Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:43 PM GMT

PDB ID	:	1YDP
Title	:	1.9A crystal structure of HLA-G
Authors	:	Clements, C.S.; Kjer-nielsen, L.; Kostenko, L.; Hoare, H.L.; Dunstone, M.A.; Moses, E.; Freed, K.; Brooks, A.G.; Rossjohn, J.; Mccluskey, J.
Deposited on	:	2004-12-25
Resolution	:	1.90 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:

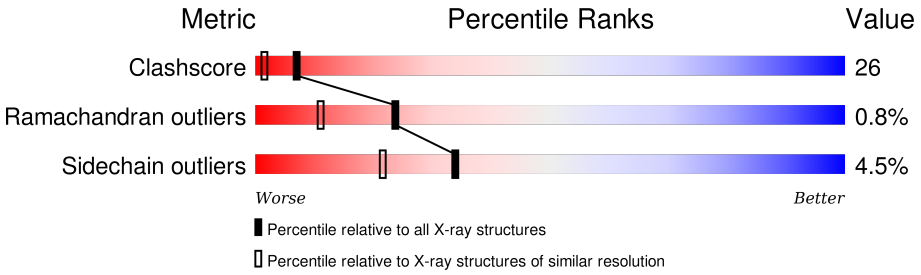
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
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 1.90 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	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)

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	275	67% 28%
2	B	100	61% 31% 5%
3	P	9	78% 11% 11%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3590 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 MHC class I antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	S	0	0	0
			2242	1397	404	429	12			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	42	SER	CYS	ENGINEERED	UNP P17693

- 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			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1A	MET	-	INITIATING METHIONINE	UNP P61769

- Molecule 3 is a protein called histone 2a peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	9	Total	C	N	O	0	0	0
			81	52	18	11			

- Molecule 4 is COBALT (II) ION (three-letter code: CO) (formula: Co).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Co	0	0
			1	1		

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	2	Total Cl 2 2	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	311	Total O 311 311	0	0
6	B	105	Total O 105 105	0	0
6	P	11	Total O 11 11	0	0

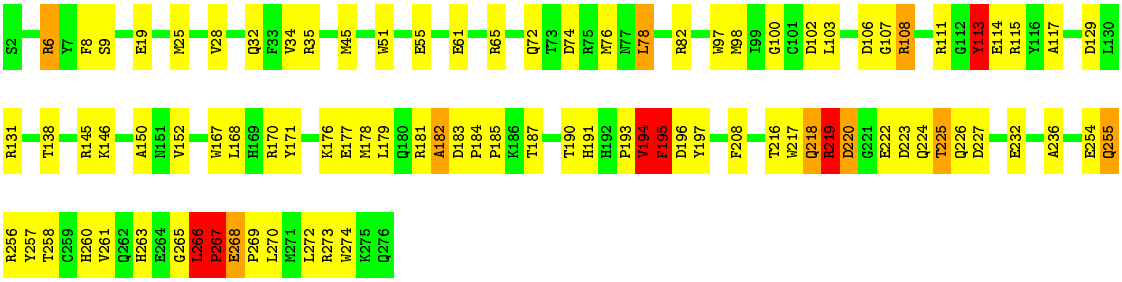
3 Residue-property plots (i)

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.

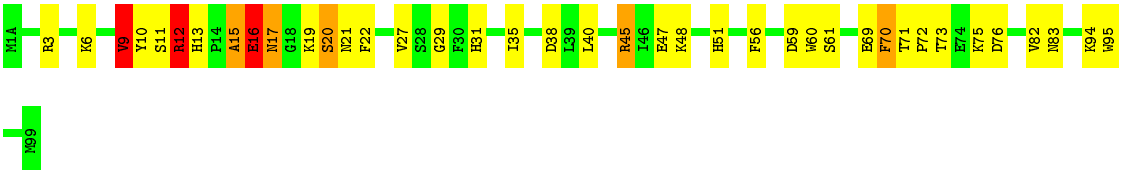
- Molecule 1: MHC class I antigen

Chain A:



- Molecule 2: Beta-2-microglobulin

Chain B:



- Molecule 3: histone 2a peptide

Chain P:

