



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 12:53 am GMT

PDB ID : 4I0P  
Title : HLA-DO in complex with HLA-DM  
Authors : Guce, A.I.; Mortimer, S.E.; Stern, L.J.  
Deposited on : 2012-11-18  
Resolution : 3.20 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : recalc28949

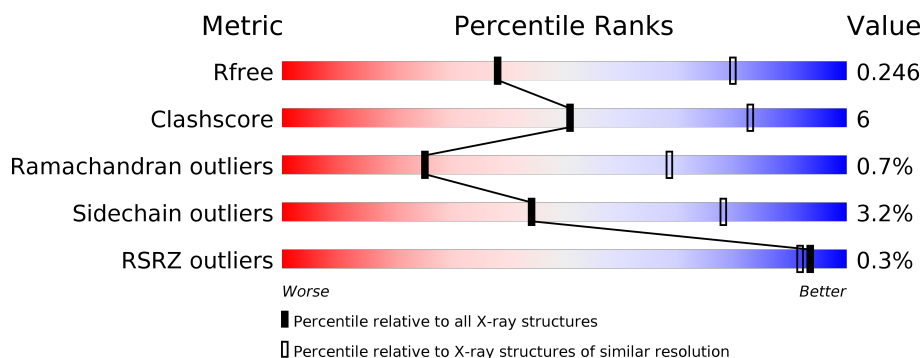
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	100719	1015 (3.22-3.18)
Clashscore	112137	1009 (3.20-3.20)
Ramachandran outliers	110173	1118 (3.22-3.18)
Sidechain outliers	110143	1117 (3.22-3.18)
RSRZ outliers	101464	1020 (3.22-3.18)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	188	<div> <div>%</div> <div> <div></div> <div>83%</div> <div>17%</div> </div> </div>
1	E	188	<div> <div>%</div> <div> <div></div> <div>85%</div> <div>15%</div> </div> </div>
2	B	191	<div> <div></div> <div> <div></div> <div>82%</div> <div>15%</div> <div>••</div> </div> </div>
2	F	191	<div> <div>%</div> <div> <div></div> <div>83%</div> <div>15%</div> <div>••</div> </div> </div>
3	C	181	<div> <div></div> <div> <div></div> <div>86%</div> <div>13%</div> <div>•</div> </div> </div>
3	G	181	<div> <div></div> <div> <div></div> <div>85%</div> <div>14%</div> <div>•</div> </div> </div>

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Mol	Chain	Length	Quality of chain
4	D	189	<div> <div>%</div> <div> </div> <div>78%17%</div> <div>• •</div> </div>
4	H	189	<div> <div> </div> <div>84%14%</div> <div>•</div> </div>

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
5	NAG	A	301	-	-	-	X

## 2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 13370 atoms, of which 1306 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-DMA protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	188	Total	C	H	N	O	S	0	0	0
			1667	984	159	239	279	6			
1	E	188	Total	C	H	N	O	S	0	0	0
			1667	984	159	239	279	6			

- Molecule 2 is a protein called HLA class II histocompatibility antigen, DM beta chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	186	Total	C	H	N	O	S	0	0	0
			1630	938	155	252	274	11			
2	F	187	Total	C	H	N	O	S	0	0	0
			1638	944	155	253	275	11			

- Molecule 3 is a protein called HLA class II histocompatibility antigen, DO alpha chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	181	Total	C	H	N	O	S	0	0	0
			1609	939	152	251	264	3			
3	G	181	Total	C	H	N	O	S	0	0	0
			1609	939	152	251	264	3			

- Molecule 4 is a protein called HLA class II histocompatibility antigen, DO beta chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
4	D	188	Total	C	H	N	O	S	0	0	0
			1704	968	187	263	278	8			
4	H	189	Total	C	H	N	O	S	0	0	0
			1710	971	187	264	280	8			

