



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:56 AM GMT

PDB ID : 3QEH  
Title : Crystal structure of human N12-i15, an ADCC and non-neutralizing anti-HIV-1 Env antibody  
Authors : Guan, Y.; DeVico, A.L.; Lewis, G.K.; Pazgier, M.  
Deposited on : 2011-01-20  
Resolution : 2.59 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.

---

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) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : trunk26865

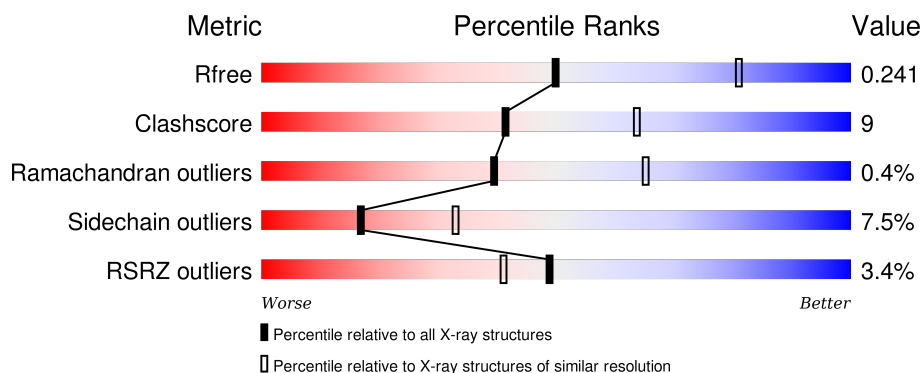
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.59 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))
$R_{free}$	91344	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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	232	<div> <div>75%</div> <div>14%</div> <div>7%</div> </div>
1	C	232	<div> <div>74%</div> <div>14%</div> <div>8%</div> </div>
1	E	232	<div> <div>3%</div> <div>74%</div> <div>13%</div> <div>8%</div> </div>
1	G	232	<div> <div>9%</div> <div>66%</div> <div>16%</div> <div>17%</div> </div>
2	B	218	<div> <div>75%</div> <div>21%</div> </div>

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Length	Quality of chain
2	D	218	<div> <div>%</div> <div> </div> <div>81%15% . .</div> </div>
2	F	218	<div> <div>4%</div> <div> </div> <div>80%16% . .</div> </div>
2	H	218	<div> <div>5%</div> <div> </div> <div>62%22%5%11%</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
4	CL	B	215	-	-	-	X
5	GOL	D	214	-	-	-	X
5	GOL	E	222	-	-	-	X
5	GOL	F	214	-	-	-	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 13446 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 Fab fragment of human anti-HIV-1 Env antibody N12-i15, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	216	Total	C	N	O	S	0	0	0
			1649	1043	281	318	7			
1	C	213	Total	C	N	O	S	0	0	0
			1630	1031	278	314	7			
1	E	214	Total	C	N	O	S	0	0	0
			1638	1037	279	315	7			
1	G	192	Total	C	N	O	S	0	0	0
			1475	939	250	279	7			

- Molecule 2 is a protein called Fab fragment of human anti-HIV-1 Env antibody N12-i15, light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	215	Total	C	N	O	S	0	0	0
			1655	1038	280	331	6			
2	D	215	Total	C	N	O	S	0	0	0
			1655	1038	280	331	6			
2	F	212	Total	C	N	O	S	0	0	0
			1631	1023	276	326	6			
2	H	193	Total	C	N	O	S	0	0	0
			1498	941	254	297	6			

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	G	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	G	1	Total	Cl	0	0
			1	1		
4	D	1	Total	Cl	0	0
			1	1		
4	H	1	Total	Cl	0	0
			1	1		
4	B	3	Total	Cl	0	0
			3	3		
4	C	2	Total	Cl	0	0
			2	2		
4	A	1	Total	Cl	0	0
			1	1		
4	F	1	Total	Cl	0	0
			1	1		

- Molecule 5 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
5	D	1	Total	C	O	0	0
			6	3	3		
5	E	1	Total	C	O	0	0
			6	3	3		
5	F	1	Total	C	O	0	0
			6	3	3		
5	H	1	Total	C	O	0	0
			6	3	3		

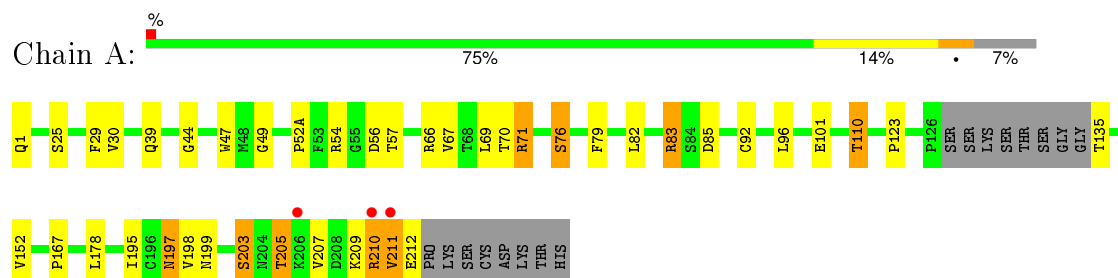
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	104	Total	O	0	0
			104	104		
6	B	78	Total	O	0	0
			78	78		
6	C	75	Total	O	0	0
			75	75		
6	D	82	Total	O	0	0
			82	82		
6	E	64	Total	O	0	0
			64	64		
6	F	63	Total	O	0	0
			63	63		
6	G	63	Total	O	0	0
			63	63		
6	H	37	Total	O	0	0
			37	37		

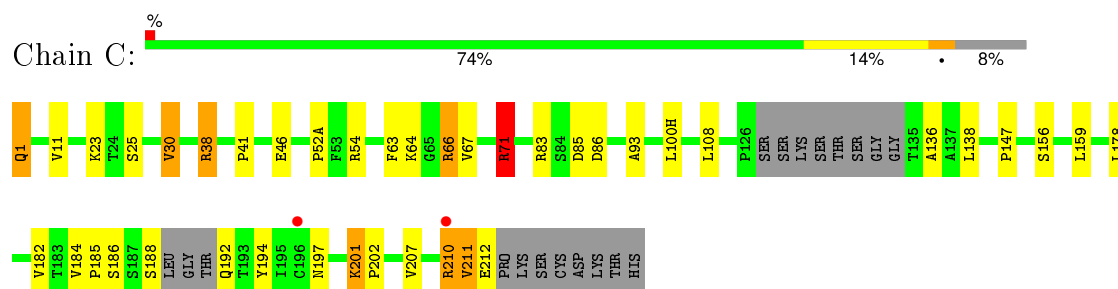
### 3 Residue-property plots

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 ( $\text{RSRZ} > 2$ ). 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.

