



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 28, 2022 – 10:13 AM JST

PDB ID : 7WPH
Title : SARS-CoV2 RBD bound to Fab06
Authors : Lin, J.Q.; El Sahili, A.; Lescar, J.
Deposited on : 2022-01-23
Resolution : 2.89 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.31.2
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) : 2.31.2

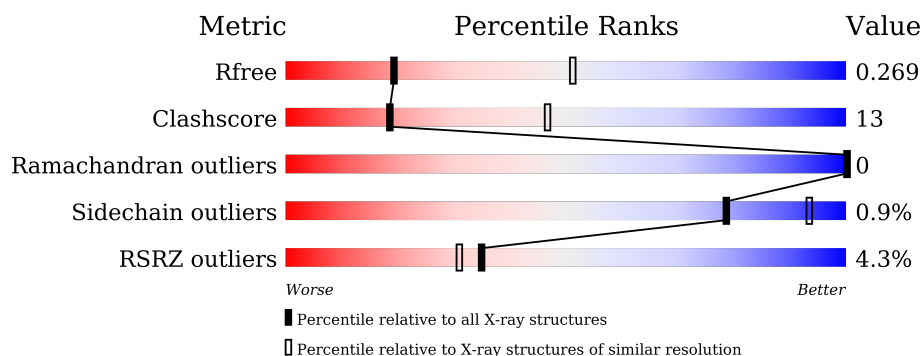
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.89 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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	A	276	<div> <div></div> <div>63%8%29%</div> </div>
1	B	276	<div> <div>6%</div> <div>59%11%29%</div> </div>
2	D	215	<div> <div>10%</div> <div>75%23%</div> </div>
2	L	215	<div> <div>%</div> <div>78%20%</div> </div>
3	E	228	<div> <div>4%</div> <div>67%25%8%</div> </div>
3	H	228	<div> <div>%</div> <div>64%32%</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9534 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 Spike protein S1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	196	Total	C	N	O	S	0	0	0
			1552	995	259	290	8			
1	B	195	Total	C	N	O	S	0	0	0
			1543	989	257	289	8			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	316	SER	-	expression tag	UNP P0DTC2
A	317	ALA	-	expression tag	UNP P0DTC2
A	318	ALA	-	expression tag	UNP P0DTC2
B	316	SER	-	expression tag	UNP P0DTC2
B	317	ALA	-	expression tag	UNP P0DTC2
B	318	ALA	-	expression tag	UNP P0DTC2

- Molecule 2 is a protein called FAB06 light chain.

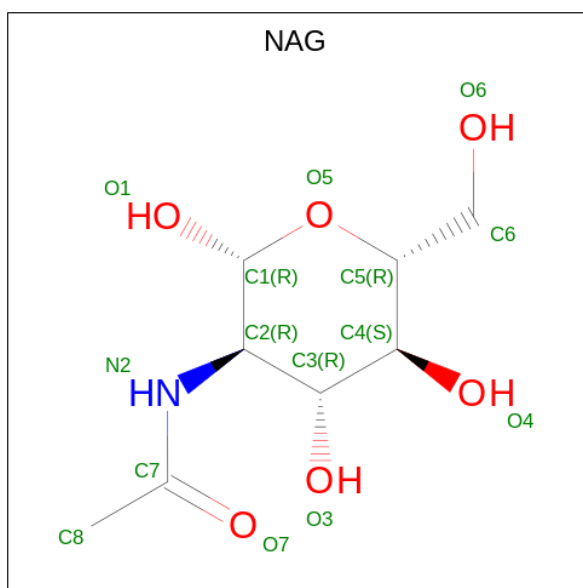
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	212	Total	C	N	O	S	0	0	0
			1570	978	262	325	5			
2	L	211	Total	C	N	O	S	0	0	0
			1561	973	260	323	5			

- Molecule 3 is a protein called Fab06 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	210	Total	C	N	O	S	0	0	0
			1563	987	263	308	5			
3	H	218	Total	C	N	O	S	0	0	0
			1622	1022	275	320	5			

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:

C₈H₁₅NO₆).



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

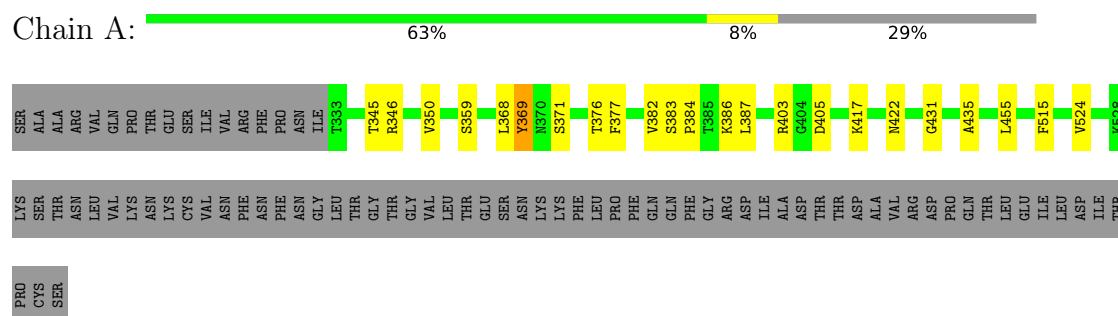
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	33	Total	O	0	0
			33	33		
5	B	16	Total	O	0	0
			16	16		
5	D	16	Total	O	0	0
			16	16		
5	E	8	Total	O	0	0
			8	8		
5	H	7	Total	O	0	0
			7	7		
5	L	15	Total	O	0	0
			15	15		

