



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2023 – 07:48 AM EDT

PDB ID : 7M3I
Title : Structure of SARS-CoV-2 spike protein receptor binding domain in complex with a neutralizing antibody, CV2-75 Fab
Authors : Hurlburt, N.K.; Pancera, M.
Deposited on : 2021-03-18
Resolution : 2.80 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 : 2.35.1
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) : 2.35.1

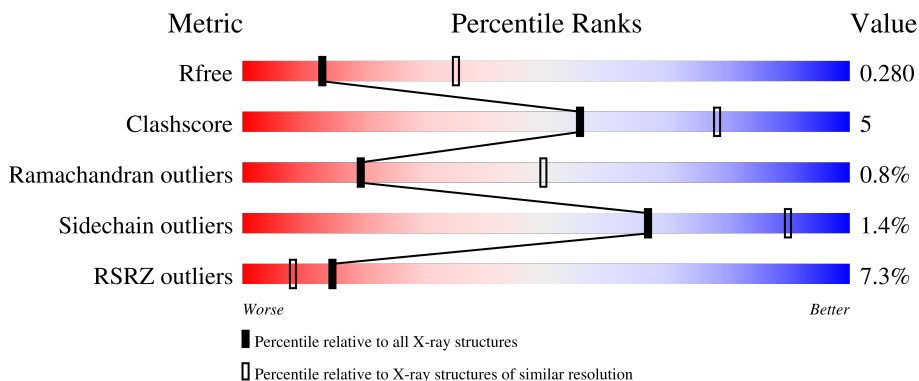
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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.



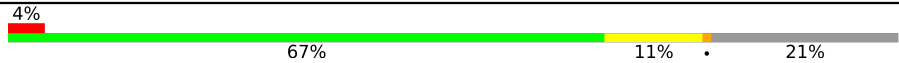
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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	227	
1	H	227	
2	B	214	
2	L	214	
3	C	273	

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Mol	Chain	Length	Quality of chain
3	R	273	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into four segments: a small red segment at the beginning labeled '4%', a large green segment labeled '67%', a small yellow segment labeled '11%', and a grey segment at the end labeled '21%'. The segments are separated by thin white lines.</p>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 9872 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 CV2-75 Fab Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	221	Total 1658	C 1054	N 266	O 334	S 4	0	0	0
1	H	205	Total 1561	C 996	N 249	O 312	S 4	0	0	0

- Molecule 2 is a protein called CV2-75 Fab Light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	211	Total 1585	C 990	N 267	O 324	S 4	0	0	0
2	L	211	Total 1585	C 990	N 267	O 324	S 4	0	0	0

- Molecule 3 is a protein called Spike protein S1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	217	Total 1718	C 1099	N 290	O 321	S 8	0	0	0
3	R	217	Total 1721	C 1102	N 290	O 321	S 8	0	0	0

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



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

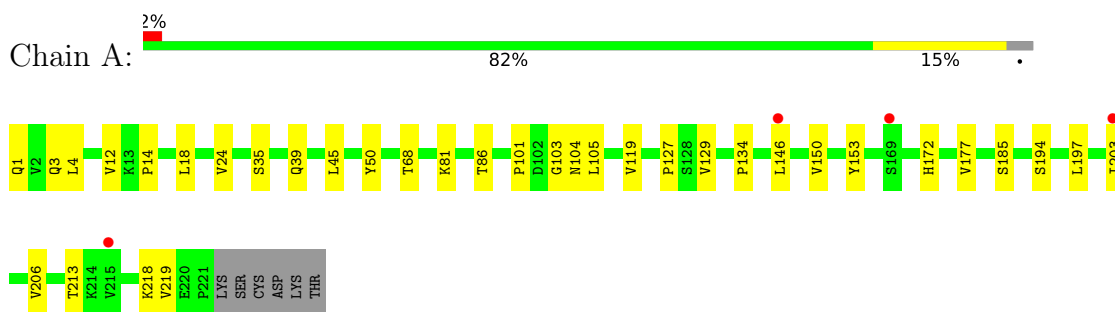
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	O	0	0
			1	1		
5	C	7	Total	O	0	0
			7	7		
5	H	2	Total	O	0	0
			2	2		
5	L	2	Total	O	0	0
			2	2		
5	R	4	Total	O	0	0
			4	4		