4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α, β, γ	77.15Å 77.15Å 151.72Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 1.90	Depositor
% Data completeness (in resolution range)	(Not available) (50.00-1.90)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.235 , 0.264	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3590	wwPDB-VP
Average B, all atoms (Å^2)	48.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CO, 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	0.85	12/2303 (0.5%)	1.30	27/3124 (0.9%)
2	B	1.21	7/860 (0.8%)	1.08	8/1162 (0.7%)
3	P	0.38	0/82	0.82	0/108
All	All	0.95	19/3245 (0.6%)	1.23	35/4394 (0.8%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4
2	B	0	4
All	All	0	8

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	12	ARG	CD-NE	-16.82	1.17	1.46
1	A	268	GLU	CB-CG	-16.16	1.21	1.52
2	B	9	VAL	C-N	-16.00	0.97	1.34
2	B	15	ALA	C-N	13.59	1.65	1.34
1	A	182	ALA	C-N	12.37	1.62	1.34
2	B	16	GLU	C-N	-11.29	1.08	1.34
2	B	82	VAL	C-N	11.15	1.59	1.34
1	A	267	PRO	N-CD	11.14	1.63	1.47
1	A	266	LEU	CB-CG	10.53	1.83	1.52
1	A	217	TRP	C-N	9.71	1.56	1.34
2	B	83	ASN	C-N	-9.59	1.11	1.34
1	A	190	THR	C-N	-8.85	1.13	1.34
1	A	176	LYS	C-N	-8.68	1.14	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	267	PRO	CB-CG	8.19	1.90	1.50
2	B	17	ASN	CB-CG	-8.02	1.32	1.51
1	A	268	GLU	C-N	-7.02	1.21	1.34
1	A	219	ARG	CA-CB	-6.87	1.38	1.53
1	A	219	ARG	CD-NE	-5.78	1.36	1.46
1	A	220	ASP	C-N	-5.30	1.23	1.33

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	219	ARG	CD-NE-CZ	28.68	163.75	123.60
1	A	220	ASP	CB-CG-OD2	21.32	137.49	118.30
1	A	220	ASP	CB-CG-OD1	-20.91	99.48	118.30
2	B	9	VAL	O-C-N	-19.23	91.93	122.70
1	A	195	PHE	CB-CG-CD1	-15.11	110.22	120.80
1	A	195	PHE	CB-CG-CD2	14.37	130.86	120.80
2	B	9	VAL	CA-C-N	13.58	147.07	117.20
1	A	266	LEU	CB-CG-CD1	10.88	129.50	111.00
1	A	268	GLU	CA-CB-CG	10.51	136.52	113.40
1	A	176	LYS	O-C-N	-10.28	106.26	122.70
1	A	220	ASP	CA-CB-CG	9.12	133.47	113.40
1	A	194	VAL	C-N-CA	-8.86	99.55	121.70
1	A	190	THR	O-C-N	-8.49	109.12	122.70
1	A	219	ARG	CB-CA-C	-7.83	94.75	110.40
2	B	83	ASN	O-C-N	-7.55	110.62	122.70
2	B	15	ALA	C-N-CA	-7.49	102.99	121.70
1	A	176	LYS	CA-C-N	7.45	133.60	117.20
1	A	113	TYR	CB-CA-C	-7.41	95.57	110.40
1	A	195	PHE	CA-CB-CG	6.87	130.38	113.90
1	A	113	TYR	CB-CG-CD2	-6.86	116.89	121.00
2	B	82	VAL	C-N-CA	-6.85	104.58	121.70
1	A	182	ALA	C-N-CA	-6.83	104.62	121.70
1	A	78	LEU	C-N-CA	-6.83	104.63	121.70
1	A	6	ARG	CB-CA-C	-6.54	97.31	110.40
1	A	267	PRO	N-CA-CB	6.35	110.92	103.30
1	A	190	THR	CA-C-N	6.29	131.04	117.20
2	B	83	ASN	CA-C-N	6.23	130.90	117.20
2	B	16	GLU	O-C-N	-5.86	113.32	122.70
2	B	9	VAL	C-N-CA	5.81	136.22	121.70
1	A	138	THR	O-C-N	-5.74	113.52	122.70
1	A	267	PRO	CA-CB-CG	-5.59	93.39	104.00
1	A	218	GLN	CB-CG-CD	-5.52	97.24	111.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	113	TYR	CB-CG-CD1	5.42	124.25	121.00
1	A	267	PRO	CA-N-CD	-5.37	103.98	111.50
1	A	220	ASP	O-C-N	5.09	131.85	123.20

