



# Full wwPDB X-ray Structure Validation Report ⓘ

Dec 13, 2023 – 03:57 pm GMT

PDB ID : 4U1K  
Title : HLA class I micropolymorphisms determine peptide-HLA landscape and dictate differential HIV-1 escape through identical epitopes  
Authors : Rizkallah, P.J.; Cole, D.K.; Fuller, A.; Sewell, A.K.  
Deposited on : 2014-07-15  
Resolution : 2.09 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](mailto: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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467  
Mogul : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
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) : 2.36

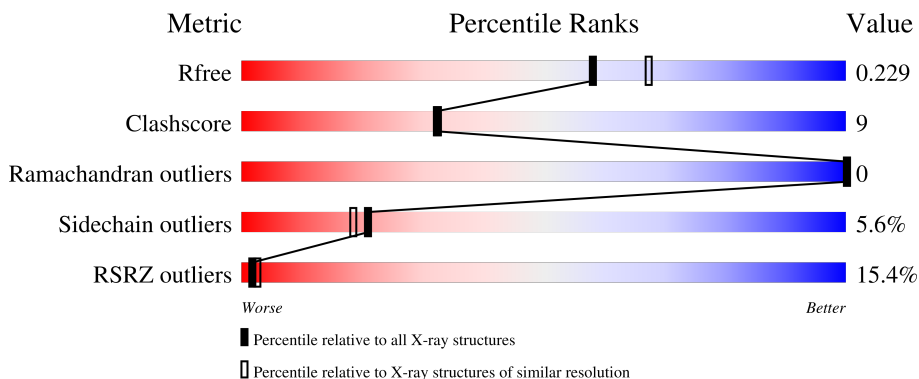
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.09 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	277	
1	D	277	
2	B	100	
2	E	100	
3	C	9	

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Mol	Chain	Length	Quality of chain
3	F	9	 89% 11%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	GOL	A	302	-	X	X	-
4	GOL	A	303	-	X	-	-
4	GOL	D	303	-	X	X	-
4	GOL	D	305	-	-	X	-

## 2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 6989 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 HLA class I histocompatibility antigen, B-7 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	277	Total	C	N	O	S	0	2	0
			2287	1418	417	445	7			
1	D	277	Total	C	N	O	S	0	4	0
			2302	1426	420	449	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP P01889
D	0	MET	-	initiating methionine	UNP P01889

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
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 Nef.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	9	Total	C	N	O	S	0	0	0
			76	48	16	11	1			
3	F	9	Total	C	N	O	S	0	0	0
			76	48	16	11	1			

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 6 3 3	0	0
4	A	1	Total C O 6 3 3	0	0
4	A	1	Total C O 6 3 3	0	0
4	D	1	Total C O 6 3 3	0	0
4	D	1	Total C O 6 3 3	0	0
4	D	1	Total C O 6 3 3	0	0
4	D	1	Total C O 6 3 3	0	0
4	D	1	Total C O 6 3 3	0	0

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $C_2H_6O_2$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	D	1	Total C O 4 2 2	0	0
5	D	1	Total C O 4 2 2	0	0
5	D	1	Total C O 4 2 2	0	0

- Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	O S	0	0
			5	4 1		

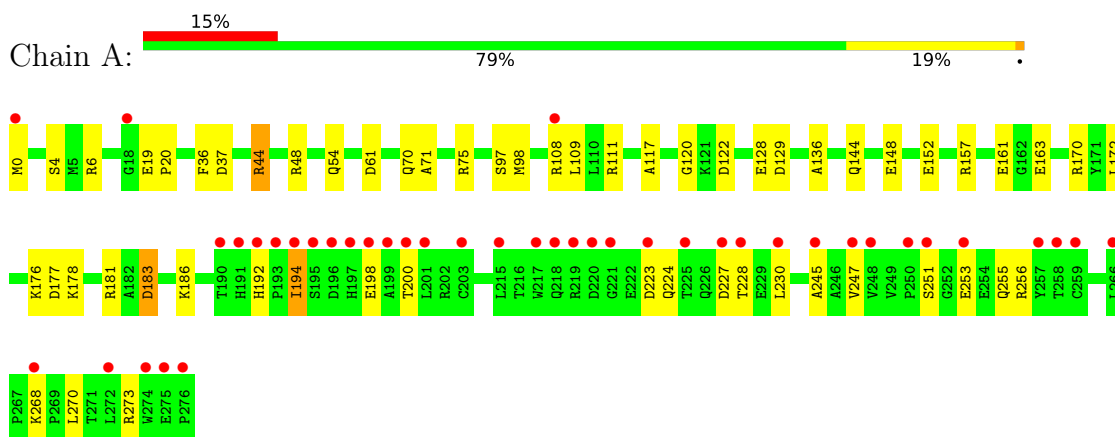
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	206	Total	O	0	0
			206	206		
7	B	31	Total	O	0	0
			31	31		
7	C	9	Total	O	0	0
			9	9		
7	D	205	Total	O	0	0
			205	205		
7	E	32	Total	O	0	0
			32	32		
7	F	14	Total	O	0	0
			14	14		