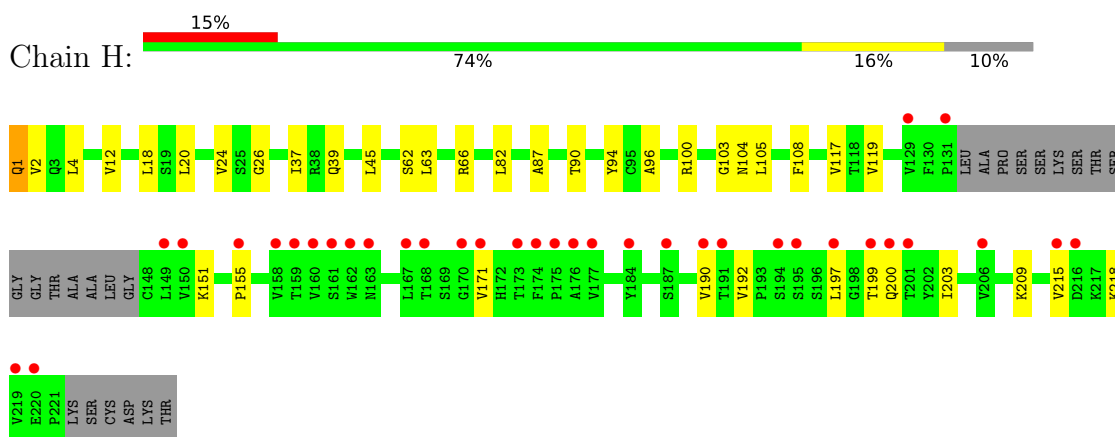
3 Residue-property plots [i](#)

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.