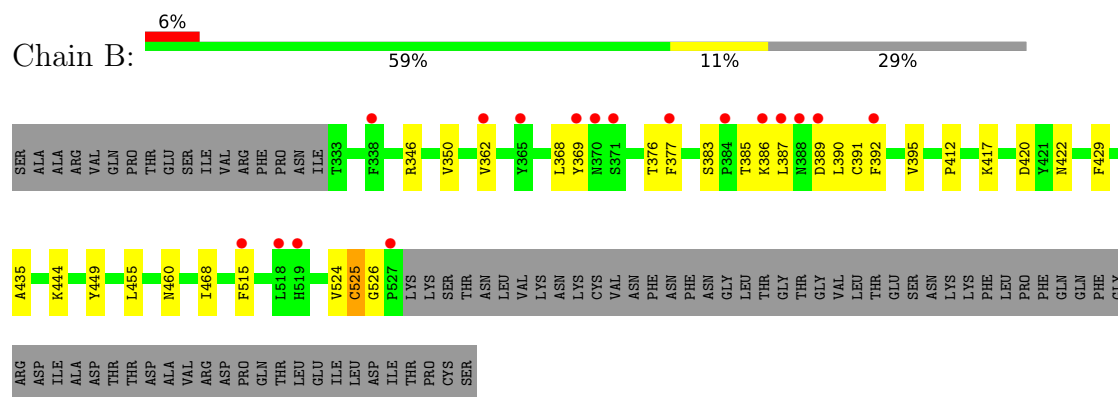
3 Residue-property plots [i](#)

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.