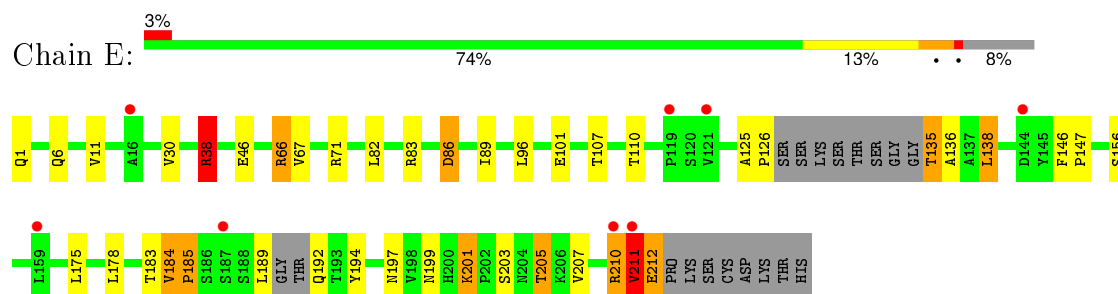
- Molecule 1: Fab fragment of human anti-HIV-1 Env antibody N12-i15, heavy chain



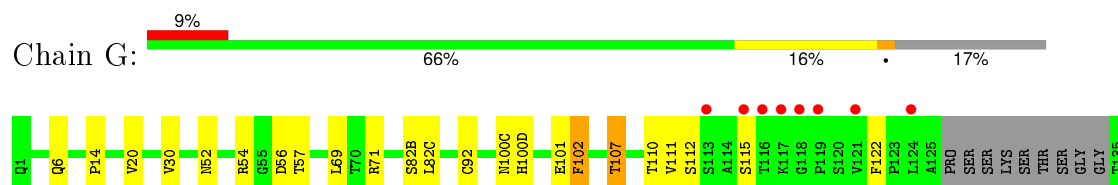
- Molecule 1: Fab fragment of human anti-HIV-1 Env antibody N12-i15, heavy chain

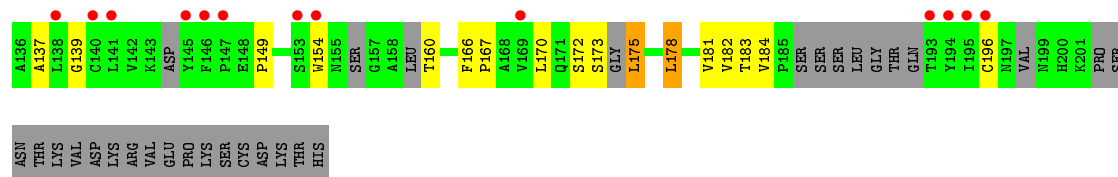


- Molecule 1: Fab fragment of human anti-HIV-1 Env antibody N12-i15, heavy chain



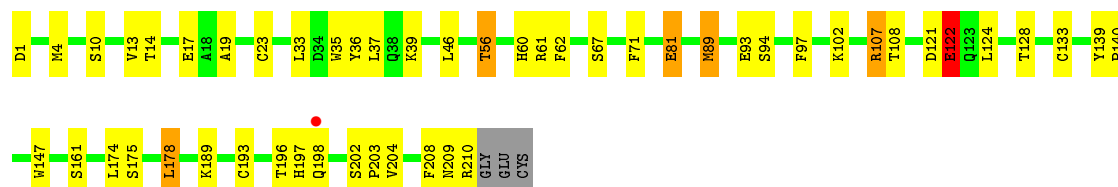
- Molecule 1: Fab fragment of human anti-HIV-1 Env antibody N12-i15, heavy chain





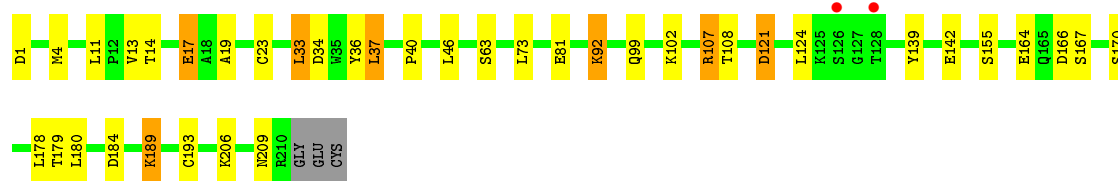
- Molecule 2: Fab fragment of human anti-HIV-1 Env antibody N12-i15,light chain

Chain B: 75% 21%



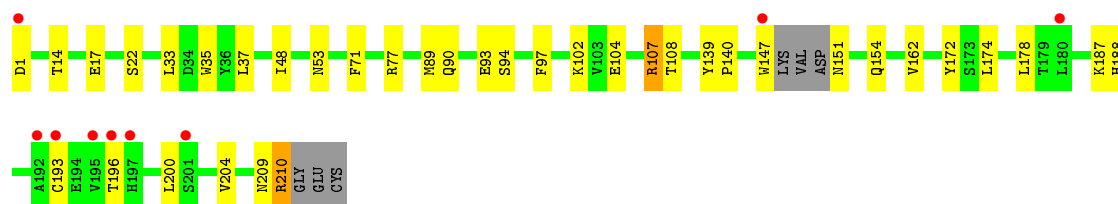
- Molecule 2: Fab fragment of human anti-HIV-1 Env antibody N12-i15,light chain

