



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 08:08 PM BST

PDB ID : 1KTD  
Title : CRYSTAL STRUCTURE OF CLASS II MHC MOLECULE IEK BOUND TO PIGEON CYTOCHROME C PEPTIDE  
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Deposited on : 2002-01-15  
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](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.

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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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.13.1  
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) : 2.13.1

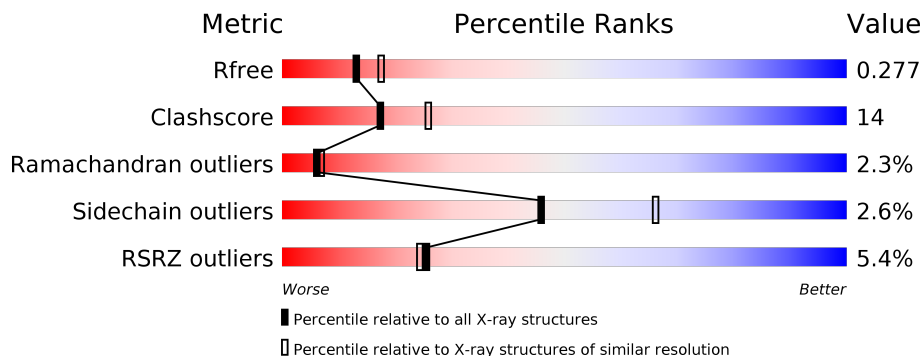
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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))
$R_{free}$	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (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  $\geq 3$ , 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	182	 70% 29%
1	C	182	 70% 29%
2	B	215	 73% 26%
2	D	215	 68% 29%

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6814 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 H-2 class II histocompatibility antigen, E-D alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	182	1497	963	243	287	4	0	0	0
1	C	182	1497	963	243	287	4	0	0	0

- Molecule 2 is a protein called Fusion protein consisting of cytochrome C peptide, glycine rich linker, and MHC E-beta-k.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	215	1721	1088	298	329	6	84	0	0
2	D	215	1721	1088	298	329	6	69	0	0

There are 34 discrepancies between the modelled and reference sequences:

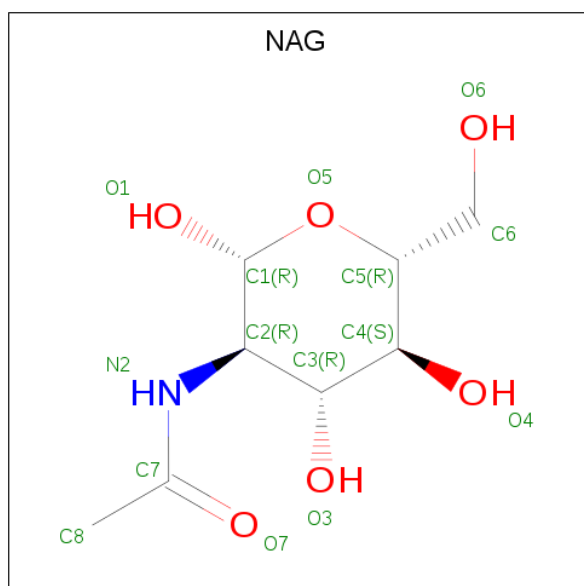
Chain	Residue	Modelled	Actual	Comment	Reference
B	1	ALA	ARG	conflict	UNP P00021
B	12	SER	THR	engineered mutation	UNP P00021
B	15	GLY	-	linker	UNP P00021
B	16	GLY	-	linker	UNP P00021
B	17	GLY	-	linker	UNP P00021
B	18	GLY	-	linker	UNP P00021
B	19	SER	-	linker	UNP P00021
B	20	LEU	-	linker	UNP P00021
B	21	VAL	-	linker	UNP P00021
B	22	GLY	-	linker	UNP P00021
B	23	GLY	-	linker	UNP P00021
B	24	GLY	-	linker	UNP P00021
B	25	SER	-	linker	UNP P00021
B	26	GLY	-	linker	UNP P00021
B	27	GLY	-	linker	UNP P00021