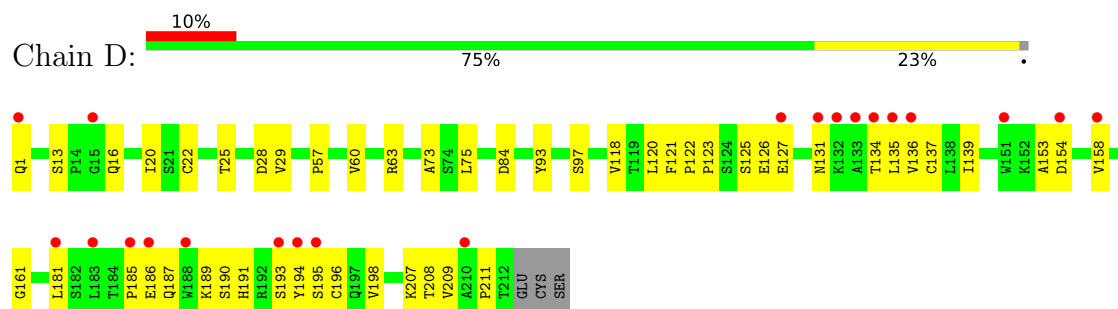
- Molecule 1: Spike protein S1



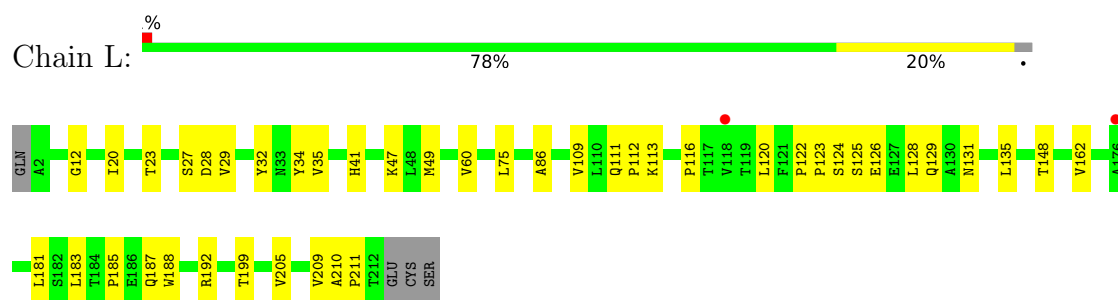
- Molecule 1: Spike protein S1



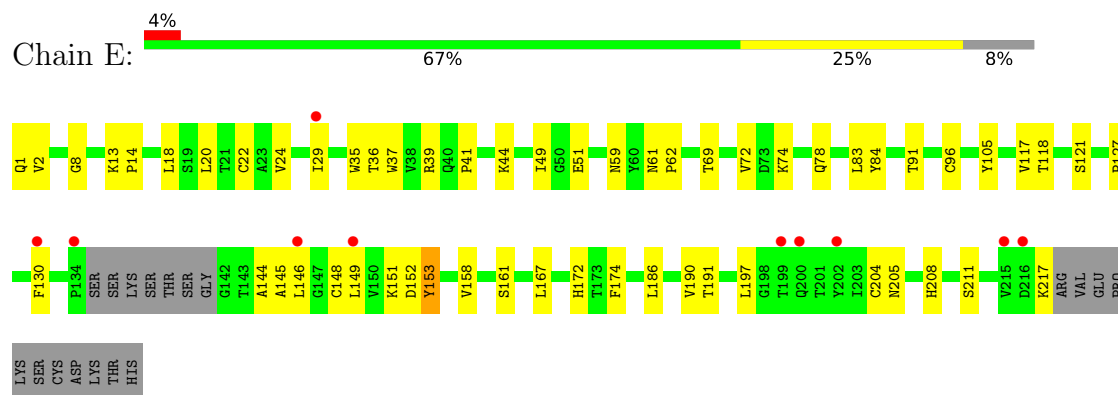
- Molecule 2: FAB06 light chain



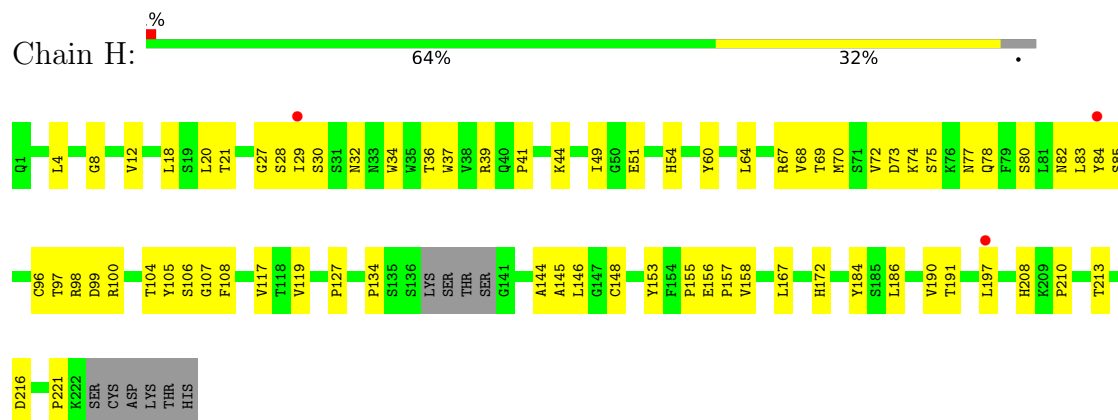
- Molecule 2: FAB06 light chain



• Molecule 3: Fab06 heavy chain



• Molecule 3: Fab06 heavy chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	50.17Å 266.35Å 112.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.15 – 2.89 51.86 – 2.89	Depositor EDS
% Data completeness (in resolution range)	99.3 (48.15-2.89) 99.3 (51.86-2.89)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.14 (at 2.91Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.229 , 0.269 0.229 , 0.269	Depositor DCC
R_{free} test set	1742 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	63.9	Xtriage
Anisotropy	0.495	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 52.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9534	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.37% 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

5.1 Standard geometry

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	A	0.49	0/1596	0.66	0/2172
1	B	0.43	0/1587	0.62	0/2161
2	D	0.49	0/1607	0.65	0/2194
2	L	0.45	0/1598	0.64	0/2182
3	E	0.45	0/1604	0.61	2/2192 (0.1%)
3	H	0.47	0/1664	0.62	0/2272
All	All	0.46	0/9656	0.63	2/13173 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	152	ASP	N-CA-C	5.53	125.92	111.00
3	E	153	TYR	N-CA-CB	-5.40	100.87	110.60

There are no chirality outliers.

There are no planarity outliers.

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	1552	0	1472	19	0
1	B	1543	0	1459	29	1
2	D	1570	0	1520	60	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	L	1561	0	1509	47	0
3	E	1563	0	1519	44	1
3	H	1622	0	1580	56	0
4	A	14	0	13	1	0
4	B	14	0	13	0	0
5	A	33	0	0	1	0
5	B	16	0	0	0	0
5	D	16	0	0	0	0
5	E	8	0	0	0	0
5	H	7	0	0	0	0
5	L	15	0	0	1	0
All	All	9534	0	9085	244	2

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 13.