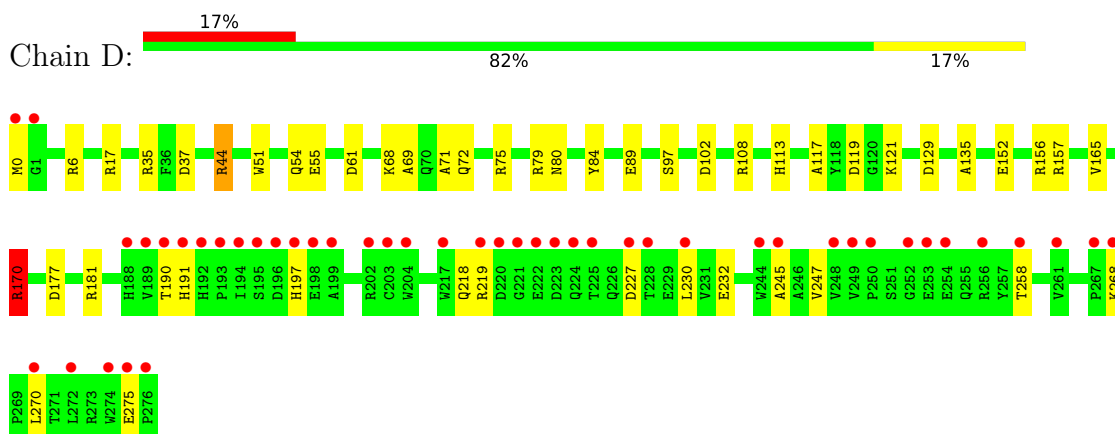
### 3 Residue-property plots [i](#)

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.

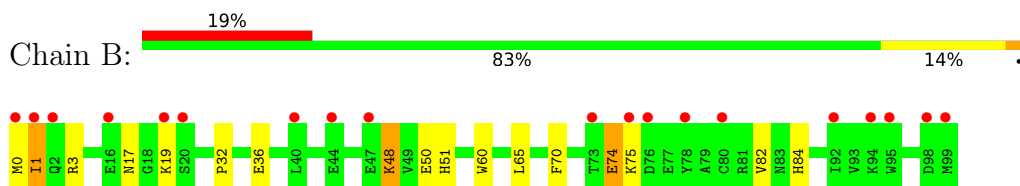
- Molecule 1: HLA class I histocompatibility antigen, B-7 alpha chain



- Molecule 1: HLA class I histocompatibility antigen, B-7 alpha chain

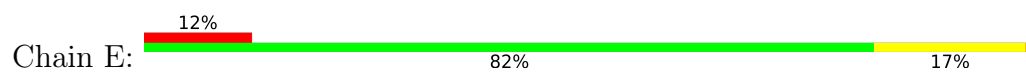


- Molecule 2: Beta-2-microglobulin

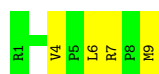


- Molecule 2: Beta-2-microglobulin

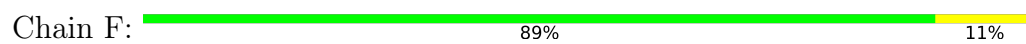




- Molecule 3: Protein Nef



- Molecule 3: Protein Nef



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	46.27Å 49.13Å 107.46Å 91.37° 93.57° 95.75°	Depositor
Resolution (Å)	48.86 – 2.09 48.86 – 2.09	Depositor EDS
% Data completeness (in resolution range)	97.5 (48.86-2.09) 97.5 (48.86-2.09)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.11 (at 2.08Å)	Xtrriage
Refinement program	REFMAC 5.8.0049	Depositor
R, $R_{free}$	0.176 , 0.226 0.183 , 0.229	Depositor DCC
$R_{free}$ test set	2757 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	28.0	Xtrriage
Anisotropy	0.508	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 50.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6989	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.33% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, SO4, EDO

