



Full wwPDB X-ray Structure Validation Report ⓘ

May 10, 2025 – 07:10 pm BST

PDB ID : 9F90 / pdb_00009f90
Title : Crystal structure of a designed three-motif Respiratory Syncytial Virus immunogen in complex with motavizumab fab
Authors : Castro, K.M.; Correia, B.E.
Deposited on : 2024-05-07
Resolution : 2.91 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<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>.

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

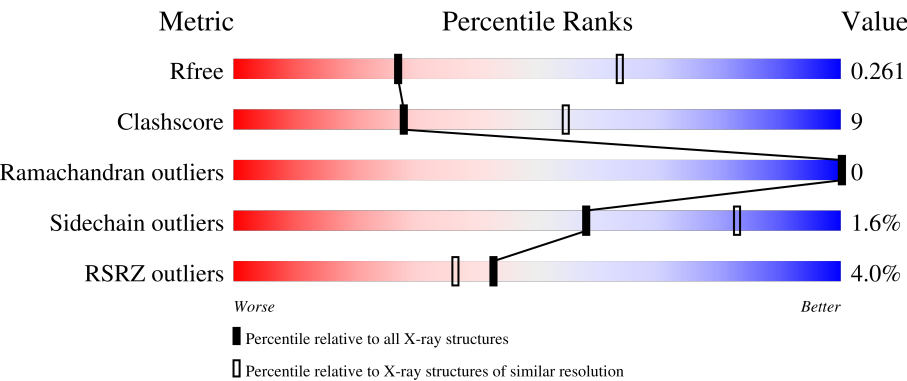
MolProbity	:	4-5-2 with Phenix2.0rc1
Xtriage (Phenix)	:	2.0rc1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.43.1

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.91 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}	164625	2797 (2.94-2.90)
Clashscore	180529	3049 (2.94-2.90)
Ramachandran outliers	177936	2981 (2.94-2.90)
Sidechain outliers	177891	2983 (2.94-2.90)
RSRZ outliers	164620	2799 (2.94-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	B	213	<div><div>2%</div><div></div><div>78%</div><div>21%</div><div></div></div>
1	C	213	<div><div>2%</div><div></div><div>78%</div><div>21%</div><div></div></div>
2	A	225	<div><div>2%</div><div></div><div>79%</div><div>15%</div><div>6%</div></div>
2	D	225	<div><div>2%</div><div></div><div>73%</div><div>20%</div><div>6%</div></div>
3	G	134	<div><div>10%</div><div></div><div>60%</div><div>25%</div><div>14%</div></div>

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Mol	Chain	Length	Quality of chain
3	H	134	<div><div></div><div>7%</div><div>63%</div><div>25%</div><div>13%</div></div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8287 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 Motavizumab Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	211	Total	C	N	O	S	0	0	0
			1607	1010	268	323	6			
1	C	210	Total	C	N	O	S	0	0	0
			1604	1009	265	324	6			

- Molecule 2 is a protein called Motavizumab Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	212	Total	C	N	O	S	0	0	0
			1611	1032	260	312	7			
2	D	212	Total	C	N	O	S	0	0	0
			1608	1029	260	312	7			

- Molecule 3 is a protein called RSVF-multi-epitope designed scaffold.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	117	Total	C	N	O	S	0	0	0
			928	601	146	179	2			
3	G	115	Total	C	N	O	S	0	0	0
			880	573	133	173	1			

- Molecule 4 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	K	0	0
			1	1		
4	A	1	Total	K	0	0
			1	1		
4	C	1	Total	K	0	0
			1	1		
4	H	1	Total	K	0	0
			1	1		

