



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 05:13 PM BST

PDB ID : 6FG2
Title : CRYSTAL STRUCTURE OF FAB OF NATALIZUMAB IN COMPLEX WITH FAB OF NAA84.
Authors : Bertrand, T.; Pouzieux, S.
Deposited on : 2018-01-09
Resolution : 2.79 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
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

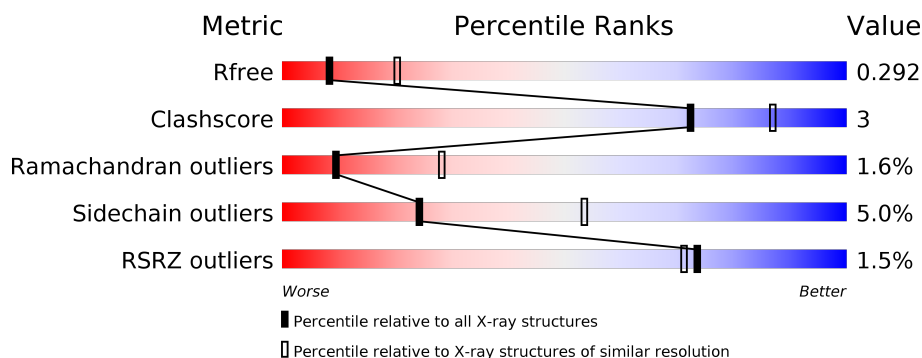
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.79 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	4107 (2.80-2.76)
Clashscore	141614	4575 (2.80-2.76)
Ramachandran outliers	138981	4487 (2.80-2.76)
Sidechain outliers	138945	4489 (2.80-2.76)
RSRZ outliers	127900	4027 (2.80-2.76)

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	D	240	<div> <div>2%</div> <div> <div></div> <div>80%</div> <div>12%</div> <div>6%</div> </div> </div>
1	F	240	<div> <div>%</div> <div> <div></div> <div>84%</div> <div>10%</div> <div>6%</div> </div> </div>
2	E	215	<div> <div>%</div> <div> <div></div> <div>86%</div> <div>10%</div> <div>•</div> </div> </div>
2	G	215	<div> <div></div> <div> <div></div> <div>85%</div> <div>10%</div> <div>•</div> </div> </div>
3	H	234	<div> <div>4%</div> <div> <div></div> <div>77%</div> <div>12%</div> <div>•</div> <div>9%</div> </div> </div>
3	I	234	<div> <div>2%</div> <div> <div></div> <div>83%</div> <div>10%</div> <div>7%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
4	L	211	<div><div>%</div><div><div></div><div>84%</div><div>16%</div></div></div>
4	M	211	<div><div></div><div><div>86%</div><div>14%</div></div></div>
5	A	2	<div><div></div><div><div>100%</div></div></div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13124 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 HEAVY CHAIN FAB NAA84.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	225	Total	C	N	O	S	0	0	0
			1690	1067	281	335	7			
1	F	225	Total	C	N	O	S	0	0	0
			1690	1067	281	335	7			

- Molecule 2 is a protein called LIGHT CHAIN FAB NAA84.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	211	Total	C	N	O	S	0	0	0
			1581	984	268	325	4			
2	G	211	Total	C	N	O	S	0	0	0
			1581	984	268	325	4			

- Molecule 3 is a protein called HEAVY CHAIN FAB NATALIZUMAB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	212	Total	C	N	O	S	0	0	0
			1616	1021	269	318	8			
3	I	218	Total	C	N	O	S	0	0	0
			1656	1045	276	327	8			

- Molecule 4 is a protein called LIGHT CHAIN FAB NATALIZUMAB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	L	211	Total	C	N	O	S	0	0	0
			1641	1029	276	330	6			
4	M	211	Total	C	N	O	S	0	0	0
			1641	1029	276	330	6			

- Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

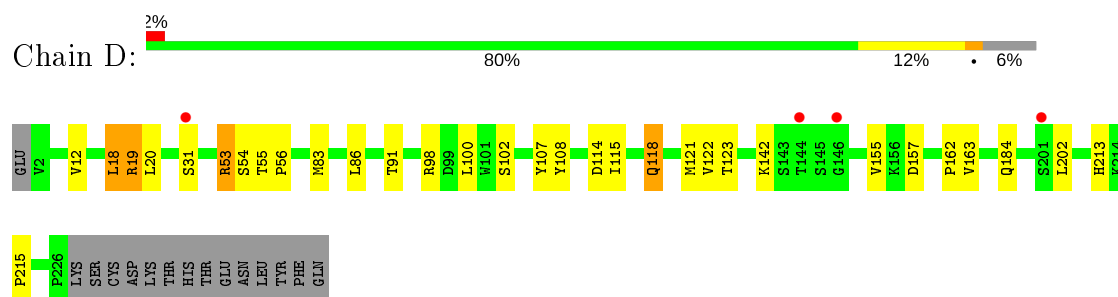


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
5	A	2	28	16	2	10	0	0	0

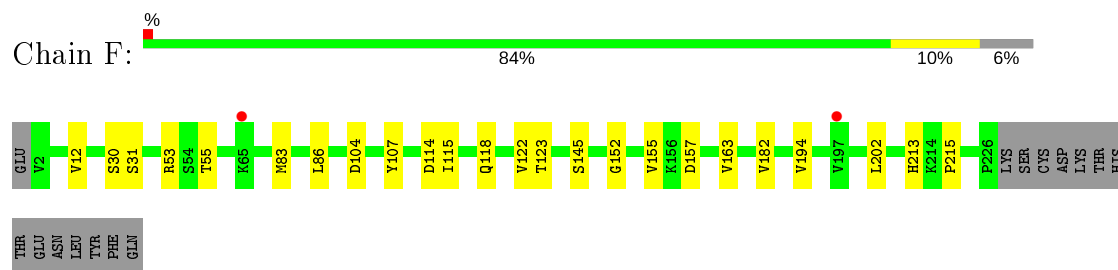
3 Residue-property plots [i](#)

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.