- Molecule 5 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



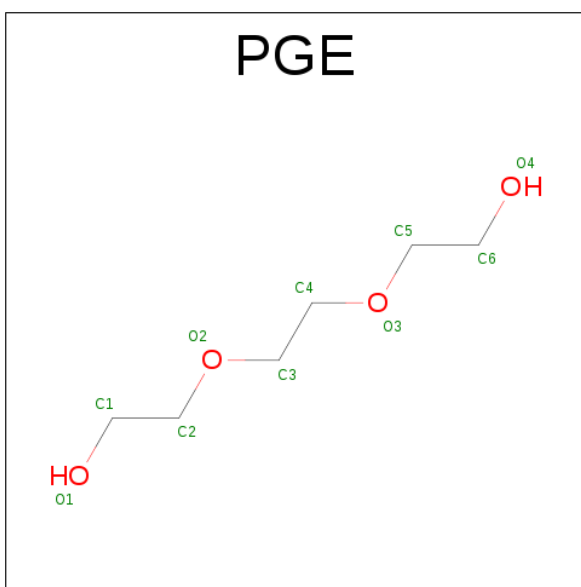
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	C	1	Total	C	N	O	0	0
			14	8	1	5		
5	D	1	Total	C	N	O	0	0
			14	8	1	5		
5	E	1	Total	C	N	O	0	0
			14	8	1	5		
5	G	1	Total	C	N	O	0	0
			14	8	1	5		
5	G	1	Total	C	N	O	0	0
			14	8	1	5		
5	H	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



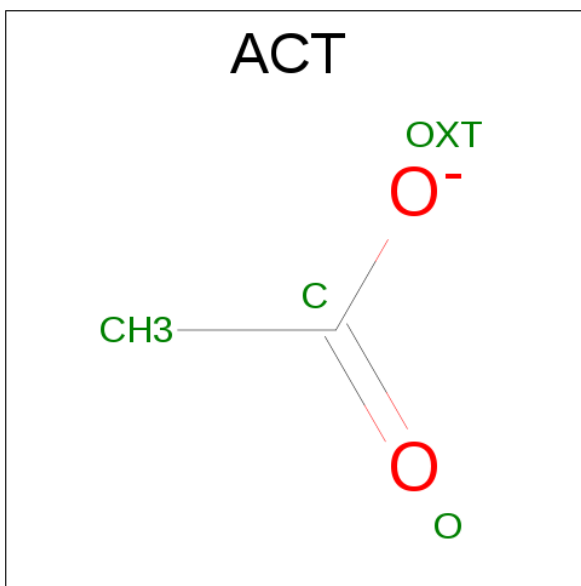
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	D	1	Total	C	O	0	0
			6	3	3		
6	F	1	Total	C	O	0	0
			6	3	3		
6	H	1	Total	C	O	0	0
			6	3	3		

- Molecule 7 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula:  $C_6H_{14}O_4$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	H	1	Total	C	O	0	0
			10	6	4		

- Molecule 8 is ACETATE ION (three-letter code: ACT) (formula:  $C_2H_3O_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	H	1	Total	C	O	0	0
			4	2	2		

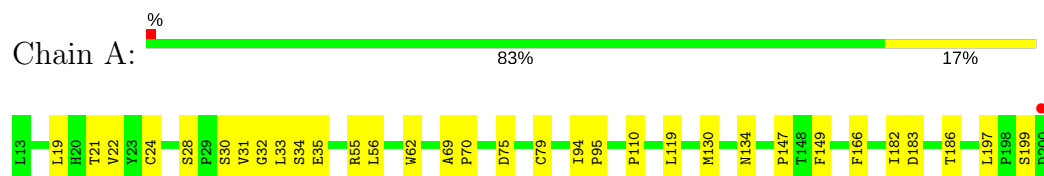
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	B	2	Total	O	0	0
			2	2		
9	C	1	Total	O	0	0
			1	1		
9	F	1	Total	O	0	0
			1	1		
9	G	2	Total	O	0	0
			2	2		

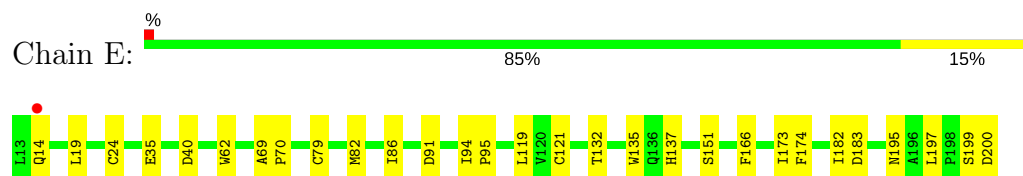
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