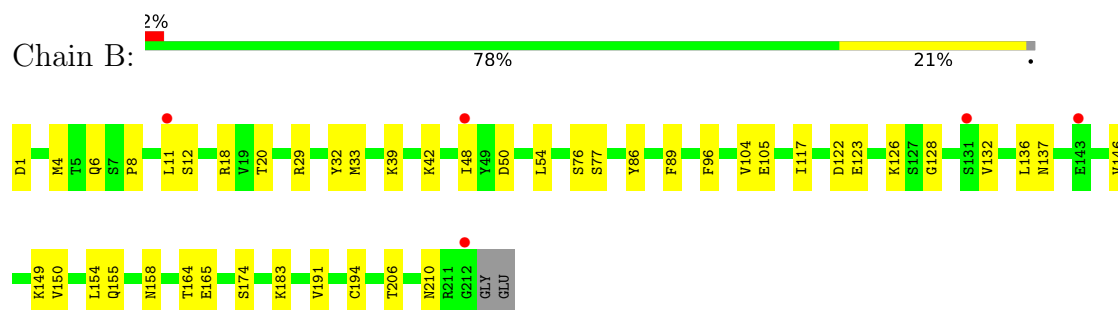
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	8	Total 8	O 8	0	0
5	A	10	Total 10	O 10	0	0
5	C	11	Total 11	O 11	0	0
5	D	7	Total 7	O 7	0	0
5	H	6	Total 6	O 6	0	0
5	G	3	Total 3	O 3	0	0

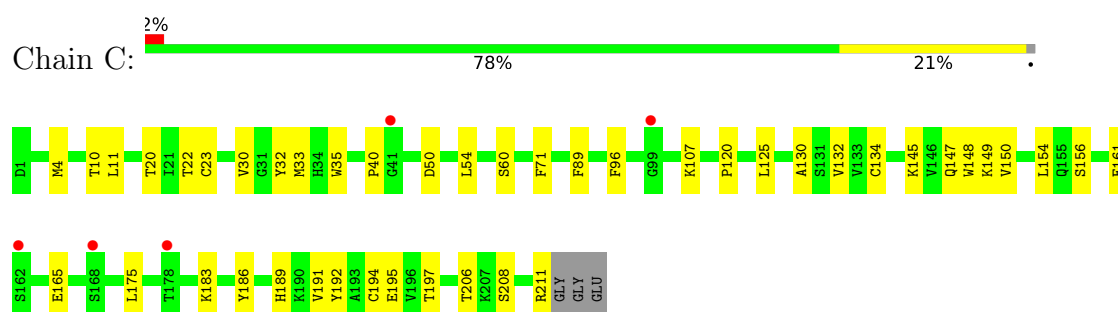
3 Residue-property plots

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.

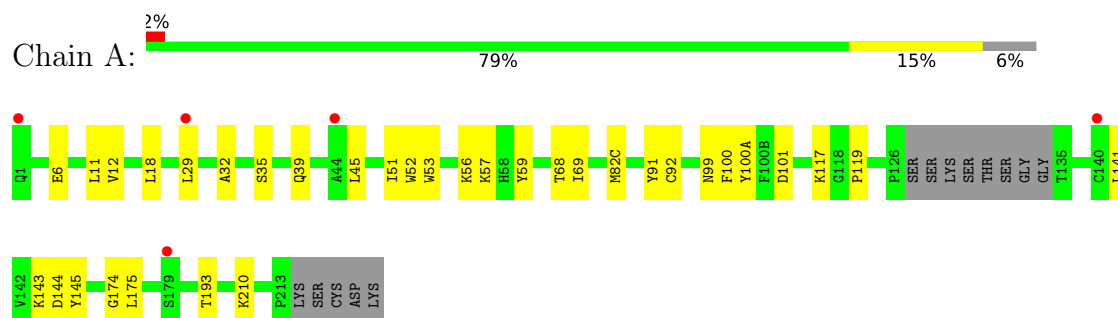
• Molecule 1: Motavizumab Fab light chain



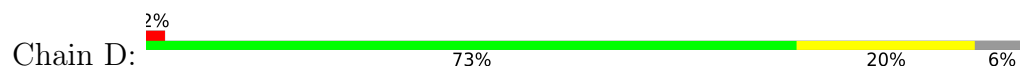
• Molecule 1: Motavizumab Fab light chain