All (244) 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:B:392:PHE:CE1	1:B:515:PHE:HB3	1.77	1.18
3:E:172:HIS:HD2	3:E:174:PHE:CE1	1.68	1.10
3:E:174:PHE:O	3:E:186:LEU:CD1	2.03	1.06
2:L:120:LEU:CD2	2:L:209:VAL:HG23	1.84	1.06
2:L:120:LEU:HD23	2:L:209:VAL:HG23	1.10	1.05
3:E:174:PHE:O	3:E:186:LEU:HD11	1.60	1.01
3:E:172:HIS:CD2	3:E:174:PHE:CE1	2.50	0.98
2:D:154:ASP:HA	2:D:193:SER:HB3	1.42	0.98
2:L:120:LEU:HG	2:L:209:VAL:HG21	1.51	0.91
1:B:392:PHE:CD1	1:B:515:PHE:HB3	2.05	0.91
3:H:158:VAL:HG22	3:H:208:HIS:HD2	1.38	0.85
3:E:41:PRO:HG2	3:E:44:LYS:HB2	1.57	0.85
3:H:41:PRO:HG2	3:H:44:LYS:HB2	1.56	0.85
2:D:120:LEU:HD12	2:D:207:LYS:HB3	1.58	0.85
3:E:172:HIS:CD2	3:E:174:PHE:CZ	2.65	0.85
3:E:172:HIS:HD2	3:E:174:PHE:CZ	1.94	0.85
2:L:120:LEU:HD23	2:L:209:VAL:CG2	2.03	0.84
2:L:148:THR:OG1	2:L:199:THR:CG2	2.26	0.83
2:D:120:LEU:HD12	2:D:207:LYS:CB	2.09	0.82
2:L:28:ASP:OD1	2:L:29:VAL:N	2.13	0.82
3:H:29:ILE:HD11	3:H:74:LYS:HA	1.62	0.81
3:H:156:GLU:HG2	3:H:157:PRO:HA	1.62	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:29:ILE:HD12	3:H:72:VAL:HG13	1.63	0.80
2:L:120:LEU:CD2	2:L:209:VAL:CG2	2.59	0.80
1:B:369:TYR:HD1	1:B:377:PHE:CE2	2.01	0.78
2:D:154:ASP:CA	2:D:193:SER:HB3	2.12	0.78
1:B:392:PHE:CE1	1:B:515:PHE:CB	2.65	0.78
3:E:29:ILE:HD12	3:E:72:VAL:CG1	2.13	0.77
2:D:120:LEU:CD1	2:D:207:LYS:CB	2.63	0.76
3:E:29:ILE:HD12	3:E:72:VAL:HG13	1.67	0.76
1:B:362:VAL:HG13	1:B:526:GLY:HA3	1.67	0.76
2:L:28:ASP:CG	2:L:29:VAL:H	1.90	0.75
3:E:174:PHE:O	3:E:186:LEU:HD12	1.88	0.73
2:D:120:LEU:HD12	2:D:207:LYS:C	2.08	0.73
1:B:391:CYS:HA	1:B:525:CYS:HB3	1.70	0.72
3:E:29:ILE:HD11	3:E:74:LYS:HA	1.69	0.72
1:A:382:VAL:HG11	1:A:387:LEU:HD21	1.71	0.71
2:D:120:LEU:HD21	2:D:137:CYS:SG	2.31	0.71
3:E:204:CYS:HB3	3:E:217:LYS:HB2	1.71	0.71
3:E:158:VAL:HG13	3:E:186:LEU:HD23	1.73	0.69
1:B:346:ARG:NH1	3:H:105:TYR:O	2.26	0.68
2:L:120:LEU:HG	2:L:209:VAL:CG2	2.22	0.68
1:B:368:LEU:HD12	1:B:368:LEU:H	1.60	0.67
3:E:18:LEU:HD13	3:E:117:VAL:HG11	1.77	0.66
2:D:120:LEU:CD1	2:D:207:LYS:HB2	2.25	0.66
2:L:148:THR:OG1	2:L:199:THR:HG22	1.93	0.66
2:D:97:SER:O	3:E:59:ASN:ND2	2.29	0.65
1:B:392:PHE:CZ	1:B:515:PHE:HB3	2.30	0.65
2:D:189:LYS:HE2	2:D:211:PRO:HB2	1.79	0.65
3:H:158:VAL:HG22	3:H:208:HIS:CD2	2.27	0.65
2:D:122:PRO:HB3	2:D:209:VAL:CG1	2.27	0.64
3:H:36:THR:HG22	3:H:51:GLU:HB3	1.79	0.64
2:L:41:HIS:CE1	2:L:86:ALA:HB2	2.32	0.64
1:A:377:PHE:HE2	1:A:384:PRO:HG3	1.63	0.63
2:L:135:LEU:HD11	2:L:183:LEU:HD12	1.81	0.63
2:D:135:LEU:HB2	2:D:181:LEU:HB3	1.81	0.62
2:L:120:LEU:CG	2:L:209:VAL:CG2	2.77	0.62
2:L:122:PRO:HB3	2:L:209:VAL:CG1	2.29	0.62
3:H:155:PRO:HD2	3:H:210:PRO:HB2	1.82	0.62
2:L:148:THR:OG1	2:L:199:THR:HG23	2.00	0.60
2:D:121:PHE:HB2	2:D:136:VAL:HG22	1.84	0.60
2:D:120:LEU:CD1	2:D:207:LYS:O	2.49	0.60
2:D:25:THR:OG1	2:D:28:ASP:OD1	2.19	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:67:ARG:HD2	3:H:84:TYR:O	2.01	0.59
3:H:153:TYR:HE1	3:H:156:GLU:HG3	1.67	0.59
2:D:120:LEU:CD1	2:D:207:LYS:C	2.71	0.59
2:D:120:LEU:HD23	2:D:137:CYS:HA	1.86	0.58
3:H:158:VAL:CG2	3:H:208:HIS:HD2	2.13	0.58
2:D:118:VAL:HG21	2:D:198:VAL:HG21	1.85	0.58
2:D:154:ASP:HB2	2:D:191:HIS:HB3	1.85	0.58
2:L:29:VAL:HG13	2:L:35:VAL:HG21	1.84	0.58
2:L:123:PRO:HD2	2:L:188:TRP:CZ2	2.39	0.57
2:L:126:GLU:CD	2:L:126:GLU:H	2.07	0.57
2:D:161:GLY:O	2:D:181:LEU:HA	2.04	0.57
3:H:8:GLY:HA3	3:H:20:LEU:HD23	1.86	0.57
2:D:134:THR:HG21	3:E:151:LYS:CE	2.35	0.56
3:H:100:ARG:HG3	3:H:107:GLY:O	2.06	0.56
2:D:122:PRO:HB3	2:D:209:VAL:HG11	1.86	0.56
2:D:154:ASP:N	2:D:193:SER:O	2.32	0.56
2:L:113:LYS:HD3	2:L:113:LYS:H	1.71	0.55
