR:HG22	5:H:204:HOH:O	1.98	0.64
1:D:215:LEU:HD13	1:D:243:LYS:HD3	1.80	0.64
4:H:35:LEU:O	4:H:36:ILE:HD12	1.99	0.63
2:B:0:MET:HG2	2:B:1:ILE:N	2.14	0.62
1:A:204:TRP:HB3	1:A:206:LEU:HD11	1.81	0.62
1:D:226:GLN:O	1:D:227:ASP:CB	2.47	0.62
5:E:106:HOH:O	4:H:89:MET:HE2	1.98	0.62
1:A:144:LYS:O	1:A:148:GLU:HG3	2.00	0.62
1:D:178:THR:HG23	4:H:51:ILE:HD11	1.82	0.62
4:H:76:LYS:HE3	5:H:261:HOH:O	2.00	0.62
1:D:107:TRP:HZ3	1:D:172:LEU:HD13	1.65	0.61
1:D:202:ARG:HD2	1:D:204:TRP:NE1	2.15	0.61
1:D:131:ARG:HD3	1:D:157:ARG:NH2	2.14	0.61
4:H:85:ILE:O	4:H:88:TYR:HD2	1.83	0.60
2:E:12:ARG:HD2	4:H:89:MET:CE	2.31	0.60
1:D:250:PRO:HB2	1:D:253:GLN:HG3	1.83	0.60
4:H:35:LEU:C	4:H:36:ILE:HD12	2.23	0.59
2:B:4:THR:OG1	2:B:86:THR:HG21	2.02	0.59
1:A:258:THR:HG23	1:A:260:HIS:NE2	2.18	0.59
1:A:225:THR:C	1:A:227:ASP:H	2.05	0.59
1:A:207:SER:HA	1:A:240:THR:HG21	1.84	0.58
1:D:219:ARG:HE	1:D:224:GLN:CD	2.08	0.57
1:A:5:MET:HB2	1:A:168:LEU:HG	1.86	0.57
1:D:162:GLY:O	1:D:165:VAL:HG22	2.05	0.57
1:D:219:ARG:HE	1:D:224:GLN:HE22	1.46	0.57
1:A:194:VAL:HG21	1:A:198:GLU:HG3	1.86	0.56
1:A:207:SER:HA	1:A:240:THR:CG2	2.35	0.56
2:E:24:ASN:HB3	2:E:65:LEU:HD11	1.88	0.56
1:D:49:ALA:O	1:D:52:ILE:HG22	2.05	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:H:68:GLU:N	4:H:68:GLU:CD	2.59	0.55
1:D:253:GLN:HB3	1:D:256:ARG:CD	2.36	0.55
4:G:3:LYS:HB2	4:G:3:LYS:NZ	2.21	0.55
1:D:5:MET:HB2	1:D:168:LEU:HD13	1.87	0.55
1:D:250:PRO:O	1:D:252:GLY:N	2.40	0.55
1:D:117:ALA:HB2	2:E:60:TRP:CE2	2.42	0.55
1:A:236:ALA:HB3	1:A:240:THR:HG22	1.90	0.54
1:D:202:ARG:HD2	1:D:204:TRP:CE2	2.42	0.54
2:E:13:HIS:NE2	4:H:89:MET:HE2	2.22	0.54
1:A:70:HIS:O	1:A:73:THR:HG22	2.08	0.54
4:G:57:THR:OG1	4:G:59:ASP:OD1	2.23	0.53
4:H:27:LYS:HG2	4:H:47:ALA:HB2	1.90	0.53
1:A:40:ALA:HB3	5:A:379:HOH:O	2.09	0.53
4:H:23:THR:HG23	4:H:49:TYR:CE1	2.43	0.53
1:A:168:LEU:HD22	1:A:172:LEU:CD2	2.39	0.53
2:E:22:PHE:HD2	4:H:89:MET:CE	2.23	0.52
1:D:266:LEU:HD13	1:D:270:LEU:HD13	1.90	0.52
1:A:49:ALA:O	1:A:52:ILE:HG22	2.08	0.52
1:A:258:THR:HG22	5:A:374:HOH:O	2.08	0.52
1:D:191:HIS:NE2	1:D:199:ALA:HB1	2.25	0.52
1:D:177:GLU:HB3	4:H:51:ILE:HD13	1.92	0.51
1:D:191:HIS:CE1	1:D:199:ALA:CB	2.93	0.51
2:E:12:ARG:HD2	4:H:89:MET:HE3	1.92	0.51
1:D:219:ARG:HA	1:D:256:ARG:O	2.10	0.51
1:D:214:THR:HB	1:D:262:GLN:HB2	1.93	0.50
1:D:105:SER:O	1:D:106:ASP:HB2	2.11	0.50
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.46	0.50
4:H:36:ILE:HD13	4:H:65:PHE:CE2	2.46	0.50
1:D:110:LEU:HB3	1:D:111:ARG:NH1	2.27	0.49
1:D:189:MET:HE3	1:D:201:LEU:HD22	1.95	0.49