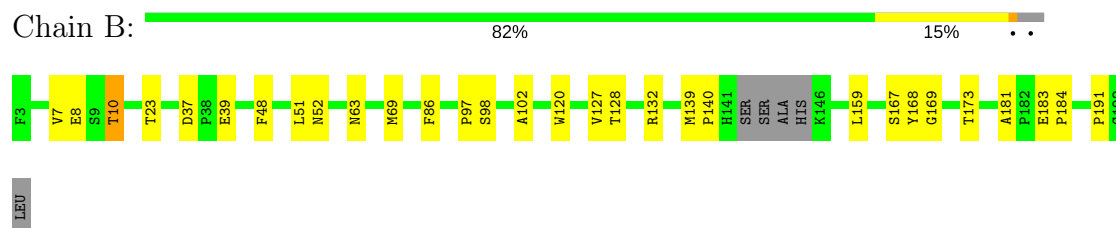
- Molecule 1: HLA-DMA protein



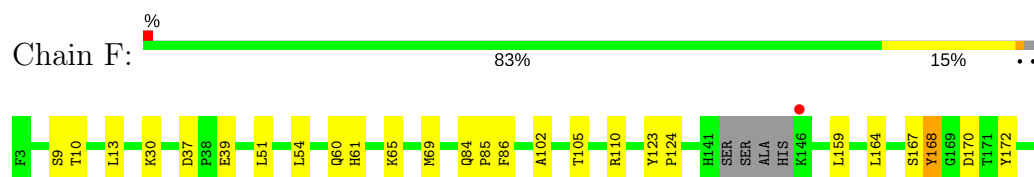
- Molecule 1: HLA-DMA protein



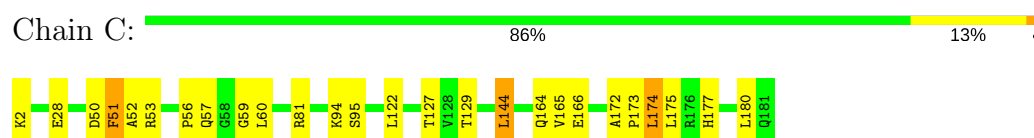
- Molecule 2: HLA class II histocompatibility antigen, DM beta chain



- Molecule 2: HLA class II histocompatibility antigen, DM beta chain




- Molecule 3: HLA class II histocompatibility antigen, DO alpha chain




- Molecule 3: HLA class II histocompatibility antigen, DO alpha chain

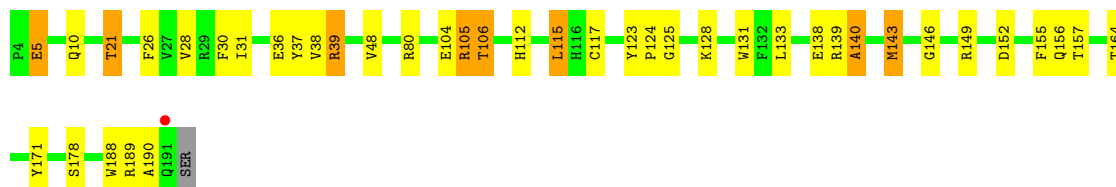


Chain G:  85% 14% .




- Molecule 4: HLA class II histocompatibility antigen, DO beta chain

Chain D:  78% 17% . .



- Molecule 4: HLA class II histocompatibility antigen, DO beta chain

Chain H:  84% 14% .



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	83.26Å 147.10Å 95.96Å 90.00° 106.49° 90.00°	Depositor
Resolution (Å)	43.91 – 3.20 43.91 – 3.20	Depositor EDS
% Data completeness (in resolution range)	94.0 (43.91-3.20) 93.9 (43.91-3.20)	Depositor EDS
$R_{merge}$	0.68	Depositor
$R_{sym}$	0.68	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.92 (at 3.19Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, $R_{free}$	0.188 , 0.242 0.190 , 0.246	Depositor DCC
$R_{free}$ test set	1855 reflections (5.40%)	DCC
Wilson B-factor (Å <sup>2</sup> )	49.0	Xtriage
Anisotropy	0.397	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 19.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	13370	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.06% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, PGE, NAG, ACT

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	0.24	0/1561	0.42	0/2135
1	E	0.24	0/1561	0.44	0/2135
2	B	0.23	0/1519	0.43	0/2079
2	F	0.23	0/1527	0.43	0/2090
3	C	0.24	0/1502	0.49	0/2046
3	G	0.24	0/1502	0.46	0/2046
4	D	0.24	0/1557	0.45	0/2116
4	H	0.24	0/1563	0.45	0/2124
All	All	0.24	0/12292	0.45	0/16771