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Chain	Residue	Modelled	Actual	Comment	Reference
B	28	GLY	-	linker	UNP P00021
B	29	GLY	-	linker	UNP P00021
D	1	ALA	ARG	conflict	UNP P00021
D	12	SER	THR	engineered mutation	UNP P00021
D	15	GLY	-	linker	UNP P00021
D	16	GLY	-	linker	UNP P00021
D	17	GLY	-	linker	UNP P00021
D	18	GLY	-	linker	UNP P00021
D	19	SER	-	linker	UNP P00021
D	20	LEU	-	linker	UNP P00021
D	21	VAL	-	linker	UNP P00021
D	22	GLY	-	linker	UNP P00021
D	23	GLY	-	linker	UNP P00021
D	24	GLY	-	linker	UNP P00021
D	25	SER	-	linker	UNP P00021
D	26	GLY	-	linker	UNP P00021
D	27	GLY	-	linker	UNP P00021
D	28	GLY	-	linker	UNP P00021
D	29	GLY	-	linker	UNP P00021

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



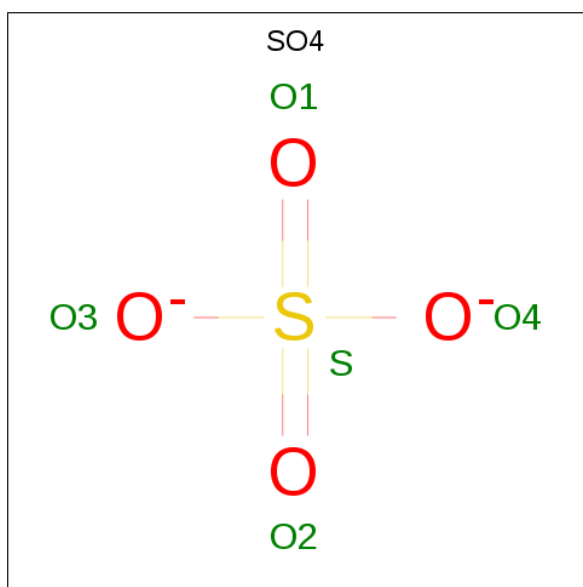
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
			Total	C	N			O
3	A	1	14	8	1	5	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		

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

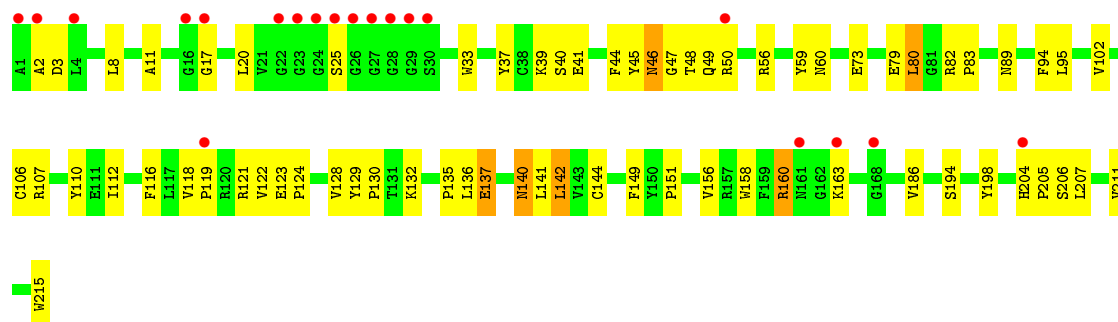


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	96	Total	O	0	0
			96	96		
5	B	99	Total	O	0	0
			99	99		
5	C	71	Total	O	0	0
			71	71		
5	D	51	Total	O	0	0
			51	51		





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	145.52Å 57.30Å 116.57Å 90.00° 94.23° 90.00°	Depositor
Resolution (Å)	10.00 – 2.40 116.25 – 2.40	Depositor EDS
% Data completeness (in resolution range)	96.3 (10.00-2.40) 96.2 (116.25-2.40)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.20 (at 2.40Å)	Xtrriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.220 , 0.279 0.219 , 0.277	Depositor DCC
$R_{free}$ test set	1810 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	44.4	Xtrriage
Anisotropy	0.408	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 59.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6814	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	48.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% 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: NAG, SO4

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	0.43	0/1539	0.65	0/2091
1	C	0.39	0/1539	0.61	0/2091
2	B	0.39	0/1766	0.63	0/2400
2	D	0.36	0/1766	0.61	0/2400
All	All	0.39	0/6610	0.63	0/8982

There are no bond length outliers.