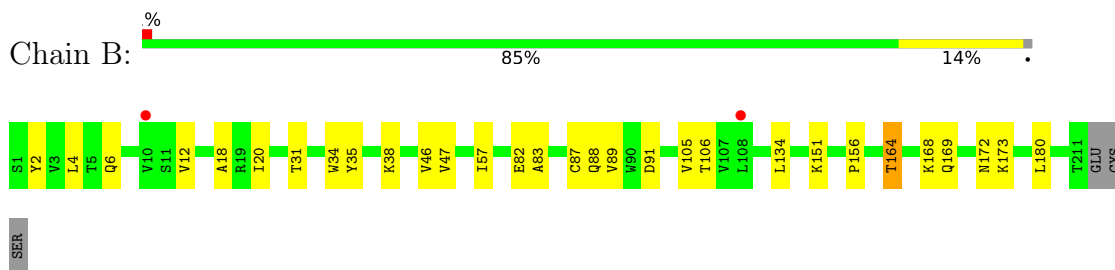
- Molecule 1: CV2-75 Fab Heavy chain



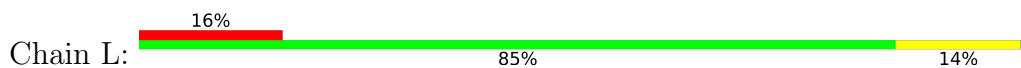
- Molecule 1: CV2-75 Fab Heavy chain

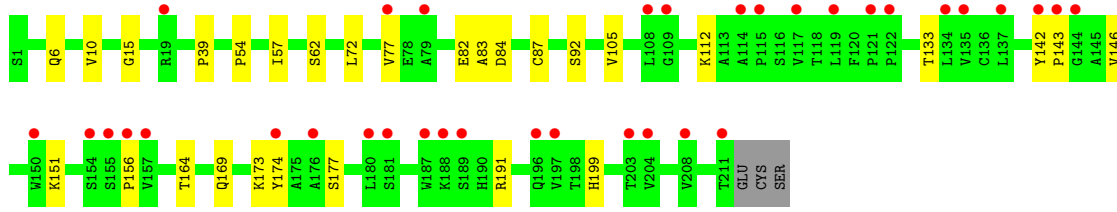


- Molecule 2: CV2-75 Fab Light chain

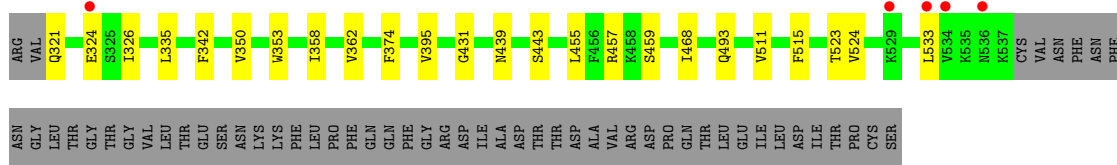


- Molecule 2: CV2-75 Fab Light chain

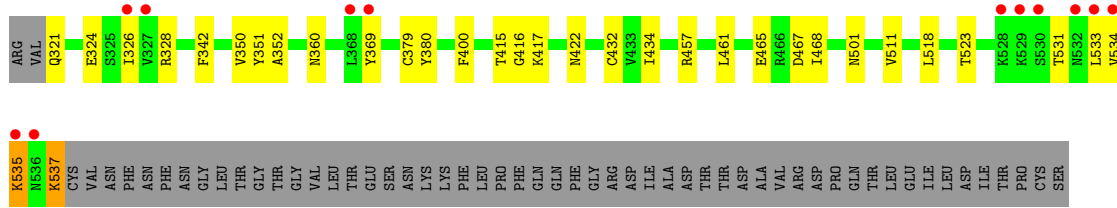




• Molecule 3: Spike protein S1



• Molecule 3: Spike protein S1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	86.37Å 127.52Å 155.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.32 – 2.80 49.32 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.5 (49.32-2.80) 99.5 (49.32-2.80)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 2.81Å)	Xtrriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.245 , 0.281 0.245 , 0.280	Depositor DCC
R_{free} test set	2092 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	79.6	Xtrriage
Anisotropy	0.433	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 61.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9872	wwPDB-VP
Average B, all atoms (Å ²)	101.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.20% 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, PCA

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	0.28	0/1696	0.49	0/2323
1	H	0.27	0/1597	0.52	1/2187 (0.0%)
2	B	0.26	0/1626	0.46	0/2227
2	L	0.26	0/1626	0.47	0/2227
3	C	0.29	0/1765	0.48	0/2401
3	R	0.28	0/1768	0.48	0/2405
All	All	0.27	0/10078	0.48	1/13770 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	H	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	1	PCA	O-C-N	-7.33	110.97	122.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	H	1	PCA	Mainchain

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	1658	0	1614	19	0
1	H	1561	0	1513	17	0
2	B	1585	0	1523	17	0
2	L	1585	0	1523	17	0
3	C	1718	0	1645	14	0
3	R	1721	0	1654	18	0
4	C	14	0	13	1	0
4	R	14	0	13	0	0
5	B	1	0	0	0	0
5	C	7	0	0	0	0
5	H	2	0	0	0	0
5	L	2	0	0	0	0
5	R	4	0	0	0	0
All	All	9872	0	9498	98	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 5.