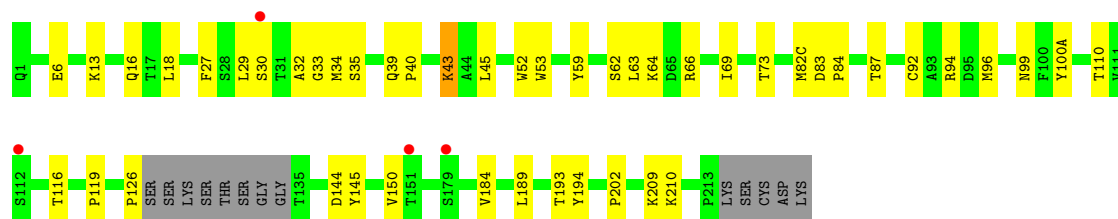


• Molecule 2: Motavizumab Fab heavy chain

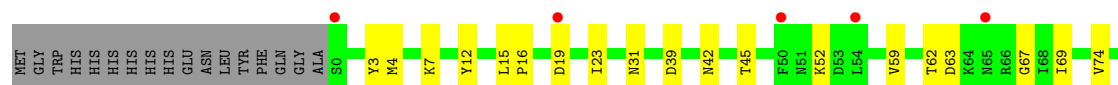


• Molecule 2: Motavizumab Fab heavy chain

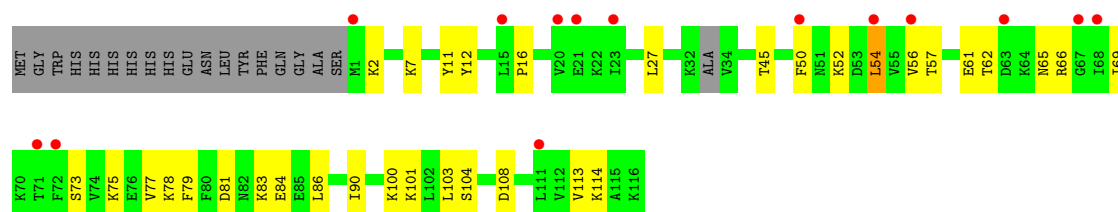




• Molecule 3: RSVF-multi-epitope designed scaffold



• Molecule 3: RSVF-multi-epitope designed scaffold



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	184.07Å 66.86Å 109.41Å 90.00° 103.87° 90.00°	Depositor
Resolution (Å)	61.50 – 2.91 61.50 – 2.91	Depositor EDS
% Data completeness (in resolution range)	99.7 (61.50-2.91) 99.8 (61.50-2.91)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.61 (at 2.91Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487, PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.224 , 0.262 0.224 , 0.261	Depositor DCC
R_{free} test set	1412 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	45.4	Xtriage
Anisotropy	1.342	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8287	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.15% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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:
K

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	B	0.49	0/1644	0.70	0/2230
1	C	0.44	0/1641	0.64	0/2226
2	A	0.48	0/1654	0.76	2/2265 (0.1%)
2	D	0.57	0/1651	0.81	0/2262
3	G	0.48	0/895	0.76	0/1215
3	H	0.48	0/943	0.79	1/1271 (0.1%)
All	All	0.49	0/8428	0.74	3/11469 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	69	ILE	N-CA-C	-6.31	106.79	111.90
2	A	174	GLY	N-CA-C	-6.09	106.36	114.95
2	A	29	LEU	N-CA-C	-5.66	106.42	113.15