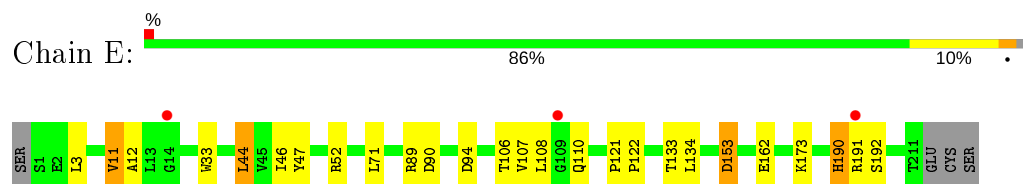
- Molecule 1: HEAVY CHAIN FAB NAA84



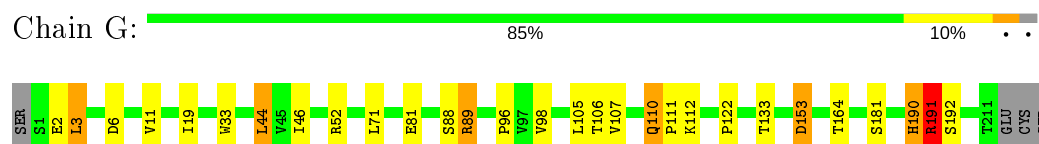
- Molecule 1: HEAVY CHAIN FAB NAA84



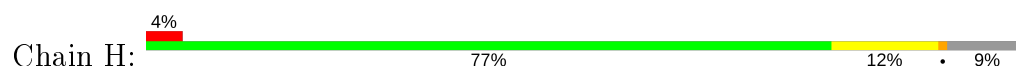
- Molecule 2: LIGHT CHAIN FAB NAA84

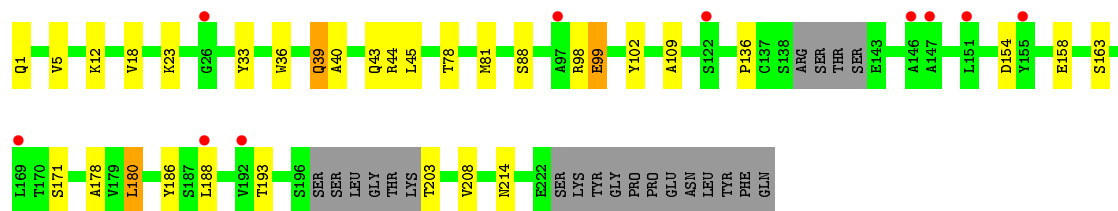


- Molecule 2: LIGHT CHAIN FAB NAA84

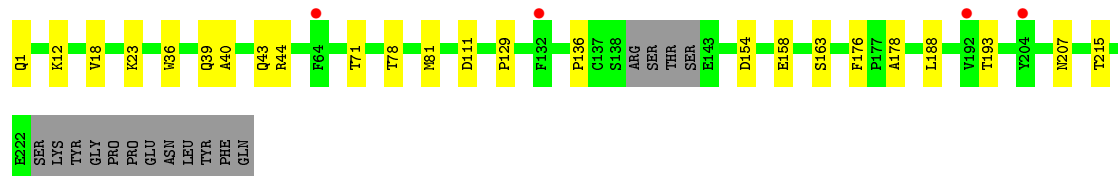
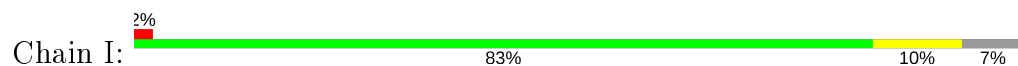


- Molecule 3: HEAVY CHAIN FAB NATALIZUMAB

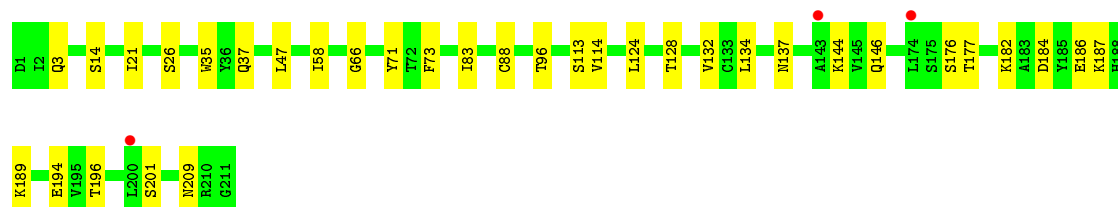
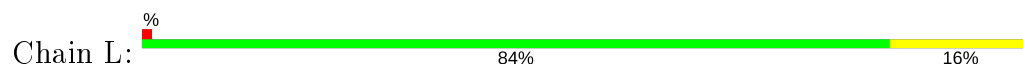




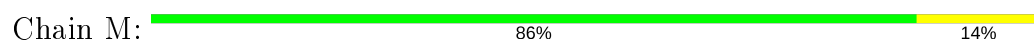
● Molecule 3: HEAVY CHAIN FAB NATALIZUMAB



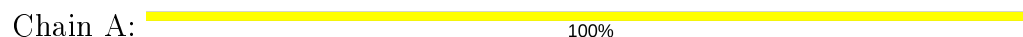
● Molecule 4: LIGHT CHAIN FAB NATALIZUMAB



● Molecule 4: LIGHT CHAIN FAB NATALIZUMAB



● Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	105.40 Å 66.62 Å 199.50 Å 90.00° 98.97° 90.00°	Depositor
Resolution (Å)	55.21 – 2.79 55.21 – 2.80	Depositor EDS
% Data completeness (in resolution range)	58.6 (55.21-2.79) 58.6 (55.21-2.80)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.14 (at 2.81 Å)	Xtriage
Refinement program	BUSTER 2.11.7	Depositor
R, R_{free}	0.238 , 0.263 0.258 , 0.292	Depositor DCC
R_{free} test set	1992 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	54.0	Xtriage
Anisotropy	0.177	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 34.4	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	0.86	EDS
Total number of atoms	13124	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.63% 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: NAG