All (98) 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:A:177:VAL:HG12	2:B:164:THR:HG23	1.63	0.80
3:R:457:ARG:NH1	3:R:467:ASP:OD2	2.23	0.72
2:L:54:PRO:HG2	2:L:57:ILE:HD13	1.75	0.69
1:A:129:VAL:HG22	1:A:150:VAL:HG12	1.75	0.67
1:A:203:ILE:HD13	1:A:218:LYS:HB3	1.77	0.66
1:A:150:VAL:HG11	1:A:206:VAL:HG21	1.79	0.64
2:B:82:GLU:HG3	2:B:106:THR:HA	1.83	0.60
1:H:203:ILE:HG22	1:H:218:LYS:HB3	1.81	0.60
4:C:600:NAG:H83	4:C:600:NAG:H3	1.85	0.59
3:C:321:GLN:HG3	3:C:324:GLU:HG3	1.85	0.59
1:H:151:LYS:HE3	2:L:133:THR:HG21	1.84	0.59
3:C:439:ASN:O	3:C:443:SER:OG	2.20	0.58
1:H:103:GLY:O	1:H:105:LEU:N	2.36	0.58
2:B:151:LYS:HA	2:B:156:PRO:HA	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:4:LEU:HD22	1:H:24:VAL:HG22	1.86	0.57
1:A:103:GLY:O	1:A:105:LEU:N	2.37	0.57
3:C:358:ILE:HB	3:C:395:VAL:HG23	1.86	0.56
3:C:468:ILE:HD12	3:R:518:LEU:HD21	1.86	0.56
2:L:151:LYS:HA	2:L:156:PRO:HA	1.88	0.56
2:L:169:GLN:NE2	2:L:173:LYS:HB3	2.20	0.55
3:R:352:ALA:HB2	3:R:468:ILE:HG22	1.89	0.55
2:B:168:LYS:HE2	2:B:172:ASN:HA	1.89	0.55
3:R:342:PHE:CE2	3:R:511:VAL:HG11	2.43	0.54
3:R:417:LYS:O	3:R:422:ASN:ND2	2.39	0.54
2:B:12:VAL:HG21	2:B:18:ALA:HB2	1.89	0.53
3:C:523:THR:HG23	3:C:524:VAL:HG23	1.90	0.53
3:R:321:GLN:HG3	3:R:324:GLU:HG3	1.89	0.53
1:A:134:PRO:HG3	1:A:197:LEU:HD11	1.89	0.53
2:L:173:LYS:HG3	2:L:174:TYR:H	1.73	0.53
2:B:34:TRP:HB2	2:B:47:VAL:HB	1.90	0.52
1:A:86:THR:O	1:A:119:VAL:HG21	2.09	0.52
2:L:112:LYS:HG3	2:L:143:PRO:HG2	1.92	0.52
1:H:2:VAL:HA	1:H:26:GLY:HA3	1.93	0.51
1:A:12:VAL:HG21	1:A:18:LEU:HD13	1.92	0.51
1:H:39:GLN:HB2	1:H:45:LEU:HD23	1.91	0.51
3:C:335:LEU:HD23	3:C:362:VAL:HG13	1.93	0.51
3:R:360:ASN:H	3:R:523:THR:HB	1.75	0.51
1:H:18:LEU:HD12	1:H:117:VAL:HG11	1.92	0.50
3:R:461:LEU:HD22	3:R:465:GLU:HB3	1.92	0.50
2:L:173:LYS:HG3	2:L:174:TYR:N	2.27	0.49
3:C:431:GLY:HA2	3:C:515:PHE:CD2	2.46	0.49
1:H:66:ARG:HH21	1:H:82:LEU:HD11	1.78	0.49
1:A:39:GLN:HB2	1:A:45:LEU:HD23	1.93	0.49
1:A:129:VAL:HG21	1:A:206:VAL:HG11	1.93	0.49
1:H:18:LEU:HD21	1:H:20:LEU:HD13	1.95	0.47
1:A:4:LEU:HD22	1:A:24:VAL:HG12	1.97	0.47
2:B:4:LEU:HD11	2:B:89:VAL:HG12	1.97	0.46
2:B:46:VAL:HA	2:B:57:ILE:HG12	1.98	0.46
3:C:431:GLY:HA2	3:C:515:PHE:HD2	1.81	0.46
2:L:39:PRO:HD3	2:L:82:GLU:OE2	2.16	0.46
1:H:96:ALA:HB1	1:H:108:PHE:HB3	1.98	0.46
1:A:127:PRO:HB3	1:A:153:TYR:HB3	1.97	0.46
1:A:146:LEU:HD13	1:A:219:VAL:HG21	1.97	0.45
