



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 06:49 pm BST

PDB ID : 2R0K
Title : Protease domain of HGFA with inhibitor Fab58
Authors : Eigenbrot, C.; Shia, S.
Deposited on : 2007-08-20
Resolution : 3.51 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
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)	:	2.11

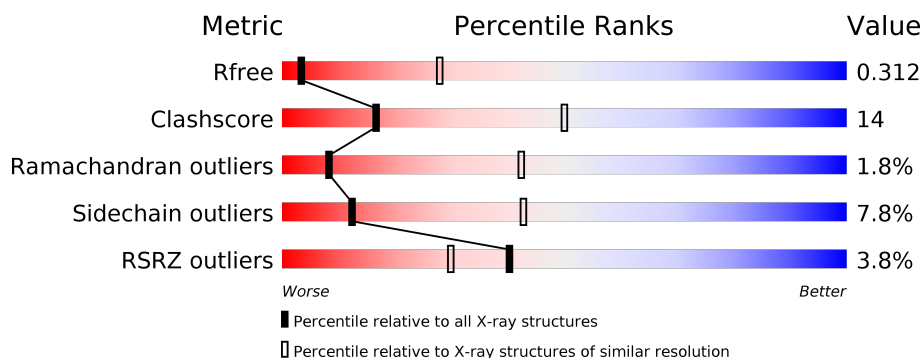
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.51 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	1161 (3.60-3.44)
Clashscore	141614	1244 (3.60-3.44)
Ramachandran outliers	138981	1206 (3.60-3.44)
Sidechain outliers	138945	1207 (3.60-3.44)
RSRZ outliers	127900	1080 (3.60-3.44)

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 ≥ 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	283	<div> <div>5%</div> <div> <div></div> <div>62%</div> <div>22%</div> <div>•</div> <div>16%</div> </div> </div>
2	L	214	<div> <div>5%</div> <div> <div></div> <div>68%</div> <div>27%</div> <div>5%</div> </div> </div>
3	H	225	<div> <div></div> <div> <div>64%</div> <div>25%</div> <div>5%</div> <div>•</div> <div>5%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5042 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 Hepatocyte growth factor activator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	239	Total	C	N	O	S	0	0	0
			1824	1156	316	339	13			

- Molecule 2 is a protein called antibody light chain of Fab58.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	214	Total	C	N	O	S	0	0	0
			1637	1024	271	336	6			

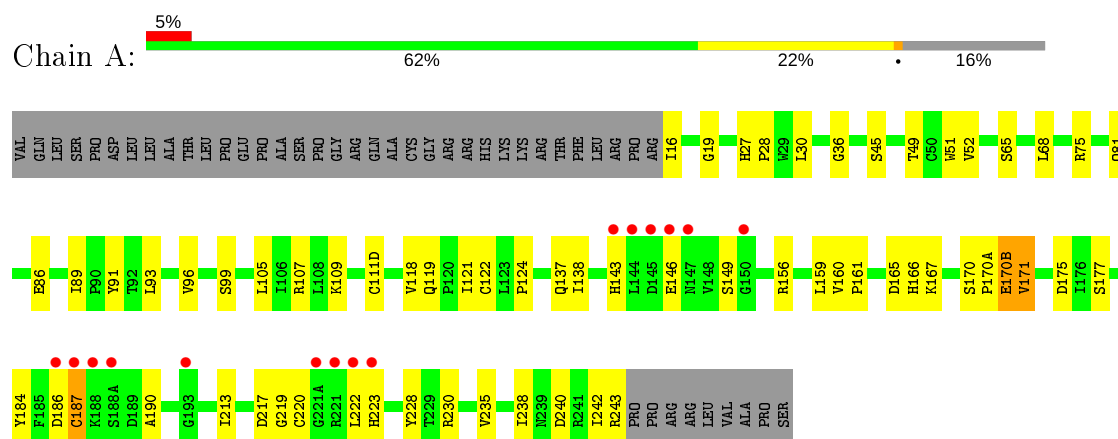
- Molecule 3 is a protein called antibody heavy chain of Fab58, Fab portion only.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	213	Total	C	N	O	S	7	0	0
			1581	1001	266	308	6			