Chain D: 81% 15%



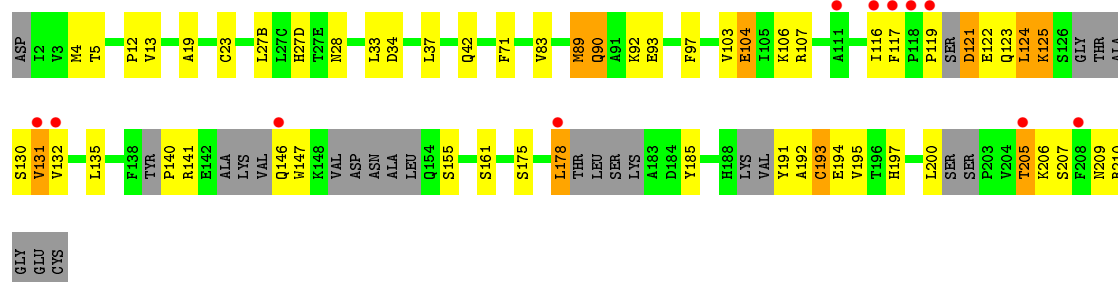
- Molecule 2: Fab fragment of human anti-HIV-1 Env antibody N12-i15,light chain

Chain F: 80% 16%



- Molecule 2: Fab fragment of human anti-HIV-1 Env antibody N12-i15,light chain

Chain H: 62% 22% 5% 11%





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	125.12Å 125.12Å 149.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.59 40.18 – 2.59	Depositor EDS
% Data completeness (in resolution range)	99.8 (20.00-2.59) 99.7 (40.18-2.59)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.39 (at 2.58Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.216 , 0.254 0.198 , 0.241	Depositor DCC
$R_{free}$ test set	3613 reflections (5.34%)	DCC
Wilson B-factor (Å <sup>2</sup> )	35.2	Xtriage
Anisotropy	0.505	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 40.0	EDS
Estimated twinning fraction	0.941 for H, K, L 0.059 for K, H, -L 0.034 for h,-k,-l	Xtriage
Reported twinning fraction	0.941 for H, K, L 0.059 for K, H, -L	Depositor
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 71531 reflections (0.001%)	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	13446	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 36.93 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 4.6355e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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, 0.375 respectively for untwinned datasets, and 0.333, 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, SO4, 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	0.95	1/1688 (0.1%)	0.87	1/2299 (0.0%)
1	C	0.94	0/1668	0.89	3/2270 (0.1%)
1	E	0.87	0/1676	0.87	5/2281 (0.2%)
1	G	0.85	2/1506 (0.1%)	0.82	0/2042
2	B	0.92	4/1691 (0.2%)	0.88	2/2295 (0.1%)
2	D	0.93	1/1691 (0.1%)	0.88	4/2295 (0.2%)
2	F	0.88	3/1666 (0.2%)	0.80	0/2260
2	H	0.82	0/1525	0.83	0/2055
All	All	0.90	11/13111 (0.1%)	0.86	15/17797 (0.1%)

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	1
1	E	0	1
All	All	0	2

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	92	CYS	CB-SG	-6.42	1.71	1.82
2	B	122	GLU	CB-CG	6.11	1.63	1.52
1	G	92	CYS	CB-SG	-6.04	1.72	1.82
2	F	71	PHE	CE1-CZ	5.79	1.48	1.37
2	B	193	CYS	CB-SG	-5.66	1.72	1.81
2	D	81	GLU	CG-CD	5.58	1.60	1.51
2	B	81	GLU	CG-CD	5.41	1.60	1.51
2	B	35	TRP	CB-CG	-5.26	1.40	1.50

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	35	TRP	CB-CG	-5.23	1.40	1.50
2	F	193	CYS	CB-SG	-5.07	1.73	1.81
1	G	102	PHE	CE2-CZ	5.03	1.47	1.37

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	ARG	NE-CZ-NH2	-7.05	116.78	120.30
1	C	71	ARG	NE-CZ-NH2	-6.17	117.22	120.30
2	D	33	LEU	CA-CB-CG	-5.92	101.68	115.30
2	D	37	LEU	CB-CG-CD2	5.89	121.01	111.00
1	E	38	ARG	NE-CZ-NH2	-5.86	117.37	120.30
1	E	66	ARG	NE-CZ-NH2	-5.84	117.38	120.30
2	B	193	CYS	CA-CB-SG	-5.80	103.55	114.00
2	D	193	CYS	CA-CB-SG	-5.76	103.63	114.00
1	E	38	ARG	NE-CZ-NH1	5.55	123.07	120.30
2	D	178	LEU	CA-CB-CG	5.54	128.04	115.30
1	C	66	ARG	NE-CZ-NH2	-5.43	117.58	120.30
1	C	71	ARG	NE-CZ-NH1	5.39	123.00	120.30
1	E	178	LEU	CA-CB-CG	5.38	127.68	115.30
1	E	86	ASP	CB-CG-OD1	5.21	122.99	118.30
2	B	178	LEU	CA-CB-CG	5.07	126.95	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	203	SER	Peptide
1	E	211	VAL	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	1649	0	1615	34	0
1	C	1630	0	1593	27	0
1	E	1638	0	1604	26	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	1475	0	1434	27	0
2	B	1655	0	1620	30	0
2	D	1655	0	1620	18	0
2	F	1631	0	1593	19	0
2	H	1498	0	1448	40	0
3	A	5	0	0	0	0
3	E	5	0	0	0	0
3	G	5	0	0	1	0
4	A	1	0	0	0	0
4	B	3	0	0	0	0
4	C	2	0	0	0	0
4	D	1	0	0	0	0
4	F	1	0	0	0	0
4	G	1	0	0	0	1
4	H	1	0	0	0	0
5	D	6	0	8	0	0
5	E	6	0	8	0	0
5	F	6	0	8	0	0
5	H	6	0	8	0	0
6	A	104	0	0	6	0
6	B	78	0	0	1	0
6	C	75	0	0	3	0
6	D	82	0	0	2	0
6	E	64	0	0	1	2
6	F	63	0	0	0	3
6	G	63	0	0	0	0
6	H	37	0	0	1	0
All	All	13446	0	12559	217	3

