



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 2, 2022 – 08:05 PM EST

PDB ID : 7S5Q
Title : Crystal structure of SARS-CoV-2 B.1.351 variant receptor binding domain in complex with neutralizing antibodies CS44 and COVA1-16
Authors : Yuan, M.; Wilson, I.A.
Deposited on : 2021-09-11
Resolution : 2.88 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	:	2.27
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)	:	2.27

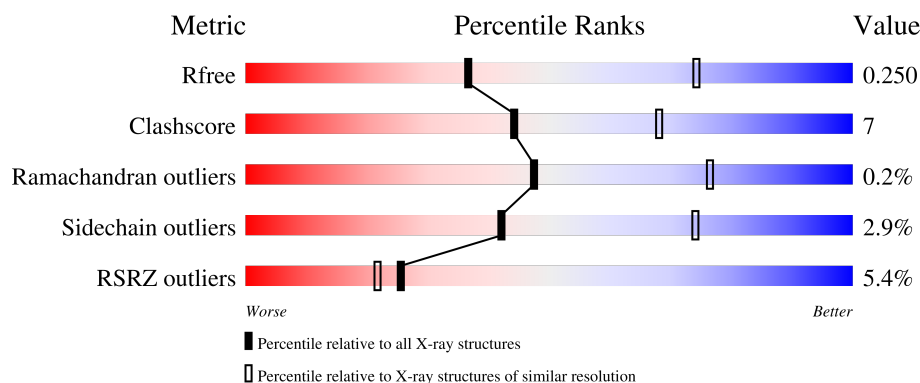
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.88 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	2691 (2.90-2.86)
Clashscore	141614	2947 (2.90-2.86)
Ramachandran outliers	138981	2868 (2.90-2.86)
Sidechain outliers	138945	2871 (2.90-2.86)
RSRZ outliers	127900	2629 (2.90-2.86)

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	231	<div> <div>%</div> <div> <div></div> <div>78%</div> <div>7%</div> <div>14%</div> </div> </div>
2	B	226	<div> <div>10%</div> <div> <div></div> <div>73%</div> <div>23%</div> <div>..</div> </div> </div>
3	C	215	<div> <div>13%</div> <div> <div></div> <div>79%</div> <div>20%</div> <div>.</div> </div> </div>
4	H	232	<div> <div>%</div> <div> <div></div> <div>82%</div> <div>14%</div> <div>..</div> </div> </div>
5	L	214	<div> <div>%</div> <div> <div></div> <div>84%</div> <div>14%</div> <div>..</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	NAG	A	601	-	-	-	X

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 8323 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	198	Total	C	N	O	S	0	0	0
			1572	1011	262	291	8			

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	417	ASN	LYS	variant	UNP P0DTC2
A	484	LYS	GLU	variant	UNP P0DTC2
A	501	TYR	ASN	variant	UNP P0DTC2
A	542	SER	-	expression tag	UNP P0DTC2
A	543	GLY	-	expression tag	UNP P0DTC2
A	544	HIS	-	expression tag	UNP P0DTC2
A	545	HIS	-	expression tag	UNP P0DTC2
A	546	HIS	-	expression tag	UNP P0DTC2
A	547	HIS	-	expression tag	UNP P0DTC2
A	548	HIS	-	expression tag	UNP P0DTC2
A	549	HIS	-	expression tag	UNP P0DTC2

- Molecule 2 is a protein called CS44 Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	219	Total	C	N	O	S	0	0	0
			1628	1022	275	322	9			

- Molecule 3 is a protein called CS44 Light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	214	Total	C	N	O	S	4	2	0
			1650	1034	278	333	5			

