



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:24 AM UTC

PDB ID : 8I5I / pdb_00008i5i
Title : Crystal structure of SARS-CoV-2 delta variant spike receptor-binding domain (RBD) in complex with NCV2SG53 Fab
Authors : Yamamoto, A.; Higashiura, A.
Deposited on : 2023-01-25
Resolution : 3.06 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	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

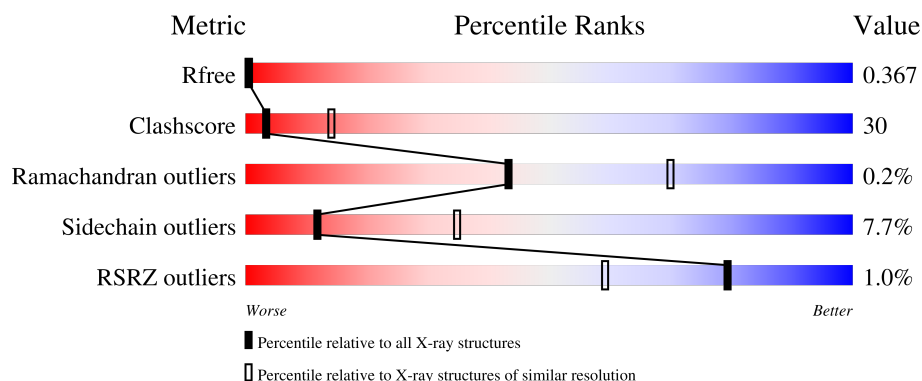
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

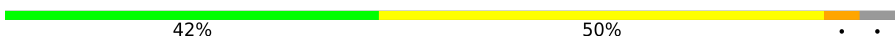
The reported resolution of this entry is 3.06 Å.

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}	180053	2469 (3.10-3.02)
Clashscore	190562	2569 (3.10-3.02)
Ramachandran outliers	187476	2424 (3.10-3.02)
Sidechain outliers	187428	2423 (3.10-3.02)
RSRZ outliers	180081	2469 (3.10-3.02)

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	R	224	
2	H	227	
3	L	214	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4780 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	R	191	Total	C	N	O	S	0	0	0
			1522	976	256	282	8			

There are 10 discrepancies between the modelled and reference sequences:

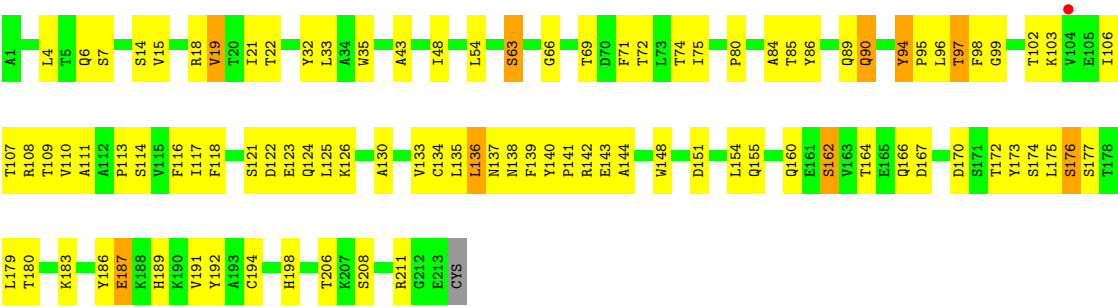
Chain	Residue	Modelled	Actual	Comment	Reference
R	452	ARG	LEU	variant	UNP P0DTC2
R	478	LYS	THR	variant	UNP P0DTC2
R	537	THR	-	expression tag	UNP P0DTC2
R	538	GLY	-	expression tag	UNP P0DTC2
R	539	HIS	-	expression tag	UNP P0DTC2
R	540	HIS	-	expression tag	UNP P0DTC2
R	541	HIS	-	expression tag	UNP P0DTC2
R	542	HIS	-	expression tag	UNP P0DTC2
R	543	HIS	-	expression tag	UNP P0DTC2
R	544	HIS	-	expression tag	UNP P0DTC2