2:D:25:THR:N	2:D:28:ASP:OD2	2.39	0.55
2:D:134:THR:HG21	3:E:151:LYS:HE3	1.88	0.55
3:E:127:PRO:HB3	3:E:153:TYR:HB3	1.88	0.55
1:A:377:PHE:HZ	1:A:384:PRO:HB3	1.71	0.55
3:E:24:VAL:HG21	3:E:29:ILE:HG22	1.87	0.55
1:B:369:TYR:HD1	1:B:377:PHE:CD2	2.24	0.55
1:B:377:PHE:CD2	1:B:377:PHE:O	2.60	0.55
2:L:183:LEU:HD22	2:L:187:GLN:HG3	1.88	0.55
3:E:69:THR:HG23	3:E:84:TYR:HE2	1.72	0.54
2:D:28:ASP:OD2	2:D:29:VAL:N	2.39	0.54
1:B:468:ILE:HG12	1:B:468:ILE:O	2.08	0.54
2:L:49:MET:HE2	2:L:49:MET:HA	1.89	0.54
2:D:120:LEU:CD1	2:D:207:LYS:HB3	2.30	0.54
3:H:39:ARG:HB3	3:H:49:ILE:HD11	1.90	0.54
3:E:8:GLY:HA3	3:E:20:LEU:HD23	1.88	0.54
1:B:389:ASP:O	1:B:390:LEU:HG	2.07	0.53
2:D:25:THR:OG1	2:D:28:ASP:CG	2.47	0.53
3:H:156:GLU:CG	3:H:157:PRO:HA	2.37	0.53
3:E:1:GLN:HG3	3:E:2:VAL:N	2.23	0.53
2:D:154:ASP:OD1	2:D:154:ASP:O	2.27	0.53
1:B:449:TYR:HB2	3:H:32:ASN:OD1	2.09	0.53
2:L:125:SER:HA	2:L:128:LEU:HD12	1.90	0.53
1:A:377:PHE:CE2	1:A:384:PRO:HG3	2.42	0.53
2:L:28:ASP:CG	2:L:29:VAL:N	2.58	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:39:ARG:HB3	3:E:49:ILE:HD11	1.90	0.53
1:B:449:TYR:C	3:H:32:ASN:HD21	2.13	0.52
2:D:120:LEU:HD11	2:D:207:LYS:HB2	1.90	0.52
3:H:29:ILE:O	3:H:54:HIS:HB3	2.09	0.52
1:B:387:LEU:O	1:B:390:LEU:HD12	2.10	0.52
3:E:158:VAL:CG1	3:E:186:LEU:HD23	2.37	0.52
2:L:47:LYS:HE2	2:L:49:MET:HE1	1.92	0.52
1:A:377:PHE:CZ	1:A:384:PRO:HB3	2.45	0.51
3:E:61:ASN:OD1	3:E:62:PRO:HD2	2.10	0.51
1:B:444:LYS:HE2	3:H:34:TRP:CE2	2.45	0.51
2:L:116:PRO:HG2	2:L:205:VAL:HG21	1.93	0.51
2:D:122:PRO:HB3	2:D:209:VAL:HG12	1.92	0.51
2:D:195:SER:CB	2:D:208:THR:HG22	2.40	0.51
2:D:93:TYR:HB2	3:E:105:TYR:HB3	1.93	0.51
3:H:145:ALA:HB2	3:H:191:THR:HG22	1.93	0.51
1:A:382:VAL:HG11	1:A:387:LEU:CD2	2.39	0.51
3:E:22:CYS:O	3:E:78:GLN:HB2	2.11	0.51
1:B:391:CYS:CA	1:B:525:CYS:HB3	2.36	0.51
3:E:144:ALA:HB3	3:E:197:LEU:HD21	1.93	0.51
3:H:134:PRO:HD2	3:H:221:PRO:HA	1.94	0.50
2:L:131:ASN:HA	2:L:185:PRO:HG2	1.92	0.50
2:L:124:SER:HB3	2:L:126:GLU:OE2	2.11	0.50
4:A:601:NAG:H61	5:A:724:HOH:O	2.11	0.50
3:H:144:ALA:HB3	3:H:197:LEU:HD21	1.94	0.49
1:A:376:THR:HB	1:A:435:ALA:HB3	1.95	0.49
3:E:91:THR:HG23	3:E:118:THR:HA	1.94	0.49
1:B:420:ASP:HB3	1:B:460:ASN:OD1	2.13	0.49
2:L:111:GLN:CG	2:L:112:PRO:HD2	2.43	0.49
2:D:186:GLU:O	2:D:190:SER:HB3	2.11	0.49
2:D:158:VAL:HG11	2:D:181:LEU:HD11	1.94	0.48
3:H:21:THR:HG22	3:H:80:SER:OG	2.13	0.48
3:H:153:TYR:CE1	3:H:156:GLU:HG3	2.47	0.48
2:D:187:GLN:HA	2:D:190:SER:OG	2.13	0.48
1:B:350:VAL:HG22	1:B:422:ASN:HB3	1.95	0.48
2:D:122:PRO:HD3	2:D:209:VAL:HG11	1.94	0.48
3:E:145:ALA:HB2	3:E:191:THR:HG22	1.93	0.48
1:A:359:SER:HA	1:A:524:VAL:HG22	1.94	0.48
1:B:395:VAL:HG23	1:B:524:VAL:HG21	1.95	0.48
2:D:120:LEU:HD21	2:D:196:CYS:CB	2.44	0.47
2:L:135:LEU:N	2:L:135:LEU:HD12	2.29	0.47
3:H:99:ASP:OD1	3:H:106:SER:HB3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:417:LYS:HE2	1:B:455:LEU:HD12	1.95	0.47
3:E:18:LEU:HB3	3:E:83:LEU:HB3	1.96	0.47
1:B:376:THR:HB	1:B:435:ALA:HB3	1.95	0.47
3:E:161:SER:HB3	3:E:205:ASN:HB2	1.97	0.47
3:H:27:GLY:O	3:H:77:ASN:ND2	2.41	0.47
3:H:18:LEU:HB3	3:H:83:LEU:HB3	1.96	0.46
1:A:377:PHE:HZ	1:A:384:PRO:CB	2.28	0.46
2:D:131:ASN:HA	2:D:185:PRO:HG2	1.97	0.46
2:D:139:ILE:HG12	2:D:198:VAL:HG11	1.97	0.46
2:L:122:PRO:HB3	2:L:209:VAL:HG11	1.98	0.46
1:A:350:VAL:HG22	1:A:422:ASN:HB3	1.97	0.46
2:D:123:PRO:HA	2:D:127:GLU:OE1	2.15	0.46
2:D:120:LEU:HD21	2:D:196:CYS:HB2	1.98	0.46
1:A:369:TYR:N	1:A:369:TYR:CD1	2.83	0.46
3:H:104:THR:O	3:H:105:TYR:C	2.53	0.46
2:L:49:MET:HE2	2:L:60:VAL:HG22	1.97	0.46