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 (217) 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:E:210:ARG:O	1:E:210:ARG:HD3	1.41	1.20
2:H:141:ARG:HG2	2:H:141:ARG:O	1.50	1.08
1:E:199:ASN:OD1	1:E:205:THR:HG22	1.66	0.95
1:G:166:PHE:O	1:G:178:LEU:HD21	1.67	0.95
1:G:170:LEU:HD11	1:G:175:LEU:HD13	1.49	0.94
1:C:138:LEU:HD13	1:C:210:ARG:HD2	1.52	0.90

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:ARG:HA	1:A:210:ARG:NE	1.88	0.88
1:E:210:ARG:CD	1:E:210:ARG:O	2.25	0.83
1:A:123:PRO:HB3	1:A:210:ARG:HD3	1.61	0.81
1:A:123:PRO:CB	1:A:210:ARG:HD3	2.12	0.79
2:H:141:ARG:CG	2:H:141:ARG:O	2.30	0.79
2:D:11:LEU:HD21	2:D:19:ALA:HB1	1.65	0.78
1:A:199:ASN:OD1	1:A:205:THR:HG22	1.83	0.78
2:B:189:LYS:HZ2	2:B:210:ARG:HA	1.49	0.77
2:D:107:ARG:HD3	2:D:108:THR:O	1.86	0.75
1:G:173:SER:C	1:G:175:LEU:HD23	2.07	0.75
1:C:1:GLN:N	1:C:1:GLN:CD	2.42	0.73
2:H:147:TRP:CD2	2:H:178:LEU:HD11	2.24	0.73
2:D:99:GLN:HG2	6:D:419:HOH:O	1.87	0.73
1:E:6:GLN:HE21	1:E:107:THR:HG23	1.54	0.72
2:F:151:ASN:OD1	2:F:151:ASN:O	2.08	0.72
2:B:189:LYS:NZ	2:B:210:ARG:HA	2.04	0.72
1:A:83:ARG:HG3	1:A:85:ASP:OD1	1.90	0.71
1:E:38:ARG:HD2	1:E:46:GLU:OE1	1.91	0.71
2:H:178:LEU:H	2:H:178:LEU:HD12	1.53	0.71
1:C:83:ARG:HG3	1:C:85:ASP:OD1	1.91	0.70
1:A:110:THR:CG2	6:A:245:HOH:O	2.40	0.68
1:E:6:GLN:NE2	1:E:107:THR:HG23	2.08	0.68
1:C:52(A):PRO:O	1:C:71:ARG:HD2	1.95	0.67
2:H:140:PRO:HD2	2:H:197:HIS:CE1	2.29	0.67
2:D:189:LYS:HD2	2:D:209:ASN:HB3	1.77	0.67
1:A:211:VAL:O	1:A:212:GLU:HB2	1.95	0.66
2:B:89:MET:HE2	2:B:97:PHE:CZ	2.30	0.66
1:A:167:PRO:HG2	2:B:161:SER:HB2	1.77	0.66
2:B:107:ARG:HD3	2:B:108:THR:O	1.95	0.66
1:A:25:SER:HB3	1:C:25:SER:HB3	1.77	0.66
2:D:166:ASP:O	2:D:170:SER:HA	1.96	0.65
1:A:70:THR:HG22	1:A:79:PHE:HB2	1.79	0.64
1:A:210:ARG:HA	1:A:210:ARG:CZ	2.28	0.64
1:E:96:LEU:HD12	1:E:101:GLU:OE1	1.98	0.64
1:A:123:PRO:HB3	1:A:210:ARG:CD	2.27	0.63
1:C:1:GLN:CD	1:C:1:GLN:H1	2.01	0.62
1:C:66:ARG:NH2	1:C:86:ASP:OD2	2.34	0.61
2:B:89:MET:CE	2:B:97:PHE:CZ	2.84	0.61
2:F:107:ARG:CD	2:F:108:THR:O	2.49	0.61
1:A:39:GLN:HG3	1:A:44:GLY:O	2.01	0.60
1:E:210:ARG:O	1:E:210:ARG:NH1	2.30	0.60