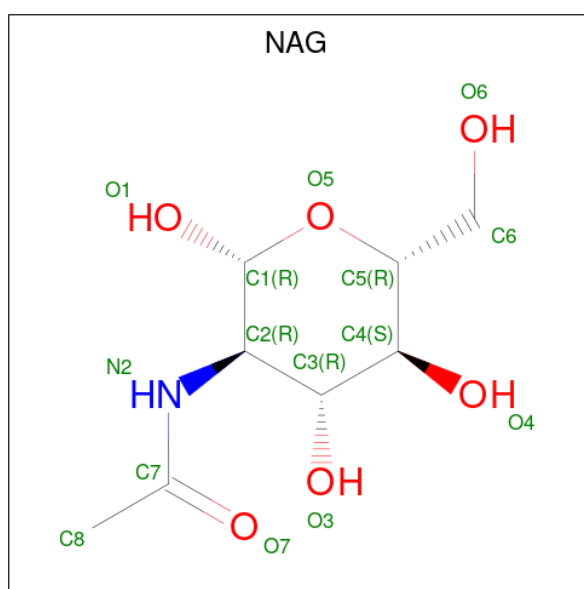
- Molecule 4 is a protein called COVA1-16 Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	225	Total	C	N	O	S	0	0	0
			1715	1081	292	334	8			

- Molecule 5 is a protein called COVA1-16 Light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	L	212	Total	C	N	O	S	0	0	0
			1631	1019	272	336	4			

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



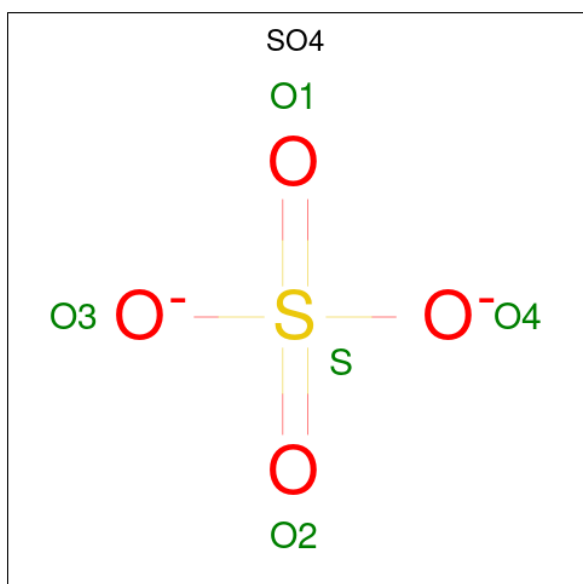
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 7 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

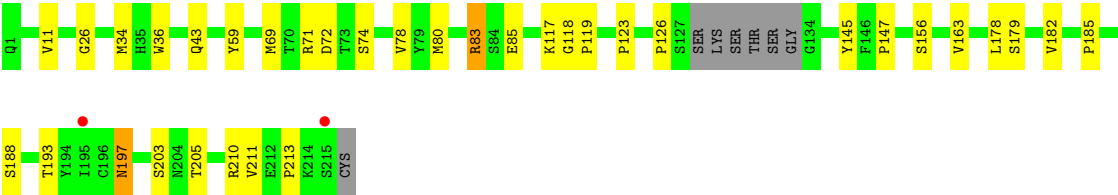
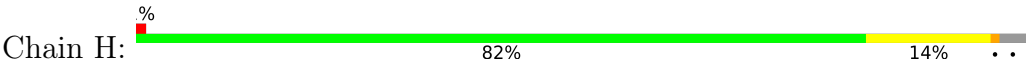
- Molecule 8 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



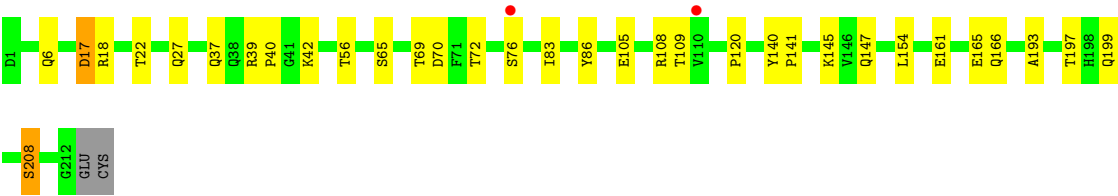
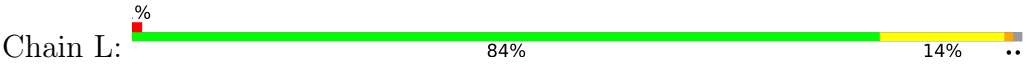
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total O S 5 4 1	0	0
8	A	1	Total O S 5 4 1	0	0
8	A	1	Total O S 5 4 1	0	0
8	A	1	Total O S 5 4 1	0	0
8	A	1	Total O S 5 4 1	0	0
8	A	1	Total O S 5 4 1	0	0
8	B	1	Total O S 5 4 1	0	0
8	H	1	Total O S 5 4 1	0	0
8	H	1	Total O S 5 4 1	0	0
8	H	1	Total O S 5 4 1	0	0
8	L	1	Total O S 5 4 1	0	0