2:L:120:LEU:CG	2:L:209:VAL:HG21	2.30	0.46
2:L:120:LEU:CG	2:L:209:VAL:HG23	2.39	0.46
1:A:368:LEU:O	1:A:371:SER:HB2	2.16	0.46
3:H:64:LEU:O	3:H:68:VAL:HG22	2.15	0.46
1:A:417:LYS:HE2	1:A:455:LEU:HD12	1.97	0.45
3:E:146:LEU:C	3:E:146:LEU:HD12	2.37	0.45
2:D:13:SER:O	2:D:16:GLN:HB2	2.16	0.45
3:H:36:THR:OG1	3:H:97:THR:HG23	2.16	0.45
3:H:18:LEU:HD13	3:H:117:VAL:HG11	1.97	0.45
2:L:123:PRO:HD3	2:L:135:LEU:HG	1.98	0.45
2:D:153:ALA:HA	2:D:193:SER:O	2.16	0.45
1:A:382:VAL:HG12	1:A:383:SER:N	2.31	0.45
3:H:28:SER:C	3:H:30:SER:H	2.19	0.45
2:D:195:SER:HB2	2:D:208:THR:HG22	1.98	0.45
3:E:36:THR:HG22	3:E:51:GLU:HB3	1.98	0.45
3:H:153:TYR:CE1	3:H:184:TYR:HB2	2.52	0.45
2:D:118:VAL:CG2	2:D:198:VAL:HG21	2.46	0.45
2:D:57:PRO:HD2	2:D:60:VAL:HG21	1.97	0.45
2:L:41:HIS:ND1	2:L:86:ALA:HB2	2.32	0.45
2:L:188:TRP:CZ2	2:L:211:PRO:HA	2.52	0.45
1:B:449:TYR:O	3:H:32:ASN:ND2	2.50	0.44
3:H:4:LEU:HB3	3:H:96:CYS:SG	2.57	0.44
3:E:146:LEU:HD12	3:E:146:LEU:O	2.18	0.44
3:H:155:PRO:HD2	3:H:210:PRO:CB	2.48	0.44
3:H:12:VAL:O	3:H:119:VAL:HA	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:189:LYS:HG2	2:D:211:PRO:HG3	2.00	0.43
2:L:123:PRO:HD2	2:L:188:TRP:CH2	2.52	0.43
3:H:37:TRP:HD1	3:H:70:MET:CE	2.32	0.43
2:L:47:LYS:CE	2:L:49:MET:HE1	2.47	0.43
2:L:162:VAL:HG22	2:L:181:LEU:HD13	2.00	0.43
2:D:120:LEU:CD2	2:D:137:CYS:SG	3.04	0.43
1:A:345:THR:HG22	1:A:346:ARG:HG2	1.99	0.43
3:H:28:SER:C	3:H:30:SER:N	2.72	0.43
3:H:29:ILE:HD12	3:H:72:VAL:CG1	2.43	0.43
3:H:98:ARG:O	3:H:108:PHE:HA	2.18	0.43
2:D:125:SER:OG	2:D:126:GLU:OE1	2.37	0.43
3:E:37:TRP:CZ3	3:E:96:CYS:HB3	2.54	0.43
3:H:146:LEU:HD12	3:H:146:LEU:C	2.38	0.43
2:L:111:GLN:HG3	2:L:112:PRO:HD2	2.00	0.43
2:D:122:PRO:CB	2:D:209:VAL:HG11	2.49	0.43
2:D:194:TYR:O	2:D:208:THR:HA	2.19	0.43
3:E:13:LYS:HG3	3:E:14:PRO:HD2	2.01	0.43
2:D:118:VAL:HG21	2:D:198:VAL:CG2	2.48	0.42
2:L:32:TYR:HB3	2:L:34:TYR:CD2	2.55	0.42
1:B:369:TYR:CD1	1:B:377:PHE:CE2	2.93	0.42
2:D:122:PRO:CD	2:D:209:VAL:HG11	2.49	0.42
2:L:113:LYS:HD3	2:L:113:LYS:N	2.33	0.42
3:H:127:PRO:HD2	3:H:213:THR:HG21	2.01	0.42
2:L:20:ILE:HD12	2:L:75:LEU:HD23	2.01	0.42
2:D:20:ILE:HD12	2:D:75:LEU:HD23	2.01	0.42
2:D:63:ARG:HH22	2:D:84:ASP:CG	2.23	0.42
3:H:69:THR:HG23	3:H:84:TYR:HE2	1.84	0.42
1:A:403:ARG:NH2	1:A:405:ASP:OD2	2.53	0.42
2:L:12:GLY:O	2:L:109:VAL:HA	2.20	0.42
2:L:125:SER:O	2:L:129:GLN:HG2	2.20	0.42
3:H:67:ARG:HD2	3:H:85:SER:HB3	2.02	0.42
3:H:67:ARG:CD	3:H:85:SER:HB3	2.50	0.42
3:E:167:LEU:HD21	3:E:190:VAL:HG21	2.01	0.41
2:D:136:VAL:HG11	3:E:149:LEU:HD13	2.02	0.41
3:E:14:PRO:HG2	3:E:121:SER:OG	2.19	0.41
1:B:386:LYS:N	1:B:386:LYS:HD2	2.34	0.41
2:D:127:GLU:OE2	3:E:149:LEU:CD2	2.68	0.41
3:H:96:CYS:SG	3:H:96:CYS:O	2.78	0.41
1:A:386:LYS:HA	1:A:386:LYS:HD2	1.92	0.41
3:E:208:HIS:CE1	3:E:211:SER:HG	2.39	0.41
3:H:54:HIS:HB2	3:H:74:LYS:HE2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:60:TYR:HE1	3:H:70:MET:HG3	1.86	0.41
3:E:29:ILE:HA	3:E:35:TRP:CZ2	2.55	0.41
3:H:69:THR:OG1	3:H:82:ASN:HB2	2.21	0.41
1:A:382:VAL:HG12	1:A:383:SER:O	2.21	0.41
1:A:431:GLY:HA2	1:A:515:PHE:CD2	2.56	0.41
3:H:146:LEU:HD12	3:H:146:LEU:O	2.20	0.41
2:L:192:ARG:NH1	5:L:301:HOH:O	2.53	0.41
2:L:210:ALA:HA	2:L:211:PRO:HD3	1.95	0.41
1:B:369:TYR:CD1	1:B:377:PHE:CD2	3.08	0.41
3:H:73:ASP:OD1	3:H:75:SER:OG	2.36	0.41
3:H:97:THR:OG1	3:H:98:ARG:N	2.55	0.40
1:B:383:SER:OG	1:B:385:THR:HG22	2.21	0.40
3:H:167:LEU:HD21	3:H:190:VAL:HG21	2.03	0.40
2:D:127:GLU:HG2	3:E:130:PHE:CG	2.57	0.40
1:B:412:PRO:HG3	1:B:429:PHE:HB3	2.03	0.40
2:D:22:CYS:HB3	2:D:73:ALA:HB3	2.02	0.40
3:H:29:ILE:HD13	3:H:78:GLN:N	2.35	0.40