- Molecule 2 is a protein called Fab Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	218	Total	C	N	O	S	0	0	0
			1639	1039	273	318	9			

- Molecule 3 is a protein called Fab Light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	213	Total	C	N	O	S	0	0	0
			1619	1016	270	329	4			



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	77.61Å 77.61Å 519.15Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	67.21 – 3.06 67.21 – 3.06	Depositor EDS
% Data completeness (in resolution range)	39.3 (67.21-3.06) 39.3 (67.21-3.06)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.56 (at 3.07Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.228 , 0.372 0.232 , 0.367	Depositor DCC
R_{free} test set	386 reflections (2.03%)	wwPDB-VP
Wilson B-factor (Å ²)	77.2	Xtriage
Anisotropy	0.233	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 112.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4780	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.61% 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](#)

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	R	0.43	0/1565	0.65	0/2127
2	H	0.44	0/1678	0.63	0/2281
3	L	0.43	0/1653	0.64	0/2246
All	All	0.43	0/4896	0.64	0/6654

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	R	1522	0	1445	119	0
2	H	1639	0	1617	97	0
3	L	1619	0	1588	77	0
All	All	4780	0	4650	285	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 30.

All (285) 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:R:452:ARG:HG2	1:R:494:SER:HA	1.40	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:33:TYR:HB2	2:H:100:PHE:HB2	1.49	0.93
1:R:383:SER:HB2	1:R:386:LYS:HD3	1.52	0.92
1:R:360:ASN:H	1:R:523:THR:HB	1.35	0.91
2:H:134:ALA:HB1	2:H:222:PRO:HA	1.53	0.90
1:R:418:ILE:HA	1:R:422:ASN:HD22	1.37	0.90
2:H:148:GLY:HA3	2:H:190:VAL:HG12	1.58	0.86
1:R:354:ASN:HD21	1:R:356:LYS:HE2	1.40	0.86
2:H:172:VAL:HA	2:H:191:VAL:HG22	1.58	0.84
3:L:90:GLN:NE2	3:L:97:THR:OG1	2.09	0.84
1:R:349:SER:HB3	1:R:452:ARG:H	1.45	0.82
2:H:135:PRO:HG2	2:H:222:PRO:HG3	1.62	0.81
2:H:39:GLN:HB2	2:H:45:LEU:HG	1.61	0.80
2:H:51:ILE:HB	2:H:70:ILE:HG12	1.66	0.77
1:R:392:PHE:HA	1:R:517:LEU:HD13	1.65	0.76
2:H:90:ASP:O	2:H:94:TYR:OH	2.04	0.76
2:H:145:ALA:H	2:H:195:SER:HB3	1.51	0.76
2:H:150:LEU:HA	2:H:188:SER:HB3	1.68	0.74
2:H:62:ASP:OD1	2:H:65:LYS:NZ	2.20	0.74