1:H:171:VAL:HA	1:H:190:VAL:HG13	1.99	0.45
3:R:537:LYS:HA	3:R:537:LYS:HD2	1.63	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:146:VAL:HG12	2:L:199:HIS:HB2	1.99	0.45
2:L:39:PRO:HD3	2:L:82:GLU:OE1	2.16	0.45
3:R:379:CYS:HA	3:R:432:CYS:HA	1.98	0.45
1:A:127:PRO:HD2	1:A:213:THR:HG21	1.99	0.45
3:C:350:VAL:O	3:C:353:TRP:HD1	1.99	0.44
1:H:100:ARG:NH1	1:H:105:LEU:HB3	2.33	0.44
3:C:326:ILE:HD12	3:C:326:ILE:HA	1.88	0.44
1:H:87:ALA:O	1:H:90:THR:HG22	2.16	0.44
1:H:18:LEU:HB3	1:H:82:LEU:HB3	1.99	0.43
2:B:35:TYR:HE1	2:B:88:GLN:HB3	1.83	0.43
2:B:20:ILE:HD11	2:B:105:VAL:HG11	2.01	0.42
2:B:169:GLN:HG2	2:B:173:LYS:O	2.19	0.42
3:R:535:LYS:HB2	3:R:535:LYS:HE2	1.72	0.42
1:H:12:VAL:O	1:H:119:VAL:HA	2.20	0.42
2:L:10:VAL:O	2:L:105:VAL:HA	2.19	0.42
2:L:62:SER:O	2:L:72:LEU:HD12	2.19	0.42
3:R:415:THR:HG22	3:R:416:GLY:H	1.85	0.42
3:C:457:ARG:NH1	3:C:459:SER:O	2.51	0.42
2:L:15:GLY:N	2:L:77:VAL:O	2.52	0.42
2:B:38:LYS:HG2	2:B:83:ALA:HB2	2.02	0.42
1:A:68:THR:HB	1:A:81:LYS:HB2	2.00	0.42
2:B:2:TYR:HB2	2:B:91:ASP:OD2	2.20	0.42
2:L:92:SER:HB2	3:R:380:TYR:CD2	2.55	0.42
3:R:350:VAL:HA	3:R:400:PHE:HB2	2.01	0.42
3:R:434:ILE:HB	3:R:511:VAL:HG13	2.00	0.42
1:H:37:ILE:O	1:H:94:TYR:N	2.42	0.41
3:R:533:LEU:HD23	3:R:533:LEU:HA	1.95	0.41
2:L:6:GLN:NE2	2:L:87:CYS:SG	2.93	0.41
1:H:62:SER:HB2	1:H:63:LEU:HD12	2.02	0.41
2:B:31:THR:O	2:B:89:VAL:HG23	2.20	0.41
3:C:455:LEU:HD22	3:C:493:GLN:HG3	2.02	0.41
1:A:18:LEU:O	1:A:81:LYS:HA	2.21	0.41
2:B:6:GLN:NE2	2:B:87:CYS:SG	2.94	0.41
2:L:83:ALA:HB2	2:L:105:VAL:O	2.21	0.41
3:R:326:ILE:HD12	3:R:326:ILE:HA	1.85	0.41
1:A:35:SER:HA	1:A:50:TYR:HA	2.03	0.41
2:B:12:VAL:HG11	2:B:18:ALA:HA	2.02	0.41
2:B:134:LEU:HB2	2:B:180:LEU:HB3	2.02	0.41
1:A:177:VAL:HG22	1:A:185:SER:O	2.20	0.41
3:C:342:PHE:CE2	3:C:511:VAL:HG11	2.55	0.41
1:A:101:PRO:HB3	3:C:374:PHE:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:164:THR:HG22	2:L:177:SER:H	1.86	0.40
3:R:501:ASN:HD22	3:R:501:ASN:HA	1.76	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	219/227 (96%)	204 (93%)	12 (6%)	3 (1%)	11	34
1	H	201/227 (88%)	187 (93%)	11 (6%)	3 (2%)	10	33
2	B	209/214 (98%)	198 (95%)	11 (5%)	0	100	100
2	L	209/214 (98%)	189 (90%)	17 (8%)	3 (1%)	11	34
3	C	215/273 (79%)	198 (92%)	16 (7%)	1 (0%)	29	61
3	R	215/273 (79%)	202 (94%)	13 (6%)	0	100	100
All	All	1268/1428 (89%)	1178 (93%)	80 (6%)	10 (1%)	19	49