There are no bond angle outliers.

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	1497	0	1420	42	0
1	C	1497	0	1422	42	0
2	B	1721	0	1637	47	0
2	D	1721	0	1638	56	0
3	A	28	0	26	3	0
3	B	14	0	13	0	0
3	C	14	0	13	0	0
4	B	5	0	0	0	0
5	A	96	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	99	0	0	2	0
5	C	71	0	0	0	0
5	D	51	0	0	1	0
All	All	6814	0	6169	171	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 14.

All (171) 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)
2:D:160:ARG:HB3	2:D:160:ARG:HH11	1.30	0.96
1:A:51:PHE:O	2:B:2:ALA:HA	1.76	0.86
2:B:73:GLU:HG3	2:B:75:ARG:HH12	1.41	0.86
1:C:123:ARG:HE	1:C:161:TYR:HE2	1.26	0.83
2:D:140:ASN:HD22	2:D:141:LEU:N	1.88	0.72
2:B:73:GLU:HG3	2:B:75:ARG:NH1	2.05	0.72
2:D:102:VAL:O	2:D:106:CYS:HB2	1.91	0.70
2:B:11:ALA:HB2	2:B:94:PHE:CZ	2.29	0.67
2:B:65:LEU:HD12	2:B:75:ARG:O	1.94	0.67
1:C:8:ILE:HG12	2:D:41:GLU:HG2	1.77	0.67
2:D:156:VAL:HB	2:D:186:VAL:HG21	1.77	0.67
2:B:141:LEU:HD22	2:B:187:MET:HB3	1.76	0.66
2:D:82:ARG:HB2	2:D:83:PRO:HD3	1.78	0.66
2:D:207:LEU:HD13	2:D:211:VAL:HG23	1.78	0.65
1:A:109:ILE:HG22	1:A:112:PHE:CE1	2.33	0.64
2:D:160:ARG:CB	2:D:160:ARG:HH11	2.09	0.64
2:D:11:ALA:HB2	2:D:94:PHE:CZ	2.34	0.63
1:C:55:GLU:OE1	1:C:57:GLN:HB2	1.99	0.62
2:D:140:ASN:HD22	2:D:141:LEU:H	1.46	0.62
1:A:59:ALA:O	1:A:63:ILE:HG12	1.99	0.62
2:B:167:THR:HA	2:D:17:GLY:HA2	1.80	0.62
2:D:50:ARG:HG2	2:D:50:ARG:HH11	1.64	0.61
1:C:147:LYS:NZ	1:C:149:HIS:HE1	1.97	0.61
2:B:143:VAL:HG22	2:B:187:MET:HG2	1.81	0.61
1:C:124:ASN:HA	1:C:160:PHE:CZ	2.36	0.61
3:A:602:NAG:O6	2:B:21:VAL:HG13	2.02	0.60
1:A:3:GLU:HG3	5:B:813:HOH:O	2.01	0.59
2:B:157:ARG:HG2	5:B:847:HOH:O	2.03	0.59
2:B:79:GLU:OE2	2:B:82:ARG:NH2	2.34	0.59
2:D:20:LEU:HD11	2:D:79:GLU:HB3	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:111:LYS:HG2	1:A:140:ARG:CZ	2.33	0.58
2:B:97:GLN:O	2:B:101:GLU:HG3	2.02	0.58
2:B:23:GLY:O	2:B:25:SER:N	2.37	0.58
2:B:124:PRO:HB3	2:B:149:PHE:HB3	1.85	0.57
2:B:73:GLU:HB2	2:B:89:ASN:OD1	2.04	0.57
