



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:28 AM GMT

PDB ID : 2AJF
Title : Structure of SARS coronavirus spike receptor-binding domain complexed with its receptor
Authors : Li, F.; Li, W.; Farzan, M.; Harrison, S.C.
Deposited on : 2005-08-01
Resolution : 2.90 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

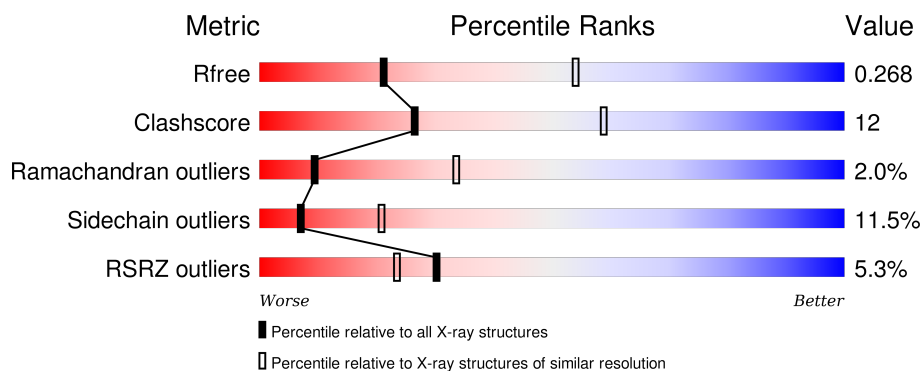
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	597	<div> <div>3%</div> <div>66%</div> <div>30%</div> <div>.</div> </div>
1	B	597	<div> <div>6%</div> <div>67%</div> <div>28%</div> <div>5%</div> </div>
2	E	180	<div> <div>9%</div> <div>72%</div> <div>21%</div> <div>...</div> </div>
2	F	180	<div> <div>8%</div> <div>67%</div> <div>24%</div> <div>5%</div> <div>..</div> </div>

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 12777 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 Angiotensin-converting enzyme-Related Carboxypeptidase (Ace2).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	597	Total	C	N	O	S	0	0	0
			4870	3115	806	920	29			
1	B	597	Total	C	N	O	S	0	0	0
			4870	3115	806	920	29			

- Molecule 2 is a protein called SARS-coronavirus spike protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	174	Total	C	N	O	S	0	0	0
			1403	909	228	259	7			
2	F	174	Total	C	N	O	S	0	0	0
			1403	909	228	259	7			

- Molecule 3 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	3	Total	C	N	O	0	0
			39	22	2	15		
3	B	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 4 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 5 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Zn	0	0
			1	1		
5	A	1	Total	Zn	0	0
			1	1		

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Cl	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Cl	0	0
			1	1		

