Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 07:52 PM GMT

PDB ID	:	1HOC
Title	:	THE THREE-DIMENSIONAL STRUCTURE OF H-2DB AT 2.4 ANGSTROMS RESOLUTION: IMPLICATIONS FOR ANTIGEN-DETERMINANT SELECTION
Authors	:	Young, A.C.M.; Zhang, W.; Sacchettini, J.C.
Deposited on	:	1994-01-02
Resolution	:	2.40 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 (i) symbol.

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

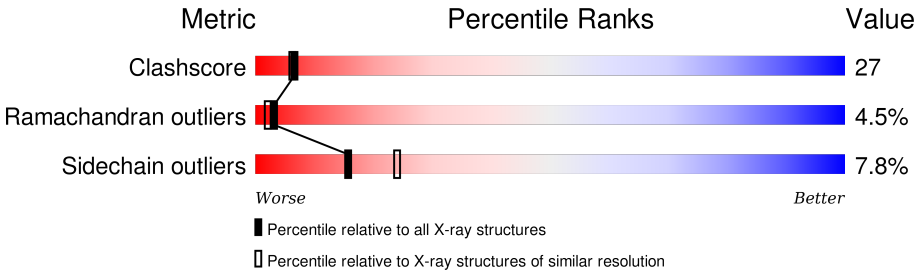
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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))
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)

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 <=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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	272	44% 41% 14%
2	B	99	46% 47% 5%
3	C	9	44% 44% 11%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3155 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 CLASS I HISTOCOMPATIBILITY ANTIGEN (H2-DB) (ALPHA CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	272	2224	1403	392	420	9	0	0	0

- Molecule 2 is a protein called BETA 2-MICROGLOBULIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	99	821	524	138	152	7	0	0	0

- Molecule 3 is a protein called 9-RESIDUE PEPTIDE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	9	69	38	11	18	2	0	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	22	22	22	0	0
4	B	17	17	17	0	0
4	C	2	2	2	0	0

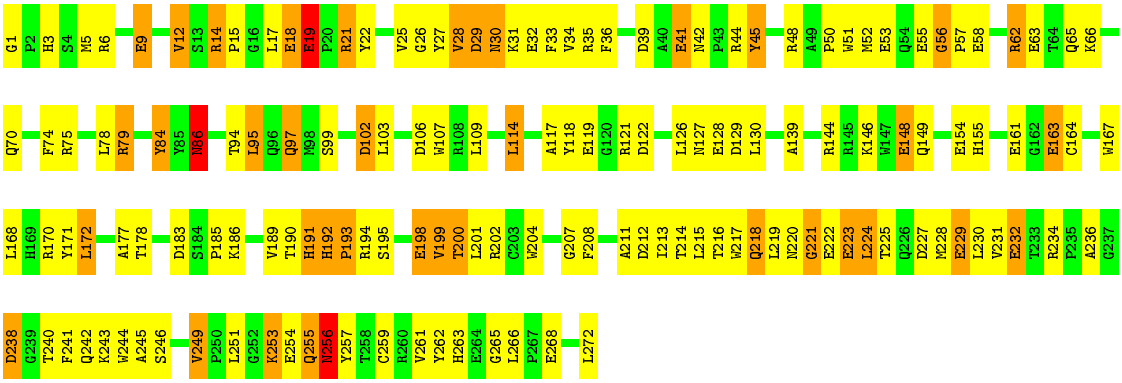
3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

Note EDS was not executed.

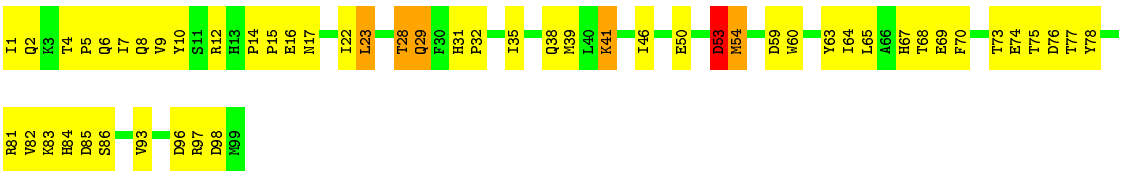
- Molecule 1: CLASS I HISTOCOMPATIBILITY ANTIGEN (H2-DB) (ALPHA CHAIN)

Chain A:



- Molecule 2: BETA 2-MICROGLOBULIN

Chain B:



- Molecule 3: 9-RESIDUE PEPTIDE

Chain C:

4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α, β, γ	92.92Å 94.54Å 133.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.40	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-2.40)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT, X-PLOR	Depositor
R, R_{free}	0.198 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3155	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

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	1.17	19/2288 (0.8%)	1.43	29/3105 (0.9%)
2	B	1.11	4/847 (0.5%)	1.45	11/1148 (1.0%)
3	C	1.47	2/68 (2.9%)	1.34	0/88
All	All	1.16	25/3203 (0.8%)	1.44	40/4341 (0.9%)

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	19	GLU	CD-OE1	7.68	1.34	1.25
1	A	53	GLU	CD-OE2	7.19	1.33	1.25
1	A	229	GLU	CD-OE1	7.16	1.33	1.25
1	A	148	GLU	CD-OE2	6.98	1.33	1.25
1	A	161	GLU	CD-OE1	6.42	1.32	1.25
3	C	7	GLU	CD-OE2	6.13	1.32	1.25
1	A	128	GLU	CD-OE2	6.05	1.32	1.25
1	A	163	GLU	CD-OE1	6.03	1.32	1.25
1	A	222	GLU	CD-OE1	6.00	1.32	1.25
1	A	55	GLU	CD-OE1	5.96	1.32	1.25
1	A	86	ASN	C-O	-5.87	1.12	1.23
1	A	63	GLU	CD-OE2	-5.82	1.19	1.25
1	A	223	GLU	CD-OE2	5.75	1.31	1.25
3	C	4	GLU	CD-OE2	5.71	1.31	1.25
1	A	198	GLU	CD-OE1	5.66	1.31	1.25
1	A	58	GLU	CD-OE1	5.51	1.31	1.25
2	B	50	GLU	CD-OE2	5.42	1.31	1.25
1	A	86	ASN	CG-OD1	-5.42	1.12	1.24
1	A	232	GLU	CD-OE1	5.37	1.31	1.25
2	B	69	GLU	CD-OE2	5.36	1.31	1.25
1	A	154	GLU	CD-OE2	5.28	1.31	1.25
1	A	9	GLU	CD-OE1	5.26	1.31	1.25
2	B	74	GLU	CD-OE1	5.23	1.31	1.25
1	A	63	GLU	CD-OE1	5.17	1.31	1.25
2	B	16	GLU	CD-OE1	5.13	1.31	1.25

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	54	MET	N-CA-CB	8.72	126.30	110.60
2	B	96	ASP	CB-CG-OD2	-8.43	110.71	118.30
1	A	56	GLY	C-N-CD	-7.80	103.43	120.60
2	B	53	ASP	CB-CG-OD2	-7.10	111.91	118.30
1	A	29	ASP	CB-CG-OD2	-7.01	111.99	118.30
1	A	106	ASP	CB-CG-OD2	-6.63	112.33	118.30
1	A	84	TYR	CB-CG-CD2	-6.62	117.03	121.00
2	B	96	ASP	CB-CG-OD1	6.54	124.19	118.30
1	A	45	TYR	CB-CG-CD1	6.47	124.88	121.00
1	A	118	TYR	CB-CG-CD1	-6.45	117.13	121.00
1	A	45	TYR	CB-CG-CD2	-6.41	117.15	121.00
2	B	98	ASP	CB-CG-OD1	-6.34	112.59	118.30
1	A	129	ASP	CB-CG-OD2	-6.34	112.59	118.30
2	B	59	ASP	CB-CG-OD2	-6.27	112.66	118.30
1	A	86	ASN	CB-CG-ND2	6.25	131.71	116.70
1	A	212	ASP	CB-CG-OD1	-6.03	112.88	118.30
1	A	102	ASP	CB-CG-OD2	-5.98	112.92	118.30
1	A	122	ASP	CB-CG-OD1	5.92	123.63	118.30
2	B	53	ASP	CB-CG-OD1	5.84	123.56	118.30
1	A	238	ASP	CB-CG-OD1	5.82	123.53	118.30
2	B	85	ASP	CB-CG-OD1	-5.79	113.09	118.30
2	B	53	ASP	CB-CA-C	-5.77	98.86	110.40
1	A	14	ARG	N-CA-C	-5.77	95.43	111.00
1	A	86	ASN	OD1-CG-ND2	-5.76	108.65	121.90
1	A	29	ASP	CB-CG-OD1	5.70	123.43	118.30
1	A	249	VAL	C-N-CD	-5.68	108.11	120.60
1	A	251	LEU	N-CA-CB	-5.64	99.13	110.40
1	A	238	ASP	CB-CG-OD2	-5.61	113.25	118.30
1	A	39	ASP	CB-CG-OD1	-5.44	113.40	118.30
1	A	218	GLN	CB-CA-C	-5.40	99.59	110.40
2	B	76	ASP	CB-CG-OD2	-5.34	113.50	118.30
1	A	227	ASP	CB-CG-OD1	-5.26	113.56	118.30
1	A	122	ASP	CB-CG-OD2	-5.25	113.58	118.30
2	B	8	GLN	N-CA-CB	5.14	119.85	110.60
1	A	129	ASP	CB-CG-OD1	5.11	122.90	118.30
1	A	97	GLN	N-CA-CB	5.10	119.78	110.60
1	A	29	ASP	N-CA-C	-5.08	97.29	111.00
1	A	256	ASN	CB-CA-C	5.03	120.46	110.40
1	A	21	ARG	NE-CZ-NH1	5.03	122.81	120.30
1	A	62	ARG	NE-CZ-NH1	5.01	122.80	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	A	2224	0	2100	122	1
2	B	821	0	796	46	0
3	C	69	0	61	8	0
4	A	22	0	0	6	1
4	B	17	0	0	0	0
4	C	2	0	0	1	0
All	All	3155	0	2957	165	1