1:D:191:HIS:HD2	1:D:201:LEU:HD21	1.77	0.49
1:D:253:GLN:HB3	1:D:256:ARG:HD2	1.93	0.49
1:D:17:ARG:HH11	1:D:17:ARG:HG3	1.77	0.48
1:A:249:VAL:HG13	1:A:257:TYR:CE2	2.48	0.48
4:H:59:ASP:OD1	4:H:76:LYS:HG3	2.13	0.48
1:D:202:ARG:NH1	1:D:244:TRP:CZ3	2.81	0.48
1:A:214:THR:HB	1:A:262:GLN:HB2	1.95	0.48
1:D:177:GLU:CB	4:H:51:ILE:HD13	2.43	0.48
2:B:5:PRO:HB3	2:B:30:PHE:HB3	1.96	0.48
1:D:138:MET:O	1:D:138:MET:HE2	2.12	0.48
1:D:224:GLN:NE2	1:D:224:GLN:HA	2.28	0.48
1:D:104:GLY:O	1:D:106:ASP:O	2.32	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:97:ARG:HG3	2:B:97:ARG:O	2.14	0.48
1:A:204:TRP:HB3	1:A:206:LEU:CD1	2.43	0.47
4:H:82:MET:HA	4:H:85:ILE:HD12	1.96	0.47
1:A:73:THR:HG21	5:A:310:HOH:O	2.14	0.47
1:D:97:ARG:NH2	5:D:361:HOH:O	2.39	0.47
1:A:266:LEU:HD22	1:A:270:LEU:HD11	1.95	0.47
4:G:3:LYS:HG3	4:G:4:VAL:N	2.30	0.47
2:B:73:THR:HG22	2:B:74:GLU:N	2.29	0.46
2:E:4:THR:HG22	5:E:103:HOH:O	2.15	0.46
4:H:2:LYS:O	4:H:2:LYS:HG2	2.16	0.46
2:E:22:PHE:HD2	4:H:89:MET:HE1	1.81	0.46
1:D:177:GLU:OE2	4:H:42:LYS:HE3	2.16	0.46
1:D:193:ALA:C	1:D:195:SER:H	2.19	0.46
1:D:274:TRP:O	1:D:275:GLU:HB2	2.15	0.46
1:D:107:TRP:CZ3	1:D:172:LEU:HD13	2.49	0.46
1:D:189:MET:HE1	1:D:274:TRP:HB2	1.98	0.46
1:D:202:ARG:HG3	1:D:202:ARG:HH11	1.80	0.46
2:E:51:HIS:HA	2:E:65:LEU:O	2.16	0.45
1:A:194:VAL:HG22	1:A:198:GLU:O	2.16	0.45
1:D:266:LEU:HD22	1:D:270:LEU:CD1	2.45	0.45
1:D:221:GLY:O	1:D:222:GLU:HB2	2.16	0.45
2:B:51:HIS:HA	2:B:65:LEU:O	2.17	0.45
1:A:81:LEU:HD13	1:A:118:TYR:CD1	2.51	0.45
1:D:253:GLN:HB3	1:D:256:ARG:NE	2.31	0.45
1:D:107:TRP:CD1	1:D:169:ARG:NH2	2.85	0.45
1:D:218:GLN:O	1:D:257:TYR:HA	2.17	0.45
1:D:191:HIS:HD2	1:D:201:LEU:CD2	2.30	0.45
1:A:11:SER:HB3	1:A:95:VAL:CG1	2.47	0.45
1:D:81:LEU:HD13	1:D:118:TYR:CD1	2.52	0.45
4:H:35:LEU:C	4:H:35:LEU:HD12	2.38	0.44
1:A:258:THR:OG1	1:A:271:THR:HG23	2.18	0.44
1:D:224:GLN:O	1:D:226:GLN:N	2.50	0.44
2:B:29:GLY:HA2	2:B:61:SER:HB2	1.99	0.44
2:B:9:VAL:HG13	5:B:114:HOH:O	2.17	0.44
1:A:272:LEU:N	1:A:272:LEU:HD22	2.32	0.44
1:D:248:VAL:HG13	1:D:248:VAL:O	2.18	0.44
2:B:73:THR:HG22	2:B:75:LYS:N	2.25	0.44
1:A:214:THR:HG22	5:A:339:HOH:O	2.17	0.44
2:B:31:HIS:ND1	2:B:32:PRO:HA	2.33	0.44
1:A:204:TRP:CZ3	2:B:99:MET:HB3	2.53	0.43
1:D:206:LEU:HD23	1:D:242:GLN:HB3	1.99	0.43
1:D:219:ARG:CG	1:D:224:GLN:OE1	2.67	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:E:36:GLU:HB2	2:E:83:ASN:HB3	2.01	0.43
1:A:42:SER:O	1:A:43:GLN:HB2	2.19	0.43
1:D:51:TRP:CZ2	1:D:179:LEU:HD11	2.54	0.43
1:A:225:THR:C	1:A:227:ASP:N	2.72	0.43
1:D:14:ARG:HG2	5:D:368:HOH:O	2.17	0.43