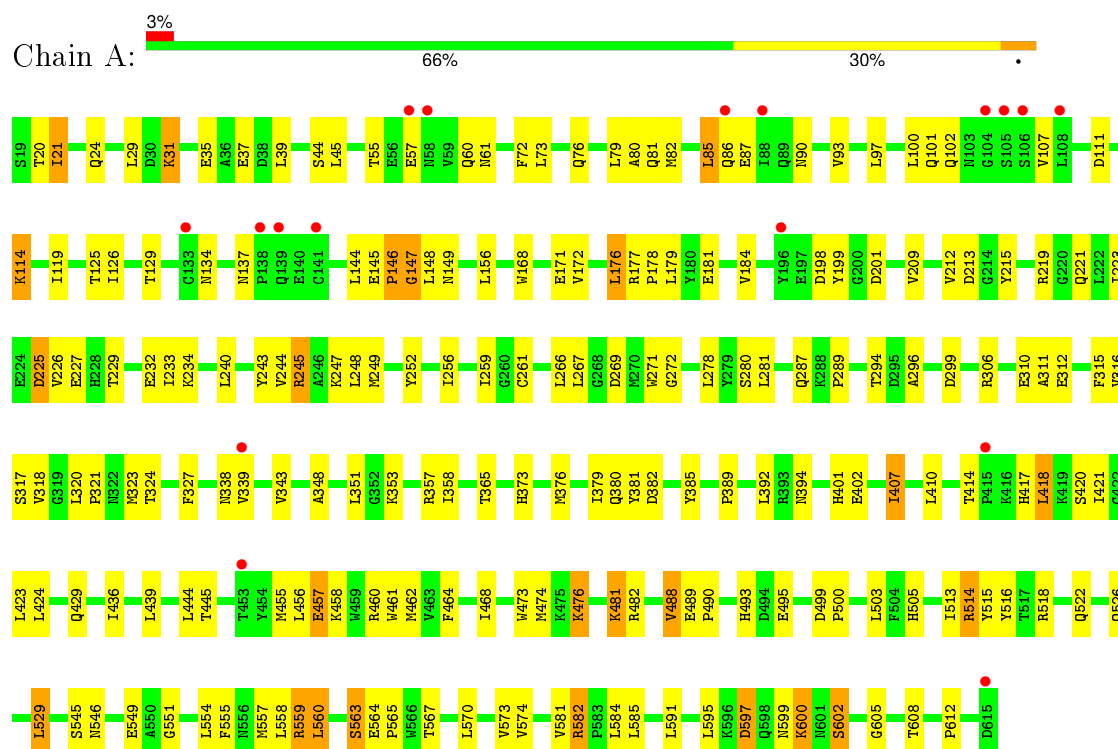
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	36	Total	O	0	0
			36	36		
7	B	24	Total	O	0	0
			24	24		
7	E	4	Total	O	0	0
			4	4		
7	F	1	Total	O	0	0
			1	1		

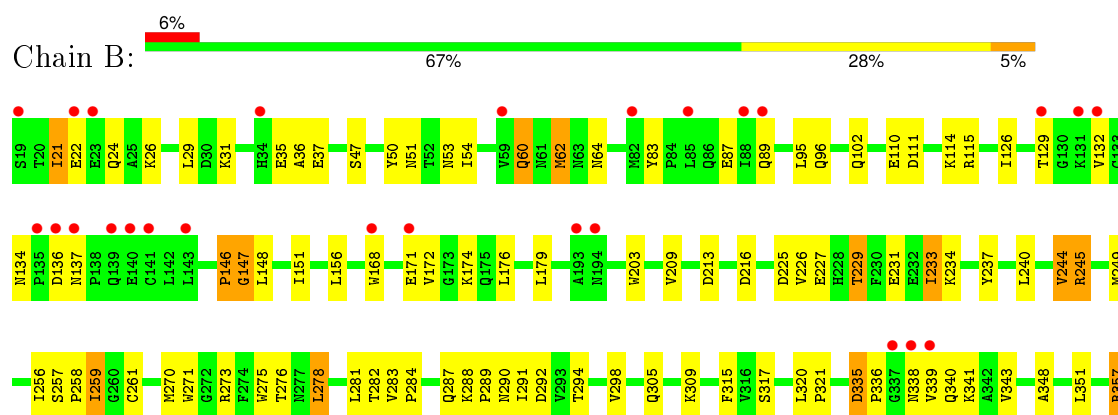
3 Residue-property plots

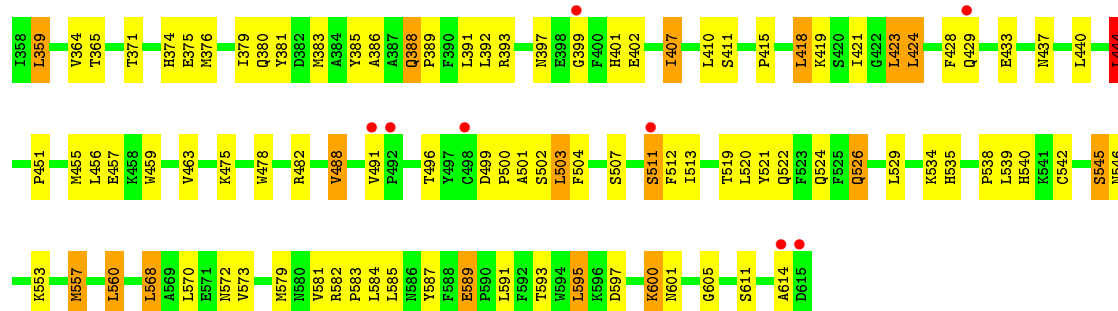
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: Angiotensin-converting enzyme-Related Carboxypeptidase (Ace2)