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	183	ASP	Mainchain
1	A	194	VAL	Peptide
1	A	219	ARG	Sidechain
1	A	268	GLU	Mainchain
2	B	12	ARG	Sidechain
2	B	15	ALA	Mainchain
2	B	16	GLU	Mainchain
2	B	9	VAL	Mainchain

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	2242	0	2094	121	1
2	B	837	0	800	45	0
3	P	81	0	94	2	0
4	B	1	0	0	0	0
5	B	2	0	0	0	0
6	A	311	0	0	28	1
6	B	105	0	0	5	0
6	P	11	0	0	1	0
All	All	3590	0	2988	163	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 26.

All (163) 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:266:LEU:CG	1:A:266:LEU:CB	1.83	1.53
1:A:266:LEU:CD1	1:A:270:LEU:HD22	1.43	1.47
1:A:267:PRO:CG	1:A:267:PRO:CB	1.90	1.39
1:A:266:LEU:CB	1:A:266:LEU:HG	1.50	1.30
1:A:108:ARG:CD	6:A:385:HOH:O	1.93	1.17
1:A:108:ARG:HD2	6:A:385:HOH:O	1.43	1.16
1:A:6:ARG:NH1	1:A:113:TYR:HE1	1.42	1.15
1:A:6:ARG:NH1	1:A:113:TYR:CE1	2.17	1.12
1:A:266:LEU:HD12	1:A:270:LEU:HD22	1.18	1.11
1:A:266:LEU:HD12	1:A:270:LEU:CD2	1.86	1.05
1:A:266:LEU:CD1	1:A:270:LEU:CD2	2.35	1.04
1:A:102:ASP:OD2	1:A:111:ARG:NE	1.98	0.96
2:B:16:GLU:OE2	2:B:19:LYS:HB2	1.68	0.94
1:A:266:LEU:HD11	1:A:270:LEU:HD22	1.46	0.92
1:A:222:GLU:HB2	6:A:522:HOH:O	1.71	0.89
1:A:194:VAL:HG12	1:A:197:TYR:CE2	2.10	0.87
1:A:218:GLN:HB2	1:A:258:THR:OG1	1.77	0.85
1:A:34:VAL:CG2	1:A:45:MET:HG3	2.07	0.84
1:A:28:VAL:HG11	1:A:179:LEU:HD13	1.60	0.84
1:A:111:ARG:NE	6:A:462:HOH:O	2.11	0.84
1:A:103:LEU:HD13	1:A:168:LEU:HD23	1.59	0.83
2:B:16:GLU:OE2	2:B:19:LYS:CB	2.29	0.80
2:B:16:GLU:O	2:B:17:ASN:C	2.17	0.78
2:B:19:LYS:HG3	6:B:1093:HOH:O	1.84	0.78
1:A:263:HIS:O	1:A:266:LEU:HB2	1.85	0.77
1:A:111:ARG:CZ	6:A:462:HOH:O	2.32	0.76
1:A:270:LEU:H	1:A:270:LEU:HD23	1.53	0.73
1:A:187:THR:HB	1:A:272:LEU:HD21	1.71	0.73
1:A:218:GLN:O	1:A:257:TYR:HA	1.91	0.71
1:A:34:VAL:HG22	1:A:45:MET:HG3	1.73	0.71
2:B:16:GLU:CG	2:B:19:LYS:HB2	2.20	0.71
1:A:194:VAL:O	1:A:195:PHE:C	2.26	0.71
1:A:107:GLY:HA3	6:A:579:HOH:O	1.91	0.71
2:B:16:GLU:OE2	2:B:19:LYS:CG	2.39	0.70
1:A:76:MET:HE3	1:A:76:MET:O	1.91	0.70
1:A:266:LEU:CG	1:A:266:LEU:CA	2.68	0.69
2:B:11:SER:HB2	2:B:21:ASN:ND2	2.08	0.69
2:B:3:ARG:HH21	2:B:61:SER:HB3	1.58	0.68
1:A:194:VAL:O	1:A:195:PHE:O	2.12	0.67
1:A:266:LEU:HD13	1:A:270:LEU:HD22	1.66	0.67
2:B:19:LYS:O	2:B:72:PRO:HD2	1.94	0.67
1:A:98:MET:HG3	6:B:1096:HOH:O	1.94	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:22:PHE:CE1	2:B:69:GLU:HG2	2.30	0.66
1:A:261:VAL:HB	1:A:270:LEU:HD21	1.77	0.66
1:A:218:GLN:HE21	1:A:260:HIS:HB2	1.61	0.66
1:A:194:VAL:CG1	1:A:197:TYR:CE2	2.80	0.65