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	B	1607	0	1560	32	0
1	C	1604	0	1556	31	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	1611	0	1582	23	0
2	D	1608	0	1576	29	0
3	G	880	0	805	19	0
3	H	928	0	913	18	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	H	1	0	0	0	0
5	A	10	0	0	0	0
5	B	8	0	0	0	0
5	C	11	0	0	0	0
5	D	7	0	0	0	0
5	G	3	0	0	0	0
5	H	6	0	0	1	0
All	All	8287	0	7992	143	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 (143) 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:34:MET:HE1	2:D:94:ARG:HD3	1.54	0.87
1:C:132:VAL:HG12	1:C:148:TRP:CH2	2.19	0.78
3:G:27:LEU:HD13	3:G:52:LYS:HA	1.66	0.77
1:B:154:LEU:HD11	1:C:156:SER:HB3	1.69	0.74
3:H:106:ASN:HB2	3:H:111:LEU:HD11	1.70	0.74
3:H:90:ILE:HA	3:H:93:LEU:HD12	1.70	0.70
2:D:99:ASN:HB2	2:D:100(A):TYR:CE2	2.27	0.69
1:B:12:SER:HB3	1:C:10:THR:HG22	1.75	0.69
2:D:119:PRO:HB3	2:D:145:TYR:HB3	1.76	0.68
1:C:120:PRO:HD3	1:C:132:VAL:HG22	1.74	0.68
2:D:193:THR:HG23	2:D:210:LYS:HE2	1.75	0.67
3:H:91:ASN:O	3:H:100:LYS:NZ	2.30	0.65
2:A:119:PRO:HB3	2:A:145:TYR:HB3	1.79	0.64
3:H:78:LYS:NZ	5:H:301:HOH:O	2.31	0.63
1:B:39:LYS:HB2	1:B:42:LYS:HD3	1.82	0.62
1:B:89:PHE:HE1	1:B:96:PHE:HB3	1.66	0.61
2:A:39:GLN:HB2	2:A:45:LEU:HD23	1.82	0.61
1:C:4:MET:SD	1:C:33:MET:HE1	2.39	0.61
2:D:40:PRO:HB2	2:D:43:LYS:HG3	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:19:ASP:O	3:H:23:ILE:HG13	2.01	0.60
1:B:164:THR:HG22	1:B:174:SER:H	1.68	0.59
2:D:99:ASN:HA	3:G:101:LYS:HE3	1.86	0.58
1:B:89:PHE:CE1	1:B:96:PHE:HB3	2.40	0.57
2:A:56:LYS:HD3	2:A:57:LYS:H	1.70	0.56
2:D:87:THR:HG23	2:D:110:THR:HA	1.87	0.56
1:C:147:GLN:HB3	1:C:195:GLU:HB3	1.88	0.56
1:B:123:GLU:OE1	1:B:123:GLU:N	2.25	0.56
3:G:2:LYS:HD3	3:G:81:ASP:HA	1.87	0.56
2:A:143:LYS:HG2	2:A:144:ASP:OD1	2.07	0.55
2:A:141:LEU:CD2	2:A:143:LYS:HB2	2.37	0.55
2:D:34:MET:CE	2:D:94:ARG:HD3	2.32	0.54
1:C:89:PHE:CE1	1:C:96:PHE:HB3	2.43	0.54
1:B:1:ASP:OD1	1:B:1:ASP:N	2.37	0.54
3:G:16:PRO:HB3	3:G:108:ASP:HA	1.88	0.54
3:G:86:LEU:O	3:G:90:ILE:HG13	2.08	0.54
3:H:3:TYR:HB2	3:H:80:PHE:HB2	1.90	0.53
3:H:39:ASP:OD2	3:H:42:ASN:ND2	2.40	0.53
1:B:122:ASP:O	1:B:126:LYS:HG3	2.08	0.53
1:B:6:GLN:NE2	1:B:86:TYR:O	2.41	0.53
2:A:56:LYS:HD3	2:A:57:LYS:N	2.24	0.53
1:B:154:LEU:HB2	1:C:154:LEU:HB2	1.91	0.53
2:A:99:ASN:HB2	2:A:100(A):TYR:CE2	2.43	0.52
1:B:32:TYR:HB3	1:B:50:ASP:HA	1.91	0.52
1:B:150:VAL:HG12	1:B:155:GLN:HE21	1.73	0.52