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 27.

All (165) 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:A:215:LEU:HD12	1:A:261:VAL:HG22	1.55	0.88
2:B:54:MET:HE2	2:B:64:ILE:HD11	1.58	0.85
2:B:54:MET:CE	2:B:64:ILE:HD11	2.09	0.82
1:A:249:VAL:HG22	1:A:257:TYR:CZ	2.16	0.79
1:A:99:SER:HB3	1:A:114:LEU:HD22	1.64	0.79
1:A:199:VAL:HG21	1:A:254:GLU:OE2	1.87	0.74
1:A:35:ARG:HH22	2:B:53:ASP:HA	1.53	0.73
1:A:229:GLU:HB2	1:A:246:SER:OG	1.90	0.71
4:A:273:HOH:O	2:B:1:ILE:HG23	1.89	0.71
2:B:9:VAL:CG2	2:B:93:VAL:HG23	2.21	0.71
3:C:3:ASN:ND2	3:C:4:GLU:H	1.89	0.70
1:A:35:ARG:HH22	2:B:53:ASP:CA	2.05	0.69
1:A:185:PRO:HD2	1:A:266:LEU:HD13	1.73	0.69
1:A:218:GLN:HG2	1:A:223:GLU:OE1	1.91	0.68
1:A:228:MET:HE3	1:A:230:LEU:HD11	1.75	0.68
2:B:17:ASN:OD1	2:B:73:THR:HA	1.94	0.68
1:A:228:MET:CE	1:A:230:LEU:HD11	2.24	0.68
2:B:2:GLN:HG2	2:B:32:PRO:HD3	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:ASP:O	1:A:31:LYS:N	2.27	0.66
1:A:103:LEU:CD2	1:A:168:LEU:HD23	2.25	0.66
1:A:219:LEU:HB2	1:A:224:LEU:HD23	1.79	0.65
1:A:103:LEU:HD21	1:A:168:LEU:HD23	1.77	0.64
2:B:29:GLN:HA	2:B:29:GLN:HE21	1.62	0.64
2:B:6:GLN:HB2	2:B:28:THR:HG22	1.79	0.64
1:A:144:ARG:O	1:A:148:GLU:HG3	1.97	0.64
1:A:5:MET:HB2	1:A:168:LEU:HD13	1.79	0.63
1:A:202:ARG:HD3	1:A:246:SER:HB3	1.79	0.63
1:A:219:LEU:HB2	1:A:224:LEU:CD2	2.30	0.61
1:A:78:LEU:HD13	1:A:95:LEU:HB2	1.81	0.61
1:A:42:ASN:HD21	1:A:44:ARG:NH2	1.99	0.60
1:A:190:THR:HG22	1:A:191:HIS:N	2.16	0.60
1:A:41:GLU:H	1:A:41:GLU:CD	2.03	0.60
1:A:190:THR:HG22	1:A:191:HIS:H	1.66	0.60
2:B:9:VAL:HG23	2:B:93:VAL:HG23	1.84	0.60
1:A:9:GLU:HG3	4:A:279:HOH:O	2.01	0.60
1:A:200:THR:HG21	4:A:289:HOH:O	2.01	0.59
1:A:78:LEU:CD1	1:A:95:LEU:HB2	2.32	0.59
2:B:38:GLN:HB2	2:B:81:ARG:HB3	1.83	0.59
2:B:9:VAL:HG21	2:B:93:VAL:HG23	1.82	0.59
1:A:42:ASN:HD21	1:A:44:ARG:HH21	1.51	0.59
1:A:216:THR:O	1:A:259:CYS:HA	2.03	0.58
1:A:79:ARG:HB3	1:A:79:ARG:NH1	2.19	0.57
1:A:35:ARG:CB	1:A:48:ARG:HD2	2.34	0.57
1:A:220:ASN:HA	1:A:256:ASN:HD21	1.67	0.57
1:A:190:THR:O	1:A:191:HIS:HB2	2.04	0.57
3:C:3:ASN:HB3	4:C:300:HOH:O	2.03	0.57
1:A:35:ARG:HH22	2:B:53:ASP:C	2.08	0.56
1:A:18:GLU:O	1:A:19:GLU:HB2	2.06	0.55
1:A:249:VAL:HG22	1:A:257:TYR:CE1	2.42	0.55
2:B:41:LYS:HA	2:B:77:THR:O	2.07	0.55
1:A:201:LEU:HD21	1:A:254:GLU:HG2	1.88	0.55
1:A:12:VAL:HG12	1:A:21:ARG:HB3	1.88	0.55
2:B:9:VAL:HG23	2:B:93:VAL:CG2	2.36	0.55
1:A:62:ARG:O	1:A:65:GLN:HG2	2.06	0.55