1:R:417:LYS:O	1:R:422:ASN:ND2	2.21	0.73
1:R:454:ARG:NH1	1:R:456:PHE:O	2.22	0.72
2:H:83:MET:HB3	2:H:86:LEU:HD21	1.73	0.71
2:H:155:PHE:HB2	2:H:184:LEU:HD23	1.72	0.70
2:H:209:HIS:HD1	2:H:212:SER:HG	1.39	0.70
2:H:87:ARG:HD2	2:H:89:GLU:HG2	1.73	0.70
3:L:110:VAL:HG22	3:L:141:PRO:HD3	1.73	0.69
2:H:12:VAL:HG11	2:H:18:LEU:HB2	1.75	0.68
3:L:80:PRO:HA	3:L:106:ILE:HG12	1.75	0.68
1:R:353:TRP:CD1	1:R:466:ARG:HD3	2.29	0.68
1:R:357:ARG:HG3	1:R:396:TYR:HE1	1.60	0.67
3:L:192:TYR:O	3:L:208:SER:OG	2.13	0.67
1:R:359:SER:HA	1:R:524:VAL:HG22	1.78	0.66
1:R:441:LEU:HD23	1:R:509:ARG:HH12	1.61	0.66
3:L:151:ASP:HA	3:L:191:VAL:HB	1.76	0.66
2:H:83:MET:HE1	2:H:118:VAL:HG21	1.78	0.65
2:H:117:LEU:HD11	2:H:119:THR:HG23	1.77	0.65
1:R:409:GLN:HA	1:R:414:GLN:HG2	1.79	0.65
2:H:70:ILE:HD11	2:H:79:LEU:HD11	1.78	0.65
1:R:454:ARG:HH21	1:R:469:SER:HB3	1.62	0.64
3:L:63:SER:HB2	3:L:74:THR:HB	1.78	0.64
1:R:392:PHE:O	1:R:523:THR:OG1	2.15	0.62
1:R:401:VAL:HG22	1:R:509:ARG:HD2	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:49:SER:HB2	2:H:60:TYR:HA	1.81	0.62
3:L:6:GLN:HB2	3:L:99:GLY:HA3	1.82	0.62
3:L:108:ARG:HG3	3:L:109:THR:O	1.99	0.62
2:H:128:PRO:HD2	2:H:214:THR:HB	1.82	0.61
3:L:111:ALA:HB3	3:L:140:TYR:N	2.15	0.61
1:R:393:THR:HB	1:R:517:LEU:HA	1.82	0.61
3:L:135:LEU:HD11	3:L:137:ASN:HB2	1.82	0.61
3:L:18:ARG:HG3	3:L:75:ILE:O	2.00	0.61
1:R:357:ARG:NH1	1:R:359:SER:OG	2.34	0.61
2:H:144:THR:HG23	2:H:195:SER:H	1.67	0.60
3:L:151:ASP:N	3:L:191:VAL:O	2.29	0.60
1:R:497:PHE:CE2	1:R:507:PRO:HB3	2.37	0.60
2:H:163:TRP:HZ3	2:H:203:TYR:HB3	1.64	0.60
3:L:139:PHE:HE2	3:L:142:ARG:HA	1.67	0.60
1:R:338:PHE:O	1:R:340:GLU:N	2.35	0.59
1:R:461:LEU:HB2	1:R:465:GLU:HB3	1.84	0.59
2:H:177:ALA:HA	2:H:187:LEU:HB3	1.84	0.59
3:L:186:TYR:O	3:L:192:TYR:OH	2.14	0.59
3:L:187:GLU:CD	3:L:211:ARG:HD2	2.27	0.59
3:L:167:ASP:HB2	3:L:172:THR:O	2.03	0.59
2:H:128:PRO:HB3	2:H:154:TYR:HB3	1.86	0.58
3:L:162:SER:O	3:L:175:LEU:HG	2.03	0.58
3:L:116:PHE:H	3:L:135:LEU:HD23	1.68	0.58
1:R:419:ALA:O	1:R:424:LYS:HD2	2.04	0.58
1:R:476:GLY:N	1:R:487:ASN:HB3	2.18	0.58
2:H:73:ASP:OD1	2:H:76:LYS:N	2.37	0.58
1:R:397:ALA:HA	1:R:512:VAL:O	2.03	0.57
2:H:37:ILE:HG13	2:H:47:TRP:HA	1.86	0.57
3:L:113:PRO:HA	3:L:139:PHE:HB3	1.86	0.57
1:R:474:GLN:HG2	1:R:476:GLY:O	2.05	0.57
3:L:85:THR:HG22	3:L:103:LYS:HG3	1.87	0.57