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:209:LYS:O	1:A:210:ARG:HD2	2.02	0.59
1:C:1:GLN:N	1:C:1:GLN:OE1	2.34	0.59
2:F:107:ARG:HD3	2:F:108:THR:O	2.01	0.59
2:H:178:LEU:N	2:H:178:LEU:HD12	2.18	0.59
1:A:210:ARG:CA	1:A:210:ARG:NE	2.65	0.58
2:H:147:TRP:CE2	2:H:178:LEU:HD11	2.39	0.58
2:B:14:THR:HB	2:B:17:GLU:HG3	1.84	0.58
2:F:89:MET:HE2	2:F:97:PHE:CZ	2.39	0.58
2:F:209:ASN:O	2:F:210:ARG:NE	2.29	0.58
2:F:147:TRP:HB2	2:F:154:GLN:HB2	1.86	0.57
2:H:27(B):LEU:HD22	2:H:90:GLN:HG2	1.86	0.57
2:B:121:ASP:HB2	2:B:122:GLU:OE1	2.05	0.57
2:H:12:PRO:HB2	2:H:106:LYS:HB2	1.87	0.57
1:E:194:TYR:HB2	1:E:210:ARG:HD2	1.87	0.57
2:B:33:LEU:HD22	2:B:71:PHE:CG	2.40	0.57
2:H:33:LEU:HD22	2:H:71:PHE:CG	2.40	0.57
2:H:119:PRO:HB3	2:H:130:SER:N	2.19	0.56
1:G:173:SER:C	1:G:175:LEU:CD2	2.72	0.56
2:H:4:MET:SD	2:H:90:GLN:HB2	2.46	0.56
1:C:211:VAL:O	1:C:212:GLU:HB2	2.05	0.56
1:G:54:ARG:HB2	1:G:56:ASP:OD1	2.06	0.56
1:C:136:ALA:HB2	1:C:186:SER:HB3	1.88	0.56
2:H:209:ASN:O	2:H:210:ARG:HB2	2.07	0.55
2:B:174:LEU:C	2:B:174:LEU:HD23	2.26	0.55
2:F:14:THR:HB	2:F:17:GLU:OE1	2.06	0.55
2:D:33:LEU:HG	2:D:34:ASP:N	2.22	0.55
2:H:119:PRO:O	2:H:121:ASP:N	2.39	0.55
2:H:124:LEU:HD22	2:H:125:LYS:HB3	1.88	0.54
1:A:70:THR:CG2	1:A:79:PHE:HB2	2.38	0.54
1:C:194:TYR:HB2	1:C:210:ARG:HH11	1.73	0.54
1:C:210:ARG:NH2	1:C:212:GLU:O	2.41	0.53
1:G:167:PRO:HG2	2:H:161:SER:HB2	1.91	0.53
2:D:13:VAL:HG21	2:D:19:ALA:HB2	1.90	0.53
1:G:6:GLN:HE21	1:G:107:THR:CG2	2.21	0.53
1:A:71:ARG:O	1:A:71:ARG:HG2	2.08	0.53
1:C:210:ARG:CZ	1:C:210:ARG:O	2.56	0.53
2:H:140:PRO:HD2	2:H:197:HIS:HE1	1.71	0.53
2:H:146:GLN:HB2	2:H:194:GLU:HB3	1.90	0.52
2:H:12:PRO:HB3	2:H:104:GLU:OE2	2.10	0.52
1:G:139:GLY:HA2	1:G:154:TRP:CH2	2.45	0.52
2:B:93:GLU:O	2:B:94:SER:C	2.48	0.52

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:185:TYR:HA	2:H:191:TYR:OH	2.09	0.52
1:A:96:LEU:HD12	1:A:101:GLU:OE1	2.10	0.51
1:G:6:GLN:HE21	1:G:107:THR:HG23	1.75	0.51
2:B:202:SER:HB2	2:B:203:PRO:HD2	1.92	0.51
1:E:210:ARG:C	1:E:210:ARG:HD3	2.23	0.51
2:H:193:CYS:N	2:H:206:LYS:O	2.37	0.51
1:A:57:THR:HG21	1:A:69:LEU:HB2	1.92	0.51
1:C:38:ARG:NH1	1:C:46:GLU:OE1	2.44	0.51
1:E:184:VAL:HG22	1:E:185:PRO:HD2	1.93	0.51
2:D:36:TYR:CE2	2:D:46:LEU:HD13	2.46	0.51
1:A:67:VAL:HG22	1:A:82:LEU:HD13	1.93	0.50
1:E:11:VAL:HG21	1:E:147:PRO:HG3	1.94	0.50
1:A:152:VAL:HG22	1:A:198:VAL:HG22	1.94	0.50
1:E:66:ARG:NH2	1:E:86:ASP:OD2	2.40	0.50
2:D:11:LEU:CD2	2:D:19:ALA:HB1	2.39	0.50
2:B:36:TYR:CE2	2:B:46:LEU:HD13	2.47	0.50
2:F:93:GLU:O	2:F:94:SER:C	2.50	0.50
2:F:89:MET:HE3	2:F:97:PHE:CE2	2.47	0.49
1:A:210:ARG:CZ	1:A:211:VAL:H	2.25	0.49
2:H:192:ALA:HA	2:H:207:SER:HB3	1.94	0.49
1:G:82(C):LEU:HD13	1:G:111:VAL:HG22	1.93	0.49
2:H:83:VAL:HG13	2:H:103:VAL:O	2.13	0.49
1:G:52:ASN:ND2	3:G:221:SO4:O2	2.44	0.49
2:B:39:LYS:NZ	2:B:81:GLU:OE1	2.35	0.49
1:C:41:PRO:HA	6:C:371:HOH:O	2.13	0.49
2:D:92:LYS:NZ	6:D:348:HOH:O	2.45	0.49
1:A:110:THR:HG23	6:A:245:HOH:O	2.07	0.49
1:C:93:ALA:HB1	1:C:100(H):LEU:HB3	1.95	0.48
2:H:124:LEU:HA	2:H:125:LYS:HA	1.49	0.48
2:B:210:ARG:HB2	6:B:392:HOH:O	2.14	0.48
1:A:52(A):PRO:O	1:A:71:ARG:HD2	2.12	0.48
1:E:197:ASN:HD22	1:E:207:VAL:HG22	1.79	0.48
2:B:67:SER:HA	2:B:71:PHE:CE1	2.49	0.48
2:F:107:ARG:HD2	2:F:108:THR:O	2.14	0.47
1:C:63:PHE:O	1:C:67:VAL:HG12	2.13	0.47
1:C:11:VAL:HG21	1:C:147:PRO:HG3	1.96	0.47
1:E:6:GLN:HB3	1:E:107:THR:CG2	2.44	0.47
2:H:197:HIS:H	2:H:200:LEU:HD12	1.80	0.47
1:C:184:VAL:HG11	1:C:194:TYR:OH	2.14	0.47
2:D:107:ARG:HD2	2:D:139:TYR:HB3	1.97	0.47
1:C:66:ARG:HH22	1:C:86:ASP:CG	2.17	0.47

