



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 10, 2021 – 10:07 AM EDT

PDB ID : 7KLH
Title : SARS-CoV-2 RBD in complex with Fab 15033-7
Authors : Li, Z.; Rini, J.M.
Deposited on : 2020-10-30
Resolution : 3.00 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 : 2.23.1
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) : 2.23.1

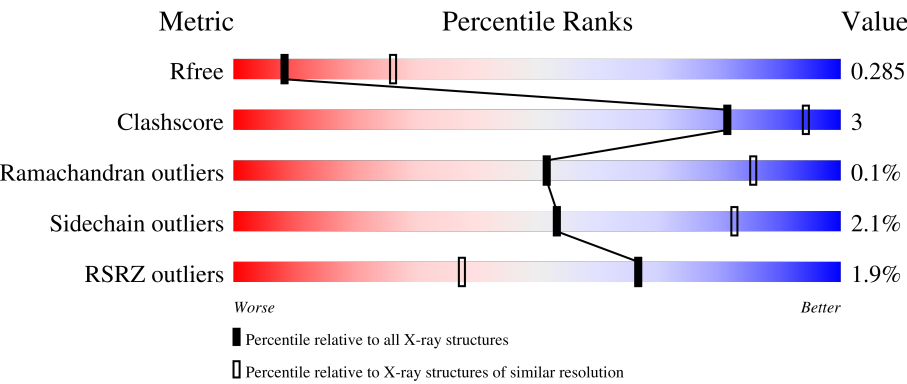
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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.



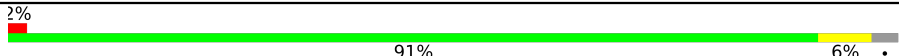
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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	H	225	<div><div>3%</div><div><div></div><div>90%</div><div>9%</div><div>.</div></div></div>
1	I	225	<div><div>3%</div><div><div></div><div>90%</div><div>9%</div><div>.</div></div></div>
2	L	214	<div><div></div><div><div></div><div>90%</div><div>9%</div><div>.</div></div></div>
2	M	214	<div><div></div><div><div></div><div>94%</div><div>6%</div><div>.</div></div></div>
3	A	201	<div><div>2%</div><div><div></div><div>92%</div><div>8%</div><div>.</div></div></div>

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Mol	Chain	Length	Quality of chain
3	B	201	 A horizontal bar chart showing the quality of chain B. The bar is divided into three segments: a small red segment at the beginning labeled '2%', a large green segment in the middle labeled '91%', and a small yellow segment at the end labeled '6%'. A small grey dot is visible at the far right end of the bar.

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9705 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Fab 15033-7 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	223	Total	C	N	O	S	0	0	0
			1635	1031	272	324	8			
1	I	223	Total	C	N	O	S	0	0	0
			1635	1031	272	324	8			

- Molecule 2 is a protein called Fab 15033-7 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	214	Total	C	N	O	S	0	0	0
			1642	1024	276	336	6			
2	M	214	Total	C	N	O	S	0	0	0
			1642	1024	276	336	6			

- Molecule 3 is a protein called Spike glycoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	B	195	Total	C	N	O	S	0	0	0
			1543	989	257	289	8			
3	A	200	Total	C	N	O	S	0	0	0
			1580	1013	264	295	8			

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

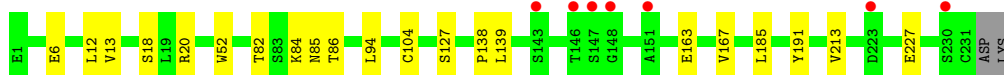
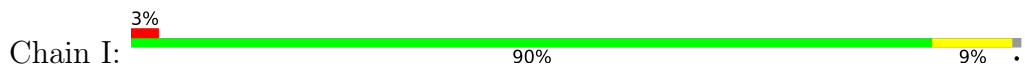
3 Residue-property plots [i](#)

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.