1:A:65:GLN:HG2	1:A:66:LYS:N	2.22	0.54
1:A:94:THR:HG22	4:A:281:HOH:O	2.07	0.54
1:A:130:LEU:HD12	1:A:130:LEU:N	2.21	0.54
1:A:121:ARG:NH2	2:B:1:ILE:HG21	2.23	0.54
3:C:6:MET:CE	3:C:6:MET:HA	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:ILE:HG13	1:A:262:TYR:O	2.07	0.54
1:A:208:PHE:HE1	1:A:211:ALA:HA	1.72	0.54
2:B:63:TYR:O	2:B:64:ILE:HG13	2.08	0.53
2:B:29:GLN:NE2	2:B:29:GLN:N	2.57	0.53
2:B:15:PRO:HG2	2:B:97:ARG:HB2	1.91	0.53
1:A:190:THR:HG22	1:A:192:HIS:ND1	2.23	0.53
3:C:6:MET:HE2	3:C:6:MET:HA	1.90	0.53
1:A:35:ARG:HB3	1:A:48:ARG:HD2	1.91	0.52
2:B:29:GLN:NE2	2:B:29:GLN:CA	2.73	0.52
1:A:183:ASP:O	1:A:208:PHE:HA	2.09	0.52
2:B:28:THR:HG23	2:B:29:GLN:NE2	2.24	0.52
1:A:202:ARG:HH11	1:A:246:SER:HB3	1.74	0.52
1:A:220:ASN:OD1	1:A:221:GLY:N	2.42	0.52
1:A:70:GLN:O	1:A:74:PHE:HD1	1.93	0.52
1:A:97:GLN:NE2	3:C:5:ASN:OD1	2.43	0.52
1:A:146:LYS:O	1:A:149:GLN:HB3	2.09	0.52
1:A:32:GLU:OE2	1:A:48:ARG:HD3	2.11	0.51
1:A:126:LEU:HD12	1:A:127:ASN:H	1.76	0.51
1:A:33:PHE:CD2	1:A:34:VAL:HG13	2.46	0.51
1:A:234:ARG:NH1	2:B:10:TYR:CB	2.74	0.51
1:A:193:PRO:HG2	1:A:194:ARG:HH21	1.76	0.51
2:B:46:ILE:HG21	2:B:68:THR:HG21	1.93	0.51
1:A:25:VAL:HG12	1:A:27:TYR:CE1	2.46	0.50
1:A:191:HIS:O	1:A:200:THR:O	2.29	0.50
1:A:74:PHE:O	1:A:78:LEU:HB2	2.11	0.50
1:A:219:LEU:O	1:A:220:ASN:OD1	2.29	0.50
1:A:190:THR:HB	1:A:202:ARG:HB3	1.94	0.50
2:B:82:VAL:HG12	2:B:83:LYS:N	2.26	0.50
2:B:12:ARG:NH2	2:B:22:ILE:HD13	2.26	0.49
1:A:35:ARG:HB2	1:A:48:ARG:HD2	1.94	0.49
1:A:102:ASP:O	1:A:109:LEU:HD12	2.12	0.49
1:A:202:ARG:HD2	1:A:244:TRP:CE3	2.47	0.49
1:A:191:HIS:NE2	1:A:199:VAL:HG22	2.28	0.49
2:B:29:GLN:CA	2:B:29:GLN:HE21	2.24	0.49
1:A:208:PHE:CE1	1:A:241:PHE:HB2	2.48	0.49
1:A:75:ARG:HG2	4:A:284:HOH:O	2.13	0.49
3:C:3:ASN:ND2	3:C:4:GLU:N	2.58	0.48
2:B:41:LYS:HB2	2:B:78:TYR:CD1	2.48	0.48
1:A:79:ARG:HH11	1:A:79:ARG:HB3	1.77	0.48
1:A:130:LEU:CD1	1:A:130:LEU:N	2.76	0.48
1:A:198:GLU:HG2	1:A:199:VAL:N	2.27	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:TYR:HB3	1:A:139:ALA:HB1	1.94	0.48
1:A:229:GLU:O	1:A:245:ALA:HA	2.13	0.48
1:A:253:LYS:NZ	1:A:253:LYS:HB3	2.29	0.48
1:A:191:HIS:NE2	1:A:199:VAL:CG2	2.77	0.47
1:A:35:ARG:NH2	2:B:53:ASP:HA	2.25	0.47
1:A:62:ARG:HG3	1:A:65:GLN:OE1	2.14	0.47
1:A:236:ALA:HB1	2:B:12:ARG:HG3	1.97	0.46
1:A:231:VAL:HG12	1:A:244:TRP:O	2.14	0.46
1:A:263:HIS:HD2	1:A:265:GLY:H	1.62	0.46
2:B:35:ILE:HD11	2:B:82:VAL:CG1	2.46	0.46
1:A:190:THR:OG1	1:A:204:TRP:NE1	2.49	0.46