- Molecule 9 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	16	Total O 16 16	0	0
9	B	1	Total O 1 1	0	0
9	C	1	Total O 1 1	0	0
9	H	13	Total O 13 13	0	0
9	L	7	Total O 7 7	0	0



● Molecule 5: COVA1-16 Light chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	89.64Å 120.97Å 137.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.56 – 2.88 40.56 – 2.88	Depositor EDS
% Data completeness (in resolution range)	98.8 (40.56-2.88) 98.7 (40.56-2.88)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 2.90Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.216 , 0.257 0.211 , 0.250	Depositor DCC
R_{free} test set	1687 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	56.9	Xtriage
Anisotropy	0.125	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 46.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8323	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.90% 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: SO4, EDO, 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	0.31	0/1617	0.56	1/2200 (0.0%)
2	B	0.32	0/1665	0.64	0/2271
3	C	0.30	0/1693	0.63	1/2300 (0.0%)
4	H	0.31	0/1759	0.59	1/2395 (0.0%)
5	L	0.31	0/1666	0.60	0/2266
All	All	0.31	0/8400	0.60	3/11432 (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	A	466	ARG	CG-CD-NE	5.79	123.96	111.80
4	H	83	ARG	CB-CG-CD	-5.14	98.23	111.60
3	C	135	LEU	CB-CG-CD1	-5.04	102.43	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1572	0	1499	10	0
2	B	1628	0	1587	34	0
3	C	1650	0	1605	28	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	1715	0	1658	19	0
5	L	1631	0	1575	22	0
6	A	14	0	13	1	0
7	A	8	0	12	0	0
7	H	4	0	6	1	0
7	L	8	0	12	2	0
8	A	30	0	0	0	0
8	B	5	0	0	0	0
8	H	15	0	0	0	0
8	L	5	0	0	0	0
9	A	16	0	0	0	0
9	B	1	0	0	0	0
9	C	1	0	0	0	0
9	H	13	0	0	0	0
9	L	7	0	0	1	0
All	All	8323	0	7967	107	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 7.