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
3	C	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
3	C	51	PHE	Peptide

## 5.2 Too-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 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	1508	159	1418	21	0
1	E	1508	159	1418	18	0
2	B	1475	155	1418	21	0
2	F	1483	155	1429	20	0
3	C	1457	152	1401	16	0
3	G	1457	152	1400	21	0
4	D	1517	187	1469	33	0
4	H	1523	187	1474	24	0
5	A	14	0	13	0	0
5	C	14	0	13	0	0
5	D	14	0	13	0	0
5	E	14	0	13	1	0
5	G	28	0	26	0	0
5	H	14	0	13	0	0
6	D	6	0	8	0	0
6	F	6	0	8	1	0
6	H	6	0	8	0	0
7	H	10	0	14	0	0
8	H	4	0	3	0	0
9	B	2	0	0	0	0
9	C	1	0	0	0	0
9	F	1	0	0	0	0
9	G	2	0	0	0	0
All	All	12064	1306	11559	153	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 6.

All (153) 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)
3:G:77:SER:O	3:G:79:ARG:N	2.12	0.82
4:D:189:ARG:HB2	4:D:190:ALA:O	1.91	0.70
2:F:110:ARG:NH1	4:H:185:SER:OG	2.30	0.65
2:F:167:SER:O	2:F:168:TYR:CB	2.45	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:115:LEU:HD13	4:H:131:TRP:CH2	2.33	0.64
4:D:188:TRP:CZ3	4:D:190:ALA:HB2	2.34	0.63
4:D:189:ARG:HB2	4:D:190:ALA:HA	1.82	0.60
2:B:63:ASN:HA	2:B:69:MET:HE3	1.84	0.59
4:D:104:GLU:O	4:D:105:ARG:C	2.40	0.59
4:D:189:ARG:HB2	4:D:190:ALA:CA	2.32	0.59
4:H:139:ARG:O	4:H:140:ALA:HB2	2.03	0.58
3:C:52:ALA:O	3:C:53:ARG:HB2	2.04	0.57
1:E:182:ILE:O	1:E:183:ASP:CG	2.43	0.57
4:D:139:ARG:O	4:D:140:ALA:HB2	2.05	0.57
4:H:37:TYR:CD2	4:H:38:VAL:HG12	2.41	0.56
4:D:21:THR:O	4:D:80:ARG:NH1	2.38	0.56
3:G:52:ALA:O	3:G:53:ARG:HB2	2.06	0.56
3:C:164:GLN:HG3	3:C:164:GLN:O	2.05	0.56
1:E:86:ILE:HD13	2:F:54:LEU:HD11	1.88	0.55
1:A:183:ASP:OD1	1:A:183:ASP:C	2.45	0.55
1:A:182:ILE:HG13	1:A:182:ILE:O	2.06	0.55
2:B:167:SER:O	2:B:169:GLY:N	2.39	0.54
3:C:174:LEU:HD12	3:C:174:LEU:C	2.28	0.54
1:E:183:ASP:OD1	1:E:183:ASP:C	2.46	0.53
2:B:183:GLU:HG3	2:B:184:PRO:HD2	1.90	0.53
3:G:77:SER:O	3:G:78:ASN:C	2.46	0.53
4:D:37:TYR:CD2	4:D:38:VAL:HG12	2.43	0.52
2:F:192:GLY:O	2:F:193:LEU:HB3	2.09	0.52
1:A:182:ILE:O	1:A:183:ASP:CG	2.48	0.52
3:G:164:GLN:O	3:G:164:GLN:HG3	2.10	0.52
1:A:94:ILE:CD1	2:B:51:LEU:HD21	2.40	0.52
4:H:189:ARG:O	4:H:190:ALA:HB3	2.10	0.52
1:A:75:ASP:HB3	2:B:7:VAL:HG21	1.91	0.52
4:D:189:ARG:CB	4:D:190:ALA:O	2.56	0.52
2:B:10:THR:HG23	2:B:23:THR:OG1	2.10	0.51
3:C:180:LEU:C	3:C:180:LEU:HD12	2.31	0.51
4:H:37:TYR:CE2	4:H:38:VAL:HG12	2.45	0.51
4:H:139:ARG:O	4:H:140:ALA:CB	2.57	0.51
1:E:182:ILE:HG13	1:E:182:ILE:O	2.10	0.50
1:E:24:CYS:SG	1:E:79:CYS:C	2.90	0.50
3:C:59:GLY:O	3:C:60:LEU:HB2	2.10	0.50
2:B:97:PRO:HD2	2:B:181:ALA:HB2	1.93	0.50