1:R:503:VAL:HA	1:R:506:GLN:HG3	1.86	0.56
1:R:367:VAL:HG13	1:R:370:ASN:HD21	1.70	0.56
2:H:98:ARG:O	2:H:110:PRO:HA	2.06	0.56
2:H:148:GLY:HA2	2:H:163:TRP:CZ2	2.39	0.56
2:H:148:GLY:HA2	2:H:163:TRP:HZ2	1.71	0.56
1:R:350:VAL:HA	1:R:400:PHE:HB2	1.87	0.56
1:R:418:ILE:HD13	1:R:422:ASN:ND2	2.19	0.56
1:R:439:ASN:HB2	1:R:506:GLN:HB3	1.88	0.56
2:H:137:SER:O	2:H:137:SER:OG	2.18	0.56
1:R:431:GLY:HA3	1:R:513:LEU:O	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:175:PHE:HB2	2:H:188:SER:O	2.07	0.55
3:L:166:GLN:HG3	3:L:173:TYR:CE1	2.41	0.55
3:L:137:ASN:OD1	3:L:138:ASN:N	2.39	0.55
1:R:424:LYS:NZ	1:R:425:LEU:H	2.05	0.54
2:H:52:SER:O	2:H:72:ARG:NH1	2.39	0.54
3:L:144:ALA:HB2	3:L:198:HIS:HD2	1.71	0.54
1:R:391:CYS:HA	1:R:524:VAL:O	2.08	0.54
2:H:14:PRO:HG3	2:H:120:VAL:HG12	1.88	0.54
2:H:105:ILE:HD12	2:H:106:ILE:HG23	1.87	0.54
2:H:128:PRO:HG3	2:H:209:HIS:HB2	1.90	0.54
1:R:441:LEU:CD2	1:R:509:ARG:HH12	2.21	0.54
2:H:180:GLN:OE1	2:H:186:SER:HB3	2.08	0.53
3:L:139:PHE:CE2	3:L:142:ARG:HA	2.43	0.53
1:R:349:SER:OG	1:R:451:TYR:HA	2.09	0.53
2:H:49:SER:OG	2:H:50:TYR:N	2.41	0.53
2:H:158:PRO:O	2:H:209:HIS:HD2	1.91	0.53
1:R:383:SER:OG	1:R:386:LYS:HB2	2.09	0.52
2:H:13:ARG:HG2	2:H:122:SER:HA	1.91	0.52
1:R:361:CYS:SG	1:R:362:VAL:N	2.82	0.52
2:H:149:CYS:O	2:H:188:SER:HB2	2.10	0.52
2:H:162:SER:HB3	2:H:206:ASN:HB2	1.91	0.52
1:R:357:ARG:HG3	1:R:396:TYR:CE1	2.44	0.52
2:H:32:TYR:HB2	2:H:34:MET:HE2	1.90	0.52
2:H:99:GLU:OE1	3:L:94:TYR:OH	2.26	0.52
2:H:148:GLY:HA3	2:H:190:VAL:HA	1.91	0.52
1:R:390:LEU:O	1:R:525:CYS:HA	2.10	0.52
3:L:143:GLU:N	3:L:143:GLU:OE1	2.39	0.52
1:R:438:SER:HB2	1:R:509:ARG:HG3	1.92	0.51
2:H:60:TYR:HD2	2:H:66:GLY:H	1.58	0.51
1:R:494:SER:H	2:H:56:SER:HB3	1.75	0.51
3:L:148:TRP:CZ3	3:L:194:CYS:HB2	2.45	0.51
2:H:204:ILE:HD13	2:H:219:LYS:HA	1.91	0.51
1:R:336:CYS:HB3	1:R:362:VAL:O	2.10	0.51
1:R:347:PHE:HB2	1:R:401:VAL:HG23	1.93	0.51
1:R:385:THR:O	1:R:388:ASN:ND2	2.44	0.50
3:L:148:TRP:CE2	3:L:179:LEU:HB2	2.46	0.50
1:R:497:PHE:CD2	1:R:507:PRO:HB3	2.46	0.50
2:H:191:VAL:HG12	2:H:192:THR:O	2.12	0.50
1:R:403:ARG:HG2	1:R:497:PHE:HE1	1.76	0.50
3:L:125:LEU:HD23	3:L:130:ALA:HB2	1.94	0.50
2:H:178:VAL:HG21	3:L:160:GLN:HB3	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:125:LEU:HD21	3:L:183:LYS:HG3	1.94	0.50
1:R:338:PHE:HE2	1:R:368:LEU:HD13	1.77	0.49