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	0.94	0/2350	0.97	8/3192 (0.3%)
1	D	0.95	2/2365 (0.1%)	0.93	5/3212 (0.2%)
2	B	0.62	0/860	0.77	0/1162
2	E	0.68	0/860	0.79	1/1162 (0.1%)
3	C	1.21	0/78	1.25	1/104 (1.0%)
3	F	1.10	0/78	1.31	1/104 (1.0%)
All	All	0.88	2/6591 (0.0%)	0.92	16/8936 (0.2%)

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
2	B	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	97	SER	CB-OG	5.75	1.49	1.42
1	D	165	VAL	CB-CG1	-5.04	1.42	1.52

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	44	ARG	NE-CZ-NH2	7.39	123.99	120.30
1	D	170	ARG	NE-CZ-NH2	-6.89	116.85	120.30
1	A	122	ASP	CB-CG-OD1	6.88	124.50	118.30
1	D	157	ARG	NE-CZ-NH2	6.87	123.73	120.30
2	E	59	ASP	CB-CG-OD1	6.40	124.06	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	129	ASP	CB-CG-OD1	6.23	123.90	118.30
1	A	98	MET	CG-SD-CE	-5.77	90.97	100.20
3	F	7	ARG	NE-CZ-NH2	-5.75	117.43	120.30
3	C	9	MET	CA-CB-CG	-5.54	103.89	113.30
1	A	157	ARG	NE-CZ-NH2	5.51	123.05	120.30
1	D	37	ASP	CB-CG-OD1	5.24	123.02	118.30
1	A	48	ARG	NE-CZ-NH1	5.18	122.89	120.30
1	D	6	ARG	NE-CZ-NH1	5.09	122.84	120.30
1	A	98	MET	CA-CB-CG	5.04	121.88	113.30
1	A	170	ARG	NE-CZ-NH1	-5.03	117.78	120.30
1	D	129	ASP	CB-CG-OD1	5.03	122.82	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	0	MET	Peptide

## 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	2287	0	2122	46	0
1	D	2302	0	2133	37	0
2	B	837	0	803	10	0
2	E	837	0	803	8	0
3	C	76	0	86	11	0
3	F	76	0	86	0	0
4	A	18	0	23	16	0
4	D	30	0	39	18	0
5	A	12	0	18	3	0
5	D	12	0	18	1	0
6	A	5	0	0	0	0
7	A	206	0	0	22	0
7	B	31	0	0	1	0
7	C	9	0	0	0	0
7	D	205	0	0	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	E	32	0	0	0	0
7	F	14	0	0	0	0
All	All	6989	0	6131	111	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 9.