1:C:33:MET:HG3	1:C:71:PHE:CD2	2.44	0.52
3:G:84:GLU:N	3:G:84:GLU:OE1	2.42	0.52
2:A:6:GLU:OE2	2:A:91:TYR:HA	2.10	0.51
1:C:54:LEU:HD11	1:C:60:SER:HA	1.91	0.51
2:D:35:SER:HB2	2:D:52:TRP:CE3	2.45	0.51
1:C:132:VAL:HG12	1:C:148:TRP:HH2	1.72	0.51
2:D:13:LYS:O	2:D:16:GLN:HG3	2.11	0.50
1:C:189:HIS:O	1:C:211:ARG:HD3	2.11	0.50
2:A:141:LEU:HD23	2:A:143:LYS:HB2	1.93	0.50
1:B:8:PRO:HB3	1:C:11:LEU:HD13	1.94	0.50
2:A:144:ASP:HB3	2:A:175:LEU:HD13	1.93	0.50
2:D:30:SER:OG	2:D:73:THR:HG21	2.11	0.50
3:G:54:LEU:HB3	3:G:79:PHE:HZ	1.76	0.50
1:C:186:TYR:HA	1:C:192:TYR:OH	2.12	0.49
1:B:117:ILE:HD12	1:B:194:CYS:HB2	1.95	0.48
1:C:32:TYR:HB3	1:C:50:ASP:HA	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:120:PRO:HG3	1:C:130:ALA:HB1	1.96	0.48
1:C:161:GLU:HB2	1:C:175:LEU:HD21	1.96	0.48
3:G:65:ASN:O	3:G:66:ARG:HD3	2.13	0.48
3:H:59:VAL:HG23	3:H:74:VAL:HG22	1.96	0.48
1:C:150:VAL:HG22	1:C:192:TYR:CD2	2.49	0.47
2:D:39:GLN:HB2	2:D:45:LEU:HD23	1.95	0.47
2:A:143:LYS:HG2	2:A:144:ASP:CG	2.39	0.47
2:D:6:GLU:HG3	2:D:92:CYS:SG	2.54	0.47
2:A:18:LEU:HD22	2:A:82(C):MET:HE2	1.97	0.47
2:D:119:PRO:HA	2:D:144:ASP:O	2.15	0.47
3:G:56:VAL:HG23	3:G:77:VAL:HG21	1.97	0.47
3:H:4:MET:HB2	3:H:15:LEU:HB2	1.97	0.46
3:H:62:THR:HA	3:H:67:GLY:O	2.16	0.46
2:A:143:LYS:HG2	2:A:144:ASP:N	2.31	0.46
2:D:126:PRO:HG2	2:D:189:LEU:HD22	1.98	0.46
1:B:164:THR:HG23	1:B:165:GLU:O	2.16	0.46
1:C:194:CYS:O	1:C:206:THR:HA	2.16	0.45
1:B:18:ARG:HD3	1:C:22:THR:HG21	1.99	0.45
2:D:209:LYS:HA	2:D:209:LYS:HD2	1.81	0.45
2:A:35:SER:HB2	2:A:52:TRP:CE3	2.52	0.45
2:D:184:VAL:HG11	2:D:194:TYR:CE1	2.51	0.45
1:B:149:LYS:HE2	1:B:154:LEU:HD21	1.98	0.45
2:D:27:PHE:CD2	2:D:34:MET:HE2	2.51	0.45
1:B:154:LEU:HB2	1:C:154:LEU:CB	2.46	0.45
1:B:76:SER:OG	1:B:77:SER:N	2.50	0.45
3:G:7:LYS:HD3	3:G:12:TYR:CE1	2.52	0.45
1:C:149:LYS:HE2	1:C:154:LEU:HD11	1.99	0.45
2:D:83:ASP:OD1	2:D:84:PRO:HD2	2.17	0.45
3:G:11:TYR:CE1	3:G:114:LYS:HD3	2.52	0.45
3:H:7:LYS:HD2	3:H:12:TYR:CZ	2.52	0.45
3:G:61:GLU:HG2	3:G:69:ILE:HD12	1.99	0.45
3:H:86:LEU:O	3:H:90:ILE:HG13	2.16	0.44
2:A:100:PHE:CE2	3:H:97:ASN:HB3	2.52	0.44
3:G:50:PHE:HE1	3:G:56:VAL:HG22	1.82	0.44
1:B:194:CYS:O	1:B:206:THR:HA	2.17	0.44
2:D:116:THR:HG21	2:D:202:PRO:O	2.18	0.44
3:G:54:LEU:HD13	3:G:54:LEU:HA	1.72	0.44
1:B:20:THR:HG21	1:C:20:THR:HG21	2.00	0.44
1:B:164:THR:HG22	1:B:174:SER:N	2.30	0.44
2:D:29:LEU:HD23	2:D:29:LEU:HA	1.91	0.44