2:B:5:PRO:HA	2:B:28:THR:O	2.15	0.46
1:A:6:ARG:NH2	1:A:102:ASP:OD1	2.48	0.46
1:A:230:LEU:HD22	1:A:243:LYS:NZ	2.31	0.45
1:A:198:GLU:CG	1:A:199:VAL:N	2.79	0.45
1:A:14:ARG:HG2	1:A:15:PRO:HD2	1.98	0.45
2:B:12:ARG:HB2	2:B:22:ILE:HG22	1.98	0.45
1:A:9:GLU:OE1	1:A:22:TYR:OH	2.31	0.45
1:A:193:PRO:HB2	1:A:194:ARG:H	1.27	0.45
2:B:84:HIS:HD2	2:B:86:SER:OG	1.98	0.45
1:A:28:VAL:HG21	1:A:51:TRP:CZ2	2.51	0.45
1:A:208:PHE:HB2	1:A:263:HIS:CE1	2.52	0.45
1:A:189:VAL:HA	1:A:202:ARG:O	2.17	0.45
1:A:155:HIS:NE2	3:C:4:GLU:O	2.48	0.45
1:A:30:ASN:ND2	1:A:30:ASN:H	2.15	0.45
1:A:249:VAL:HG13	1:A:257:TYR:CE2	2.51	0.45
2:B:29:GLN:NE2	2:B:29:GLN:H	2.13	0.45
1:A:242:GLN:O	1:A:243:LYS:HB2	2.16	0.44
1:A:253:LYS:NZ	1:A:253:LYS:O	2.50	0.44
1:A:249:VAL:CG1	1:A:257:TYR:CE2	3.00	0.44
1:A:238:ASP:OD1	1:A:240:THR:OG1	2.34	0.44
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.51	0.44
1:A:119:GLU:HA	1:A:119:GLU:OE1	2.17	0.44
1:A:217:TRP:HA	1:A:259:CYS:HA	1.99	0.44
1:A:1:GLY:O	1:A:3:HIS:NE2	2.50	0.44
1:A:163:GLU:O	1:A:167:TRP:HB2	2.17	0.44
2:B:12:ARG:CB	2:B:22:ILE:HG22	2.48	0.44
3:C:3:ASN:HD22	3:C:4:GLU:H	1.62	0.43
1:A:30:ASN:HD22	1:A:30:ASN:H	1.66	0.43
1:A:36:PHE:HB2	1:A:45:TYR:CD1	2.54	0.43
1:A:26:GLY:O	1:A:32:GLU:HA	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:148:GLU:HA	4:A:288:HOH:O	2.18	0.43
1:A:6:ARG:HB2	1:A:27:TYR:HB2	2.00	0.43
1:A:207:GLY:O	1:A:240:THR:HG21	2.18	0.43
1:A:126:LEU:HD12	1:A:127:ASN:N	2.34	0.43
1:A:214:THR:O	1:A:262:TYR:HD1	2.01	0.43
1:A:208:PHE:CE1	1:A:211:ALA:HA	2.52	0.43
1:A:228:MET:HE2	1:A:230:LEU:HD11	2.00	0.43
1:A:199:VAL:HG12	1:A:249:VAL:HB	2.01	0.42
2:B:14:PRO:HA	2:B:15:PRO:HD3	1.94	0.42
2:B:7:ILE:HD13	2:B:82:VAL:CG2	2.49	0.42
1:A:52:MET:HE2	1:A:171:TYR:CE1	2.54	0.42
1:A:35:ARG:HH12	2:B:53:ASP:HA	1.85	0.42
2:B:2:GLN:HA	2:B:31:HIS:O	2.19	0.42
1:A:229:GLU:HB3	1:A:244:TRP:HZ3	1.85	0.42
2:B:54:MET:HE1	2:B:64:ILE:HD11	1.99	0.41
2:B:41:LYS:HB2	2:B:78:TYR:CE1	2.55	0.41
2:B:39:MET:HB3	2:B:46:ILE:HD12	2.01	0.41
1:A:259:CYS:SG	1:A:272:LEU:HB2	2.60	0.41
2:B:23:LEU:O	2:B:67:HIS:HA	2.20	0.41
1:A:103:LEU:HD23	1:A:168:LEU:HD23	2.00	0.41
1:A:172:LEU:HA	1:A:172:LEU:HD12	1.84	0.41
2:B:75:THR:O	2:B:75:THR:HG22	2.21	0.41
1:A:5:MET:HE3	1:A:164:CYS:SG	2.61	0.40
1:A:25:VAL:HG12	1:A:27:TYR:HE1	1.86	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	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:ASN:O	4:A:293:HOH:O[2_575]	2.13	0.07