All (111) 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:D:71:ALA:HB1	4:D:305:GOL:H12	1.40	1.01
4:A:302:GOL:H12	3:C:4:VAL:HG13	1.41	0.99
4:A:302:GOL:C1	3:C:4:VAL:HG13	1.92	0.98
1:A:109:LEU:HD22	1:A:161[B]:GLU:HB3	1.62	0.81
4:A:302:GOL:H11	3:C:4:VAL:N	1.97	0.79
1:D:71:ALA:HB1	4:D:305:GOL:C1	2.15	0.75
1:A:70:GLN:HE22	4:A:301:GOL:H12	1.52	0.74
1:A:163:GLU:OE2	4:A:302:GOL:H2	1.88	0.74
1:A:6:ARG:HG2	7:A:406:HOH:O	1.87	0.74
1:D:75:ARG:CZ	4:D:305:GOL:H11	2.18	0.73
4:A:302:GOL:H11	3:C:4:VAL:H	1.53	0.73
1:A:71:ALA:O	1:A:75:ARG:HD3	1.90	0.71
1:D:170:ARG:NH1	7:D:401:HOH:O	2.23	0.68
1:A:20:PRO:HD2	1:A:75:ARG:HG2	1.76	0.67
1:D:72:GLN:CD	7:D:403:HOH:O	2.31	0.67
1:D:75:ARG:NE	4:D:305:GOL:H11	2.10	0.66
1:A:36:PHE:CD1	7:A:411:HOH:O	2.48	0.65
1:A:163:GLU:OE2	4:A:302:GOL:C2	2.45	0.64
4:A:302:GOL:H11	3:C:4:VAL:HG13	1.77	0.63
1:D:230:LEU:HD23	1:D:245:ALA:HB2	1.80	0.62
5:A:306:EDO:H11	7:A:456:HOH:O	2.00	0.62
1:D:75:ARG:NH2	4:D:305:GOL:H11	2.16	0.60
2:E:22:PHE:CE2	2:E:69:GLU:HG3	2.38	0.59
2:B:48:LYS:O	2:B:48:LYS:HG3	2.02	0.59
1:D:69:ALA:HA	4:D:304:GOL:H2	1.84	0.59
1:D:135:ALA:HB3	5:D:308:EDO:H21	1.85	0.58
1:D:80:ASN:HB3	4:D:303:GOL:H32	1.86	0.58
1:A:4:SER:HB2	7:A:406:HOH:O	2.03	0.58
7:A:566:HOH:O	2:B:1:ILE:HG21	2.04	0.57
4:A:302:GOL:H12	3:C:4:VAL:CG1	2.26	0.57
4:A:302:GOL:C1	3:C:4:VAL:H	2.18	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:PHE:CE1	7:A:411:HOH:O	2.58	0.56
1:A:172:LEU:O	1:A:176:LYS:HB3	2.06	0.56
1:A:136:ALA:HB2	5:A:305:EDO:H11	1.88	0.56
1:A:4:SER:CB	7:A:406:HOH:O	2.53	0.56
4:A:302:GOL:H32	3:C:4:VAL:HG13	1.89	0.55
1:A:36:PHE:CG	7:A:411:HOH:O	2.60	0.55
5:A:306:EDO:C1	7:A:456:HOH:O	2.54	0.55
2:B:48:LYS:O	2:B:48:LYS:CG	2.54	0.55
4:A:303:GOL:H2	7:D:408:HOH:O	2.05	0.55
1:A:6:ARG:NH1	7:A:406:HOH:O	2.38	0.54
1:D:69:ALA:CA	4:D:304:GOL:H2	2.36	0.54
1:A:37:ASP:N	7:A:411:HOH:O	2.40	0.54
1:D:80:ASN:HA	4:D:303:GOL:H31	1.90	0.54
1:D:89:GLU:CD	7:D:478:HOH:O	2.45	0.54
1:A:71:ALA:O	1:A:75:ARG:CD	2.56	0.54
1:A:97:SER:HB2	7:A:442:HOH:O	2.09	0.53
1:A:54:GLN:NE2	7:A:413:HOH:O	2.41	0.53
4:A:302:GOL:H32	3:C:4:VAL:CG1	2.38	0.53
1:D:68:LYS:NZ	7:D:402:HOH:O	2.25	0.53
1:A:61[B]:ASP:OD1	4:A:303:GOL:H32	2.08	0.53
1:A:253:GLU:OE2	1:A:256:ARG:NH1	2.42	0.53
2:E:32:PRO:O	2:E:84:HIS:HE1	1.92	0.53
1:A:75:ARG:NH2	7:A:412:HOH:O	2.41	0.52
1:D:170:ARG:HH11	1:D:170:ARG:HG3	1.76	0.51
1:A:44:ARG:NH2	7:A:414:HOH:O	2.43	0.51
1:D:51:TRP:O	1:D:54:GLN:HG2	2.11	0.50
4:A:303:GOL:O2	7:A:401:HOH:O	2.19	0.50
2:B:50:GLU:HG3	7:B:125:HOH:O	2.11	0.50
1:A:227:ASP:O	1:A:247:VAL:HG23	2.10	0.50
1:A:192:HIS:HB2	1:A:200:THR:HB	1.93	0.50
1:D:55:GLU:OE2	1:D:170:ARG:NE	2.45	0.49
2:B:36:GLU:O	2:B:82:VAL:HA	2.12	0.49