All (107) 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)
5:L:39:ARG:HH11	5:L:42:LYS:HZ1	1.11	0.90
2:B:181:VAL:HG21	3:C:135:LEU:HD11	1.66	0.77
5:L:140:TYR:CD1	5:L:141:PRO:HA	2.21	0.74
4:H:34:MET:HG2	4:H:78:VAL:HG21	1.71	0.73
4:H:11:VAL:HB	4:H:147:PRO:HG3	1.69	0.72
3:C:39:LYS:NZ	3:C:81:GLU:O	2.23	0.72
2:B:171:GLN:O	2:B:173:SER:N	2.25	0.70
4:H:26:GLY:H	7:H:301:EDO:H21	1.56	0.69
3:C:35:TRP:HB2	3:C:48:ILE:HB	1.74	0.69
5:L:22:THR:HG22	5:L:72:THR:HG22	1.74	0.69
2:B:55:GLY:HA3	2:B:71:ARG:HH21	1.58	0.68
5:L:83:ILE:HD11	5:L:105:GLU:HA	1.77	0.67
5:L:40:PRO:HB3	5:L:165:GLU:HG3	1.76	0.66
3:C:140:TYR:CD1	3:C:141:PRO:HA	2.31	0.66
5:L:17:ASP:N	5:L:17:ASP:OD1	2.30	0.64
4:H:156:SER:H	4:H:197:ASN:HD21	1.46	0.64
2:B:13:LYS:O	2:B:16:THR:OG1	2.16	0.63
3:C:47:LEU:HD23	3:C:58:ILE:HD12	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:121:VAL:HG21	2:B:207:VAL:HG21	1.82	0.61
2:B:151:THR:HG22	2:B:199:ASN:HB3	1.83	0.60
5:L:6:GLN:NE2	5:L:86:TYR:O	2.34	0.60
5:L:145:LYS:HB3	5:L:197:THR:HB	1.83	0.60
5:L:39:ARG:NH1	5:L:42:LYS:HZ1	1.93	0.59
2:B:123:PRO:HD3	2:B:209:LYS:HD3	1.84	0.58
3:C:8:PRO:O	3:C:102:THR:HG23	2.04	0.58
3:C:120:PRO:HD3	3:C:132:VAL:HG22	1.85	0.58
2:B:27:ILE:HG23	2:B:97:CYS:H	1.69	0.57
2:B:119:PRO:HB3	2:B:145:TYR:HB3	1.85	0.57
3:C:108:ARG:HB3	3:C:140:TYR:CD2	2.41	0.56
4:H:72:ASP:OD1	4:H:74:SER:OG	2.23	0.56
4:H:119:PRO:HB3	4:H:145:TYR:HB3	1.89	0.55
2:B:38:ARG:HB3	2:B:48:ILE:HD11	1.89	0.54
2:B:85:GLU:OE2	2:B:85:GLU:HA	2.07	0.54
3:C:31:SER:HA	3:C:50:GLY:HA2	1.89	0.54
2:B:144:ASP:OD1	2:B:171:GLN:NE2	2.38	0.54
4:H:178:LEU:HA	7:L:301:EDO:H22	1.89	0.53
5:L:39:ARG:HB2	5:L:42:LYS:HE2	1.90	0.53
2:B:112:SER:HG	2:B:146:PHE:HZ	1.54	0.53
3:C:117:ILE:HG12	3:C:209:PHE:HD1	1.71	0.53
2:B:36:TRP:CE2	2:B:80:MET:HB2	2.44	0.52
2:B:55:GLY:CA	2:B:71:ARG:HH21	2.23	0.52
2:B:164:HIS:CE1	3:C:174:SER:HG	2.27	0.52
3:C:128:GLY:HA2	3:C:183:LYS:HE3	1.90	0.52
2:B:2:VAL:HG22	2:B:26:GLY:HA3	1.92	0.52
4:H:59:TYR:HE1	4:H:69:MET:HB2	1.75	0.51
5:L:108:ARG:NH2	5:L:109:THR:OG1	2.43	0.51
4:H:83:ARG:HG3	4:H:85:GLU:H	1.75	0.51
4:H:83:ARG:HG3	4:H:85:GLU:OE1	2.11	0.51
5:L:105:GLU:OE2	5:L:166:GLN:OE1	2.29	0.50
4:H:185:PRO:O	4:H:188:SER:OG	2.28	0.50
3:C:81:GLU:H	3:C:81:GLU:CD	2.14	0.50
1:A:431:GLY:HA2	1:A:515:PHE:CE2	2.47	0.50
5:L:193:ALA:HB2	5:L:208:SER:HB3	1.92	0.50
2:B:181:VAL:HG11	3:C:135:LEU:HD12	1.92	0.50
4:H:36:TRP:CE2	4:H:80:MET:HB2	2.46	0.50
3:C:61:ARG:HB2	3:C:76:SER:O	2.12	0.50
2:B:46:GLU:OE2	2:B:62:LYS:HE3	2.13	0.49