2:B:21:ASN:ND2	2:B:22:PHE:H	1.95	0.65
2:B:73:THR:HG22	2:B:75:LYS:H	1.61	0.65
2:B:40:LEU:HD23	2:B:45:ARG:HA	1.78	0.64
1:A:194:VAL:HG11	1:A:197:TYR:CZ	2.34	0.62
1:A:178:MET:HG2	6:A:458:HOH:O	2.00	0.62
1:A:266:LEU:HB2	1:A:266:LEU:HG	1.73	0.61
1:A:102:ASP:HB2	1:A:111:ARG:HG2	1.82	0.61
1:A:236:ALA:O	2:B:12:ARG:HD3	2.00	0.61
2:B:20:SER:HA	2:B:70:PHE:O	2.01	0.61
2:B:13:HIS:HB2	2:B:21:ASN:OD1	2.01	0.61
1:A:72:GLN:HB3	6:A:545:HOH:O	2.01	0.61
2:B:16:GLU:CD	2:B:19:LYS:HB2	2.22	0.60
1:A:6:ARG:HH11	1:A:113:TYR:HE1	1.42	0.60
1:A:187:THR:CB	1:A:272:LEU:HD21	2.32	0.59
1:A:185:PRO:HD2	1:A:266:LEU:HD23	1.83	0.59
1:A:194:VAL:HG11	1:A:197:TYR:OH	2.03	0.59
2:B:13:HIS:H	2:B:21:ASN:HD21	1.49	0.59
1:A:6:ARG:NH1	1:A:113:TYR:CZ	2.70	0.58
1:A:19:GLU:HB2	6:A:557:HOH:O	2.03	0.58
1:A:108:ARG:CG	6:A:385:HOH:O	2.43	0.57
1:A:181:ARG:HB3	6:A:563:HOH:O	2.04	0.57
2:B:27:VAL:HG11	2:B:35:ILE:CD1	2.36	0.56
1:A:191:HIS:HB2	1:A:274:TRP:NE1	2.19	0.56
1:A:261:VAL:HB	1:A:270:LEU:CD2	2.36	0.56
2:B:16:GLU:HG3	2:B:16:GLU:O	2.06	0.56
1:A:34:VAL:HG21	1:A:45:MET:HG3	1.86	0.56
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.40	0.55
1:A:108:ARG:CZ	1:A:108:ARG:HB2	2.27	0.55
2:B:6:LYS:HE2	2:B:29:GLY:HA3	1.87	0.55
1:A:65:ARG:NH1	6:A:460:HOH:O	2.40	0.55
1:A:273:ARG:HH11	1:A:273:ARG:HB2	1.71	0.54
2:B:21:ASN:HB3	2:B:70:PHE:CE1	2.43	0.54
1:A:6:ARG:CZ	1:A:113:TYR:CE1	2.89	0.54
1:A:260:HIS:CE1	1:A:269:PRO:HB2	2.43	0.54
1:A:218:GLN:O	1:A:258:THR:N	2.36	0.53
1:A:167:TRP:CE2	3:P:1:ARG:HG3	2.43	0.53
1:A:61:GLU:HG2	6:A:449:HOH:O	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:273:ARG:NH1	1:A:273:ARG:HB2	2.24	0.53
1:A:194:VAL:CG1	1:A:197:TYR:CZ	2.92	0.53
1:A:224:GLN:HA	6:A:468:HOH:O	2.09	0.53
1:A:115:ARG:HD3	6:A:560:HOH:O	2.09	0.53
1:A:216:THR:OG1	1:A:260:HIS:HB3	2.09	0.53
1:A:265:GLY:HA2	6:A:569:HOH:O	2.09	0.52
1:A:191:HIS:CD2	1:A:193:PRO:HG3	2.45	0.51
2:B:3:ARG:NH2	2:B:61:SER:HB3	2.25	0.51
2:B:73:THR:HB	2:B:76:ASP:HB2	1.92	0.51
1:A:184:PRO:HB2	1:A:266:LEU:CD2	2.41	0.50
2:B:20:SER:OG	2:B:71:THR:HG22	2.12	0.50
2:B:11:SER:HB2	2:B:21:ASN:HD22	1.74	0.50
1:A:196:ASP:OD1	6:A:454:HOH:O	2.20	0.50
2:B:31:HIS:CD2	6:B:1072:HOH:O	2.65	0.49
2:B:16:GLU:OE2	2:B:19:LYS:HG2	2.11	0.49
2:B:10:TYR:N	2:B:10:TYR:CD1	2.80	0.49
1:A:227:ASP:HB2	6:A:541:HOH:O	2.12	0.49
1:A:150:ALA:HB3	1:A:152:VAL:HG13	1.92	0.49
1:A:270:LEU:HA	6:A:529:HOH:O	2.12	0.49
1:A:266:LEU:HD11	1:A:270:LEU:CD2	2.25	0.49
1:A:6:ARG:HD3	1:A:100:GLY:HA3	1.94	0.49
1:A:102:ASP:CG	1:A:113:TYR:OH	2.52	0.48