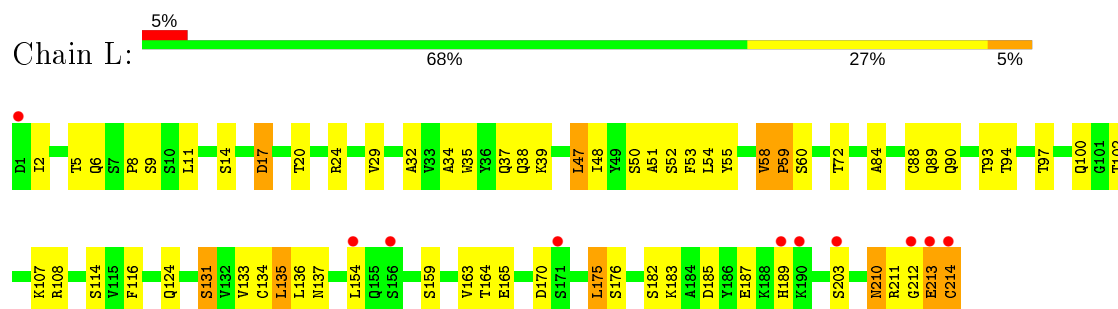
3 Residue-property plots [i](#)

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.

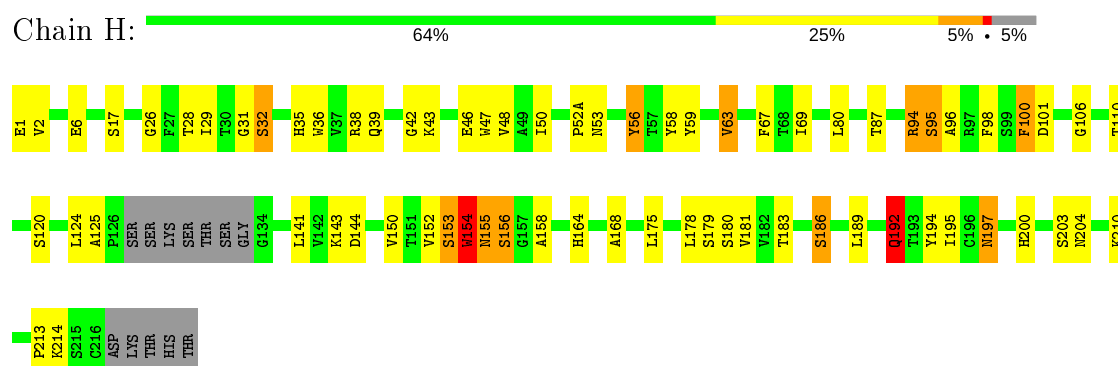
- Molecule 1: Hepatocyte growth factor activator



- Molecule 2: antibody light chain of Fab58



- Molecule 3: antibody heavy chain of Fab58, Fab portion only



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	188.66Å 75.61Å 69.14Å 90.00° 92.67° 90.00°	Depositor
Resolution (Å)	50.00 – 3.51 49.66 – 3.51	Depositor EDS
% Data completeness (in resolution range)	87.7 (50.00-3.51) 87.8 (49.66-3.51)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.55 (at 3.48Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.251 , 0.312 0.258 , 0.312	Depositor DCC
R_{free} test set	526 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	47.6	Xtriage
Anisotropy	0.225	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 31.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.030 for -h,-k,l	Xtriage
F_o, F_c correlation	0.80	EDS
Total number of atoms	5042	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.99% 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](#)

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	0.40	0/1878	0.60	1/2558 (0.0%)
2	L	0.41	0/1673	0.59	1/2273 (0.0%)
3	H	0.67	2/1619 (0.1%)	0.65	1/2208 (0.0%)
All	All	0.50	2/5170 (0.0%)	0.61	3/7039 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	1	GLU	N-CA	-18.31	1.09	1.46
3	H	192	GLN	CB-CG	8.69	1.76	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	175	LEU	CA-CB-CG	5.43	127.78	115.30
1	A	36	GLY	O-C-N	5.25	131.10	122.70
3	H	156	SER	N-CA-C	-5.02	97.44	111.00