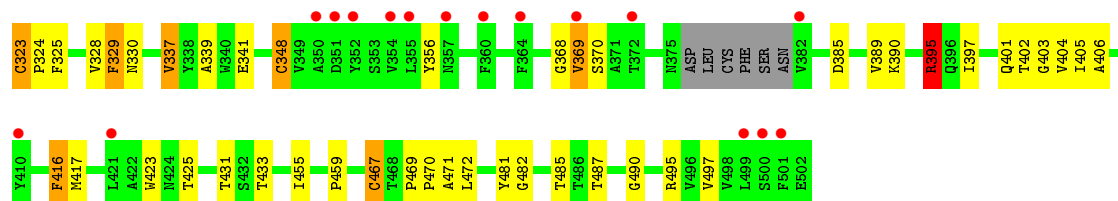


- Molecule 1: Angiotensin-converting enzyme-Related Carboxypeptidase (Ace2)

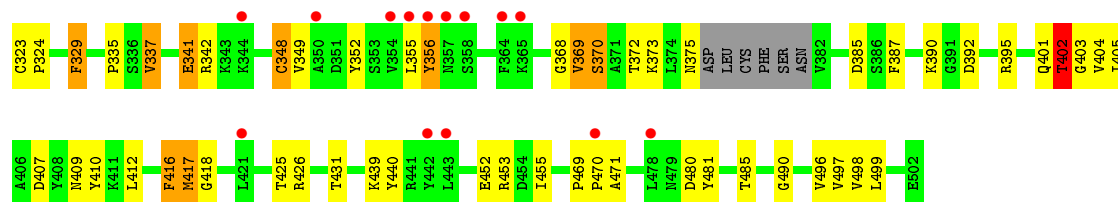




• Molecule 2: SARS-coronavirus spike protein



• Molecule 2: SARS-coronavirus spike protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.30Å 119.43Å 113.24Å 90.00° 91.97° 90.00°	Depositor
Resolution (Å)	47.40 – 2.90 47.39 – 2.81	Depositor EDS
% Data completeness (in resolution range)	92.4 (47.40-2.90) 85.9 (47.39-2.81)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.95 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.218 , 0.275 0.212 , 0.268	Depositor DCC
R_{free} test set	2311 reflections (5.52%)	DCC
Wilson B-factor (Å ²)	73.2	Xtriage
Anisotropy	0.075	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 81.2	EDS
Estimated twinning fraction	0.038 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 50791 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12777	wwPDB-VP
Average B, all atoms (Å ²)	90.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.38% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: ZN, BMA, NAG, CL

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	0.47	0/5007	0.62	0/6803
1	B	0.45	0/5007	0.58	2/6803 (0.0%)
2	E	0.47	0/1448	0.62	0/1972
2	F	0.44	0/1448	0.60	0/1972
All	All	0.46	0/12910	0.60	2/17550 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	444	LEU	CA-CB-CG	5.95	128.99	115.30
1	B	423	LEU	N-CA-C	-5.10	97.22	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	4870	0	4639	124	0
1	B	4870	0	4641	124	0
2	E	1403	0	1326	24	0
2	F	1403	0	1325	31	0
3	A	39	0	34	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	39	0	34	1	0
4	A	42	0	39	2	0
4	B	14	0	13	0	0
4	E	14	0	13	0	0
4	F	14	0	13	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	1	0	0	1	0
6	B	1	0	0	0	0
7	A	36	0	0	3	0
7	B	24	0	0	3	0
7	E	4	0	0	0	0
7	F	1	0	0	0	0
All	All	12777	0	12077	302	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 12.