1:D:108:ARG:CZ	1:D:108:ARG:HB2	2.43	0.49
2:B:32:PRO:O	2:B:84:HIS:HE1	1.97	0.48
1:A:152:GLU:OE1	3:C:6:LEU:HB3	2.14	0.48
1:A:255:GLN:O	1:A:273:ARG:NH1	2.47	0.48
1:A:111:ARG:HD2	1:A:128:GLU:HG3	1.96	0.47
1:A:36:PHE:CZ	7:A:411:HOH:O	2.68	0.47
1:A:111:ARG:NH1	7:A:402:HOH:O	2.23	0.47
2:E:24:ASN:HB3	2:E:65:LEU:HD11	1.97	0.47
4:D:305:GOL:H31	7:D:421:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:80:ASN:OD1	4:D:303:GOL:H31	2.15	0.46
1:D:71:ALA:CB	4:D:305:GOL:H12	2.29	0.46
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.50	0.46
1:A:161[B]:GLU:OE1	1:A:161[B]:GLU:C	2.53	0.46
1:D:119:ASP:OD2	4:D:302:GOL:O1	2.28	0.45
1:A:37:ASP:C	7:A:411:HOH:O	2.56	0.45
1:A:44:ARG:NH1	1:D:75:ARG:NH1	2.65	0.45
1:D:152[A]:GLU:O	1:D:156:ARG:HG2	2.18	0.44
1:D:177:ASP:O	1:D:181:ARG:NH1	2.50	0.44
4:D:305:GOL:H32	7:D:462:HOH:O	2.18	0.44
1:D:102:ASP:OD2	1:D:113:HIS:HE1	2.02	0.43
1:A:70:GLN:HE22	4:A:301:GOL:C1	2.27	0.43
1:A:181:ARG:NH2	1:A:183:ASP:OD2	2.52	0.43
1:D:80:ASN:HA	4:D:303:GOL:C3	2.49	0.43
1:D:84:TYR:CE1	4:D:303:GOL:H32	2.53	0.43
2:E:37:VAL:HG21	2:E:66:TYR:CE1	2.52	0.43
1:A:152:GLU:OE2	3:C:7:ARG:NH1	2.46	0.43
1:D:80:ASN:CB	4:D:303:GOL:H32	2.47	0.43
1:D:69:ALA:N	4:D:304:GOL:H2	2.33	0.43
1:A:224:GLN:O	1:A:228:THR:OG1	2.32	0.42
7:A:414:HOH:O	1:D:79:ARG:NH1	2.52	0.42
1:A:194:ILE:HG12	1:A:198:GLU:HB2	2.01	0.42
1:D:230:LEU:CD2	1:D:245:ALA:HB2	2.47	0.42
1:A:108:ARG:NE	7:A:425:HOH:O	2.52	0.42
1:D:35:ARG:HD3	2:E:53:ASP:OD2	2.20	0.42
2:B:17:ASN:HD21	2:B:74:GLU:HA	1.86	0.41
1:D:170:ARG:NH1	1:D:170:ARG:HG3	2.34	0.41
2:B:51:HIS:HA	2:B:65:LEU:O	2.19	0.41
1:A:19:GLU:CB	1:A:75:ARG:NH2	2.84	0.41
1:A:36:PHE:CD2	7:A:411:HOH:O	2.72	0.41
1:A:144:GLN:O	1:A:148:GLU:HG3	2.21	0.41
1:A:230:LEU:HD23	1:A:245:ALA:HB2	2.03	0.41
1:D:113:HIS:HD2	7:D:553:HOH:O	2.03	0.41
2:E:59:ASP:O	2:E:60:TRP:HB2	2.20	0.41
1:A:120:GLY:O	2:B:3:ARG:NH1	2.51	0.41
1:A:108:ARG:HH21	1:A:108:ARG:HB3	1.86	0.40
1:D:44:ARG:NH2	1:D:61:ASP:OD1	2.46	0.40
2:E:19:LYS:O	2:E:72:PRO:HD2	2.22	0.40
1:D:117:ALA:HB2	2:E:60:TRP:CE2	2.56	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	A	277/277 (100%)	266 (96%)	11 (4%)	0	100	100
1	D	279/277 (101%)	270 (97%)	9 (3%)	0	100	100
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%)	6 (86%)	1 (14%)	0	100	100
3	F	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
All	All	766/772 (99%)	741 (97%)	25 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	237/235 (101%)	227 (96%)	10 (4%)	30	30
1	D	239/235 (102%)	222 (93%)	17 (7%)	14	11
2	B	95/95 (100%)	89 (94%)	6 (6%)	18	15
2	E	95/95 (100%)	90 (95%)	5 (5%)	22	20
3	C	9/9 (100%)	9 (100%)	0	100	100
3	F	9/9 (100%)	9 (100%)	0	100	100
All	All	684/678 (101%)	646 (94%)	38 (6%)	21	18