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	270/272 (99%)	221 (82%)	33 (12%)	16 (6%)	2 0
2	B	97/99 (98%)	91 (94%)	5 (5%)	1 (1%)	19 28
3	C	7/9 (78%)	6 (86%)	1 (14%)	0	100 100
All	All	374/380 (98%)	318 (85%)	39 (10%)	17 (4%)	3 2

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	30	ASN
1	A	178	THR
1	A	191	HIS
1	A	195	SER
2	B	53	ASP
1	A	28	VAL
1	A	177	ALA
1	A	193	PRO
1	A	19	GLU
1	A	200	THR
1	A	56	GLY
1	A	221	GLY
1	A	255	GLN
1	A	50	PRO
1	A	199	VAL
1	A	57	PRO
1	A	192	HIS

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	A	230/230 (100%)	211 (92%)	19 (8%)	14 21
2	B	94/94 (100%)	87 (93%)	7 (7%)	17 26
3	C	8/8 (100%)	8 (100%)	0	100 100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	332/332 (100%)	306 (92%)	26 (8%)	16 24

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	VAL
1	A	17	LEU
1	A	18	GLU
1	A	41	GLU
1	A	79	ARG
1	A	86	ASN
1	A	95	LEU
1	A	107	TRP
1	A	114	LEU
1	A	170	ARG
1	A	172	LEU
1	A	186	LYS
1	A	224	LEU
1	A	225	THR
1	A	232	GLU
1	A	253	LYS
1	A	255	GLN
1	A	256	ASN
1	A	268	GLU
2	B	4	THR
2	B	23	LEU
2	B	28	THR
2	B	29	GLN
2	B	41	LYS
2	B	65	LEU
2	B	70	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	ASN
1	A	54	GLN
1	A	96	GLN
1	A	97	GLN
1	A	127	ASN
1	A	226	GLN

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Mol	Chain	Res	Type
1	A	255	GLN
1	A	256	ASN
1	A	263	HIS
2	B	2	GLN
2	B	29	GLN
2	B	38	GLN
2	B	67	HIS
2	B	84	HIS
3	C	3	ASN
3	C	5	ASN

5.3.3 RNA

There are no RNA molecules 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

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.