3:C:144:LEU:N	3:C:144:LEU:HD22	2.26	0.50
1:A:69:ALA:N	1:A:70:PRO:HD2	2.27	0.50
2:B:7:VAL:O	2:B:7:VAL:HG23	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:26:PHE:CZ	4:D:28:VAL:CG2	2.95	0.50
4:H:117:CYS:HB2	4:H:131:TRP:CZ2	2.48	0.49
2:B:7:VAL:CG2	2:B:7:VAL:O	2.59	0.49
1:E:119:LEU:HG	1:E:166:PHE:CE1	2.48	0.49
1:A:110:PRO:HG2	2:B:120:TRP:CZ3	2.48	0.49
1:E:62:TRP:CD2	2:F:86:PHE:HB3	2.48	0.49
3:G:174:LEU:HD23	3:G:175:LEU:N	2.26	0.49
3:G:59:GLY:O	3:G:60:LEU:HB2	2.13	0.49
4:D:10:GLN:HB2	4:D:31:ILE:HB	1.95	0.48
3:G:89:VAL:HG21	3:G:165:VAL:HG11	1.95	0.48
4:D:189:ARG:HB2	4:D:190:ALA:C	2.34	0.48
1:E:19:LEU:O	1:E:35:GLU:HA	2.14	0.48
3:C:172:ALA:HB1	3:C:173:PRO:HD2	1.94	0.48
4:H:21:THR:O	4:H:80:ARG:NH1	2.46	0.48
4:D:115:LEU:HD13	4:D:131:TRP:CH2	2.49	0.47
2:B:37:ASP:OD1	2:B:39:GLU:HG2	2.14	0.47
4:H:104:GLU:O	4:H:105:ARG:C	2.53	0.47
3:C:144:LEU:N	3:C:144:LEU:CD2	2.77	0.47
4:D:189:ARG:CB	4:D:190:ALA:CA	2.93	0.47
4:H:103:PRO:HA	4:H:115:LEU:HA	1.96	0.47
1:A:119:LEU:HG	1:A:166:PHE:CE1	2.49	0.47
3:G:50:ASP:HB3	3:G:52:ALA:O	2.15	0.47
1:A:28:SER:O	1:A:130:MET:HG3	2.14	0.47
1:A:147:PRO:HG2	1:A:149:PHE:CZ	2.49	0.47
4:H:87:ALA:N	4:H:88:PRO:HD2	2.29	0.47
1:A:21:THR:HG23	2:B:8:GLU:HG3	1.96	0.47
2:F:167:SER:O	2:F:168:TYR:HB3	2.15	0.46
3:G:113:PHE:HB2	3:G:144:LEU:HD13	1.98	0.46
4:H:105:ARG:O	4:H:106:THR:C	2.53	0.46
4:D:112:HIS:HA	4:D:164:THR:HG22	1.98	0.46
4:D:37:TYR:CE2	4:D:38:VAL:HG12	2.51	0.46
1:A:19:LEU:O	1:A:35:GLU:HA	2.15	0.46
4:D:26:PHE:CE2	4:D:28:VAL:HG23	2.50	0.46
3:C:122:LEU:HD23	3:C:127:THR:HA	1.98	0.45
1:E:197:LEU:HD22	2:F:102:ALA:HB3	1.98	0.45
3:C:50:ASP:O	3:C:51:PHE:HB2	2.16	0.45
3:G:113:PHE:HB2	3:G:144:LEU:CD1	2.47	0.45
1:E:137:HIS:CD2	1:E:174:PHE:CZ	3.04	0.45
2:B:98:SER:HB2	2:B:120:TRP:CE2	2.52	0.45
4:D:133:LEU:HD12	4:D:171:TYR:CE2	2.51	0.45
1:E:121:CYS:HB2	1:E:135:TRP:CZ2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:82:MET:O	1:E:86:ILE:O	2.33	0.45
4:H:21:THR:CG2	4:H:21:THR:O	2.64	0.45
4:D:36:GLU:OE2	4:D:39:ARG:NH1	2.50	0.45
2:F:167:SER:O	2:F:168:TYR:HB2	2.18	0.44
3:C:94:LYS:HE2	4:D:152:ASP:OD2	2.17	0.44
2:F:30:LYS:HE3	3:G:49:GLY:O	2.17	0.44
3:G:174:LEU:C	3:G:174:LEU:HD23	2.38	0.44
4:H:21:THR:O	4:H:21:THR:HG22	2.17	0.44
1:A:95:PRO:CD	3:C:57:GLN:HA	2.47	0.44
4:D:139:ARG:O	4:D:140:ALA:CB	2.66	0.44
4:H:112:HIS:HA	4:H:164:THR:HG22	2.00	0.44
3:C:177:HIS:HB2	4:H:143:MET:SD	2.58	0.44
2:F:170:ASP:O	2:F:190:THR:HG22	2.18	0.43
1:A:55:ARG:HG3	1:A:56:LEU:CD1	2.48	0.43
1:A:94:ILE:HD13	2:B:51:LEU:HD21	2.00	0.43
2:F:84:GLN:N	2:F:85:PRO:HD2	2.33	0.43
3:G:50:ASP:O	3:G:51:PHE:HB2	2.17	0.43