All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	0	MET
1	A	177	ASP
1	A	178	LYS
1	A	183	ASP
1	A	186	LYS
1	A	194	ILE
1	A	223	ASP
1	A	251	SER
1	A	268	LYS
1	A	270	LEU
2	B	1	ILE
2	B	19	LYS
2	B	48	LYS
2	B	70	PHE
2	B	74	GLU
2	B	75	LYS
1	D	0	MET
1	D	17	ARG
1	D	44	ARG
1	D	121	LYS
1	D	170	ARG
1	D	190	THR
1	D	191	HIS
1	D	197	HIS
1	D	218	GLN
1	D	219	ARG
1	D	227	ASP
1	D	232	GLU
1	D	247	VAL
1	D	258	THR
1	D	268	LYS
1	D	270	LEU
1	D	275	GLU
2	E	1	ILE
2	E	20	SER
2	E	70	PHE
2	E	75	LYS
2	E	91	LYS

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

Mol	Chain	Res	Type
1	A	70	GLN
1	A	113	HIS
1	A	188	HIS
1	A	260	HIS
2	B	84	HIS
1	D	113	HIS
1	D	155	GLN
1	D	191	HIS
1	D	218	GLN
2	E	84	HIS

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

15 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
4	GOL	A	302	-	5,5,5	1.86	3 (60%)	5,5,5	3.85	4 (80%)
4	GOL	D	304	-	5,5,5	0.81	0	5,5,5	1.43	1 (20%)
4	GOL	A	301	-	5,5,5	1.18	0	5,5,5	1.69	1 (20%)
5	EDO	A	304	-	3,3,3	0.57	0	2,2,2	0.36	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	EDO	A	305	-	3,3,3	0.48	0	2,2,2	0.29	0
5	EDO	D	307	-	3,3,3	0.63	0	2,2,2	0.27	0
5	EDO	D	306	-	3,3,3	0.84	0	2,2,2	0.49	0
4	GOL	D	302	-	5,5,5	0.57	0	5,5,5	0.70	0
5	EDO	D	308	-	3,3,3	0.50	0	2,2,2	0.72	0
5	EDO	A	306	-	3,3,3	0.49	0	2,2,2	0.20	0
4	GOL	A	303	-	5,5,5	1.05	0	5,5,5	1.70	2 (40%)
6	SO4	A	307	-	4,4,4	0.45	0	6,6,6	0.34	0
4	GOL	D	305	-	5,5,5	1.11	0	5,5,5	1.50	1 (20%)
4	GOL	D	303	-	5,5,5	1.90	1 (20%)	5,5,5	2.37	2 (40%)
4	GOL	D	301	-	5,5,5	0.99	0	5,5,5	1.31	1 (20%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	A	302	-	-	2/4/4/4	-
4	GOL	D	304	-	-	0/4/4/4	-
4	GOL	A	301	-	-	2/4/4/4	-
5	EDO	A	304	-	-	1/1/1/1	-
5	EDO	A	305	-	-	0/1/1/1	-
5	EDO	D	307	-	-	1/1/1/1	-
5	EDO	D	306	-	-	1/1/1/1	-
4	GOL	D	302	-	-	3/4/4/4	-
5	EDO	D	308	-	-	1/1/1/1	-
5	EDO	A	306	-	-	1/1/1/1	-
4	GOL	A	303	-	-	4/4/4/4	-
4	GOL	D	305	-	-	2/4/4/4	-
4	GOL	D	303	-	-	4/4/4/4	-
4	GOL	D	301	-	-	2/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	303	GOL	O2-C2	-3.83	1.31	1.43
4	A	302	GOL	O2-C2	-2.52	1.35	1.43
4	A	302	GOL	O1-C1	2.43	1.52	1.42
4	A	302	GOL	C1-C2	2.08	1.60	1.51