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	D	0.38	0/1733	0.68	1/2362 (0.0%)
1	F	0.36	0/1733	0.64	0/2362
2	E	0.38	0/1619	0.62	1/2212 (0.0%)
2	G	0.36	0/1619	0.62	1/2212 (0.0%)
3	H	0.38	0/1655	0.62	0/2253
3	I	0.36	0/1696	0.62	0/2309
4	L	0.41	0/1678	0.61	0/2280
4	M	0.39	0/1678	0.59	0/2280
All	All	0.38	0/13411	0.63	3/18270 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	18	LEU	C-N-CA	6.39	137.68	121.70
2	G	190	HIS	C-N-CA	5.64	135.80	121.70
2	E	190	HIS	C-N-CA	5.60	135.69	121.70

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	D	1690	0	1642	13	0
1	F	1690	0	1642	8	0
2	E	1581	0	1517	13	0
2	G	1581	0	1518	16	0
3	H	1616	0	1564	10	0
3	I	1656	0	1609	9	0
4	L	1641	0	1595	14	0
4	M	1641	0	1595	12	0
5	A	28	0	25	0	0
All	All	13124	0	12707	87	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 3.

All (87) 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:G:110:GLN:HB3	2:G:111:PRO:HA	1.49	0.91
1:D:53:ARG:HG3	1:D:102:SER:HB2	1.69	0.74
2:G:190:HIS:HA	2:G:191:ARG:HB2	1.71	0.71
2:E:190:HIS:HA	2:E:191:ARG:HB2	1.79	0.65
3:H:178:ALA:HB2	3:H:188:LEU:HD23	1.81	0.62
2:G:110:GLN:HB3	2:G:111:PRO:CA	2.26	0.60
2:G:81:GLU:HB2	2:G:107:VAL:HG22	1.85	0.59
1:F:182:VAL:CG1	2:G:164:THR:HG23	2.32	0.59
2:E:44:LEU:HD21	2:E:47:TYR:HB3	1.86	0.57
4:L:37:GLN:HB2	4:L:47:LEU:HD11	1.86	0.57
4:M:37:GLN:HB2	4:M:47:LEU:HD11	1.86	0.57
3:I:163:SER:HB3	3:I:207:ASN:HB2	1.87	0.56
1:D:18:LEU:HA	1:D:19:ARG:HB2	1.88	0.56
3:H:40:ALA:HB3	3:H:43:GLN:HB2	1.88	0.55
1:F:182:VAL:HG12	2:G:164:THR:HG23	1.87	0.55
3:I:40:ALA:HB3	3:I:43:GLN:HB2	1.89	0.55
1:D:184:GLN:HG2	2:E:162:GLU:HG3	1.88	0.54
1:F:114:ASP:HA	2:G:44:LEU:HD12	1.90	0.52
1:F:213:HIS:CD2	1:F:215:PRO:HD2	2.44	0.52
2:G:110:GLN:CB	2:G:111:PRO:HA	2.33	0.52
1:D:213:HIS:CD2	1:D:215:PRO:HD2	2.44	0.52
3:H:23:LYS:HA	3:H:78:THR:HG22	1.91	0.51
1:D:114:ASP:HA	2:E:44:LEU:HD12	1.92	0.51
2:E:46:ILE:HD13	2:E:52:ARG:HB3	1.93	0.51
4:L:132:VAL:HG22	4:L:177:THR:HG23	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:M:132:VAL:HG22	4:M:177:THR:HG23	1.93	0.50
4:L:47:LEU:HD23	4:L:58:ILE:HD12	1.94	0.49
3:I:23:LYS:HA	3:I:78:THR:HG22	1.94	0.49
2:G:46:ILE:HD13	2:G:52:ARG:HB3	1.94	0.49
1:F:83:MET:HB3	1:F:86:LEU:HD21	1.95	0.48
3:I:178:ALA:HB2	3:I:188:LEU:HD23	1.95	0.48
3:H:39:GLN:HG3	3:H:45:LEU:HD23	1.96	0.48
4:M:47:LEU:HD23	4:M:58:ILE:HD12	1.95	0.47
4:M:146:GLN:HB3	4:M:194:GLU:HB3	1.95	0.47
1:D:118:GLN:HG3	1:D:118:GLN:H	1.47	0.47