1:B:128:GLY:HA2	1:B:183:LYS:HB2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:12:SER:HA	1:B:105:GLU:O	2.18	0.43
2:D:18:LEU:HD22	2:D:82(C):MET:HE2	2.00	0.43
2:D:35:SER:HB2	2:D:52:TRP:CD2	2.54	0.43
1:B:137:ASN:O	1:B:174:SER:HA	2.18	0.43
1:C:132:VAL:HG12	1:C:148:TRP:CZ3	2.53	0.43
2:A:100(A):TYR:HE1	2:A:101:ASP:HB3	1.84	0.43
3:H:4:MET:HE2	3:H:4:MET:HB3	1.70	0.43
3:G:45:THR:OG1	3:G:62:THR:O	2.37	0.43
1:B:158:ASN:OD1	1:B:158:ASN:N	2.50	0.43
2:A:193:THR:HG23	2:A:210:LYS:HE2	2.01	0.42
1:C:134:CYS:HB2	1:C:148:TRP:CZ2	2.54	0.42
1:B:29:ARG:HB2	1:B:29:ARG:NH1	2.35	0.42
3:H:31:ASN:HA	3:H:52:LYS:HG3	2.00	0.42
1:C:125:LEU:O	1:C:183:LYS:HD2	2.19	0.42
1:B:191:VAL:HG22	1:B:210:ASN:OD1	2.20	0.42
3:G:7:LYS:HG3	3:G:75:LYS:HB2	2.02	0.42
1:C:23:CYS:HB2	1:C:35:TRP:CH2	2.54	0.42
1:C:132:VAL:CG1	1:C:148:TRP:CZ3	3.03	0.41
2:A:11:LEU:HD12	2:A:11:LEU:HA	1.89	0.41
2:A:6:GLU:HG3	2:A:92:CYS:SG	2.60	0.41
2:A:12:VAL:HG11	2:A:82(C):MET:HG3	2.02	0.41
2:D:64:LYS:C	2:D:66:ARG:H	2.27	0.41
1:B:48:ILE:HD13	1:B:54:LEU:HA	2.01	0.41
2:D:32:ALA:HA	2:D:53:TRP:CD1	2.56	0.41
2:D:33:GLY:O	2:D:96:MET:HA	2.20	0.41
2:D:62:SER:O	2:D:63:LEU:C	2.60	0.41
2:D:99:ASN:HB2	2:D:100(A):TYR:CZ	2.55	0.41
2:A:59:TYR:CE1	2:A:69:ILE:HG22	2.55	0.41
1:C:191:VAL:HG23	1:C:208:SER:OG	2.21	0.41
3:H:83:LYS:O	3:H:87:LEU:HG	2.20	0.41
3:G:103:LEU:HD11	3:G:113:VAL:HG21	2.02	0.41
1:B:4:MET:HE1	1:B:33:MET:SD	2.61	0.41
2:A:51:ILE:HB	2:A:69:ILE:HD13	2.02	0.41
1:C:145:LYS:HB3	1:C:197:THR:HB	2.02	0.41
3:G:100:LYS:O	3:G:104:SER:OG	2.26	0.41
3:H:63:ASP:OD1	3:H:67:GLY:N	2.48	0.41
1:B:11:LEU:HD22	1:B:104:VAL:HG22	2.03	0.40
2:D:59:TYR:CZ	2:D:69:ILE:HG22	2.56	0.40
1:C:40:PRO:HB3	1:C:165:GLU:HG3	2.03	0.40
3:H:16:PRO:HB3	3:H:107:ILE:O	2.21	0.40
2:A:32:ALA:HA	2:A:53:TRP:CD1	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:83:LYS:HE2	3:G:83:LYS:HB2	1.74	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	B	209/213 (98%)	200 (96%)	9 (4%)	0	100	100
1	C	208/213 (98%)	202 (97%)	6 (3%)	0	100	100
2	A	208/225 (92%)	203 (98%)	5 (2%)	0	100	100
2	D	208/225 (92%)	199 (96%)	9 (4%)	0	100	100
3	G	111/134 (83%)	107 (96%)	4 (4%)	0	100	100
3	H	115/134 (86%)	111 (96%)	4 (4%)	0	100	100
All	All	1059/1144 (93%)	1022 (96%)	37 (4%)	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	B	182/184 (99%)	179 (98%)	3 (2%)	58	83
1	C	182/184 (99%)	180 (99%)	2 (1%)	70	89