3:C:166:GLU:HG2	3:C:173:PRO:HB3	2.00	0.43
1:E:94:ILE:CD1	2:F:51:LEU:HD21	2.48	0.43
1:A:22:VAL:HB	2:B:7:VAL:HG22	2.01	0.43
4:D:123:TYR:CG	4:D:124:PRO:HA	2.53	0.43
3:C:81:ARG:CZ	4:D:5:GLU:HG2	2.49	0.43
3:G:52:ALA:O	3:G:53:ARG:CB	2.66	0.43
4:H:76:ASP:HA	4:H:80:ARG:HD2	2.01	0.43
1:A:24:CYS:SG	1:A:79:CYS:C	2.97	0.43
4:D:125:GLY:CA	4:D:155:PHE:CD1	3.01	0.43
3:C:28:GLU:HB2	4:D:149:ARG:CZ	2.48	0.43
2:F:65:LYS:O	2:F:69:MET:HG2	2.18	0.43
4:D:143:MET:SD	3:G:177:HIS:HB2	2.59	0.43
1:A:197:LEU:HD22	2:B:102:ALA:HB3	2.00	0.42
2:F:37:ASP:OD1	2:F:39:GLU:HG2	2.19	0.42
4:D:105:ARG:O	4:D:106:THR:C	2.58	0.42
3:G:144:LEU:HD22	3:G:144:LEU:N	2.34	0.42
4:H:55:GLN:N	4:H:56:PRO:CD	2.82	0.42
1:A:62:TRP:CD2	2:B:86:PHE:HB3	2.54	0.42
1:E:195:ASN:OD1	2:F:105:THR:N	2.52	0.42
4:D:146:GLY:O	4:D:157:THR:HG22	2.20	0.42
3:G:10:GLY:N	3:G:10(A):PRO:CD	2.83	0.42
1:A:31:VAL:HG22	1:A:32:GLY:N	2.35	0.41
3:G:88:ARG:HG3	3:G:88:ARG:O	2.20	0.41
2:F:13:LEU:N	2:F:13:LEU:HD12	2.34	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:117:CYS:HB2	4:D:131:TRP:CZ2	2.55	0.41
2:F:123:TYR:CG	2:F:124:PRO:HA	2.55	0.41
2:B:132:ARG:HB2	2:B:173:THR:HB	2.02	0.41
1:E:94:ILE:HG23	1:E:95:PRO:HD2	2.02	0.41
4:H:189:ARG:O	4:H:190:ALA:CB	2.68	0.41
2:F:60:GLN:HG3	2:F:61:HIS:N	2.36	0.41
3:G:122:LEU:HD23	3:G:127:THR:HA	2.03	0.41
4:H:133:LEU:HD23	4:H:133:LEU:C	2.41	0.41
4:H:26:PHE:CE2	4:H:28:VAL:HG23	2.55	0.41
1:E:199:SER:HA	1:E:200:ASP:HA	1.92	0.41
2:F:164:LEU:HD22	2:F:172:TYR:CE1	2.56	0.41
4:H:21:THR:HG22	4:H:80:ARG:NH1	2.36	0.41
4:D:115:LEU:O	4:D:115:LEU:HD12	2.21	0.41
4:D:188:TRP:CZ3	4:D:190:ALA:CB	3.04	0.41
2:F:65:LYS:HE2	6:F:201:GOL:O2	2.21	0.41
2:B:139:MET:HA	2:B:140:PRO:HD3	1.94	0.40
4:D:21:THR:HG22	4:D:80:ARG:HG2	2.03	0.40
3:G:142:ASP:CG	3:G:144:LEU:HD23	2.42	0.40
4:H:115:LEU:HD12	4:H:115:LEU:O	2.20	0.40
1:A:33:LEU:HD12	1:A:34:SER:N	2.35	0.40
2:B:127:VAL:HG22	2:B:128:THR:N	2.36	0.40
2:B:48:PHE:HA	2:B:52:ASN:HB2	2.03	0.40
4:D:30:PHE:HB3	4:D:37:TYR:CZ	2.57	0.40
1:E:14:GLN:HA	5:E:301:NAG:H82	2.03	0.40
1:E:69:ALA:N	1:E:70:PRO:HD2	2.36	0.40
3:G:109:VAL:O	3:G:146:ARG:HA	2.20	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	186/188 (99%)	174 (94%)	12 (6%)	0	100	100
1	E	186/188 (99%)	174 (94%)	12 (6%)	0	100	100
2	B	182/191 (95%)	175 (96%)	5 (3%)	2 (1%)	17	58
2	F	183/191 (96%)	177 (97%)	5 (3%)	1 (0%)	32	74
3	C	179/181 (99%)	166 (93%)	12 (7%)	1 (1%)	28	72
3	G	179/181 (99%)	165 (92%)	12 (7%)	2 (1%)	17	58
4	D	186/189 (98%)	169 (91%)	15 (8%)	2 (1%)	17	58
4	H	187/189 (99%)	170 (91%)	15 (8%)	2 (1%)	17	58
All	All	1468/1498 (98%)	1370 (93%)	88 (6%)	10 (1%)	25	68