All (302) 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:320:LEU:HD13	1:A:380:GLN:HG2	1.36	1.05
1:A:261:CYS:HB2	1:A:488:VAL:CG2	1.93	0.97
1:B:245:ARG:HH21	1:B:245:ARG:HG2	1.32	0.94
1:A:261:CYS:HB2	1:A:488:VAL:HG22	1.48	0.94
2:F:335:PRO:HG3	2:F:341:GLU:HG2	1.49	0.93
1:B:589:GLU:HA	1:B:589:GLU:OE1	1.69	0.91
1:B:407:ILE:HD11	1:B:522:GLN:O	1.71	0.91
1:B:371:THR:OG1	7:B:1923:HOH:O	1.88	0.89
1:A:348:ALA:HB1	1:A:379:ILE:HD12	1.53	0.89
2:E:329:PHE:H	2:E:329:PHE:HD1	1.20	0.87
1:A:348:ALA:HB1	1:A:379:ILE:CD1	2.06	0.85
1:B:411:SER:OG	7:B:1906:HOH:O	1.94	0.84
2:F:469:PRO:HA	2:F:471:ALA:H	1.41	0.84
1:B:261:CYS:HB2	1:B:488:VAL:HG22	1.62	0.82
1:B:21:ILE:HA	1:B:24:GLN:HE21	1.45	0.80
1:A:229:THR:HG23	1:A:516:TYR:OH	1.84	0.78
2:E:329:PHE:HB3	2:E:497:VAL:HG21	1.63	0.78
1:B:320:LEU:HD13	1:B:380:GLN:HG2	1.68	0.75
1:A:493:HIS:HD2	1:A:499:ASP:OD1	1.73	0.72
1:A:310:GLU:HG2	1:A:421:ILE:HD11	1.72	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:457:GLU:HG2	1:B:512:PHE:HB3	1.72	0.71
1:B:209:VAL:HG23	1:B:216:ASP:HA	1.73	0.70
1:B:24:GLN:HG3	1:B:83:TYR:HE2	1.57	0.70
1:B:294:THR:O	1:B:298:VAL:HG23	1.94	0.68
1:B:47:SER:HA	1:B:62:MET:HG2	1.73	0.68
1:A:320:LEU:CD1	1:A:380:GLN:HG2	2.20	0.67
1:B:132:VAL:HG12	1:B:171:GLU:HG3	1.75	0.67
1:B:111:ASP:O	1:B:115:ARG:HB3	1.93	0.67
1:B:315:PHE:HB2	7:B:1907:HOH:O	1.93	0.67
1:A:529:LEU:HD21	1:A:554:LEU:HD13	1.77	0.67
1:A:126:ILE:HD11	1:A:176:LEU:HD22	1.77	0.66
1:B:168:TRP:O	1:B:172:VAL:HG22	1.94	0.66
1:B:240:LEU:O	1:B:244:VAL:HG13	1.95	0.66
1:A:248:LEU:HD21	1:A:278:LEU:HD23	1.77	0.66
1:A:462:MET:HE3	1:A:468:ILE:HD11	1.78	0.66
1:B:440:LEU:HD13	1:B:444:LEU:CD2	2.25	0.66
1:B:524:GLN:HE22	1:B:579:MET:HA	1.61	0.65
1:B:245:ARG:HH21	1:B:245:ARG:CG	2.08	0.65
1:A:597:ASP:O	1:A:600:LYS:HB2	1.95	0.65
2:F:323:CYS:N	2:F:348:CYS:SG	2.70	0.65
1:A:402:GLU:HB3	1:A:518:ARG:HD2	1.77	0.65
1:A:261:CYS:HB2	1:A:488:VAL:HG23	1.79	0.64
1:B:126:ILE:HA	1:B:129:THR:HG22	1.80	0.64
1:A:229:THR:OG1	1:A:581:VAL:HB	1.98	0.63
1:B:419:LYS:HD2	1:B:428:PHE:HB3	1.80	0.63
1:A:225:ASP:O	1:A:229:THR:HG22	1.99	0.63