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:184:VAL:HG12	1:C:185:PRO:O	2.15	0.47
1:G:137:ALA:HB2	1:G:183:THR:HG22	1.97	0.47
1:G:139:GLY:HA2	1:G:154:TRP:HH2	1.81	0.46
6:A:223:HOH:O	2:B:56:THR:HG23	2.15	0.46
2:F:104:GLU:OE2	2:F:172:TYR:OH	2.25	0.46
2:H:121:ASP:OD2	2:H:122:GLU:HG2	2.15	0.46
1:G:182:VAL:HG22	1:G:183:THR:N	2.31	0.46
1:G:139:GLY:HA3	1:G:181:VAL:HG12	1.98	0.46
1:G:122:PHE:CE1	2:H:123:GLN:HG3	2.51	0.46
2:D:121:ASP:HA	2:D:124:LEU:HD12	1.97	0.45
1:A:39:GLN:OE1	6:A:355:HOH:O	2.20	0.45
2:F:139:TYR:CG	2:F:140:PRO:HA	2.52	0.45
1:A:110:THR:HG22	6:A:245:HOH:O	2.12	0.45
1:C:23:LYS:NZ	6:C:511:HOH:O	2.49	0.45
1:G:14:PRO:HD3	1:G:112:SER:C	2.37	0.45
1:A:197:ASN:HD22	1:A:207:VAL:HG13	1.82	0.45
2:H:93:GLU:HA	6:H:435:HOH:O	2.16	0.44
1:G:100(C):ASN:OD1	1:G:100(D):HIS:N	2.50	0.44
2:B:13:VAL:HG21	2:B:19:ALA:HB2	2.00	0.44
1:E:96:LEU:HD12	1:E:101:GLU:CD	2.38	0.44
2:H:131:VAL:HB	2:H:178:LEU:HD13	2.00	0.44
1:G:57:THR:HG21	1:G:69:LEU:HB2	1.99	0.44
2:B:189:LYS:NZ	2:B:210:ARG:HG3	2.33	0.44
1:E:210:ARG:NH1	1:E:211:VAL:HA	2.33	0.44
1:A:71:ARG:NH2	6:A:320:HOH:O	2.50	0.44
1:G:82(C):LEU:HD13	1:G:111:VAL:CG2	2.47	0.44
2:B:208:PHE:CD2	2:B:208:PHE:C	2.91	0.44
2:B:124:LEU:HA	2:B:124:LEU:HD23	1.85	0.44
2:H:89:MET:CE	2:H:97:PHE:CZ	3.01	0.44
1:A:209:LYS:O	1:A:210:ARG:NH1	2.51	0.44
2:H:89:MET:HE3	2:H:97:PHE:CZ	2.52	0.44
1:G:183:THR:O	1:G:184:VAL:HG23	2.18	0.44
2:F:162:VAL:HG22	2:F:174:LEU:HD12	2.00	0.44
2:D:40:PRO:HB3	2:D:164:GLU:HG3	2.00	0.43
2:H:13:VAL:HG21	2:H:19:ALA:HB2	2.00	0.43
1:E:126:PRO:HD3	1:E:138:LEU:HB3	1.99	0.43
1:C:100(H):LEU:HD23	1:C:100(H):LEU:N	2.32	0.43
1:E:201:LYS:HG3	1:E:201:LYS:H	1.66	0.43
2:F:89:MET:CE	2:F:97:PHE:CZ	3.02	0.43
2:H:135:LEU:HD11	2:H:195:VAL:HG21	2.01	0.43
2:B:133:CYS:HB2	2:B:147:TRP:CH2	2.54	0.43

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:146:PHE:HB2	1:E:175:LEU:HD23	2.00	0.42
1:E:135:THR:N	6:E:285:HOH:O	2.51	0.42
2:D:63:SER:O	2:D:73:LEU:HD12	2.18	0.42
2:H:27(D):HIS:HB3	2:H:28:ASN:OD1	2.18	0.42
2:B:89:MET:HE3	2:B:97:PHE:CZ	2.54	0.42
2:B:4:MET:HE3	2:B:23:CYS:SG	2.59	0.42
1:G:6:GLN:NE2	1:G:107:THR:HG23	2.33	0.42
2:B:197:HIS:CG	2:B:198:GLN:H	2.37	0.42
2:H:33:LEU:HG	2:H:34:ASP:N	2.35	0.42
1:G:20:VAL:CG1	1:G:107:THR:HG21	2.50	0.42
2:H:193:CYS:O	2:H:205:THR:HA	2.20	0.42
2:B:209:ASN:O	2:B:210:ARG:CB	2.67	0.42
2:F:187:LYS:HD2	2:F:188:HIS:CE1	2.54	0.42
1:A:96:LEU:HD12	1:A:101:GLU:CD	2.40	0.42
1:A:29:PHE:CG	1:A:76:SER:CB	3.02	0.42
2:B:60:HIS:O	2:B:62:PHE:N	2.53	0.42
1:C:30:VAL:HB	6:C:240:HOH:O	2.20	0.42
2:H:117:PHE:HD2	2:H:132:VAL:HG12	1.85	0.41
2:D:4:MET:CE	2:D:23:CYS:SG	3.08	0.41
2:H:192:ALA:HB1	2:H:205:THR:HG22	2.02	0.41
2:D:14:THR:HB	2:D:17:GLU:HG3	2.02	0.41
1:E:1:GLN:N	1:E:1:GLN:CD	2.74	0.41
1:A:195:ILE:HG21	1:A:207:VAL:HG12	2.02	0.41
2:D:179:THR:O	2:D:180:LEU:HD23	2.21	0.41
1:E:125:ALA:HB1	1:E:212:GLU:H	1.85	0.41
1:C:197:ASN:HD22	1:C:207:VAL:HG22	1.85	0.41
1:C:136:ALA:CB	1:C:186:SER:HB3	2.50	0.41
2:B:139:TYR:CD2	2:B:140:PRO:HA	2.56	0.41
1:A:47:TRP:CZ3	1:A:49:GLY:HA2	2.56	0.41
1:G:167:PRO:O	1:G:178:LEU:HD23	2.21	0.41
2:B:189:LYS:O	2:B:209:ASN:HA	2.20	0.41
2:H:4:MET:HE3	2:H:23:CYS:SG	2.61	0.41
2:B:174:LEU:HD23	2:B:175:SER:N	2.35	0.41
1:G:82(C):LEU:HB3	1:G:111:VAL:HG11	2.03	0.41
2:F:139:TYR:CD2	2:F:140:PRO:HA	2.56	0.41
1:E:67:VAL:HG22	1:E:82:LEU:HD13	2.03	0.41
1:E:136:ALA:O	1:E:183:THR:HA	2.21	0.41
2:H:116:ILE:HD12	2:H:193:CYS:HB2	2.02	0.41
2:D:4:MET:HE2	2:D:23:CYS:SG	2.60	0.40
2:F:200:LEU:HD13	2:F:204:VAL:HG13	2.03	0.40
1:G:101:GLU:HG3	1:G:102:PHE:CD2	2.56	0.40