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	D	140	ALA
2	F	168	TYR
3	G	78	ASN
4	H	105	ARG
4	H	140	ALA
3	C	56	PRO
4	D	105	ARG
3	G	56	PRO
2	B	168	TYR
2	B	191	PRO

### 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	167/167 (100%)	163 (98%)	4 (2%)	54	84
1	E	167/167 (100%)	162 (97%)	5 (3%)	46	79
2	B	165/169 (98%)	163 (99%)	2 (1%)	75	91
2	F	166/169 (98%)	163 (98%)	3 (2%)	64	87

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	C	156/156 (100%)	149 (96%)	7 (4%)	32	70
3	G	156/156 (100%)	153 (98%)	3 (2%)	62	86
4	D	167/168 (99%)	156 (93%)	11 (7%)	19	57
4	H	168/168 (100%)	161 (96%)	7 (4%)	34	72
All	All	1312/1320 (99%)	1270 (97%)	42 (3%)	44	78

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	SER
1	A	134	ASN
1	A	186	THR
1	A	199	SER
2	B	10	THR
2	B	159	LEU
3	C	2	LYS
3	C	95	SER
3	C	129	THR
3	C	144	LEU
3	C	165	VAL
3	C	174	LEU
3	C	175	LEU
4	D	5	GLU
4	D	21	THR
4	D	39	ARG
4	D	48	VAL
4	D	106	THR
4	D	115	LEU
4	D	128	LYS
4	D	138	GLU
4	D	143	MET
4	D	156	GLN
4	D	178	SER
1	E	40	ASP
1	E	91	ASP
1	E	132	THR
1	E	151	SER
1	E	173	ILE
2	F	9	SER
2	F	10	THR
2	F	159	LEU

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Mol	Chain	Res	Type
3	G	2	LYS
3	G	79	ARG
3	G	165	VAL
4	H	39	ARG
4	H	48	VAL
4	H	106	THR
4	H	143	MET
4	H	157	THR
4	H	163	MET
4	H	185	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	137	HIS
3	C	167	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