1:A:320:LEU:HD13	1:A:380:GLN:CG	2.23	0.62
1:B:526:GLN:HA	1:B:526:GLN:HE21	1.64	0.62
1:A:248:LEU:HD21	1:A:278:LEU:CD2	2.28	0.62
2:E:329:PHE:N	2:E:329:PHE:HD1	1.96	0.61
1:B:553:LYS:HE3	1:B:573:VAL:O	2.00	0.61
1:A:455:MET:HE2	1:A:481:LYS:HG2	1.81	0.61
2:E:395:ARG:HH11	2:E:395:ARG:HB2	1.65	0.61
1:A:72:PHE:O	1:A:76:GLN:HG2	2.01	0.60
1:B:499:ASP:N	1:B:500:PRO:HD2	2.17	0.60
1:B:351:LEU:HD11	1:B:357:ARG:HD3	1.84	0.59
1:A:296:ALA:HA	1:A:299:ASP:HB2	1.85	0.59
1:A:269:ASP:OD2	1:A:272:GLY:N	2.31	0.59
2:E:323:CYS:N	2:E:348:CYS:SG	2.76	0.59
1:B:294:THR:HG23	1:B:365:THR:HA	1.84	0.58
2:F:469:PRO:HA	2:F:471:ALA:N	2.15	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:21:ILE:HA	1:B:24:GLN:NE2	2.18	0.58
1:A:223:ILE:O	1:A:227:GLU:HG3	2.03	0.58
1:B:51:ASN:HD22	1:B:343:VAL:HG11	1.68	0.58
1:A:233:ILE:HD11	1:A:584:LEU:HD23	1.86	0.57
2:F:403:GLY:HA2	2:F:407:ASP:CG	2.24	0.57
2:E:325:PHE:HB3	2:E:329:PHE:HE1	1.70	0.56
2:F:323:CYS:HB2	2:F:324:PRO:HD3	1.87	0.56
1:B:134:ASN:HB3	1:B:137:ASN:HB3	1.87	0.56
1:B:176:LEU:HA	1:B:179:LEU:HD12	1.87	0.56
1:A:90:ASN:HB3	1:A:93:VAL:HG22	1.87	0.56
1:B:290:ASN:OD1	1:B:292:ASP:HB3	2.06	0.56
1:A:245:ARG:NH2	7:A:1572:HOH:O	2.40	0.55
1:B:540:HIS:HA	1:B:587:TYR:CE1	2.41	0.55
1:B:288:LYS:HB3	1:B:289:PRO:HD2	1.88	0.55
1:A:418:LEU:HB3	1:A:424:LEU:HB2	1.88	0.55
2:F:418:GLY:HA3	2:F:499:LEU:O	2.07	0.55
1:B:338:ASN:C	1:B:340:GLN:H	2.09	0.54
1:B:433:GLU:O	1:B:437:ASN:OD1	2.24	0.54
1:A:146:PRO:O	1:A:148:LEU:N	2.40	0.54
1:B:348:ALA:HB1	1:B:379:ILE:HD13	1.90	0.54
1:A:144:LEU:HA	1:A:148:LEU:HB3	1.90	0.54
1:B:459:TRP:O	1:B:463:VAL:HG23	2.08	0.54
1:A:267:LEU:HD13	7:A:1551:HOH:O	2.08	0.54
1:A:181:GLU:O	1:A:184:VAL:HG22	2.08	0.54
1:A:499:ASP:N	1:A:500:PRO:HD2	2.23	0.53
1:B:557:MET:HA	1:B:560:LEU:HD22	1.89	0.53
1:A:489:GLU:O	1:A:489:GLU:HG2	2.08	0.53
1:B:597:ASP:O	1:B:600:LYS:HD2	2.08	0.53
1:B:386:ALA:HA	1:B:393:ARG:HD3	1.91	0.53
1:A:134:ASN:HB3	1:A:137:ASN:H	1.73	0.53
1:B:520:LEU:HD22	1:B:579:MET:CE	2.39	0.53
1:A:407:ILE:HD13	1:A:407:ILE:N	2.22	0.53
1:B:389:PRO:HG2	1:B:392:LEU:HB2	1.91	0.53