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	155	PRO
1	H	200	GLN
2	L	142	TYR
2	L	191	ARG
1	A	104	ASN
1	A	194	SER
1	H	104	ASN
3	C	533	LEU
2	L	84	ASP
1	A	14	PRO

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	191/197 (97%)	189 (99%)	2 (1%)	76	93
1	H	181/197 (92%)	176 (97%)	5 (3%)	43	77
2	B	179/182 (98%)	178 (99%)	1 (1%)	86	96
2	L	179/182 (98%)	179 (100%)	0	100	100
3	C	189/240 (79%)	189 (100%)	0	100	100
3	R	190/240 (79%)	183 (96%)	7 (4%)	34	68
All	All	1109/1238 (90%)	1094 (99%)	15 (1%)	67	90

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

Mol	Chain	Res	Type
1	A	3	GLN
1	A	172	HIS
2	B	164	THR
1	H	192	VAL
1	H	197	LEU
1	H	199	THR
1	H	209	LYS
1	H	215	VAL
3	R	328	ARG
3	R	351	TYR
3	R	369	TYR
3	R	531	THR
3	R	534	VAL
3	R	535	LYS
3	R	537	LYS

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](#)

2 non-standard protein/DNA/RNA residues 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
1	PCA	A	1	1	7,8,9	1.80	1 (14%)	9,10,12	2.10	5 (55%)
1	PCA	H	1	1	7,8,9	1.82	1 (14%)	9,10,12	2.17	5 (55%)

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
1	PCA	A	1	1	-	0/0/11/13	0/1/1/1
1	PCA	H	1	1	-	0/0/11/13	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	1	PCA	CD-N	4.69	1.47	1.34
1	A	1	PCA	CD-N	4.64	1.46	1.34

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	1	PCA	CA-N-CD	-3.12	102.91	113.58
1	A	1	PCA	CA-N-CD	-3.08	103.02	113.58
1	H	1	PCA	OE-CD-CG	-3.05	121.44	126.76
1	A	1	PCA	OE-CD-CG	-2.91	121.68	126.76
1	H	1	PCA	CB-CA-N	2.49	110.46	103.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1	PCA	CB-CA-N	2.49	110.45	103.30
1	H	1	PCA	CG-CD-N	2.40	114.61	108.39
1	A	1	PCA	CB-CA-C	-2.39	109.42	112.70
1	A	1	PCA	CG-CD-N	2.37	114.52	108.39
1	H	1	PCA	CB-CA-C	-2.31	109.52	112.70

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	C	600	3	14,14,15	0.51	0	17,19,21	1.28	2 (11%)
4	NAG	R	600	3	14,14,15	0.35	0	17,19,21	0.52	0

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	C	600	3	-	5/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	R	600	3	-	2/6/23/26	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(°)
4	C	600	NAG	C2-N2-C7	4.26	128.97	122.90
4	C	600	NAG	C1-C2-N2	2.31	114.44	110.49

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	600	NAG	C8-C7-N2-C2
4	C	600	NAG	O7-C7-N2-C2
4	R	600	NAG	O5-C5-C6-O6
4	C	600	NAG	O5-C5-C6-O6
4	C	600	NAG	C4-C5-C6-O6
4	R	600	NAG	C4-C5-C6-O6
4	C	600	NAG	C3-C2-N2-C7