12 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
5	NAG	A	301	1	14,14,15	0.45	0	15,19,21	1.01	1 (6%)
5	NAG	C	201	3	14,14,15	0.54	0	15,19,21	1.07	1 (6%)
5	NAG	D	201	4	14,14,15	0.62	0	15,19,21	1.37	2 (13%)
6	GOL	D	202	-	5,5,5	0.34	0	5,5,5	0.23	0
5	NAG	E	301	1	14,14,15	0.48	0	15,19,21	0.91	1 (6%)
6	GOL	F	201	-	5,5,5	0.40	0	5,5,5	0.09	0
5	NAG	G	201	3	14,14,15	0.50	0	15,19,21	0.82	0
5	NAG	G	202	3	14,14,15	0.45	0	15,19,21	1.13	1 (6%)
5	NAG	H	201	4	14,14,15	0.58	0	15,19,21	0.87	0
7	PGE	H	202	-	9,9,9	0.65	0	8,8,8	0.70	0
8	ACT	H	203	-	1,3,3	1.31	0	0,3,3	0.00	-
6	GOL	H	204	-	5,5,5	0.34	0	5,5,5	0.25	0

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
5	NAG	A	301	1	-	0/6/23/26	0/1/1/1
5	NAG	C	201	3	-	0/6/23/26	0/1/1/1
5	NAG	D	201	4	-	0/6/23/26	0/1/1/1
6	GOL	D	202	-	-	0/4/4/4	0/0/0/0
5	NAG	E	301	1	-	0/6/23/26	0/1/1/1
6	GOL	F	201	-	-	0/4/4/4	0/0/0/0
5	NAG	G	201	3	-	0/6/23/26	0/1/1/1
5	NAG	G	202	3	-	0/6/23/26	0/1/1/1
5	NAG	H	201	4	-	0/6/23/26	0/1/1/1
7	PGE	H	202	-	-	0/7/7/7	0/0/0/0
8	ACT	H	203	-	-	0/0/0/0	0/0/0/0
6	GOL	H	204	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	201	NAG	C4-C3-C2	2.22	114.27	111.02
5	E	301	NAG	C1-O5-C5	2.50	115.61	112.17
5	C	201	NAG	C3-C4-C5	2.74	115.05	110.22
5	A	301	NAG	C1-O5-C5	3.05	116.37	112.17
5	D	201	NAG	C3-C4-C5	3.59	116.54	110.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	G	202	NAG	C1-O5-C5	3.73	117.31	112.17

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	301	NAG	1	0
6	F	201	GOL	1	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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	188/188 (100%)	-0.40	1 (0%) 90 85	12, 28, 54, 73	0
1	E	188/188 (100%)	-0.46	1 (0%) 90 85	8, 27, 51, 73	0
2	B	186/191 (97%)	-0.37	0 100 100	13, 30, 54, 71	0
2	F	187/191 (97%)	-0.36	1 (0%) 90 85	13, 29, 56, 89	0
3	C	181/181 (100%)	-0.47	0 100 100	11, 22, 41, 61	0
3	G	181/181 (100%)	-0.52	0 100 100	9, 19, 34, 48	0
4	D	188/189 (99%)	-0.28	1 (0%) 90 85	8, 24, 54, 76	0
4	H	189/189 (100%)	-0.38	0 100 100	11, 26, 52, 65	0
All	All	1488/1498 (99%)	-0.40	4 (0%) 93 92	8, 26, 52, 89	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	191	GLN	3.6
1	A	200	ASP	2.5
1	E	14	GLN	2.5
2	F	146	LYS	2.1

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
5	NAG	A	301	14/15	0.92	0.36	2.09	55,64,69,78	0
6	GOL	H	204	6/6	0.95	0.19	1.19	13,19,34,42	0
8	ACT	H	203	4/4	0.86	0.19	0.19	27,41,41,47	0
5	NAG	C	201	14/15	0.89	0.20	-0.35	29,45,57,61	0
7	PGE	H	202	10/10	0.91	0.17	-0.65	12,16,20,25	0
5	NAG	G	202	14/15	0.90	0.14	-0.70	14,29,46,55	0
6	GOL	D	202	6/6	0.94	0.20	-	21,25,31,47	0
5	NAG	H	201	14/15	0.80	0.39	-	25,34,44,46	0
6	GOL	F	201	6/6	0.81	0.28	-	29,41,48,53	0
5	NAG	G	201	14/15	0.78	0.34	-	63,81,86,86	0
5	NAG	D	201	14/15	0.81	0.31	-	66,85,95,98	0
5	NAG	E	301	14/15	0.93	0.30	-	40,53,61,63	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.