2:H:201:GLN:HB3	2:H:203:TYR:CE1	2.46	0.49
2:H:178:VAL:CG2	3:L:160:GLN:HB3	2.42	0.49
2:H:31:ASP:HB3	2:H:101:ASP:OD1	2.12	0.49
3:L:187:GLU:OE2	3:L:211:ARG:HD2	2.13	0.49
1:R:438:SER:HB2	1:R:509:ARG:CG	2.43	0.49
2:H:103:THR:HB	2:H:104:LYS:HZ3	1.78	0.49
1:R:384:PRO:HD2	1:R:386:LYS:NZ	2.28	0.49
1:R:455:LEU:HD22	1:R:493:GLN:HG3	1.93	0.49
1:R:454:ARG:HA	1:R:492:LEU:HD23	1.95	0.49
2:H:204:ILE:CD1	2:H:219:LYS:HA	2.41	0.49
1:R:402:ILE:HD12	1:R:406:GLU:O	2.12	0.48
2:H:32:TYR:O	2:H:72:ARG:NH1	2.37	0.48
1:R:424:LYS:HZ3	1:R:425:LEU:H	1.61	0.48
3:L:148:TRP:CD1	3:L:179:LEU:HD13	2.48	0.48
1:R:387:LEU:HD23	1:R:392:PHE:HZ	1.78	0.48
3:L:66:GLY:HA3	3:L:71:PHE:CD1	2.48	0.48
1:R:359:SER:HB3	1:R:523:THR:HB	1.95	0.48
3:L:122:ASP:OD1	3:L:123:GLU:N	2.47	0.48
2:H:132:PRO:HG3	2:H:218:LYS:HG3	1.95	0.48
3:L:194:CYS:O	3:L:206:THR:HG23	2.14	0.48
1:R:361:CYS:HB3	1:R:524:VAL:HG13	1.95	0.48
3:L:136:LEU:HB3	3:L:139:PHE:CE1	2.49	0.48
2:H:58:MET:HB2	2:H:70:ILE:HG21	1.94	0.48
1:R:345:THR:O	1:R:509:ARG:NH2	2.47	0.48
1:R:438:SER:HB3	1:R:507:PRO:O	2.13	0.48
3:L:189:HIS:O	3:L:211:ARG:NE	2.42	0.47
2:H:181:SER:C	2:H:183:GLY:H	2.22	0.47
1:R:384:PRO:O	1:R:386:LYS:HD2	2.14	0.47
2:H:154:TYR:HB2	2:H:209:HIS:CE1	2.49	0.47
2:H:101:ASP:OD2	2:H:104:LYS:NZ	2.45	0.47
3:L:111:ALA:N	3:L:140:TYR:O	2.46	0.47
1:R:396:TYR:O	1:R:513:LEU:HA	2.15	0.47
1:R:445:VAL:O	1:R:498:GLN:HG2	2.15	0.47
1:R:357:ARG:HH22	1:R:359:SER:HB2	1.80	0.46
1:R:403:ARG:HB3	1:R:504:GLY:O	2.15	0.46
3:L:89:GLN:HG2	3:L:90:GLN:O	2.15	0.46
1:R:376:THR:CG2	1:R:435:ALA:HB3	2.46	0.46
1:R:487:ASN:HA	1:R:489:TYR:HE1	1.81	0.46
2:H:78:SER:HB2	2:H:80:PHE:CZ	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:138:ASN:N	3:L:173:TYR:O	2.49	0.46
1:R:425:LEU:HD23	1:R:425:LEU:HA	1.65	0.46
2:H:203:TYR:O	2:H:204:ILE:HD13	2.14	0.46
3:L:118:PHE:HB2	3:L:133:VAL:HG23	1.98	0.46
1:R:433:VAL:HG12	1:R:512:VAL:HG22	1.97	0.46
1:R:468:ILE:H	1:R:468:ILE:HG13	1.47	0.46
1:R:338:PHE:CE1	1:R:363:ALA:HA	2.50	0.46
1:R:346:ARG:HH21	1:R:450:ASN:HB3	1.81	0.46
3:L:19:VAL:H	3:L:75:ILE:HG22	1.80	0.46
1:R:347:PHE:CE2	1:R:509:ARG:HB3	2.50	0.46
2:H:12:VAL:HG23	2:H:120:VAL:HG22	1.98	0.46
3:L:111:ALA:HB3	3:L:140:TYR:H	1.81	0.46
1:R:346:ARG:HG2	1:R:451:TYR:OH	2.17	0.45
2:H:35:SER:OG	2:H:49:SER:O	2.23	0.45
1:R:378:LYS:HG2	1:R:379:CYS:H	1.79	0.45
3:L:84:ALA:HB3	3:L:86:TYR:CE1	2.51	0.45