1:A:178:THR:CG2	4:G:49:TYR:OH	2.67	0.43
1:D:189:MET:CE	1:D:201:LEU:HD22	2.49	0.43
1:A:250:PRO:HG2	1:A:253:GLN:HB2	2.00	0.43
1:D:189:MET:HE2	1:D:274:TRP:HE3	1.83	0.42
1:A:73:THR:OG1	3:C:7:VAL:O	2.24	0.42
1:A:186:LYS:HE2	4:G:95:LEU:HD22	2.02	0.42
4:H:39:HIS:HB3	4:H:48:VAL:HG11	2.01	0.42
1:D:219:ARG:HG3	1:D:224:GLN:OE1	2.19	0.42
2:E:22:PHE:CD2	4:H:89:MET:HE1	2.54	0.42
1:D:177:GLU:OE2	4:H:50:ALA:HA	2.20	0.42
4:G:32:HIS:HB3	4:G:65:PHE:O	2.20	0.42
4:G:35:LEU:C	4:G:35:LEU:HD12	2.39	0.42
1:A:267:PRO:HB2	1:A:268:LYS:HE3	2.00	0.42
1:D:33:PHE:CD1	1:D:34:VAL:HG13	2.55	0.42
1:D:189:MET:CE	1:D:274:TRP:HB2	2.50	0.41
1:A:206:LEU:HD12	1:A:206:LEU:N	2.35	0.41
1:A:73:THR:HG23	5:A:377:HOH:O	2.19	0.41
1:D:181:ARG:NH1	1:D:183:ASP:OD2	2.45	0.41
1:D:191:HIS:NE2	1:D:199:ALA:CB	2.84	0.41
1:D:196:ASP:HB3	1:D:197:HIS:H	1.58	0.41
1:D:215:LEU:N	1:D:215:LEU:HD12	2.35	0.41
4:H:33:GLU:CD	4:H:67:GLY:H	2.24	0.41
1:D:193:ALA:HA	1:D:199:ALA:HA	2.02	0.41
4:H:23:THR:CG2	4:H:49:TYR:HE1	2.34	0.41
1:D:253:GLN:NE2	1:D:256:ARG:NH1	2.69	0.40
1:D:2:SER:HB2	1:D:103:VAL:O	2.21	0.40
1:A:88:SER:N	2:E:36:GLU:OE2	2.35	0.40
4:H:88:TYR:CD1	4:H:89:MET:N	2.90	0.40
1:D:111:ARG:HD2	1:D:113:TYR:OH	2.21	0.40

All (1) 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	Distance(Å)	Clash(Å)
4:H:85:ILE:CD1	4:H:85:ILE:CD1[8_554]	2.18	0.02

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	273/275 (99%)	268 (98%)	5 (2%)	0	100	100
1	D	273/275 (99%)	257 (94%)	11 (4%)	5 (2%)	13	3
2	B	98/100 (98%)	96 (98%)	2 (2%)	0	100	100
2	E	98/100 (98%)	97 (99%)	1 (1%)	0	100	100
3	C	7/9 (78%)	7 (100%)	0	0	100	100
3	F	7/9 (78%)	7 (100%)	0	0	100	100
4	G	97/100 (97%)	94 (97%)	3 (3%)	0	100	100
4	H	95/100 (95%)	94 (99%)	1 (1%)	0	100	100
All	All	948/968 (98%)	920 (97%)	23 (2%)	5 (0%)	38	23

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	225	THR
1	D	251	SER
1	D	194	VAL
1	D	222	GLU
1	D	252	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	231/231 (100%)	220 (95%)	11 (5%)	35	17
1	D	231/231 (100%)	220 (95%)	11 (5%)	35	17
2	B	95/95 (100%)	95 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	E	95/95 (100%)	94 (99%)	1 (1%)	84	81
3	C	8/8 (100%)	7 (88%)	1 (12%)	7	1
3	F	8/8 (100%)	7 (88%)	1 (12%)	7	1
4	G	92/92 (100%)	89 (97%)	3 (3%)	50	35
4	H	90/92 (98%)	88 (98%)	2 (2%)	64	55
All	All	850/852 (100%)	820 (96%)	30 (4%)	48	32

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	THR
1	A	35	ARG
1	A	67	VAL
1	A	98	MET
1	A	106	ASP
1	A	110	LEU
1	A	168	LEU
1	A	172	LEU
1	A	178	THR