5:L:161:GLU:HG3	9:L:401:HOH:O	2.13	0.49
2:B:61:GLN:HE22	2:B:64:GLN:HE22	1.62	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:96:TYR:HB2	2:B:102:ILE:HD11	1.96	0.47
4:H:126:PRO:HD2	4:H:213:PRO:HA	1.97	0.47
5:L:108:ARG:HG3	5:L:140:TYR:CD2	2.49	0.47
2:B:87:THR:HG23	2:B:110:THR:HA	1.97	0.47
4:H:193:THR:HG23	4:H:210:ARG:NE	2.30	0.47
1:A:431:GLY:HA2	1:A:515:PHE:CD2	2.49	0.46
2:B:181:VAL:HG21	3:C:135:LEU:CD1	2.38	0.46
3:C:189:HIS:O	3:C:211:ARG:NE	2.48	0.46
3:C:186:TYR:O	3:C:192:TYR:OH	2.31	0.46
2:B:48:ILE:HG21	2:B:80:MET:HE2	1.98	0.46
3:C:79:GLU:HB3	3:C:81:GLU:OE1	2.17	0.45
2:B:41:ARG:HE	2:B:41:ARG:HB3	1.60	0.45
5:L:154:LEU:H	5:L:154:LEU:HD12	1.81	0.45
2:B:44:ARG:NH2	3:C:100:GLN:HG3	2.32	0.45
2:B:63:PHE:O	2:B:67:VAL:HG12	2.17	0.44
5:L:108:ARG:NE	5:L:109:THR:O	2.47	0.44
2:B:46:GLU:OE2	2:B:62:LYS:CE	2.66	0.44
2:B:151:THR:CG2	2:B:199:ASN:HB3	2.48	0.44
5:L:69:THR:HG22	5:L:70:ASP:OD1	2.19	0.43
2:B:170:LEU:HD22	2:B:175:LEU:O	2.18	0.43
3:C:24:ARG:HG2	3:C:69:THR:O	2.18	0.43
5:L:40:PRO:CB	5:L:165:GLU:HG3	2.47	0.43
1:A:403:ARG:HD3	1:A:495:TYR:CE1	2.52	0.43
3:C:21:LEU:HD23	3:C:102:THR:HB	2.01	0.43
5:L:18:ARG:HG2	5:L:76:SER:HA	2.00	0.43
3:C:135:LEU:HD21	3:C:137:ASN:HB2	2.01	0.42
3:C:189:HIS:O	3:C:211:ARG:NH2	2.50	0.42
4:H:117:LYS:HD3	4:H:118:GLY:N	2.33	0.42
3:C:54:ARG:NH1	3:C:62:PHE:O	2.51	0.42
3:C:85:VAL:HG22	3:C:103:LYS:HG2	2.01	0.42
2:B:199:ASN:OD1	2:B:206:LYS:HG2	2.18	0.42
4:H:163:VAL:HG22	4:H:182:VAL:HB	2.02	0.42
5:L:37:GLN:HB2	5:L:86:TYR:CE2	2.54	0.42
1:A:415:THR:O	5:L:56:THR:HG21	2.20	0.42
1:A:376:THR:HB	1:A:435:ALA:HB3	2.01	0.41
1:A:444:LYS:HG3	1:A:448:ASN:HB2	2.02	0.41
3:C:158:ASN:OD1	3:C:158:ASN:N	2.52	0.41
4:H:203:SER:OG	4:H:205:THR:OG1	2.27	0.41
1:A:342:PHE:HB2	6:A:601:NAG:H82	2.03	0.41
1:A:439:ASN:O	1:A:443:SER:HB2	2.20	0.41
1:A:517:LEU:HD12	1:A:517:LEU:HA	1.83	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:146:PHE:HA	2:B:147:PRO:HA	1.90	0.41
2:B:87:THR:HA	2:B:109:VAL:O	2.21	0.41
1:A:387:LEU:HD23	1:A:387:LEU:HA	1.97	0.40
2:B:170:LEU:HB2	2:B:176:TYR:CE2	2.57	0.40
3:C:24:ARG:HE	3:C:24:ARG:HB3	1.38	0.40
4:H:123:PRO:HB3	4:H:211:VAL:HG12	2.04	0.40
4:H:179:SER:OG	7:L:301:EDO:H21	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	A	196/231 (85%)	190 (97%)	6 (3%)	0	100	100
2	B	215/226 (95%)	208 (97%)	5 (2%)	2 (1%)	17	45
3	C	214/215 (100%)	208 (97%)	6 (3%)	0	100	100
4	H	221/232 (95%)	216 (98%)	5 (2%)	0	100	100
5	L	210/214 (98%)	202 (96%)	8 (4%)	0	100	100
All	All	1056/1118 (94%)	1024 (97%)	30 (3%)	2 (0%)	47	76