1:B:611:SER:HB2	1:B:614:ALA:HB3	1.89	0.53
1:A:457:GLU:HA	1:A:457:GLU:OE1	2.08	0.53
1:A:245:ARG:NH1	1:A:605:GLY:O	2.42	0.52
1:B:315:PHE:CE2	1:B:376:MET:HG2	2.44	0.52
1:A:490:PRO:HA	1:A:612:PRO:HG2	1.92	0.52
1:A:455:MET:CE	1:A:481:LYS:HG2	2.39	0.52
1:A:420:SER:HB2	4:A:1546:NAG:H61	1.91	0.52
2:F:455:ILE:HG22	2:F:455:ILE:O	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:538:PRO:HB2	1:B:540:HIS:CE1	2.45	0.52
1:B:146:PRO:O	1:B:148:LEU:N	2.43	0.52
2:F:385:ASP:HB2	2:F:498:VAL:HB	1.92	0.51
2:E:425:THR:HG21	2:E:495:ARG:HD2	1.92	0.51
2:E:403:GLY:O	2:E:405:ILE:N	2.43	0.51
1:B:407:ILE:HG23	1:B:526:GLN:NE2	2.25	0.51
1:B:521:TYR:O	1:B:522:GLN:C	2.49	0.50
1:B:440:LEU:HD13	1:B:444:LEU:HD22	1.93	0.50
1:A:379:ILE:O	1:A:382:ASP:HB2	2.12	0.50
1:B:51:ASN:ND2	1:B:343:VAL:HG11	2.26	0.50
2:F:403:GLY:O	2:F:405:ILE:N	2.45	0.50
1:A:573:VAL:HG23	1:A:574:VAL:HG13	1.92	0.50
1:A:85:LEU:HB3	1:A:101:GLN:HE22	1.77	0.50
1:A:560:LEU:O	1:A:563:SER:HB3	2.11	0.50
1:A:201:ASP:O	1:A:219:ARG:HD2	2.12	0.50
1:A:229:THR:HG23	1:A:516:TYR:HH	1.74	0.50
1:A:414:THR:O	1:A:418:LEU:HD22	2.12	0.50
2:E:337:VAL:HG22	2:E:389:VAL:HG12	1.93	0.50
1:B:407:ILE:HG23	1:B:526:GLN:HE22	1.76	0.49
1:A:21:ILE:HA	1:A:24:GLN:HE21	1.77	0.49
3:B:1092:BMA:H4	2:F:402:THR:HG23	1.94	0.49
1:B:21:ILE:HA	1:B:24:GLN:HG2	1.93	0.49
1:B:261:CYS:CB	1:B:488:VAL:HG22	2.40	0.49
1:B:524:GLN:HG2	1:B:583:PRO:HG2	1.94	0.49
1:A:233:ILE:HD11	1:A:584:LEU:CD2	2.42	0.49
1:A:402:GLU:HA	1:A:402:GLU:OE2	2.12	0.49
1:A:526:GLN:HE21	1:A:526:GLN:HA	1.77	0.49
2:E:329:PHE:N	2:E:329:PHE:CD1	2.68	0.49
1:B:348:ALA:HB1	1:B:379:ILE:CD1	2.43	0.49
1:A:505:HIS:H	1:A:505:HIS:CD2	2.31	0.49
1:B:225:ASP:O	1:B:229:THR:HB	2.13	0.49
1:B:31:LYS:HE3	1:B:35:GLU:OE1	2.13	0.48
2:E:337:VAL:CG2	2:E:389:VAL:HG12	2.43	0.48
1:B:320:LEU:HD13	1:B:380:GLN:CG	2.39	0.48
1:B:203:TRP:CZ3	1:B:511:SER:HA	2.49	0.48
1:B:388:GLN:HA	1:B:388:GLN:HE21	1.77	0.48
2:F:342:ARG:HG3	2:F:385:ASP:OD1	2.14	0.48
1:A:338:ASN:HD22	1:A:339:VAL:HG23	1.77	0.48
2:E:459:PRO:HB2	2:E:467:CYS:HB3	1.96	0.48
2:F:323:CYS:HB2	2:F:324:PRO:CD	2.44	0.48