1:R:395:VAL:HG13	1:R:515:PHE:CE1	2.51	0.45
2:H:164:ASN:CB	2:H:168:LEU:HD13	2.47	0.45
1:R:350:VAL:HG23	1:R:400:PHE:CD1	2.52	0.45
1:R:387:LEU:HD23	1:R:392:PHE:CZ	2.52	0.45
2:H:152:LYS:HB3	2:H:152:LYS:HE3	1.54	0.45
3:L:116:PHE:H	3:L:135:LEU:HB3	1.82	0.45
2:H:6:GLU:OE2	2:H:95:TYR:HA	2.17	0.45
3:L:170:ASP:O	3:L:172:THR:HG23	2.17	0.45
2:H:29:PHE:CD1	2:H:34:MET:HE3	2.52	0.45
2:H:32:TYR:HA	2:H:100:PHE:O	2.16	0.45
3:L:32:TYR:N	3:L:32:TYR:CD1	2.85	0.45
1:R:377:PHE:HD1	1:R:433:VAL:O	1.99	0.44
1:R:494:SER:N	2:H:56:SER:HB3	2.32	0.44
1:R:520:ALA:HB1	1:R:521:PRO:HD2	1.99	0.44
2:H:34:MET:HB2	2:H:51:ILE:HG22	1.99	0.44
2:H:97:ALA:HA	2:H:111:TYR:O	2.17	0.44
1:R:398:ASP:OD1	1:R:398:ASP:N	2.50	0.44
3:L:164:THR:CG2	3:L:174:SER:HB2	2.47	0.44
1:R:347:PHE:HB2	1:R:401:VAL:CG2	2.48	0.44
3:L:122:ASP:O	3:L:126:LYS:HG2	2.18	0.44
3:L:179:LEU:HD12	3:L:180:THR:H	1.82	0.44
1:R:390:LEU:HD12	1:R:391:CYS:H	1.83	0.44
2:H:98:ARG:HE	2:H:111:TYR:HB2	1.83	0.44
1:R:448:ASN:C	1:R:450:ASN:H	2.26	0.44
2:H:153:ASP:HA	2:H:184:LEU:HB3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:136:LEU:HB2	3:L:175:LEU:HB3	2.00	0.44
1:R:494:SER:HB2	2:H:56:SER:HB3	1.99	0.43
3:L:7:SER:HB2	3:L:22:THR:HB	1.99	0.43
2:H:166:GLY:C	2:H:168:LEU:H	2.26	0.43
1:R:405:ASP:OD1	1:R:405:ASP:N	2.51	0.43
1:R:434:ILE:HD11	1:R:511:VAL:HB	2.01	0.43
2:H:128:PRO:CB	2:H:154:TYR:HB3	2.49	0.43
3:L:114:SER:OG	3:L:137:ASN:HB3	2.19	0.43
1:R:366:SER:O	1:R:370:ASN:N	2.44	0.43
1:R:480:CYS:O	1:R:483:VAL:HG12	2.19	0.43
3:L:21:ILE:O	3:L:72:THR:HA	2.19	0.43
3:L:136:LEU:HD13	3:L:136:LEU:HA	1.62	0.43
1:R:484:GLU:OE2	2:H:33:TYR:OH	2.14	0.43
2:H:130:VAL:HG21	2:H:216:VAL:HG23	2.01	0.43
3:L:96:LEU:HA	3:L:96:LEU:HD23	1.76	0.43
1:R:368:LEU:O	1:R:371:SER:OG	2.30	0.43
2:H:5:VAL:O	2:H:22:CYS:HA	2.19	0.43
3:L:33:LEU:HD12	3:L:89:GLN:O	2.19	0.43
3:L:108:ARG:CZ	3:L:109:THR:HG22	2.48	0.43
2:H:127:GLY:HA2	2:H:212:SER:OG	2.19	0.43
2:H:148:GLY:CA	2:H:190:VAL:HG12	2.40	0.43
3:L:32:TYR:N	3:L:32:TYR:HD1	2.16	0.43
1:R:360:ASN:N	1:R:523:THR:HB	2.18	0.43
1:R:412:PRO:HG3	1:R:429:PHE:HB3	2.01	0.43
1:R:501:ASN:HB3	1:R:505:TYR:HB2	2.00	0.43
1:R:395:VAL:HG22	1:R:515:PHE:CD1	2.54	0.42
3:L:4:LEU:HB2	3:L:98:PHE:O	2.19	0.42
3:L:140:TYR:O	3:L:198:HIS:HE1	2.02	0.42
1:R:442:ASP:O	1:R:507:PRO:HG3	2.19	0.42
3:L:33:LEU:HD21	3:L:35:TRP:NE1	2.35	0.42