All (2) 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)
1:B:369:TYR:CE2	1:B:369:TYR:CE2[2_655]	2.06	0.14
2:D:131:ASN:ND2	3:E:13:LYS:CD[4_456]	2.13	0.07

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	194/276 (70%)	186 (96%)	8 (4%)	0	100	100
1	B	193/276 (70%)	184 (95%)	9 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	210/215 (98%)	199 (95%)	11 (5%)	0	100	100
2	L	209/215 (97%)	197 (94%)	12 (6%)	0	100	100
3	E	206/228 (90%)	200 (97%)	6 (3%)	0	100	100
3	H	214/228 (94%)	204 (95%)	10 (5%)	0	100	100
All	All	1226/1438 (85%)	1170 (95%)	56 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	169/241 (70%)	168 (99%)	1 (1%)	86	96
1	B	168/241 (70%)	167 (99%)	1 (1%)	86	96
2	D	179/182 (98%)	178 (99%)	1 (1%)	86	96
2	L	178/182 (98%)	176 (99%)	2 (1%)	73	92
3	E	177/194 (91%)	176 (99%)	1 (1%)	86	96
3	H	184/194 (95%)	180 (98%)	4 (2%)	52	81
All	All	1055/1234 (86%)	1045 (99%)	10 (1%)	78	93