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	1824	0	1739	40	0
2	L	1637	0	1586	48	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	H	1581	0	1550	63	0
All	All	5042	0	4875	141	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 (141) 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:H:192:GLN:CG	3:H:192:GLN:CB	1.76	1.64
3:H:153:SER:HA	3:H:154:TRP:HB2	1.41	1.03
1:A:222:LEU:HG	1:A:223:HIS:CD2	1.97	0.99
1:A:217:ASP:HB2	3:H:28:THR:HG21	1.47	0.96
2:L:37:GLN:HB2	2:L:47:LEU:HD21	1.57	0.87
1:A:217:ASP:HB2	3:H:28:THR:CG2	2.07	0.83
2:L:213:GLU:O	2:L:214:CYS:OXT	1.96	0.83
3:H:192:GLN:CG	3:H:192:GLN:CA	2.58	0.81
3:H:35:HIS:HD2	3:H:47:TRP:HE1	1.28	0.81
3:H:35:HIS:CE1	3:H:95:SER:OG	2.39	0.75
3:H:153:SER:CA	3:H:154:TRP:HB2	2.15	0.75
3:H:29:ILE:CG2	3:H:52(A):PRO:HB2	2.20	0.72
2:L:55:TYR:OH	3:H:101:ASP:HB2	1.89	0.72
3:H:153:SER:HA	3:H:154:TRP:CB	2.18	0.71
3:H:192:GLN:CD	3:H:192:GLN:CB	2.58	0.71
2:L:212:GLY:C	2:L:214:CYS:H	1.96	0.70
2:L:136:LEU:HD13	2:L:175:LEU:HD12	1.74	0.69
3:H:29:ILE:HG22	3:H:52(A):PRO:HB2	1.75	0.69
2:L:35:TRP:HB2	2:L:48:ILE:HB	1.74	0.69
1:A:86:GLU:OE2	1:A:109:LYS:HG3	1.94	0.68
2:L:8:PRO:CG	2:L:11:LEU:HD23	2.24	0.68
1:A:19:GLY:O	1:A:156:ARG:NH1	2.27	0.67
3:H:124:LEU:HD21	3:H:141:LEU:HB2	1.76	0.67
2:L:163:VAL:HG22	2:L:175:LEU:HD23	1.75	0.67
2:L:9:SER:O	2:L:102:THR:HA	1.95	0.66
2:L:37:GLN:CB	2:L:47:LEU:HD21	2.25	0.66
3:H:29:ILE:HG23	3:H:52(A):PRO:HG2	1.77	0.66
3:H:87:THR:HG23	3:H:110:THR:HA	1.78	0.66
3:H:168:ALA:HB2	3:H:178:LEU:HD23	1.77	0.65
1:A:186:ASP:O	1:A:187:CYS:C	2.35	0.65
2:L:89:GLN:HG2	2:L:90:GLN:H	1.60	0.65
3:H:35:HIS:NE2	3:H:50:ILE:HD12	2.13	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:35:HIS:HE1	3:H:95:SER:OG	1.81	0.63
2:L:164:THR:HG22	2:L:165:GLU:O	1.97	0.63
2:L:212:GLY:O	2:L:214:CYS:SG	2.56	0.63
1:A:170:SER:HB2	1:A:170(A):PRO:HD2	1.81	0.62
2:L:38:GLN:HE22	3:H:39:GLN:HE22	1.47	0.61
2:L:89:GLN:HG2	2:L:90:GLN:N	2.16	0.60
2:L:211:ARG:HG2	2:L:212:GLY:N	2.17	0.60
2:L:133:VAL:HG21	3:H:124:LEU:HD22	1.84	0.60
3:H:29:ILE:CG2	3:H:52(A):PRO:CB	2.81	0.59
3:H:125:ALA:HB3	3:H:214:LYS:HZ2	1.66	0.59
1:A:28:PRO:HB2	1:A:119:GLN:H	1.68	0.58
3:H:154:TRP:CE3	3:H:154:TRP:HA	2.37	0.58