2:G:19:ILE:HD11	2:G:105:LEU:HD13	1.97	0.47
1:D:91:THR:HG23	1:D:123:THR:HA	1.97	0.47
2:G:122:PRO:HB3	2:G:133:THR:H	1.80	0.46
4:L:146:GLN:HB3	4:L:194:GLU:HB3	1.95	0.46
3:I:178:ALA:HA	3:I:188:LEU:HB3	1.96	0.46
2:G:3:LEU:HD21	2:G:98:VAL:HB	1.98	0.46
2:E:122:PRO:HB3	2:E:133:THR:H	1.80	0.45
4:M:66:GLY:HA3	4:M:71:TYR:HA	1.98	0.45
1:D:12:VAL:HG11	1:D:86:LEU:HD13	1.99	0.45
1:D:155:VAL:HG11	1:D:163:VAL:HG21	1.98	0.45
2:E:190:HIS:CA	2:E:191:ARG:HB2	2.47	0.44
3:I:12:LYS:HG3	3:I:18:VAL:HB	2.00	0.44
2:G:110:GLN:HG2	2:G:112:LYS:HG3	1.99	0.44
1:F:182:VAL:HG13	2:G:164:THR:HG23	2.00	0.44
3:I:176:PHE:CD2	4:M:163:THR:HG23	2.53	0.44
2:E:190:HIS:HA	2:E:191:ARG:CB	2.47	0.43
4:L:184:ASP:HA	4:L:187:LYS:HE3	2.00	0.43
4:L:66:GLY:HA3	4:L:71:TYR:HA	1.99	0.43
4:M:184:ASP:HA	4:M:187:LYS:HE3	2.00	0.43
4:M:189:LYS:HE2	4:M:209:ASN:HB2	1.99	0.43
4:L:21:ILE:HD12	4:L:73:PHE:HD2	1.83	0.43
1:D:108:TYR:O	2:E:89:ARG:HG3	2.18	0.43
3:H:12:LYS:HG3	3:H:18:VAL:HB	2.01	0.42
3:I:36:TRP:CE2	3:I:81:MET:HB2	2.54	0.42
4:L:189:LYS:HE2	4:L:209:ASN:HB2	2.01	0.42
4:L:3:GLN:HB2	4:L:26:SER:HB3	2.01	0.42
1:D:83:MET:HB2	1:D:86:LEU:HD21	2.01	0.42
4:L:35:TRP:CZ3	4:L:88:CYS:HB3	2.55	0.42
1:F:155:VAL:HG11	1:F:163:VAL:HG21	2.00	0.42
4:M:35:TRP:CZ3	4:M:88:CYS:HB3	2.55	0.42
3:H:36:TRP:CE2	3:H:81:MET:HB2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:98:ARG:HH11	1:D:115:ILE:HD13	1.85	0.41
2:E:121:PRO:HA	2:E:134:LEU:HD23	2.03	0.41
4:L:144:LYS:HB3	4:L:196:THR:HB	2.02	0.41
4:M:3:GLN:HB2	4:M:26:SER:HB3	2.03	0.41
4:M:47:LEU:HA	4:M:58:ILE:HG13	2.02	0.41
2:E:11:VAL:HG12	2:E:12:ALA:H	1.86	0.41
2:E:108:LEU:HD23	2:E:110:GLN:HB2	2.02	0.41
3:H:178:ALA:HA	3:H:188:LEU:HB3	2.02	0.41
3:I:129:PRO:HD2	3:I:215:THR:HG21	2.00	0.41
4:M:21:ILE:HD12	4:M:73:PHE:HD2	1.85	0.41
3:H:33:TYR:HB2	3:H:99:GLU:HB3	2.02	0.41
1:D:121:MET:HG2	1:D:162:PRO:HD3	2.03	0.41
4:L:47:LEU:HA	4:L:58:ILE:HG13	2.02	0.41
2:G:89:ARG:HA	2:G:96:PRO:O	2.21	0.41
3:H:102:TYR:HE2	3:H:109:ALA:HB3	1.85	0.41
4:L:124:LEU:HB3	4:L:182:LYS:HE3	2.04	0.40
2:G:33:TRP:CE2	2:G:71:LEU:HB2	2.57	0.40
3:H:180:LEU:HA	3:H:186:TYR:HA	2.03	0.40
4:L:114:VAL:HA	4:L:134:LEU:O	2.22	0.40
2:E:33:TRP:CE2	2:E:71:LEU:HB2	2.56	0.40
1:F:152:GLY:HA3	1:F:194:VAL:HG12	2.03	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	D	223/240 (93%)	206 (92%)	11 (5%)	6 (3%)	5	15
1	F	223/240 (93%)	204 (92%)	14 (6%)	5 (2%)	6	20
2	E	209/215 (97%)	194 (93%)	11 (5%)	4 (2%)	8	23
2	G	209/215 (97%)	197 (94%)	7 (3%)	5 (2%)	6	18