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:201:LYS:H	1:C:201:LYS:HG3	1.75	0.40
2:F:48:ILE:HA	2:F:53:ASN:O	2.22	0.40
1:G:20:VAL:HG13	1:G:107:THR:HG21	2.03	0.40
1:A:54:ARG:HB2	1:A:56:ASP:OD1	2.21	0.40
1:E:210:ARG:CD	1:E:210:ARG:C	2.88	0.40
2:F:209:ASN:O	2:F:210:ARG:HB2	2.19	0.40

All (3) 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)
6:E:226:HOH:O	6:F:565:HOH:O[4_555]	1.41	0.79
6:E:558:HOH:O	6:F:384:HOH:O[4_555]	1.57	0.63
4:G:222:CL:CL	6:F:446:HOH:O[3_554]	1.75	0.45

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	212/232 (91%)	198 (93%)	13 (6%)	1 (0%)	34	60
1	C	207/232 (89%)	194 (94%)	11 (5%)	2 (1%)	19	39
1	E	208/232 (90%)	198 (95%)	9 (4%)	1 (0%)	34	60
1	G	176/232 (76%)	165 (94%)	10 (6%)	1 (1%)	30	56
2	B	213/218 (98%)	208 (98%)	4 (2%)	1 (0%)	34	60
2	D	213/218 (98%)	202 (95%)	11 (5%)	0	100	100
2	F	208/218 (95%)	201 (97%)	7 (3%)	0	100	100
2	H	175/218 (80%)	170 (97%)	5 (3%)	0	100	100
All	All	1612/1800 (90%)	1536 (95%)	70 (4%)	6 (0%)	39	65

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	185	PRO
2	B	61	ARG
1	C	156	SER
1	C	202	PRO
1	A	76	SER
1	G	149	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	186/200 (93%)	174 (94%)	12 (6%)	21	42
1	C	184/200 (92%)	169 (92%)	15 (8%)	14	27
1	E	185/200 (92%)	167 (90%)	18 (10%)	10	19
1	G	164/200 (82%)	153 (93%)	11 (7%)	20	40
2	B	189/191 (99%)	177 (94%)	12 (6%)	22	44
2	D	189/191 (99%)	176 (93%)	13 (7%)	19	38
2	F	186/191 (97%)	175 (94%)	11 (6%)	24	47
2	H	171/191 (90%)	154 (90%)	17 (10%)	10	18
All	All	1454/1564 (93%)	1345 (92%)	109 (8%)	17	33

All (109) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	GLN
1	A	30	VAL
1	A	71	ARG
1	A	83	ARG
1	A	110	THR
1	A	135	THR
1	A	178	LEU
1	A	197	ASN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	203	SER
1	A	205	THR
1	A	210	ARG
1	A	211	VAL
2	B	1	ASP
2	B	10	SER
2	B	37	LEU
2	B	56	THR
2	B	89	MET
2	B	102	LYS
2	B	107	ARG
2	B	122	GLU
2	B	128	THR
2	B	178	LEU
2	B	196	THR
2	B	204	VAL
1	C	1	GLN
1	C	30	VAL
1	C	38	ARG
1	C	54	ARG
1	C	64	LYS
1	C	71	ARG
1	C	108	LEU
1	C	159	LEU
1	C	178	LEU
1	C	182	VAL
1	C	188	SER
1	C	192	GLN
1	C	201	LYS
1	C	210	ARG
1	C	211	VAL
2	D	1	ASP
2	D	17	GLU
2	D	37	LEU
2	D	92	LYS
2	D	102	LYS
2	D	107	ARG
2	D	121	ASP
2	D	142	GLU
2	D	155	SER
2	D	167	SER
2	D	184	ASP

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
2	D	189	LYS
2	D	206	LYS
1	E	30	VAL
1	E	38	ARG
1	E	71	ARG
1	E	83	ARG
1	E	89	ILE
1	E	110	THR
1	E	135	THR
1	E	138	LEU
1	E	156	SER
1	E	184	VAL
1	E	189	LEU
1	E	192	GLN
1	E	201	LYS
1	E	203	SER
1	E	205	THR
1	E	210	ARG
1	E	211	VAL
1	E	212	GLU
2	F	1	ASP
2	F	22	SER
2	F	33	LEU
2	F	37	LEU
2	F	77	ARG
2	F	90	GLN
2	F	102	LYS
2	F	107	ARG
2	F	178	LEU
2	F	196	THR
2	F	210	ARG
1	G	30	VAL
1	G	71	ARG
1	G	82(B)	SER
1	G	107	THR
1	G	110	THR
1	G	115	SER
1	G	160	THR
1	G	172	SER
1	G	175	LEU
1	G	178	LEU
1	G	196	CYS

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
2	H	5	THR
2	H	37	LEU
2	H	42	GLN
2	H	89	MET
2	H	90	GLN
2	H	92	LYS
2	H	104	GLU
2	H	107	ARG
2	H	121	ASP
2	H	124	LEU
2	H	125	LYS
2	H	131	VAL
2	H	155	SER
2	H	175	SER
2	H	178	LEU
2	H	193	CYS
2	H	205	THR

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

Mol	Chain	Res	Type
1	E	164	HIS
2	F	123	GLN

### 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](#)

Of 17 ligands modelled in this entry, 10 are monoatomic - leaving 7 for Mogul analysis.