2:F:417:MET:HA	2:F:417:MET:CE	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:320:LEU:HB3	1:B:321:PRO:HD2	1.95	0.48
2:F:409:ASN:OD1	2:F:440:TYR:HB2	2.14	0.47
2:F:329:PHE:HZ	2:F:355:LEU:HD23	1.79	0.47
1:A:351:LEU:HD12	1:A:351:LEU:N	2.28	0.47
1:A:351:LEU:HD11	1:A:357:ARG:HG2	1.96	0.47
2:E:341:GLU:O	2:E:385:ASP:HA	2.15	0.47
1:B:520:LEU:HD21	1:B:581:VAL:HG23	1.96	0.47
2:E:395:ARG:CB	2:E:395:ARG:HH11	2.27	0.47
2:F:329:PHE:HB3	2:F:497:VAL:HG21	1.96	0.47
1:A:55:THR:HG22	1:A:57:GLU:H	1.79	0.47
1:A:312:GLU:N	1:A:376:MET:HE1	2.29	0.47
1:B:375:GLU:OE1	1:B:375:GLU:HA	2.13	0.47
1:A:599:ASN:HA	1:A:602:SER:HB2	1.97	0.47
1:B:231:GLU:OE1	1:B:234:LYS:HE2	2.15	0.47
1:A:245:ARG:HG2	7:A:1572:HOH:O	2.13	0.47
1:B:305:GLN:O	1:B:309:LYS:HB2	2.14	0.47
1:B:440:LEU:O	1:B:444:LEU:HD22	2.15	0.47
1:B:245:ARG:NH2	1:B:245:ARG:HG2	2.13	0.47
2:F:356:TYR:OH	2:F:370:SER:O	2.33	0.46
1:B:415:PRO:O	1:B:419:LYS:HB2	2.15	0.46
1:A:324:THR:O	1:A:327:PHE:HB3	2.16	0.46
2:F:390:LYS:HG2	2:F:490:GLY:O	2.15	0.46
1:B:50:TYR:HE1	1:B:54:ILE:HG23	1.81	0.46
1:A:389:PRO:HG2	1:A:392:LEU:HD22	1.98	0.46
1:B:47:SER:HA	1:B:62:MET:CG	2.41	0.46
1:B:501:ALA:O	1:B:507:SER:HB3	2.16	0.46
1:A:457:GLU:HG3	1:A:513:ILE:HB	1.98	0.45
2:E:390:LYS:HG2	2:E:490:GLY:O	2.16	0.45
1:A:199:TYR:HB3	1:A:464:PHE:CD1	2.51	0.45
1:B:237:TYR:CE2	1:B:451:PRO:HG2	2.52	0.45
1:A:407:ILE:HD11	1:A:522:GLN:O	2.15	0.45
1:A:423:LEU:HA	1:A:423:LEU:HD23	1.76	0.45
1:B:245:ARG:NH1	1:B:605:GLY:O	2.49	0.45
1:B:53:ASN:HD22	1:B:340:GLN:HG3	1.81	0.45
1:A:456:LEU:C	1:A:456:LEU:HD13	2.36	0.45
1:A:597:ASP:HA	1:A:600:LYS:NZ	2.30	0.45
1:A:323:MET:HE3	1:A:376:MET:SD	2.57	0.45
2:F:392:ASP:O	2:F:395:ARG:HD3	2.17	0.45
1:B:60:GLN:O	1:B:64:ASN:HB2	2.16	0.45
1:A:252:TYR:CD2	1:A:266:LEU:HD13	2.51	0.45
1:A:168:TRP:O	1:A:172:VAL:HG22	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:385:ASP:O	2:E:497:VAL:HA	2.17	0.44
2:E:324:PRO:HD3	2:E:348:CYS:SG	2.57	0.44
1:B:336:PRO:HB2	1:B:340:GLN:HB3	1.99	0.44
1:B:275:TRP:O	1:B:278:LEU:HB2	2.17	0.44
1:A:111:ASP:HA	1:A:114:LYS:HB2	1.97	0.44