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	600	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	220/227 (96%)	0.19	4 (1%) 68 61	59, 94, 157, 229	0
1	H	204/227 (89%)	0.77	35 (17%) 1 1	58, 103, 178, 188	0
2	B	211/214 (98%)	0.02	2 (0%) 84 80	62, 100, 138, 157	0
2	L	211/214 (98%)	0.78	35 (16%) 1 1	66, 136, 178, 196	0
3	C	217/273 (79%)	0.13	5 (2%) 60 51	47, 77, 134, 160	0
3	R	217/273 (79%)	0.10	12 (5%) 25 16	53, 79, 150, 199	0
All	All	1280/1428 (89%)	0.33	93 (7%) 15 8	47, 94, 168, 229	0

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	155	SER	7.8
2	L	117	VAL	6.6
1	H	191	THR	6.2
2	L	176	ALA	5.6
1	H	162	TRP	5.5
3	C	534	VAL	5.2
3	R	533	LEU	4.7
1	H	160	VAL	4.4
3	R	536	ASN	4.4
1	H	195	SER	4.3
2	L	156	PRO	4.3
1	H	206	VAL	4.3
2	L	157	VAL	4.3
1	H	219	VAL	4.3
2	L	196	GLN	4.1
1	H	187	SER	4.1
2	L	144	GLY	3.9
3	R	530	SER	3.9
1	H	159	THR	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	169	SER	3.8
2	L	137	LEU	3.8
1	H	167	LEU	3.7
2	L	154	SER	3.7
1	H	201	THR	3.7
3	R	532	ASN	3.6
1	H	158	VAL	3.6
3	R	327	VAL	3.6
1	H	184	TYR	3.6
1	H	131	PRO	3.6
1	H	149	LEU	3.6
2	B	108	LEU	3.5
2	L	208	VAL	3.4
2	L	211	THR	3.4
2	L	150	TRP	3.4
1	H	220	GLU	3.3
2	L	122	PRO	3.2
2	L	203	THR	3.2
1	H	176	ALA	3.2
2	L	143	PRO	3.2
1	H	170	GLY	3.1
1	H	197	LEU	3.0
1	H	200	GLN	3.0
1	H	175	PRO	3.0
2	L	142	TYR	3.0
2	L	134	LEU	3.0
1	H	216	ASP	3.0
2	L	197	VAL	3.0
3	R	534	VAL	3.0
1	H	215	VAL	2.9
2	L	108	LEU	2.8
1	H	199	THR	2.8
2	L	109	GLY	2.8
2	L	174	TYR	2.8
1	H	190	VAL	2.8
3	R	529	LYS	2.8
2	L	114	ALA	2.8
2	L	135	VAL	2.7
2	L	181	SER	2.7
1	A	215	VAL	2.7
2	L	119	LEU	2.6
1	H	129	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	H	171	VAL	2.5
3	R	535	LYS	2.5
2	L	79	ALA	2.5
1	H	155	PRO	2.5
2	L	115	PRO	2.5
1	H	168	THR	2.5
1	H	163	ASN	2.4
3	R	326	ILE	2.4
2	L	187	TRP	2.4
2	L	180	LEU	2.4
2	L	189	SER	2.4
1	H	174	PHE	2.3
3	C	533	LEU	2.3
1	H	150	VAL	2.3
3	C	529	LYS	2.3
3	R	369	TYR	2.3
1	A	203	ILE	2.3
1	H	194	SER	2.3
2	L	19	ARG	2.3
3	R	368	LEU	2.3
2	L	188	LYS	2.3
1	A	146	LEU	2.3
3	C	536	ASN	2.2
3	R	528	LYS	2.2
2	L	204	VAL	2.2
2	L	77	VAL	2.2
2	L	121	PRO	2.1
3	C	324	GLU	2.1
1	H	173	THR	2.0
1	H	177	VAL	2.0
1	H	161	SER	2.0
2	B	10	VAL	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	PCA	A	1	8/9	0.74	0.28	128,145,147,148	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
1	PCA	H	1	8/9	0.81	0.20	123,127,132,141	0

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	R	600	14/15	0.87	0.18	90,108,118,118	0
4	NAG	C	600	14/15	0.89	0.21	68,94,108,111	0

6.5 Other polymers [i](#)

There are no such residues in this entry.