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	A	183/197 (93%)	181 (99%)	2 (1%)	70	89
2	D	183/197 (93%)	181 (99%)	2 (1%)	70	89
3	G	88/126 (70%)	84 (96%)	4 (4%)	23	54
3	H	100/126 (79%)	98 (98%)	2 (2%)	50	78
All	All	918/1014 (90%)	903 (98%)	15 (2%)	58	83

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

Mol	Chain	Res	Type
1	B	132	VAL
1	B	136	LEU
1	B	146	VAL
2	A	68	THR
2	A	117	LYS
1	C	30	VAL
1	C	107	LYS
2	D	43	LYS
2	D	150	VAL
3	H	45	THR
3	H	108	ASP
3	G	54	LEU
3	G	57	THR
3	G	73	SER
3	G	78	LYS

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

Mol	Chain	Res	Type
2	A	58	HIS
3	H	106	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [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, 95th 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 ²)	Q<0.9
1	B	211/213 (99%)	0.43	5 (2%) 59 53	28, 44, 58, 71	0
1	C	210/213 (98%)	0.53	5 (2%) 59 53	30, 47, 61, 70	0
2	A	212/225 (94%)	0.49	5 (2%) 59 53	25, 45, 64, 79	0
2	D	212/225 (94%)	0.57	4 (1%) 66 60	33, 50, 68, 82	0
3	G	115/134 (85%)	1.08	14 (12%) 10 9	48, 72, 97, 103	0
3	H	117/134 (87%)	0.92	10 (8%) 18 16	36, 52, 77, 94	0
All	All	1077/1144 (94%)	0.61	43 (3%) 43 36	25, 48, 77, 103	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	G	20	VAL	3.5
2	D	179	SER	3.5
3	G	68	ILE	3.4
3	H	107	ILE	3.4
3	H	65	ASN	3.3
3	H	50	PHE	3.3
3	G	50	PHE	3.2
3	H	19	ASP	3.1
1	C	99	GLY	2.9
3	G	1	MET	2.9
2	D	30	SER	2.8
2	A	179	SER	2.8
2	A	1	GLN	2.8
2	D	112	SER	2.7
3	H	0	SER	2.7
1	B	48	ILE	2.7
3	G	15	LEU	2.6
1	C	178	THR	2.6
3	H	54	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
3	G	67	GLY	2.5
3	G	23	ILE	2.5
3	G	56	VAL	2.5
1	C	162	SER	2.5
2	A	29	LEU	2.5
1	B	11	LEU	2.4
3	G	111	LEU	2.3
3	G	63	ASP	2.3
1	C	168	SER	2.3
3	H	81	ASP	2.3
3	G	21	GLU	2.2
3	G	71	THR	2.2
3	H	82	ASN	2.2
2	A	140	CYS	2.2
1	B	212	GLY	2.1
3	G	54	LEU	2.1
1	B	143	GLU	2.1
1	B	131	SER	2.1
2	A	44	ALA	2.1
3	H	109	GLU	2.1
3	G	72	PHE	2.1
2	D	151	THR	2.1
3	H	108	ASP	2.0
1	C	41	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, 95th 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(\AA^2)	Q<0.9
4	K	H	201	1/1	0.74	0.11	95,95,95,95	0
4	K	C	301	1/1	0.80	0.13	79,79,79,79	0
4	K	B	301	1/1	0.80	0.14	81,81,81,81	0
4	K	A	301	1/1	0.90	0.10	49,49,49,49	1

6.5 Other polymers [i](#)

There are no such residues in this entry.