2:D:3:ASP:HB2	2:D:112:ILE:HG12	1.86	0.57
2:B:21:VAL:HG12	2:B:22:GLY:N	2.19	0.56
2:D:37:TYR:OH	2:D:39:LYS:HD3	2.06	0.56
1:A:167:HIS:CD2	1:A:169:GLY:H	2.23	0.56
1:C:26:PHE:HB2	1:C:31:ILE:HD11	1.86	0.56
2:D:8:LEU:HD22	2:D:40:SER:HB3	1.87	0.56
1:A:117:VAL:HG12	1:A:118:ASN:N	2.21	0.56
1:C:87:PRO:HB3	1:C:112:PHE:HB3	1.88	0.55
1:A:73:MET:CE	2:B:80:LEU:HG	2.36	0.55
1:A:50:LYS:HE2	1:A:51:PHE:CE1	2.41	0.55
1:A:1:ILE:N	1:A:1:ILE:HD13	2.22	0.55
1:A:95:SER:HB2	1:A:96:PRO:HD2	1.89	0.55
2:B:90:SER:O	2:B:92:PRO:HD3	2.06	0.54
1:C:103:ASN:HB3	1:C:153:PHE:CE1	2.42	0.54
2:D:41:GLU:OE2	2:D:56:ARG:HD3	2.07	0.54
1:C:94:ARG:HH11	1:C:94:ARG:HG2	1.73	0.54
1:A:114:PRO:HB3	2:B:33:TRP:CE3	2.42	0.54
1:C:147:LYS:HZ3	1:C:149:HIS:HE1	1.55	0.54
1:C:162:ASP:OD1	1:C:177:THR:HA	2.07	0.53
1:A:87:PRO:HB3	1:A:112:PHE:HB3	1.90	0.53
2:D:124:PRO:HB3	2:D:149:PHE:HB3	1.91	0.53
2:B:10:GLN:HE21	2:B:11:ALA:H	1.56	0.53
2:D:121:ARG:HD3	2:D:151:PRO:HD3	1.91	0.53
1:A:109:ILE:N	1:A:109:ILE:HD12	2.24	0.52
2:D:122:VAL:O	2:D:149:PHE:HA	2.10	0.52
1:C:2:LYS:HA	2:D:46:ASN:OD1	2.10	0.51
2:D:48:THR:HB	2:D:107:ARG:HG2	1.92	0.51
2:D:89:ASN:HA	2:D:95:LEU:HD11	1.93	0.51
1:A:83:ALA:HA	2:B:27:GLY:O	2.11	0.51
1:C:114:PRO:HB3	2:D:33:TRP:CE3	2.46	0.51
1:C:97:VAL:C	1:C:98:ASN:HD22	2.14	0.50
1:C:74:LYS:HE2	1:C:79:ASN:OD1	2.12	0.50
1:C:147:LYS:HZ3	1:C:149:HIS:CE1	2.29	0.50
1:C:126:ARG:HG2	1:C:126:ARG:HH11	1.77	0.50
2:D:163:LYS:NZ	2:D:163:LYS:HB2	2.26	0.50
1:A:123:ARG:HG3	1:A:161:TYR:CE2	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:204:HIS:CG	2:D:205:PRO:HD2	2.47	0.49
1:C:99:LEU:HA	1:C:155:PRO:HB2	1.93	0.49
1:A:89:VAL:HG23	1:A:174:LEU:HD23	1.95	0.49
2:D:59:TYR:CE2	2:D:60:ASN:ND2	2.81	0.49
1:A:82:ASP:HA	2:B:25:SER:HB3	1.95	0.49
2:D:8:LEU:HD13	2:D:40:SER:CB	2.43	0.49
1:C:122:LEU:HA	1:C:126:ARG:O	2.13	0.48
2:D:3:ASP:HB2	2:D:112:ILE:CD1	2.42	0.48
2:B:79:GLU:CD	2:B:82:ARG:HE	2.16	0.48
1:C:167:HIS:CD2	1:C:169:GLY:H	2.31	0.48
1:A:95:SER:HB2	1:A:96:PRO:CD	2.44	0.48
2:D:20:LEU:HD12	2:D:80:LEU:HD13	1.96	0.48