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	H	206/234 (88%)	195 (95%)	8 (4%)	3 (2%)	10	30
3	I	214/234 (92%)	201 (94%)	11 (5%)	2 (1%)	17	44
4	L	209/211 (99%)	192 (92%)	16 (8%)	1 (0%)	29	58
4	M	209/211 (99%)	192 (92%)	16 (8%)	1 (0%)	29	58
All	All	1702/1800 (95%)	1581 (93%)	94 (6%)	27 (2%)	9	28

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	19	ARG
1	D	157	ASP
1	F	31	SER
1	F	145	SER
1	F	157	ASP
1	D	31	SER
1	D	53	ARG
1	D	54	SER
2	E	90	ASP
2	E	153	ASP
2	G	110	GLN
3	H	136	PRO
4	L	137	ASN
1	F	55	THR
2	G	153	ASP
3	H	154	ASP
3	I	154	ASP
4	M	137	ASN
1	D	56	PRO
2	E	173	LYS
1	F	12	VAL
2	G	2	GLU
2	G	191	ARG
2	G	192	SER
3	H	171	SER
2	E	107	VAL
3	I	136	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	D	191/206 (93%)	183 (96%)	8 (4%)	30	60
1	F	191/206 (93%)	182 (95%)	9 (5%)	26	56
2	E	176/180 (98%)	169 (96%)	7 (4%)	31	62
2	G	176/180 (98%)	166 (94%)	10 (6%)	20	48
3	H	178/198 (90%)	164 (92%)	14 (8%)	12	31
3	I	183/198 (92%)	176 (96%)	7 (4%)	33	64
4	L	187/187 (100%)	179 (96%)	8 (4%)	29	59
4	M	187/187 (100%)	177 (95%)	10 (5%)	22	51
All	All	1469/1542 (95%)	1396 (95%)	73 (5%)	24	53

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	20	LEU
1	D	55	THR
1	D	100	LEU
1	D	107	TYR
1	D	118	GLN
1	D	122	VAL
1	D	142	LYS
1	D	202	LEU
2	E	3	LEU
2	E	11	VAL
2	E	44	LEU
2	E	94	ASP
2	E	106	THR
2	E	153	ASP
2	E	192	SER
1	F	30	SER
1	F	53	ARG
1	F	104	ASP
1	F	107	TYR