2:L:54:LEU:HD22	2:L:58:VAL:HG13	1.85	0.58
2:L:136:LEU:HB2	2:L:175:LEU:HB3	1.85	0.58
1:A:81:GLN:HE21	1:A:118:VAL:HG21	1.69	0.57
1:A:49:THR:O	1:A:111(D):CYS:HB2	2.03	0.57
3:H:164:HIS:HB2	3:H:181:VAL:HG12	1.86	0.56
1:A:186:ASP:O	1:A:187:CYS:O	2.23	0.56
3:H:67:PHE:HB3	3:H:80:LEU:HD11	1.86	0.56
1:A:121:ILE:HG13	1:A:122:CYS:H	1.70	0.56
2:L:14:SER:HB3	2:L:107:LYS:HB3	1.87	0.56
1:A:187:CYS:SG	1:A:222:LEU:N	2.78	0.55
2:L:8:PRO:HG3	2:L:11:LEU:HD23	1.87	0.55
3:H:42:GLY:H	3:H:43:LYS:NZ	2.05	0.55
2:L:8:PRO:HG2	2:L:11:LEU:HD23	1.88	0.54
3:H:192:GLN:CG	3:H:192:GLN:HA	2.35	0.54
3:H:100:PHE:N	3:H:100:PHE:CD1	2.76	0.53
2:L:212:GLY:C	2:L:214:CYS:N	2.59	0.53
2:L:2:ILE:O	2:L:97:THR:OG1	2.26	0.53
1:A:28:PRO:HB2	1:A:119:GLN:N	2.25	0.52
3:H:35:HIS:CD2	3:H:47:TRP:HE1	2.19	0.51
2:L:55:TYR:CZ	3:H:101:ASP:HB2	2.45	0.51
1:A:222:LEU:CG	1:A:223:HIS:CD2	2.83	0.50
1:A:121:ILE:HG13	1:A:122:CYS:N	2.25	0.50
1:A:161:PRO:HD3	1:A:184:TYR:CZ	2.46	0.50
3:H:195:ILE:HG22	3:H:197:ASN:ND2	2.27	0.50
3:H:63:VAL:HG22	3:H:67:PHE:CG	2.47	0.50
1:A:45:SER:O	1:A:52:VAL:HA	2.12	0.50
3:H:155:ASN:OD1	3:H:194:TYR:HA	2.12	0.50
2:L:124:GLN:OE1	2:L:131:SER:HB2	2.12	0.49
2:L:50:SER:O	2:L:52:SER:N	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:6:GLU:OE1	3:H:106:GLY:HA2	2.11	0.49
3:H:2:VAL:HA	3:H:26:GLY:HA3	1.94	0.49
3:H:143:LYS:HG2	3:H:144:ASP:N	2.28	0.49
3:H:200:HIS:HD2	3:H:203:SER:OG	1.95	0.49
2:L:35:TRP:CZ3	2:L:88:CYS:HB3	2.48	0.49
3:H:195:ILE:HG13	3:H:210:LYS:HA	1.95	0.49
2:L:47:LEU:HB3	2:L:58:VAL:HG21	1.95	0.49
2:L:6:GLN:O	2:L:100:GLN:NE2	2.46	0.49
1:A:124:PRO:O	1:A:235:VAL:HG21	2.13	0.48
3:H:47:TRP:HZ2	3:H:50:ILE:HG22	1.78	0.48
1:A:99:SER:OG	3:H:31:GLY:HA2	2.13	0.48
2:L:14:SER:O	2:L:17:ASP:HB2	2.14	0.48
3:H:100:PHE:N	3:H:100:PHE:HD1	2.12	0.47
3:H:94:ARG:O	3:H:100:PHE:HA	2.14	0.47
2:L:50:SER:HB2	2:L:53:PHE:CD2	2.50	0.47
3:H:43:LYS:N	3:H:43:LYS:HD3	2.29	0.47
3:H:192:GLN:HG2	3:H:192:GLN:HA	1.96	0.47
1:A:51:TRP:CH2	1:A:107:ARG:HB2	2.50	0.47
2:L:135:LEU:HD22	2:L:137:ASN:HB2	1.97	0.47
3:H:195:ILE:HG22	3:H:197:ASN:HD21	1.78	0.47
3:H:59:TYR:CZ	3:H:69:ILE:HG22	2.50	0.47
3:H:168:ALA:HA	3:H:178:LEU:HB3	1.97	0.46