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	302	GOL	O1-C1-C2	5.44	136.27	110.20
4	A	302	GOL	O2-C2-C3	-4.87	87.67	109.12
4	D	303	GOL	O3-C3-C2	4.66	132.56	110.20
4	A	302	GOL	O3-C3-C2	3.96	129.18	110.20
4	D	304	GOL	O1-C1-C2	2.65	122.89	110.20
4	A	301	GOL	C3-C2-C1	2.55	121.62	111.70
4	D	301	GOL	O1-C1-C2	2.41	121.75	110.20
4	D	303	GOL	O1-C1-C2	2.39	121.65	110.20
4	D	305	GOL	O1-C1-C2	2.29	121.20	110.20
4	A	302	GOL	C3-C2-C1	2.28	120.59	111.70
4	A	303	GOL	O1-C1-C2	2.28	121.15	110.20
4	A	303	GOL	O2-C2-C3	-2.18	99.51	109.12

There are no chirality outliers.

All (24) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	301	GOL	O1-C1-C2-O2
4	A	301	GOL	O1-C1-C2-C3
4	A	302	GOL	C1-C2-C3-O3
4	A	303	GOL	C1-C2-C3-O3
4	A	303	GOL	O2-C2-C3-O3
4	D	302	GOL	O1-C1-C2-C3
4	D	303	GOL	O1-C1-C2-C3
4	D	303	GOL	C1-C2-C3-O3
4	D	305	GOL	C1-C2-C3-O3
5	A	306	EDO	O1-C1-C2-O2
4	A	303	GOL	O1-C1-C2-C3
4	D	301	GOL	O1-C1-C2-C3
4	A	302	GOL	O2-C2-C3-O3
4	A	303	GOL	O1-C1-C2-O2
4	D	301	GOL	O1-C1-C2-O2
4	D	302	GOL	O1-C1-C2-O2
4	D	303	GOL	O2-C2-C3-O3
4	D	305	GOL	O2-C2-C3-O3
5	A	304	EDO	O1-C1-C2-O2
5	D	306	EDO	O1-C1-C2-O2
4	D	303	GOL	O1-C1-C2-O2
4	D	302	GOL	C1-C2-C3-O3
5	D	308	EDO	O1-C1-C2-O2
5	D	307	EDO	O1-C1-C2-O2

There are no ring outliers.

10 monomers are involved in 38 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	302	GOL	11	0
4	D	304	GOL	3	0
4	A	301	GOL	2	0
5	A	305	EDO	1	0
4	D	302	GOL	1	0
5	D	308	EDO	1	0
5	A	306	EDO	2	0
4	A	303	GOL	3	0
4	D	305	GOL	8	0
4	D	303	GOL	6	0

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	277/277 (100%)	0.90	42 (15%) <b>2</b> <b>3</b>	16, 36, 124, 139	0
1	D	277/277 (100%)	0.81	46 (16%) <b>1</b> <b>2</b>	17, 35, 122, 140	0
2	B	100/100 (100%)	0.85	19 (19%) <b>1</b> <b>1</b>	25, 60, 89, 111	0
2	E	100/100 (100%)	0.60	12 (12%) <b>4</b> <b>5</b>	25, 56, 86, 111	0
3	C	9/9 (100%)	0.33	0 <b>100</b> <b>100</b>	21, 25, 33, 34	0
3	F	9/9 (100%)	0.25	0 <b>100</b> <b>100</b>	20, 25, 33, 34	0
All	All	772/772 (100%)	0.81	119 (15%) <b>2</b> <b>2</b>	16, 43, 117, 140	0

All (119) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	219	ARG	7.5
1	A	194	ILE	7.0
1	D	250	PRO	6.8
1	A	274	TRP	6.4
2	E	1	ILE	6.2
1	A	195	SER	5.8
1	A	199	ALA	5.8
1	D	225	THR	5.6
1	D	258	THR	5.6
2	E	0	MET	5.6
1	D	221	GLY	5.5
2	B	0	MET	5.5
1	A	257	TYR	5.2
1	D	228	THR	5.2
1	D	197	HIS	5.2
1	A	223	ASP	5.2
1	A	225	THR	5.2
1	A	250	PRO	5.1
1	A	198	GLU	4.8