1:R:401:VAL:CG2	1:R:509:ARG:HD2	2.48	0.42
1:R:367:VAL:HA	1:R:370:ASN:CG	2.44	0.42
2:H:152:LYS:HG2	2:H:153:ASP:CG	2.44	0.42
3:L:117:ILE:HD12	3:L:133:VAL:O	2.20	0.42
1:R:420:ASP:O	1:R:461:LEU:HG	2.20	0.42
1:R:457:ARG:HB3	1:R:459:SER:O	2.20	0.42
3:L:6:GLN:HB2	3:L:99:GLY:CA	2.49	0.42
1:R:376:THR:HG22	1:R:435:ALA:HB3	2.02	0.42
1:R:338:PHE:CD1	1:R:363:ALA:HA	2.55	0.42
2:H:2:VAL:HG13	2:H:27:PHE:HD2	1.85	0.42
1:R:349:SER:CB	1:R:452:ARG:H	2.23	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:365:TYR:CE1	1:R:388:ASN:HA	2.54	0.41
1:R:450:ASN:O	1:R:452:ARG:HG3	2.20	0.41
2:H:87:ARG:HG2	2:H:89:GLU:H	1.85	0.41
2:H:114:GLN:HA	3:L:43:ALA:HB2	2.02	0.41
2:H:145:ALA:N	2:H:195:SER:HB3	2.28	0.41
3:L:6:GLN:NE2	3:L:102:THR:OG1	2.54	0.41
3:L:162:SER:HB2	3:L:176:SER:HB3	2.02	0.41
3:L:186:TYR:O	3:L:211:ARG:HD3	2.20	0.41
1:R:339:GLY:HA2	1:R:342:PHE:CE2	2.56	0.41
1:R:358:ILE:HD13	1:R:358:ILE:HA	1.86	0.41
2:H:35:SER:HA	2:H:50:TYR:HA	2.02	0.41
2:H:159:VAL:CG2	2:H:207:VAL:HG13	2.50	0.41
3:L:48:ILE:HD13	3:L:54:LEU:HA	2.01	0.41
1:R:418:ILE:HA	1:R:418:ILE:HD13	1.92	0.41
2:H:150:LEU:HD22	2:H:187:LEU:O	2.20	0.41
1:R:424:LYS:HB3	1:R:463:PRO:HA	2.02	0.41
1:R:392:PHE:CD1	1:R:517:LEU:HD22	2.56	0.41
2:H:148:GLY:HA3	2:H:190:VAL:CG1	2.40	0.41
3:L:154:LEU:HD23	3:L:155:GLN:N	2.36	0.41
1:R:354:ASN:O	1:R:398:ASP:HA	2.21	0.41
1:R:438:SER:O	1:R:438:SER:OG	2.30	0.41
1:R:451:TYR:CB	1:R:495:TYR:HD2	2.34	0.41
2:H:34:MET:HB3	2:H:79:LEU:HD22	2.02	0.41
3:L:94:TYR:O	3:L:96:LEU:N	2.53	0.41
3:L:186:TYR:CZ	3:L:211:ARG:HG3	2.55	0.41
2:H:32:TYR:CG	2:H:98:ARG:HG3	2.55	0.41
2:H:70:ILE:HG13	2:H:71:SER:N	2.35	0.41
2:H:206:ASN:HA	2:H:217:ASP:OD1	2.21	0.41
3:L:121:SER:O	3:L:124:GLN:HB3	2.21	0.41
1:R:363:ALA:HB3	1:R:365:TYR:CZ	2.57	0.40
2:H:204:ILE:HA	2:H:218:LYS:O	2.21	0.40
3:L:14:SER:HA	3:L:107:THR:HB	2.02	0.40
1:R:353:TRP:CD1	1:R:353:TRP:H	2.38	0.40
1:R:354:ASN:ND2	1:R:356:LYS:HE2	2.22	0.40
1:R:454:ARG:NH1	1:R:457:ARG:HG3	2.37	0.40
1:R:398:ASP:OD2	1:R:423:TYR:OH	2.32	0.40
3:L:108:ARG:HD2	3:L:109:THR:H	1.87	0.40
1:R:517:LEU:O	1:R:517:LEU:HG	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	R	189/224 (84%)	154 (82%)	35 (18%)	0	100	100
2	H	214/227 (94%)	189 (88%)	25 (12%)	0	100	100
3	L	211/214 (99%)	187 (89%)	23 (11%)	1 (0%)	24	52
All	All	614/665 (92%)	530 (86%)	83 (14%)	1 (0%)	43	70