2:L:94:THR:OG1	3:H:58:TYR:CD2	2.69	0.46
1:A:175:ASP:OD1	3:H:94:ARG:NH2	2.48	0.46
2:L:213:GLU:O	2:L:214:CYS:C	2.53	0.46
1:A:16:ILE:HD13	1:A:190:ALA:HA	1.98	0.46
1:A:27:HIS:HE1	1:A:137:GLN:CD	2.19	0.46
3:H:6:GLU:OE1	3:H:106:GLY:CA	2.64	0.46
3:H:36:TRP:O	3:H:48:VAL:HG22	2.15	0.46
1:A:165:ASP:OD1	1:A:230:ARG:NH2	2.48	0.46
3:H:152:VAL:O	3:H:153:SER:HB3	2.16	0.46
2:L:108:ARG:HD2	2:L:170:ASP:O	2.16	0.46
1:A:213:ILE:HA	1:A:228:TYR:CD2	2.51	0.46
1:A:242:ILE:O	1:A:243:ARG:HB3	2.15	0.46
3:H:164:HIS:O	3:H:180:SER:HA	2.16	0.45
3:H:94:ARG:NH1	3:H:96:ALA:HB3	2.31	0.45
1:A:213:ILE:HG12	1:A:228:TYR:HE2	1.82	0.45
3:H:156:SER:H	3:H:158:ALA:H	1.64	0.44
1:A:105:LEU:HD11	1:A:238:ILE:HG23	1.99	0.44
1:A:146:GLU:HB2	1:A:220:CYS:HB2	1.99	0.44
2:L:210:ASN:HD22	2:L:210:ASN:N	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:20:THR:HG23	2:L:72:THR:HG23	1.99	0.44
2:L:187:GLU:O	2:L:211:ARG:NH1	2.51	0.44
2:L:14:SER:HA	2:L:107:LYS:H	1.81	0.44
1:A:143:HIS:CE1	1:A:149:SER:HA	2.53	0.43
2:L:189:HIS:O	2:L:211:ARG:HD3	2.17	0.43
2:L:58:VAL:HA	2:L:59:PRO:HD2	1.82	0.43
3:H:38:ARG:NE	3:H:46:GLU:OE1	2.37	0.43
2:L:213:GLU:C	2:L:214:CYS:OXT	2.55	0.43
2:L:34:ALA:HB3	2:L:89:GLN:HB3	2.01	0.43
1:A:170(B):GLU:HG2	1:A:171:VAL:HG23	2.00	0.43
2:L:32:ALA:O	2:L:90:GLN:HA	2.19	0.42
1:A:166:HIS:O	1:A:167:LYS:C	2.58	0.42
1:A:27:HIS:HE1	1:A:137:GLN:NE2	2.17	0.42
1:A:121:ILE:CG1	1:A:122:CYS:N	2.83	0.42
1:A:96:VAL:HB	3:H:56:TYR:HE2	1.84	0.42
1:A:217:ASP:O	1:A:219:GLY:C	2.58	0.42
3:H:143:LYS:CG	3:H:144:ASP:N	2.82	0.41
3:H:186:SER:HA	3:H:189:LEU:HD12	2.01	0.41
1:A:138:ILE:HG12	1:A:160:VAL:HG22	2.02	0.41
1:A:91:TYR:CE2	1:A:93:LEU:HB2	2.55	0.41
3:H:35:HIS:HD2	3:H:47:TRP:NE1	2.07	0.41
1:A:213:ILE:CG1	1:A:228:TYR:HE2	2.33	0.41
2:L:116:PHE:HB2	2:L:135:LEU:HB3	2.03	0.41
3:H:189:LEU:HD22	3:H:213:PRO:HG3	2.02	0.41
2:L:39:LYS:HG2	2:L:84:ALA:HB2	2.02	0.41
3:H:59:TYR:OH	3:H:69:ILE:N	2.40	0.40
2:L:134:CYS:O	2:L:176:SER:HA	2.21	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	237/283 (84%)	213 (90%)	22 (9%)	2 (1%)	19	59
2	L	212/214 (99%)	192 (91%)	16 (8%)	4 (2%)	8	41
3	H	209/225 (93%)	182 (87%)	21 (10%)	6 (3%)	4	33
All	All	658/722 (91%)	587 (89%)	59 (9%)	12 (2%)	8	42