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	D	199	ALA	4.7
1	D	195	SER	4.7
1	A	221	GLY	4.6
1	D	252	GLY	4.4
1	D	220	ASP	4.4
1	D	223	ASP	4.3
1	A	197	HIS	4.2
1	A	0	MET	4.1
1	D	230	LEU	4.1
1	A	268	LYS	4.1
1	D	194	ILE	4.0
1	D	245	ALA	4.0
1	A	220	ASP	3.9
1	A	219	ARG	3.9
1	D	0	MET	3.9
2	B	95	TRP	3.9
1	D	275	GLU	3.8
1	A	276	PRO	3.8
1	D	254	GLU	3.8
1	A	258	THR	3.8
2	B	1	ILE	3.7
1	D	222	GLU	3.7
1	A	191	HIS	3.6
2	E	75	LYS	3.5
1	A	230	LEU	3.5
1	D	268	LYS	3.5
1	A	253	GLU	3.4
1	D	261	VAL	3.4
1	A	200	THR	3.4
1	A	251	SER	3.4
1	A	245	ALA	3.4
1	D	192	HIS	3.3
1	D	198	GLU	3.3
1	A	201	LEU	3.3
1	D	276	PRO	3.2
1	D	272	LEU	3.2
1	D	274	TRP	3.0
1	A	227	ASP	3.0
1	D	253	GLU	3.0
1	D	1	GLY	2.9
2	B	40	LEU	2.9
1	D	203	CYS	2.9

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
2	E	15	ALA	2.9
2	B	99	MET	2.9
1	D	204	TRP	2.8
2	E	48	LYS	2.8
1	D	189	VAL	2.7
1	D	227	ASP	2.7
2	E	16	GLU	2.7
1	A	193	PRO	2.7
1	D	256	ARG	2.7
1	D	191	HIS	2.7
2	B	16	GLU	2.7
1	A	190	THR	2.7
1	A	192	HIS	2.7
2	B	76	ASP	2.6
2	E	88	SER	2.6
1	A	275	GLU	2.6
1	A	247	VAL	2.6
1	A	217	TRP	2.5
2	B	20	SER	2.5
1	A	215	LEU	2.5
1	D	270	LEU	2.5
2	E	74	GLU	2.5
1	A	203	CYS	2.5
1	D	196	ASP	2.5
1	D	249	VAL	2.4
1	D	188	HIS	2.4
1	D	217	TRP	2.4
1	D	224	GLN	2.4
1	A	18	GLY	2.4
2	B	94	LYS	2.3
1	A	259	CYS	2.3
2	B	73	THR	2.3
2	E	44	GLU	2.3
2	B	78	TYR	2.3
1	A	228	THR	2.3
2	B	80	CYS	2.3
1	A	196	ASP	2.2
1	A	272	LEU	2.2
2	E	78	TYR	2.2
2	B	2	GLN	2.2
2	B	92	ILE	2.2
2	E	18	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	248	VAL	2.1
1	A	218	GLN	2.1
1	D	267	PRO	2.1
2	B	75	LYS	2.1
1	D	244	TRP	2.1
1	A	266	LEU	2.1
1	D	190	THR	2.1
1	D	248	VAL	2.1
1	D	202	ARG	2.1
2	B	98	ASP	2.1
2	B	19	LYS	2.1
1	D	193	PRO	2.0
2	B	44	GLU	2.0
1	A	108	ARG	2.0
2	E	19	LYS	2.0
2	B	47	GLU	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, 95<sup>th</sup> 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(Å <sup>2</sup> )	Q<0.9
4	GOL	D	302	6/6	0.77	0.20	41,52,61,63	0
4	GOL	A	303	6/6	0.84	0.23	31,32,36,39	0
4	GOL	A	301	6/6	0.85	0.21	27,38,40,41	0
4	GOL	D	305	6/6	0.85	0.33	27,40,44,65	0
5	EDO	D	308	4/4	0.85	0.32	41,41,42,46	0
4	GOL	D	304	6/6	0.86	0.26	24,37,45,54	0
4	GOL	A	302	6/6	0.87	0.33	24,28,33,37	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	EDO	D	307	4/4	0.90	0.16	37,46,46,49	0
4	GOL	D	301	6/6	0.91	0.16	38,40,47,54	0
5	EDO	A	305	4/4	0.91	0.11	39,41,45,51	0
4	GOL	D	303	6/6	0.93	0.24	24,27,32,33	0
5	EDO	D	306	4/4	0.94	0.16	24,25,25,26	0
6	SO4	A	307	5/5	0.94	0.19	43,68,71,76	0
5	EDO	A	304	4/4	0.95	0.17	32,36,39,41	0
5	EDO	A	306	4/4	0.96	0.12	39,45,49,52	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.