1:A:219:ARG:HA	1:A:256:ARG:O	2.12	0.48
1:A:218:GLN:HA	1:A:222:GLU:O	2.14	0.48
1:A:194:VAL:HG12	1:A:195:PHE:N	2.28	0.47
1:A:9:SER:OG	1:A:97:TRP:HB3	2.15	0.47
1:A:145:ARG:HD3	6:A:523:HOH:O	2.15	0.47
1:A:28:VAL:HG11	1:A:179:LEU:CD1	2.39	0.46
1:A:184:PRO:CB	1:A:266:LEU:CD2	2.93	0.46
1:A:102:ASP:OD1	1:A:113:TYR:OH	2.34	0.46
1:A:182:ALA:HB1	1:A:265:GLY:HA2	1.96	0.46
1:A:218:GLN:NE2	1:A:260:HIS:HB2	2.30	0.46
2:B:47:GLU:O	2:B:48:LYS:HG2	2.15	0.46
2:B:16:GLU:O	2:B:16:GLU:CG	2.63	0.46
1:A:218:GLN:N	1:A:258:THR:O	2.42	0.46
1:A:272:LEU:N	1:A:272:LEU:HD12	2.31	0.46
1:A:270:LEU:H	1:A:270:LEU:CD2	2.24	0.46
1:A:51:TRP:CZ3	1:A:171:TYR:HB3	2.51	0.45
1:A:34:VAL:CG2	1:A:45:MET:CG	2.88	0.45
2:B:11:SER:HA	2:B:22:PHE:O	2.16	0.45
1:A:146:LYS:HD3	6:A:410:HOH:O	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:P:5:ARG:HG2	6:P:19:HOH:O	2.16	0.45
2:B:94:LYS:HE3	6:B:1081:HOH:O	2.16	0.45
2:B:20:SER:CB	2:B:71:THR:HG22	2.47	0.45
1:A:32:GLN:OE1	1:A:35:ARG:HD3	2.17	0.45
1:A:82:ARG:NH1	6:A:377:HOH:O	2.50	0.45
2:B:21:ASN:CG	2:B:22:PHE:H	2.21	0.45
2:B:59:ASP:O	2:B:60:TRP:HB2	2.17	0.44
1:A:224:GLN:O	1:A:226:GLN:N	2.51	0.44
1:A:34:VAL:HG22	1:A:45:MET:CG	2.46	0.43
1:A:185:PRO:HB3	1:A:208:PHE:HB3	1.99	0.43
2:B:38:ASP:OD1	2:B:45:ARG:NH1	2.51	0.43
1:A:191:HIS:CE1	1:A:254:GLU:HG2	2.54	0.43
1:A:182:ALA:HB3	6:A:569:HOH:O	2.17	0.43
1:A:8:PHE:HB3	2:B:56:PHE:CZ	2.53	0.43
1:A:255:GLN:HG3	1:A:273:ARG:HD3	1.99	0.43
1:A:55:GLU:OE1	1:A:170:ARG:NH2	2.52	0.43
1:A:170:ARG:NH2	6:A:546:HOH:O	2.52	0.42
1:A:232:GLU:HB2	6:A:530:HOH:O	2.19	0.42
1:A:218:GLN:NE2	6:A:339:HOH:O	2.52	0.42
1:A:273:ARG:HH11	1:A:273:ARG:CB	2.31	0.42
1:A:194:VAL:HG12	1:A:197:TYR:HE2	1.74	0.42
2:B:27:VAL:HG11	2:B:35:ILE:HD11	2.02	0.42
2:B:27:VAL:HG11	2:B:35:ILE:HD13	2.02	0.42
1:A:78:LEU:HA	1:A:78:LEU:HD23	1.89	0.42
1:A:196:ASP:CG	6:A:454:HOH:O	2.58	0.41
1:A:191:HIS:HB2	1:A:274:TRP:CE2	2.55	0.41
1:A:182:ALA:HB1	1:A:265:GLY:CA	2.51	0.41
1:A:184:PRO:CB	1:A:266:LEU:HD22	2.51	0.41
1:A:270:LEU:N	1:A:270:LEU:HD23	2.30	0.41
1:A:115:ARG:HG3	2:B:60:TRP:CH2	2.56	0.41
1:A:185:PRO:HB3	1:A:208:PHE:CD1	2.56	0.41
1:A:34:VAL:HG21	1:A:45:MET:CG	2.50	0.41
1:A:129:ASP:O	1:A:131:ARG:HG3	2.21	0.41
1:A:106:ASP:N	1:A:106:ASP:OD1	2.51	0.41
2:B:45:ARG:HG3	2:B:45:ARG:O	2.21	0.41
1:A:225:THR:HA	6:A:551:HOH:O	2.21	0.41
2:B:51:HIS:NE2	6:B:1004:HOH:O	2.37	0.40
1:A:254:GLU:H	1:A:254:GLU:CD	2.24	0.40
2:B:9:VAL:CG1	2:B:95:TRP:HB2	2.51	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:195:PHE:CZ	6:A:568:HOH:O[5_565]	2.10	0.10