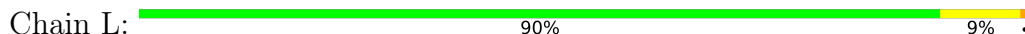
- Molecule 1: Fab 15033-7 heavy chain



- Molecule 1: Fab 15033-7 heavy chain



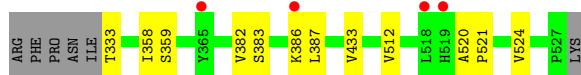
- Molecule 2: Fab 15033-7 light chain



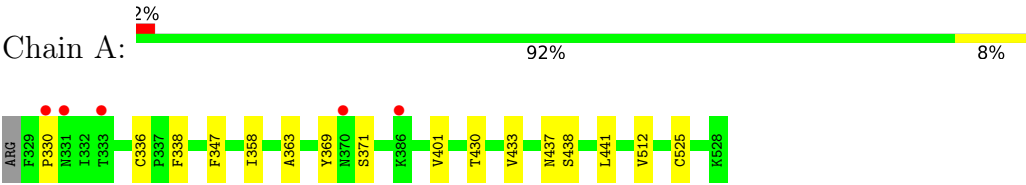
- Molecule 2: Fab 15033-7 light chain



- Molecule 3: Spike glycoprotein



- Molecule 3: Spike glycoprotein



4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, α , β , γ	197.27Å 197.27Å 211.09Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.32 – 3.00 49.32 – 2.88	Depositor EDS
% Data completeness (in resolution range)	99.6 (49.32-3.00) 87.0 (49.32-2.88)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.51 (at 2.86Å)	Xtrriage
Refinement program	PHENIX 1.16rc1_3535	Depositor
R, R_{free}	0.256 , 0.283 0.259 , 0.285	Depositor DCC
R_{free} test set	2760 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	80.5	Xtrriage
Anisotropy	0.016	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 63.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9705	wwPDB-VP
Average B, all atoms (Å ²)	103.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 55.01 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.3659e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	H	0.27	0/1675	0.47	0/2280
1	I	0.28	0/1675	0.48	0/2280
2	L	0.27	0/1677	0.47	0/2276
2	M	0.27	0/1677	0.46	0/2276
3	A	0.28	0/1625	0.43	0/2213
3	B	0.27	0/1587	0.44	0/2161
All	All	0.27	0/9916	0.46	0/13486

There are no bond length outliers.

There are no bond angle outliers.

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	H	1635	0	1586	9	0
1	I	1635	0	1586	10	0
2	L	1642	0	1594	12	0
2	M	1642	0	1594	7	0
3	A	1580	0	1498	7	0
3	B	1543	0	1459	7	0
4	A	14	0	13	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	14	0	13	1	0
All	All	9705	0	9343	50	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 (50) 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:L:156:LEU:HD21	2:L:195:LEU:HD22	1.69	0.73
3:B:433:VAL:HG22	3:B:512:VAL:HG22	1.72	0.70
1:I:82:THR:O	1:I:85:ASN:ND2	2.27	0.68
4:B:2001:NAG:O3	4:B:2001:NAG:H82	2.02	0.60
2:L:128:ARG:NH2	2:L:131:ALA:HB2	2.21	0.56
2:M:36:SER:OG	2:M:37:SER:N	2.39	0.56
2:L:36:SER:OG	2:L:37:SER:N	2.39	0.55
3:B:359:SER:HA	3:B:524:VAL:HG23	1.88	0.54
3:A:433:VAL:HG22	3:A:512:VAL:HG22	1.88	0.54
2:L:76:PHE:CD1	2:L:91:ILE:HD12	2.44	0.53
1:H:13:VAL:HG11	1:H:94:LEU:HD13	1.89	0.53
3:A:338:PHE:HE1	3:A:358:ILE:HD13	1.74	0.53
2:L:128:ARG:NE	2:L:129:THR:O	2.43	0.52
2:L:56:SER:O	2:L:57:ALA:HB3	2.11	0.50
1:I:13:VAL:HG11	1:I:94:LEU:HD13	1.95	0.48
2:M:56:SER:O	2:M:57:ALA:HB3	2.14	0.48
3:B:358:ILE:HG22	3:B:524:VAL:HG21	1.95	0.48
2:L:135:VAL:HG11	2:L:225:VAL:HG11	1.94	0.48
3:B:383:SER:O	3:B:387:LEU:HD21	2.14	0.47
3:B:520:ALA:HB1	3:B:521:PRO:HD2	1.96	0.47
1:I:185:LEU:HD13	1:I:191:TYR:CZ	2.49	0.47
2:M:135:VAL:HG21	2:M:225:VAL:HG11	1.97	0.47
1:I:12:LEU:C	1:I:12:LEU:HD13	2.35	0.47
1:I:52:TRP:CD1	2:M:116:ILE:HD11	2.50	0.47
3:A:336:CYS:HB2	3:A:363:ALA:HB2	1.97	0.47
3:A:437:ASN:OD1	3:A:438:SER:N	2.48	0.47
2:L:120:GLN:OE1	2:L:120:GLN:N	2.47	0.46
1:I:6:GLU:OE2	1:I:104:CYS:N	2.45	0.46
2:M:120:GLN:OE1	2:M:120:GLN:N	2.45	0.46
1:H:52:TRP:CG	2:L:116:ILE:HG12	2.49	0.46
2:L:76:PHE:CE1	2:L:91:ILE:HD12	2.51	0.46
2:M:187:ASP:OD1	2:M:188:SER:N	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:167:VAL:HG23	1:I:213:VAL:HG22	1.99	0.45
1:H:156:LEU:HD12	1:H:194:SER:HB3	1.98	0.45
3:A:430:THR:O	3:A:430:THR:HG23	2.17	0.45
3:B:333:THR:O	3:B:333:THR:HG22	2.17	0.44
1:H:84:LYS:O	1:H:86:THR:HG23	2.17	0.44
1:H:99:THR:HG23	1:H:99:THR:O	2.17	0.44
1:I:52:TRP:CG	2:M:116:ILE:HG12	2.53	0.43
2:L:133:PRO:HB2	2:L:156:LEU:HB2	2.01	0.43
1:H:106:ARG:O	1:H:115:MET:HA	2.18	0.42
1:H:12:LEU:C	1:H:12:LEU:HD13	2.40	0.42
3:A:347:PHE:HB3	3:A:401:VAL:HG23	2.02	0.42
1:H:2:VAL:HG13	1:H:28:PHE:CD1	2.55	0.41
1:H:157:VAL:HB	1:H:193:LEU:HD23	2.02	0.41
3:B:382:VAL:HG22	3:B:387:LEU:HD21	2.02	0.41
2:L:6:GLN:HG3	2:L:23:CYS:SG	2.61	0.41
1:I:84:LYS:O	1:I:86:THR:HG23	2.21	0.41
3:A:330:PRO:HD3	3:A:525:CYS:SG	2.61	0.41
1:I:138:PRO:O	1:I:139:LEU:HD12	2.22	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	H	221/225 (98%)	215 (97%)	6 (3%)	0	100	100
1	I	221/225 (98%)	216 (98%)	5 (2%)	0	100	100
2	L	212/214 (99%)	207 (98%)	5 (2%)	0	100	100
2	M	212/214 (99%)	207 (98%)	5 (2%)	0	100	100
3	A	198/201 (98%)	185 (93%)	13 (7%)	0	100	100
3	B	193/201 (96%)	181 (94%)	11 (6%)	1 (0%)	29	68