2:D:123:GLU:HA	2:D:206:SER:OG	2.13	0.48
2:D:79:GLU:OE2	2:D:82:ARG:NE	2.38	0.48
1:C:74:LYS:HB3	1:C:79:ASN:OD1	2.13	0.47
2:D:8:LEU:HD13	2:D:40:SER:HB2	1.95	0.47
2:D:144:CYS:HB2	2:D:158:TRP:CZ2	2.49	0.47
1:A:73:MET:HE1	2:B:80:LEU:HG	1.96	0.47
1:A:138:LEU:HB2	1:A:146:ARG:HG3	1.96	0.47
2:B:28:GLY:O	2:B:30:SER:N	2.48	0.47
1:C:8:ILE:HB	1:C:25:ASP:HB3	1.97	0.47
2:D:11:ALA:HB2	2:D:94:PHE:HZ	1.77	0.47
1:A:33:HIS:CG	1:A:136:VAL:HG11	2.49	0.47
2:B:179:ASP:O	2:B:180:TRP:HB2	2.15	0.47
2:B:22:GLY:O	2:B:24:GLY:N	2.48	0.47
1:C:124:ASN:ND2	1:C:159:ASP:OD1	2.38	0.47
2:D:82:ARG:HH11	2:D:82:ARG:HG2	1.80	0.47
2:D:80:LEU:HD22	5:D:247:HOH:O	2.15	0.47
1:C:50:LYS:HE2	1:C:51:PHE:CZ	2.50	0.46
1:C:147:LYS:NZ	1:C:149:HIS:CE1	2.80	0.46
1:C:50:LYS:HZ3	1:C:51:PHE:HE1	1.61	0.46
2:D:50:ARG:HG2	2:D:50:ARG:NH1	2.31	0.46
1:A:84:ASN:HD22	2:B:30:SER:HB2	1.81	0.45
1:C:94:ARG:NH1	1:C:94:ARG:HG2	2.31	0.45
1:A:9:GLN:HB3	2:B:40:SER:HB2	1.98	0.45
2:D:129:TYR:HA	2:D:130:PRO:HD3	1.82	0.45
1:A:11:GLU:OE1	1:A:66:ASP:OD2	2.34	0.45
2:D:128:VAL:HG11	2:D:215:TRP:HB2	1.97	0.45
1:A:74:LYS:HE2	1:A:79:ASN:OD1	2.16	0.45
2:D:118:VAL:HB	2:D:119:PRO:HD3	1.99	0.45
1:A:168:TRP:CE2	3:A:604:NAG:H83	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:94:PHE:CZ	2:B:98:LYS:HD3	2.52	0.45
2:D:3:ASP:HB2	2:D:112:ILE:CG1	2.46	0.45
2:D:47:GLY:C	2:D:49:GLN:H	2.19	0.45
1:C:122:LEU:HD12	1:C:122:LEU:N	2.32	0.45
1:C:143:HIS:H	1:C:143:HIS:CD2	2.33	0.45
2:D:118:VAL:HB	2:D:119:PRO:CD	2.46	0.45
1:C:33:HIS:CD2	1:C:136:VAL:HG11	2.52	0.45
2:D:140:ASN:ND2	2:D:141:LEU:N	2.62	0.45
2:B:21:VAL:HG12	2:B:22:GLY:H	1.81	0.45
2:B:163:LYS:CB	2:B:163:LYS:NZ	2.81	0.44
2:B:168:GLY:O	2:B:189:GLU:HG3	2.17	0.44
1:A:51:PHE:C	2:B:2:ALA:HA	2.36	0.44
1:A:135:THR:HG23	1:A:148:PHE:HB2	2.00	0.44
1:C:109:ILE:HD11	1:C:119:VAL:HG21	1.98	0.44
2:D:44:PHE:CE1	2:D:110:TYR:HB2	2.53	0.44
2:B:191:VAL:O	2:B:191:VAL:HG23	2.17	0.44
1:A:126:ARG:HH11	1:A:126:ARG:HG2	1.82	0.44
2:B:8:LEU:HD22	2:B:40:SER:HB3	1.99	0.43