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	369	TYR
1	B	525	CYS
2	D	1	GLN
3	E	148	CYS
3	H	148	CYS
3	H	172	HIS
3	H	186	LEU
3	H	216	ASP
2	L	23	THR

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Mol	Chain	Res	Type
2	L	27	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	354	ASN
2	D	71	ASN
2	D	111	GLN
2	D	191	HIS
3	E	54	HIS
3	E	59	ASN
3	E	172	HIS
3	H	78	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	A	601	1	14,14,15	0.33	0	17,19,21	1.04	2 (11%)
4	NAG	B	601	1	14,14,15	0.34	0	17,19,21	1.17	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
4	NAG	A	601	1	-	0/6/23/26	0/1/1/1
4	NAG	B	601	1	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	601	NAG	C1-O5-C5	3.99	117.60	112.19
4	A	601	NAG	C1-O5-C5	3.02	116.28	112.19
4	A	601	NAG	C1-C2-N2	-2.21	106.71	110.49

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	601	NAG	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	601	NAG	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, 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	196/276 (71%)	0.02	0 100 100	27, 43, 68, 90	0
1	B	195/276 (70%)	0.47	17 (8%) 10 7	39, 70, 117, 131	0
2	D	212/215 (98%)	0.35	21 (9%) 7 5	29, 63, 111, 125	0
2	L	211/215 (98%)	0.08	2 (0%) 84 84	47, 70, 89, 105	0
3	E	210/228 (92%)	0.31	10 (4%) 30 27	29, 63, 105, 125	0
3	H	218/228 (95%)	0.17	3 (1%) 75 75	49, 77, 108, 120	0
All	All	1242/1438 (86%)	0.23	53 (4%) 35 31	27, 65, 108, 131	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	388	ASN	5.3
1	B	519	HIS	4.4
2	D	183	LEU	4.4
2	D	188	TRP	4.3
1	B	365	TYR	4.1
3	H	197	LEU	4.1
3	E	200	GLN	4.0
1	B	338	PHE	3.9
1	B	384	PRO	3.9
2	D	132	LYS	3.7
3	E	146	LEU	3.5
2	D	195	SER	3.5
2	D	1	GLN	3.5
2	D	158	VAL	3.5
2	D	154	ASP	3.4
1	B	369	TYR	3.4
2	D	133	ALA	3.3
2	D	186	GLU	3.1
2	D	194	TYR	3.1

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Mol	Chain	Res	Type	RSRZ
3	E	130	PHE	3.1
3	E	149	LEU	3.1
1	B	370	ASN	3.1
2	D	131	ASN	3.0
1	B	392	PHE	2.9
3	E	199	THR	2.8
1	B	389	ASP	2.8
2	D	136	VAL	2.7
1	B	377	PHE	2.7
2	D	15	GLY	2.6
2	D	151	TRP	2.6
2	D	193	SER	2.6
2	L	176	ALA	2.5
3	E	134	PRO	2.5
1	B	387	LEU	2.5
2	D	210	ALA	2.5
2	D	127	GLU	2.4
2	D	134	THR	2.4
2	D	181	LEU	2.4
3	E	202	TYR	2.4
3	H	84	TYR	2.4
2	D	135	LEU	2.3
1	B	371	SER	2.2
1	B	515	PHE	2.2
2	L	118	VAL	2.2
3	E	29	ILE	2.2
1	B	518	LEU	2.2
3	E	215	VAL	2.2
3	E	216	ASP	2.2
1	B	386	LYS	2.1
2	D	185	PRO	2.1
1	B	362	VAL	2.1
1	B	527	PRO	2.1
3	H	29	ILE	2.0

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

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	NAG	B	601	14/15	0.62	0.31	86,99,107,107	0
4	NAG	A	601	14/15	0.88	0.20	59,70,79,80	0

6.5 Other polymers [i](#)

There are no such residues in this entry.