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1257/1280 (98%)	1211 (96%)	45 (4%)	1 (0%)	51 85

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	B	386	LYS

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	H	177/179 (99%)	171 (97%)	6 (3%)	37 72
1	I	177/179 (99%)	172 (97%)	5 (3%)	43 77
2	L	189/189 (100%)	184 (97%)	5 (3%)	46 78
2	M	189/189 (100%)	186 (98%)	3 (2%)	62 86
3	A	172/174 (99%)	169 (98%)	3 (2%)	60 85
3	B	168/174 (97%)	168 (100%)	0	100 100
All	All	1072/1084 (99%)	1050 (98%)	22 (2%)	53 82

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	3	GLN
1	H	18	SER
1	H	20	ARG
1	H	116	ASP
1	H	155	CYS
1	H	211	CYS
2	L	12	SER
2	L	36	SER
2	L	77	SER
2	L	128	ARG
2	L	176	SER

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Mol	Chain	Res	Type
1	I	18	SER
1	I	20	ARG
1	I	127	SER
1	I	163	GLU
1	I	227	GLU
2	M	77	SER
2	M	86	ASP
2	M	176	SER
3	A	369	TYR
3	A	371	SER
3	A	441	LEU

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

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	B	2001	3	14,14,15	0.23	0	17,19,21	0.61	0
4	NAG	A	2001	3	14,14,15	0.49	0	17,19,21	1.35	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	B	2001	3	-	4/6/23/26	0/1/1/1
4	NAG	A	2001	3	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	2001	NAG	C1-O5-C5	4.75	118.63	112.19

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	2001	NAG	C8-C7-N2-C2
4	B	2001	NAG	O7-C7-N2-C2
4	B	2001	NAG	O5-C5-C6-O6
4	A	2001	NAG	C4-C5-C6-O6
4	A	2001	NAG	O5-C5-C6-O6
4	B	2001	NAG	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	2001	NAG	1	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	H	223/225 (99%)	-0.09	6 (2%)	54	26	68, 99, 162, 192	0
1	I	223/225 (99%)	-0.03	7 (3%)	49	21	68, 95, 172, 192	0
2	L	214/214 (100%)	-0.20	1 (0%)	91	75	70, 97, 119, 179	0
2	M	214/214 (100%)	-0.07	1 (0%)	91	75	72, 104, 128, 177	0
3	A	200/201 (99%)	-0.14	5 (2%)	57	29	65, 92, 145, 183	0
3	B	195/201 (97%)	-0.10	4 (2%)	63	34	68, 96, 159, 182	0
All	All	1269/1280 (99%)	-0.10	24 (1%)	66	37	65, 99, 153, 192	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	230	SER	4.5
3	A	333	THR	4.3
3	A	330	PRO	4.2
1	I	147	SER	4.0
3	B	518	LEU	3.9
3	B	519	HIS	3.6
1	H	151	ALA	3.2
1	H	231	CYS	3.1
1	I	146	THR	3.0
1	I	151	ALA	2.9
1	I	148	GLY	2.8
3	A	331	ASN	2.7
1	H	225	LYS	2.5
3	A	386	LYS	2.5
3	A	370	ASN	2.4
3	B	386	LYS	2.3
2	L	234	CYS	2.2
1	I	143	SER	2.2
1	H	146	THR	2.2

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Mol	Chain	Res	Type	RSRZ
3	B	365	TYR	2.1
2	M	19	VAL	2.1
1	I	223	ASP	2.1
1	H	209	TYR	2.1
1	H	226	VAL	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	NAG	B	2001	14/15	0.89	0.20	92,132,139,140	0
4	NAG	A	2001	14/15	0.93	0.24	104,115,128,128	0

6.5 Other polymers [i](#)

There are no such residues in this entry.