1:C:11:GLU:OE1	1:C:66:ASP:OD2	2.35	0.43
2:D:142:LEU:HD22	2:D:158:TRP:CH2	2.53	0.43
2:B:101:GLU:HA	2:B:104:THR:OG1	2.18	0.43
1:A:140:ARG:HE	1:A:144:LEU:HB2	1.84	0.43
2:B:128:VAL:HA	2:B:143:VAL:O	2.18	0.43
1:C:110:ASP:OD1	1:C:146:ARG:HB3	2.18	0.43
1:C:137:PHE:CE1	1:C:147:LYS:HD2	2.54	0.42
2:D:204:HIS:ND1	2:D:205:PRO:HD2	2.33	0.42
5:A:690:HOH:O	2:B:16:GLY:HA2	2.18	0.42
2:B:193:GLN:HA	2:B:193:GLN:OE1	2.19	0.42
1:A:168:TRP:CD2	3:A:604:NAG:H83	2.54	0.42
1:A:167:HIS:HD2	1:A:169:GLY:H	1.67	0.42
2:B:8:LEU:HD13	2:B:40:SER:CB	2.50	0.42
1:C:76:ARG:CD	2:D:80:LEU:HD12	2.50	0.41
2:D:73:GLU:HB3	2:D:89:ASN:OD1	2.19	0.41
1:C:157:THR:HG22	1:C:180:PHE:CD2	2.55	0.41
1:C:74:LYS:O	1:C:79:ASN:N	2.45	0.41
1:C:69:ASN:O	1:C:73:MET:HG2	2.20	0.41
2:D:3:ASP:HB2	2:D:112:ILE:HD11	2.02	0.41
1:A:1:ILE:H3	1:A:1:ILE:HD13	1.85	0.41
2:D:79:GLU:CD	2:D:82:ARG:HE	2.22	0.41
2:D:160:ARG:HG3	2:D:198:TYR:CE2	2.56	0.41
1:A:16:PRO:O	1:A:18:LYS:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:THR:CG2	1:A:148:PHE:HB2	2.50	0.41
2:B:129:TYR:HA	2:B:130:PRO:HD3	1.86	0.41
2:B:165:GLU:OE1	2:B:165:GLU:HA	2.21	0.41
1:C:114:PRO:O	1:C:167:HIS:HE1	2.04	0.41
1:C:31:ILE:HG22	1:C:32:PHE:HD1	1.85	0.41
1:A:134:GLU:OE2	1:A:147:LYS:NZ	2.50	0.41
2:B:188:LEU:HG	2:B:190:THR:HG23	2.03	0.41
2:B:64:ASN:O	2:B:81:GLY:HA3	2.21	0.41
1:C:44:ARG:HD2	1:C:44:ARG:HA	1.91	0.41
2:D:45:TYR:O	2:D:46:ASN:HB2	2.21	0.41
1:A:12:PHE:O	1:A:20:GLY:HA2	2.21	0.40
1:A:91:VAL:HA	1:A:106:ILE:O	2.20	0.40
1:C:76:ARG:HD2	2:D:80:LEU:HD12	2.03	0.40
1:A:74:LYS:HB3	1:A:74:LYS:HE2	1.86	0.40
1:A:97:VAL:HG13	1:A:155:PRO:CG	2.51	0.40
1:A:9:GLN:OE1	2:B:8:LEU:HD12	2.21	0.40
2:B:163:LYS:HB3	2:B:163:LYS:NZ	2.37	0.40
2:D:48:THR:HG23	2:D:110:TYR:CD2	2.57	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	180/182 (99%)	173 (96%)	5 (3%)	2 (1%)	14	20
1	C	180/182 (99%)	169 (94%)	10 (6%)	1 (1%)	25	36
2	B	213/215 (99%)	188 (88%)	19 (9%)	6 (3%)	5	4
2	D	213/215 (99%)	186 (87%)	18 (8%)	9 (4%)	3	2
All	All	786/794 (99%)	716 (91%)	52 (7%)	18 (2%)	6	7