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	187	CYS
3	H	98	PHE
3	H	153	SER
3	H	154	TRP
1	A	171	VAL
3	H	204	ASN
2	L	213	GLU
3	H	32	SER
2	L	51	ALA
2	L	59	PRO
2	L	60	SER
3	H	100	PHE

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	199/237 (84%)	190 (96%)	9 (4%)	27	62
2	L	188/188 (100%)	170 (90%)	18 (10%)	8	35
3	H	174/185 (94%)	157 (90%)	17 (10%)	8	34
All	All	561/610 (92%)	517 (92%)	44 (8%)	12	43

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	LEU

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Mol	Chain	Res	Type
1	A	65	SER
1	A	68	LEU
1	A	75	ARG
1	A	89	ILE
1	A	159	LEU
1	A	170(B)	GLU
1	A	177	SER
1	A	240	ASP
2	L	5	THR
2	L	17	ASP
2	L	24	ARG
2	L	29	VAL
2	L	47	LEU
2	L	58	VAL
2	L	93	THR
2	L	114	SER
2	L	131	SER
2	L	135	LEU
2	L	154	LEU
2	L	159	SER
2	L	182	SER
2	L	183	LYS
2	L	185	ASP
2	L	203	SER
2	L	210	ASN
2	L	214	CYS
3	H	17	SER
3	H	32	SER
3	H	53	ASN
3	H	56	TYR
3	H	63	VAL
3	H	94	ARG
3	H	95	SER
3	H	120	SER
3	H	150	VAL
3	H	154	TRP
3	H	155	ASN
3	H	175	LEU
3	H	179	SER
3	H	183	THR
3	H	186	SER
3	H	192	GLN

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Mol	Chain	Res	Type
3	H	197	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	HIS
1	A	71	HIS
1	A	137	GLN
1	A	223	HIS
2	L	3	GLN
2	L	27	GLN
2	L	198	HIS
2	L	210	ASN
3	H	13	GLN
3	H	35	HIS
3	H	39	GLN
3	H	53	ASN
3	H	155	ASN
3	H	200	HIS

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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	A	239/283 (84%)	-0.01	15 (6%)	20	15	5, 8, 45, 63	15 (6%)
2	L	214/214 (100%)	0.22	10 (4%)	31	23	5, 17, 44, 58	7 (3%)
3	H	213/225 (94%)	-0.13	0	100	100	5, 11, 36, 62	3 (1%)
All	All	666/722 (92%)	0.03	25 (3%)	40	30	5, 12, 43, 63	25 (3%)

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	214	CYS	5.1
1	A	145	ASP	4.5
1	A	147	ASN	4.4
1	A	187	CYS	4.4
1	A	186	ASP	3.7
1	A	221	ARG	3.6
1	A	146	GLU	3.3
2	L	156	SER	3.1
2	L	189	HIS	3.0
1	A	144	LEU	3.0
2	L	212	GLY	2.8
1	A	222	LEU	2.8
1	A	143	HIS	2.7
2	L	1	ASP	2.3
2	L	190	LYS	2.3
2	L	171	SER	2.2
2	L	213	GLU	2.2
2	L	154	LEU	2.2
1	A	223	HIS	2.2
1	A	188(A)	SER	2.2
1	A	193	GLY	2.2
2	L	203	SER	2.2
1	A	221(A)	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	188	LYS	2.1
1	A	150	GLY	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.