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	SO4	A	221	-	4,4,4	0.39	0	6,6,6	0.76	0
5	GOL	D	214	-	5,5,5	0.33	0	5,5,5	0.56	0
3	SO4	E	221	-	4,4,4	0.45	0	6,6,6	0.39	0
5	GOL	E	222	-	5,5,5	0.62	0	5,5,5	0.52	0
5	GOL	F	214	-	5,5,5	0.45	0	5,5,5	1.01	0
3	SO4	G	221	-	4,4,4	0.36	0	6,6,6	0.20	0
5	GOL	H	214	-	5,5,5	0.64	0	5,5,5	0.21	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	SO4	A	221	-	-	0/0/0/0	0/0/0/0
5	GOL	D	214	-	-	0/4/4/4	0/0/0/0
3	SO4	E	221	-	-	0/0/0/0	0/0/0/0
5	GOL	E	222	-	-	0/4/4/4	0/0/0/0
5	GOL	F	214	-	-	0/4/4/4	0/0/0/0
3	SO4	G	221	-	-	0/0/0/0	0/0/0/0
5	GOL	H	214	-	-	0/4/4/4	0/0/0/0

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	221	SO4	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 ⓘ

### 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	216/232 (93%)	0.02	3 (1%) 78 74	12, 29, 62, 72	0
1	C	213/232 (91%)	0.16	2 (0%) 85 83	12, 30, 68, 78	0
1	E	214/232 (92%)	0.32	8 (3%) 45 37	13, 37, 73, 89	0
1	G	192/232 (82%)	0.53	21 (10%) 7 4	14, 45, 81, 91	0
2	B	215/218 (98%)	-0.03	1 (0%) 91 90	11, 32, 49, 59	0
2	D	215/218 (98%)	-0.09	2 (0%) 85 83	12, 33, 57, 70	0
2	F	212/218 (97%)	0.29	9 (4%) 40 32	15, 40, 84, 94	0
2	H	193/218 (88%)	0.37	11 (5%) 27 20	21, 48, 91, 103	0
All	All	1670/1800 (92%)	0.19	57 (3%) 49 41	11, 35, 77, 103	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	196	CYS	5.7
1	G	121	VAL	4.6
1	G	124	LEU	4.6
1	G	154	TRP	4.3
1	E	210	ARG	4.1
2	H	117	PHE	3.9
1	G	119	PRO	3.8
2	H	208	PHE	3.7
1	G	195	ILE	3.6
2	F	195	VAL	3.5
1	G	113	SER	3.5
1	C	210	ARG	3.4
1	A	210	ARG	3.4
1	G	140	CYS	3.2
1	G	169	VAL	3.0
1	E	121	VAL	3.0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	G	138	LEU	3.0
2	H	132	VAL	3.0
1	E	159	LEU	2.9
1	G	193	THR	2.9
1	G	118	GLY	2.8
1	G	115	SER	2.7
2	F	193	CYS	2.6
2	F	192	ALA	2.6
1	G	116	THR	2.6
2	H	205	THR	2.6
2	H	178	LEU	2.5
1	E	144	ASP	2.5
1	G	145	TYR	2.5
1	G	146	PHE	2.5
1	E	119	PRO	2.5
1	E	211	VAL	2.4
2	H	131	VAL	2.4
1	E	187	SER	2.4
2	D	128	THR	2.4
1	G	147	PRO	2.3
1	C	196	CYS	2.3
1	G	194	TYR	2.3
2	D	126	SER	2.3
2	H	118	PRO	2.3
2	H	116	ILE	2.3
2	B	198	GLN	2.2
1	A	206	LYS	2.2
2	F	180	LEU	2.2
1	E	16	ALA	2.2
1	A	211	VAL	2.2
2	F	201	SER	2.2
1	G	117	LYS	2.1
2	F	196	THR	2.1
1	G	153	SER	2.1
2	F	147	TRP	2.1
2	H	111	ALA	2.1
2	H	146	GLN	2.1
1	G	141	LEU	2.1
2	F	197	HIS	2.0
2	F	1	ASP	2.0
2	H	119	PRO	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 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	GOL	F	214	6/6	0.89	0.23	3.48	38,41,48,51	0
5	GOL	E	222	6/6	0.95	0.22	3.04	43,47,48,49	0
4	CL	B	215	1/1	0.94	0.21	2.88	65,65,65,65	0
5	GOL	D	214	6/6	0.95	0.19	2.40	50,51,51,53	0
5	GOL	H	214	6/6	0.91	0.18	0.40	53,54,54,55	0
3	SO4	E	221	5/5	0.98	0.15	0.17	53,54,55,58	0
3	SO4	A	221	5/5	0.97	0.15	-0.22	55,55,56,59	0
3	SO4	G	221	5/5	0.98	0.11	-0.71	63,63,64,65	0
4	CL	H	215	1/1	0.97	0.13	-	59,59,59,59	0
4	CL	A	222	1/1	0.97	0.36	-	46,46,46,46	0
4	CL	C	222	1/1	0.90	0.39	-	58,58,58,58	0
4	CL	F	215	1/1	0.96	0.16	-	63,63,63,63	0
4	CL	B	214	1/1	0.93	0.09	-	51,51,51,51	0
4	CL	B	216	1/1	0.84	0.14	-	50,50,50,50	0
4	CL	G	222	1/1	0.98	0.28	-	54,54,54,54	0
4	CL	D	215	1/1	0.96	0.08	-	50,50,50,50	0
4	CL	C	221	1/1	0.97	0.23	-	43,43,43,43	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.