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	2	ALA
2	D	46	ASN
2	D	135	PRO
2	D	136	LEU
2	D	194	SER
2	B	23	GLY
2	B	24	GLY
2	B	29	GLY
2	B	46	ASN
1	A	142	ASP
2	B	194	SER
2	D	132	LYS
1	A	81	PRO
1	C	157	THR
2	D	25	SER
2	D	116	PHE
2	D	137	GLU
2	B	135	PRO

### 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	165/165 (100%)	161 (98%)	4 (2%)	49	68
1	C	165/165 (100%)	161 (98%)	4 (2%)	49	68
2	B	187/187 (100%)	182 (97%)	5 (3%)	44	65
2	D	187/187 (100%)	182 (97%)	5 (3%)	44	65
All	All	704/704 (100%)	686 (97%)	18 (3%)	46	66

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	ILE
1	A	81	PRO
1	A	116	VAL
1	A	181	GLU

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Mol	Chain	Res	Type
2	B	3	ASP
2	B	4	LEU
2	B	50	ARG
2	B	140	ASN
2	B	141	LEU
1	C	24	PHE
1	C	129	THR
1	C	153	PHE
1	C	181	GLU
2	D	80	LEU
2	D	137	GLU
2	D	140	ASN
2	D	142	LEU
2	D	160	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	149	HIS
1	A	167	HIS
2	B	10	GLN
1	C	98	ASN
1	C	143	HIS
1	C	149	HIS
1	C	167	HIS
2	D	10	GLN
2	D	140	ASN
2	D	161	ASN

### 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

5 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
3	NAG	B	601	2	14,14,15	0.56	0	17,19,21	0.74	0
3	NAG	A	602	1	14,14,15	0.55	0	17,19,21	0.71	0
3	NAG	C	603	1	14,14,15	0.61	0	17,19,21	0.68	0
3	NAG	A	604	1	14,14,15	0.49	0	17,19,21	0.75	0
4	SO4	B	800	-	4,4,4	0.19	0	6,6,6	0.11	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
3	NAG	B	601	2	-	3/6/23/26	0/1/1/1
3	NAG	A	602	1	-	6/6/23/26	0/1/1/1
3	NAG	C	603	1	-	2/6/23/26	0/1/1/1
3	NAG	A	604	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	601	NAG	C3-C2-N2-C7
3	B	601	NAG	C8-C7-N2-C2
3	B	601	NAG	O7-C7-N2-C2
3	A	602	NAG	C8-C7-N2-C2
3	A	602	NAG	O7-C7-N2-C2
3	C	603	NAG	C8-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
3	C	603	NAG	O7-C7-N2-C2
3	A	604	NAG	O5-C5-C6-O6
3	A	604	NAG	C4-C5-C6-O6
3	A	602	NAG	C4-C5-C6-O6
3	A	602	NAG	C1-C2-N2-C7
3	A	602	NAG	O5-C5-C6-O6
3	A	602	NAG	C3-C2-N2-C7

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	602	NAG	1	0
3	A	604	NAG	2	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	182/182 (100%)	0.07	1 (0%) 91 89	26, 38, 59, 89	0
1	C	182/182 (100%)	0.40	7 (3%) 40 39	29, 47, 85, 125	0
2	B	205/215 (95%)	0.61	14 (6%) 17 15	29, 45, 85, 109	0
2	D	207/215 (96%)	0.71	20 (9%) 7 7	32, 53, 89, 115	0
All	All	776/794 (97%)	0.46	42 (5%) 25 24	26, 45, 83, 125	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	24	GLY	9.1
2	B	2	ALA	8.8
2	B	16	GLY	8.5
2	B	25	SER	8.3
2	D	28	GLY	8.2
1	C	1	ILE	6.8
1	C	182	GLU	6.6
2	D	26	GLY	6.1
2	D	1	ALA	6.0
2	D	23	GLY	6.0
2	D	24	GLY	5.1
2	B	1	ALA	5.0
2	B	27	GLY	5.0
2	D	163	LYS	4.6
2	D	27	GLY	4.5
2	D	29	GLY	4.1
2	B	28	GLY	4.0
1	C	2	LYS	3.9
1	A	181	GLU	3.7
2	D	2	ALA	3.7
2	B	22	GLY	3.6

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Mol	Chain	Res	Type	RSRZ
1	C	160	PHE	3.5
1	C	158	ASP	3.4
2	D	4	LEU	3.4
2	D	25	SER	3.2
2	D	161	ASN	3.0
2	B	163	LYS	2.9
2	B	192	PRO	2.7
2	D	16	GLY	2.7
2	D	30	SER	2.6
2	D	119	PRO	2.6
2	D	204	HIS	2.4
1	C	3	GLU	2.4
2	B	195	GLY	2.3
2	B	23	GLY	2.2
2	B	167	THR	2.2
2	D	22	GLY	2.2
1	C	181	GLU	2.2
2	D	50	ARG	2.2
2	D	17	GLY	2.1
2	D	168	GLY	2.1
2	B	26	GLY	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
3	NAG	B	601	14/15	0.67	0.34	67,70,72,73	0
3	NAG	A	602	14/15	0.68	0.36	82,86,91,92	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	NAG	C	603	14/15	0.76	0.28	74,76,81,81	0
3	NAG	A	604	14/15	0.90	0.19	53,56,59,60	0
4	SO4	B	800	5/5	0.98	0.21	49,50,51,52	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.