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	273/275 (99%)	261 (96%)	9 (3%)	3 (1%)	17 6
2	B	98/100 (98%)	94 (96%)	4 (4%)	0	100 100
3	P	7/9 (78%)	7 (100%)	0	0	100 100
All	All	378/384 (98%)	362 (96%)	13 (3%)	3 (1%)	24 11

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	114	GLU
1	A	195	PHE
1	A	225	THR

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	232/232 (100%)	221 (95%)	11 (5%)	32 20
2	B	95/95 (100%)	92 (97%)	3 (3%)	46 35
3	P	9/9 (100%)	8 (89%)	1 (11%)	8 2

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	336/336 (100%)	321 (96%)	15 (4%)	34 21

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	MET
1	A	74	ASP
1	A	108	ARG
1	A	113	TYR
1	A	177	GLU
1	A	195	PHE
1	A	220	ASP
1	A	223	ASP
1	A	255	GLN
1	A	266	LEU
1	A	267	PRO
2	B	20	SER
2	B	45	ARG
2	B	70	PHE
3	P	1	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	87	GLN
1	A	127	ASN
1	A	180	GLN
1	A	191	HIS
1	A	218	GLN
1	A	255	GLN
1	A	260	HIS
2	B	21	ASN
2	B	83	ASN
2	B	89	GLN

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

Of 3 ligands modelled in this entry, 3 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

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.