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	L	95	PRO

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	R	165/196 (84%)	155 (94%)	10 (6%)	17	42
2	H	185/193 (96%)	167 (90%)	18 (10%)	8	26
3	L	184/185 (100%)	171 (93%)	13 (7%)	13	37
All	All	534/574 (93%)	493 (92%)	41 (8%)	12	35

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	R	398	ASP
1	R	399	SER

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Mol	Chain	Res	Type
1	R	440	ASN
1	R	441	LEU
1	R	468	ILE
1	R	484	GLU
1	R	495	TYR
1	R	513	LEU
1	R	518	LEU
1	R	524	VAL
2	H	12	VAL
2	H	37	ILE
2	H	93	VAL
2	H	98	ARG
2	H	105	ILE
2	H	118	VAL
2	H	125	THR
2	H	129	SER
2	H	137	SER
2	H	147	LEU
2	H	150	LEU
2	H	151	VAL
2	H	161	VAL
2	H	178	VAL
2	H	197	SER
2	H	202	THR
2	H	208	ASN
2	H	223	LYS
3	L	15	VAL
3	L	19	VAL
3	L	63	SER
3	L	69	THR
3	L	90	GLN
3	L	94	TYR
3	L	97	THR
3	L	134	CYS
3	L	136	LEU
3	L	162	SER
3	L	176	SER
3	L	177	SER
3	L	187	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
2	H	59	ASN
2	H	82	GLN
2	H	180	GLN
3	L	90	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	R	191/224 (85%)	-0.21	4 (2%) 63 40	46, 65, 127, 211	0
2	H	218/227 (96%)	-0.38	1 (0%) 87 72	43, 57, 93, 127	0
3	L	213/214 (99%)	-0.25	1 (0%) 87 72	35, 66, 125, 177	0
All	All	622/665 (93%)	-0.28	6 (0%) 79 59	35, 63, 124, 211	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	163	TRP	3.1
3	L	104	VAL	2.7
1	R	342	PHE	2.2
1	R	380	TYR	2.2
1	R	407	VAL	2.2
1	R	387	LEU	2.1

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.