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	172	SER
2	B	99	GLY

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	171/203 (84%)	169 (99%)	2 (1%)	71	89
2	B	184/192 (96%)	177 (96%)	7 (4%)	33	65
3	C	187/186 (100%)	179 (96%)	8 (4%)	29	60
4	H	190/198 (96%)	187 (98%)	3 (2%)	62	85
5	L	187/189 (99%)	180 (96%)	7 (4%)	34	66
All	All	919/968 (95%)	892 (97%)	27 (3%)	42	74

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	377	PHE
1	A	493	GLN
2	B	44	ARG
2	B	84	SER
2	B	111	VAL
2	B	128	SER
2	B	177	SER
2	B	197	ASN
2	B	199	ASN
3	C	13	LEU
3	C	33	LEU
3	C	47	LEU
3	C	65	SER
3	C	83	PHE
3	C	156	SER
3	C	176	SER
3	C	197	THR
4	H	43	GLN
4	H	71	ARG
4	H	197	ASN
5	L	17	ASP
5	L	27	GLN
5	L	65	SER

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Mol	Chain	Res	Type
5	L	120	PRO
5	L	147	GLN
5	L	199	GLN
5	L	208	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	64	GLN
2	B	164	HIS
3	C	138	ASN
3	C	210	ASN
4	H	197	ASN

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](#)