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Mol	Chain	Res	Type
1	F	115	ILE
1	F	118	GLN
1	F	122	VAL
1	F	123	THR
1	F	202	LEU
2	G	3	LEU
2	G	6	ASP
2	G	11	VAL
2	G	44	LEU
2	G	88	SER
2	G	89	ARG
2	G	106	THR
2	G	153	ASP
2	G	181	SER
2	G	191	ARG
3	H	1	GLN
3	H	5	VAL
3	H	39	GLN
3	H	44	ARG
3	H	88	SER
3	H	98	ARG
3	H	99	GLU
3	H	158	GLU
3	H	163	SER
3	H	180	LEU
3	H	193	THR
3	H	203	THR
3	H	208	VAL
3	H	214	ASN
3	I	1	GLN
3	I	39	GLN
3	I	44	ARG
3	I	71	THR
3	I	111	ASP
3	I	158	GLU
3	I	193	THR
4	L	14	SER
4	L	83	ILE
4	L	96	THR
4	L	113	SER
4	L	128	THR
4	L	176	SER

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Mol	Chain	Res	Type
4	L	186	GLU
4	L	201	SER
4	M	1	ASP
4	M	14	SER
4	M	83	ILE
4	M	96	THR
4	M	116	ILE
4	M	164	GLU
4	M	176	SER
4	M	186	GLU
4	M	189	LYS
4	M	201	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 ⓘ

2 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	NAG	A	1	2,5	14,14,15	0.30	0	17,19,21	0.98	1 (5%)
5	NAG	A	2	5	14,14,15	0.33	0	17,19,21	0.67	1 (5%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	A	1	2,5	-	0/6/23/26	0/1/1/1
5	NAG	A	2	5	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1	NAG	C1-O5-C5	3.51	116.95	112.19
5	A	2	NAG	C1-O5-C5	2.04	114.95	112.19

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.6 Ligand geometry [i](#)

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	D	225/240 (93%)	0.33	4 (1%) 68 65	32, 56, 76, 92	0
1	F	225/240 (93%)	0.34	2 (0%) 84 82	32, 49, 72, 82	0
2	E	211/215 (98%)	0.37	3 (1%) 75 73	36, 59, 79, 93	0
2	G	211/215 (98%)	0.27	0 100 100	31, 46, 63, 78	0
3	H	212/234 (90%)	0.49	10 (4%) 31 25	39, 65, 82, 101	0
3	I	218/234 (93%)	0.29	4 (1%) 68 65	36, 54, 70, 96	0
4	L	211/211 (100%)	0.32	3 (1%) 75 73	36, 59, 76, 88	0
4	M	211/211 (100%)	0.26	0 100 100	39, 57, 72, 88	0
All	All	1724/1800 (95%)	0.34	26 (1%) 73 71	31, 55, 77, 101	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	144	THR	4.8
2	E	109	GLY	4.4
3	I	64	PHE	4.1
3	H	122	SER	3.9
3	H	147	ALA	3.7
3	H	151	LEU	3.7
4	L	143	ALA	3.4
3	H	26	GLY	3.1
3	I	192	VAL	2.8
1	D	201	SER	2.8
1	F	65	LYS	2.6
3	H	188	LEU	2.5
3	I	204	TYR	2.4
3	H	146	ALA	2.4
3	H	155	TYR	2.3
3	H	192	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
3	H	97	ALA	2.2
2	E	191	ARG	2.2
4	L	174	LEU	2.1
3	H	169	LEU	2.1
1	D	31	SER	2.1
1	D	146	GLY	2.1
1	F	197	VAL	2.0
2	E	14	GLY	2.0
3	I	132	PHE	2.0
4	L	200	LEU	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](#)

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(Å ²)	Q<0.9
5	NAG	A	2	14/15	0.83	0.20	88,89,91,91	0
5	NAG	A	1	14/15	0.88	0.18	70,79,82,85	0

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.