1	A	240	THR
1	A	268	LYS
3	C	2	LEU
1	D	14	ARG
1	D	67	VAL
1	D	72	GLN
1	D	128	GLU
1	D	156	LEU
1	D	169	ARG
1	D	181	ARG
1	D	197	HIS
1	D	219	ARG
1	D	223	ASP
1	D	226	GLN
2	E	70	PHE
3	F	2	LEU
4	G	3	LYS
4	G	25	LEU
4	G	68	GLU
4	H	68	GLU
4	H	74	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	141	GLN
1	A	218	GLN
1	D	72	GLN
1	D	191	HIS
1	D	218	GLN
1	D	224	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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	275/275 (100%)	0.38	20 (7%) 15 19	19, 33, 65, 96	0
1	D	275/275 (100%)	0.74	44 (16%) 3 2	17, 30, 83, 105	0
2	B	100/100 (100%)	0.40	6 (6%) 21 27	21, 33, 57, 87	1 (1%)
2	E	100/100 (100%)	0.25	4 (4%) 36 41	21, 36, 61, 81	0
3	C	9/9 (100%)	0.32	1 (11%) 6 7	25, 26, 29, 46	0
3	F	9/9 (100%)	0.78	0 100 100	20, 29, 40, 49	0
4	G	99/100 (99%)	0.21	3 (3%) 48 55	15, 24, 48, 81	2 (2%)
4	H	97/100 (97%)	0.36	4 (4%) 35 40	17, 26, 56, 100	1 (1%)
All	All	964/968 (99%)	0.45	82 (8%) 11 14	15, 30, 69, 105	4 (0%)

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	195	SER	10.3
1	D	225	THR	10.0
4	H	2	LYS	6.3
1	D	219	ARG	6.2
4	H	3	LYS	6.2
1	A	226	GLN	5.8
4	H	4	VAL	5.5
1	D	223	ASP	5.5
1	D	251	SER	5.3
1	D	275	GLU	5.1
1	D	193	ALA	5.1
1	D	194	VAL	5.0
1	D	196	ASP	4.9
1	D	226	GLN	4.8
1	D	197	HIS	4.6
1	D	268	LYS	4.0

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Mol	Chain	Res	Type	RSRZ
1	A	220	ASP	4.0
1	A	222	GLU	3.9
1	D	270	LEU	3.9
1	A	1	GLY	3.8
1	D	255	GLN	3.7
1	D	274	TRP	3.7
1	A	275	GLU	3.6
2	E	75	LYS	3.6
4	H	88	TYR	3.4
1	A	197	HIS	3.4
1	D	192	HIS	3.3
2	E	74	GLU	3.3
1	D	1	GLY	3.3
1	D	258	THR	3.3
1	D	227	ASP	3.3
1	D	105	SER	3.3
1	D	272	LEU	3.3
1	D	201	LEU	3.2
1	D	220	ASP	3.2
1	D	228	THR	3.2
1	A	268	LYS	3.1
1	D	222	GLU	3.1
1	D	257	TYR	3.1
1	A	227	ASP	3.1
2	B	48	LYS	3.0
1	D	224	GLN	3.0
2	E	0	MET	3.0
1	D	273	ARG	2.9
1	A	223	ASP	2.9
1	A	138	MET	2.9
1	D	248	VAL	2.9
1	D	267	PRO	2.9
1	D	254	GLU	2.8
1	A	225	THR	2.8
1	D	191	HIS	2.8
2	B	0	MET	2.7
1	A	108	ARG	2.6
1	A	221	GLY	2.6
1	A	256	ARG	2.5
1	A	228	THR	2.5
2	E	1	ILE	2.5
1	D	230	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
2	B	1	ILE	2.4
1	D	256	ARG	2.4
1	D	138	MET	2.4
4	G	4	VAL	2.4
1	A	18	GLY	2.4
1	D	216	THR	2.4
1	D	215	LEU	2.3
3	C	5	TYR	2.3
1	D	247	VAL	2.3
1	A	17	ARG	2.3
1	D	116	TYR	2.2
1	A	219	ARG	2.2
1	D	218	GLN	2.2
1	D	221	GLY	2.2
4	G	3	LYS	2.1
4	G	100	LYS	2.1
1	D	260	HIS	2.1
2	B	44	GLU	2.1
1	D	253	GLN	2.1
2	B	89	GLN	2.0
2	B	47	GLU	2.0
1	D	249	VAL	2.0
1	A	110	LEU	2.0
1	A	274	TRP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.