17 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
7	EDO	H	301	-	3,3,3	0.47	0	2,2,2	0.32	0
8	SO4	L	303	-	4,4,4	0.13	0	6,6,6	0.09	0
6	NAG	A	601	1	14,14,15	0.45	0	17,19,21	0.44	0
7	EDO	L	301	-	3,3,3	0.50	0	2,2,2	0.09	0
8	SO4	A	605	-	4,4,4	0.17	0	6,6,6	0.25	0
8	SO4	A	606	-	4,4,4	0.12	0	6,6,6	0.15	0
8	SO4	A	604	-	4,4,4	0.16	0	6,6,6	0.17	0
7	EDO	A	602	-	3,3,3	0.56	0	2,2,2	0.18	0
8	SO4	A	609	-	4,4,4	0.15	0	6,6,6	0.08	0
8	SO4	B	301	-	4,4,4	0.15	0	6,6,6	0.08	0
8	SO4	A	608	-	4,4,4	0.13	0	6,6,6	0.11	0
8	SO4	A	607	-	4,4,4	0.13	0	6,6,6	0.17	0
7	EDO	L	302	-	3,3,3	0.48	0	2,2,2	0.27	0
8	SO4	H	304	-	4,4,4	0.13	0	6,6,6	0.14	0
8	SO4	H	303	-	4,4,4	0.14	0	6,6,6	0.11	0
7	EDO	A	603	-	3,3,3	0.54	0	2,2,2	0.13	0
8	SO4	H	302	-	4,4,4	0.15	0	6,6,6	0.06	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
7	EDO	H	301	-	-	0/1/1/1	-
6	NAG	A	601	1	-	2/6/23/26	0/1/1/1
7	EDO	L	301	-	-	0/1/1/1	-
7	EDO	A	602	-	-	1/1/1/1	-
7	EDO	L	302	-	-	0/1/1/1	-
7	EDO	A	603	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	601	NAG	O5-C5-C6-O6
6	A	601	NAG	C4-C5-C6-O6
7	A	602	EDO	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	H	301	EDO	1	0
6	A	601	NAG	1	0
7	L	301	EDO	2	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	198/231 (85%)	-0.15	2 (1%) 82 82	19, 34, 60, 81	0
2	B	219/226 (96%)	0.62	23 (10%) 6 4	42, 78, 107, 126	0
3	C	214/215 (99%)	0.80	29 (13%) 3 2	57, 84, 106, 119	0
4	H	225/232 (96%)	-0.03	2 (0%) 84 84	23, 47, 98, 109	0
5	L	212/214 (99%)	0.07	2 (0%) 84 84	31, 53, 74, 86	0
All	All	1068/1118 (95%)	0.27	58 (5%) 25 22	19, 59, 99, 126	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	196	CYS	4.4
2	B	211	VAL	4.4
2	B	185	PRO	4.2
3	C	181	LEU	4.0
2	B	208	ASP	3.7
2	B	212	GLU	3.7
2	B	207	VAL	3.6
2	B	128	SER	3.6
3	C	182	SER	3.6
1	A	333	THR	3.6
3	C	76	SER	3.5
3	C	183	LYS	3.4
3	C	130	ALA	3.4
2	B	127	SER	3.4
2	B	134	GLY	3.1
3	C	199	GLN	3.1
2	B	193	THR	3.1
2	B	117	LYS	3.0
3	C	196	VAL	2.9
2	B	184	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	332	ILE	2.8
4	H	215	SER	2.8
2	B	186	SER	2.8
3	C	129	THR	2.8
3	C	109	THR	2.7
3	C	18	ARG	2.7
3	C	16	GLY	2.7
2	B	187	SER	2.6
3	C	4	LEU	2.5
2	B	11	VAL	2.5
3	C	154	LEU	2.5
2	B	15	GLY	2.4
3	C	20	THR	2.4
3	C	77	ARG	2.4
3	C	104	VAL	2.4
4	H	195	ILE	2.4
2	B	135	THR	2.4
3	C	127	SER	2.3
2	B	13	LYS	2.3
3	C	106	ILE	2.3
2	B	194	TYR	2.3
3	C	126	LYS	2.3
5	L	76	SER	2.2
3	C	197	THR	2.2
3	C	188	LYS	2.2
2	B	125	ALA	2.1
3	C	7	SER	2.1
3	C	24	ARG	2.1
3	C	194[A]	CYS	2.1
3	C	11	LEU	2.1
3	C	120	PRO	2.1
2	B	154	TRP	2.1
3	C	131	SER	2.1
3	C	213	GLU	2.1
2	B	155	ASN	2.0
2	B	82(C)	LEU	2.0
5	L	110	VAL	2.0
3	C	122	ASP	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 ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

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
6	NAG	A	601	14/15	0.77	0.46	77,87,97,100	0
8	SO4	A	609	5/5	0.85	0.21	80,84,93,102	5
7	EDO	L	302	4/4	0.87	0.30	52,60,63,65	0
7	EDO	A	602	4/4	0.87	0.42	39,47,47,50	0
8	SO4	A	607	5/5	0.88	0.33	45,52,66,70	5
7	EDO	L	301	4/4	0.90	0.24	46,46,58,63	0
8	SO4	A	608	5/5	0.91	0.21	56,63,68,82	5
8	SO4	A	605	5/5	0.91	0.26	29,36,41,55	5
8	SO4	B	301	5/5	0.91	0.21	53,58,61,62	5
8	SO4	H	303	5/5	0.91	0.25	69,72,93,97	5
8	SO4	H	304	5/5	0.91	0.26	42,50,56,66	5
8	SO4	L	303	5/5	0.92	0.23	82,83,98,101	5
8	SO4	A	606	5/5	0.94	0.26	37,38,39,49	5
7	EDO	A	603	4/4	0.94	0.21	38,43,44,45	0
7	EDO	H	301	4/4	0.95	0.11	46,48,49,50	0
8	SO4	H	302	5/5	0.96	0.12	31,42,54,65	5
8	SO4	A	604	5/5	0.97	0.16	37,41,51,54	0

6.5 Other polymers ⓘ

There are no such residues in this entry.