1:B:257:SER:OG	1:B:259:ILE:HD13	2.17	0.44
2:E:455:ILE:O	2:E:455:ILE:CG2	2.65	0.44
1:B:233:ILE:HD11	1:B:584:LEU:HD22	2.00	0.44
2:F:452:GLU:HG2	2:F:453:ARG:H	1.81	0.44
1:B:410:LEU:HA	1:B:410:LEU:HD12	1.90	0.44
1:A:215:TYR:HB3	1:A:567:THR:OG1	2.17	0.44
1:B:36:ALA:O	1:B:37:GLU:C	2.55	0.44
1:A:417:HIS:CE1	1:A:545:SER:OG	2.71	0.44
1:B:526:GLN:NE2	1:B:526:GLN:HA	2.32	0.44
1:A:499:ASP:HB3	6:A:902:CL:CL	2.54	0.44
1:A:461:TRP:CH2	1:A:513:ILE:HD12	2.53	0.44
2:E:469:PRO:HA	2:E:471:ALA:N	2.32	0.44
1:B:380:GLN:HA	1:B:383:MET:CE	2.48	0.44
2:E:339:ALA:HA	2:E:455:ILE:HD11	2.00	0.44
1:A:514:ARG:HG2	1:A:515:TYR:N	2.33	0.44
1:A:353:LYS:HD2	2:E:487:THR:HG21	2.00	0.44
1:A:199:TYR:HD2	1:A:464:PHE:CE1	2.35	0.44
1:A:100:LEU:HD23	1:A:100:LEU:HA	1.83	0.44
1:B:374:HIS:NE2	1:B:402:GLU:OE1	2.46	0.44
1:A:21:ILE:HD11	1:A:87:GLU:O	2.18	0.43
1:B:456:LEU:HD23	1:B:512:PHE:CE1	2.53	0.43
1:B:168:TRP:CD1	1:B:502:SER:OG	2.71	0.43
1:A:555:PHE:O	1:A:559:ARG:HG2	2.18	0.43
2:F:352:TYR:HD1	2:F:372:THR:HG23	1.83	0.43
1:B:410:LEU:HD22	1:B:522:GLN:HE21	1.83	0.43
1:B:539:LEU:HD23	1:B:587:TYR:HB2	2.01	0.43
1:B:600:LYS:HG2	1:B:601:ASN:ND2	2.33	0.43
2:F:410:TYR:HE2	2:F:412:LEU:HD13	1.83	0.43
1:A:177:ARG:NH1	1:A:495:GLU:HB3	2.34	0.43
2:F:349:VAL:HA	2:F:375:ASN:C	2.38	0.43
1:A:474:MET:CE	1:A:499:ASP:H	2.31	0.43
1:A:327:PHE:HE2	1:A:358:ILE:HG13	1.82	0.43
1:B:568:LEU:HD22	1:B:572:ASN:HD21	1.84	0.43
1:B:249:MET:HG2	1:B:256:ILE:HB	1.99	0.43
2:E:324:PRO:O	2:E:328:VAL:HG23	2.19	0.43
1:A:243:TYR:CZ	1:A:247:LYS:HE3	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:PRO:HB2	1:A:147:GLY:H	1.63	0.43
1:A:21:ILE:HA	1:A:24:GLN:NE2	2.34	0.43
1:A:80:ALA:C	1:A:82:MET:H	2.22	0.43
1:B:513:ILE:HD12	1:B:513:ILE:HA	1.92	0.43
1:B:457:GLU:CG	1:B:512:PHE:HB3	2.44	0.43
1:A:226:VAL:HG21	1:A:513:ILE:HD11	2.01	0.43
1:B:397:ASN:C	1:B:399:GLY:H	2.22	0.43
1:A:97:LEU:O	1:A:101:GLN:HG2	2.18	0.43
1:B:291:ILE:CG2	1:B:291:ILE:O	2.67	0.43
1:A:31:LYS:NZ	1:A:76:GLN:OE1	2.51	0.42
2:E:403:GLY:O	2:E:406:ALA:N	2.51	0.42
2:F:329:PHE:N	2:F:329:PHE:CD1	2.87	0.42
1:A:318:VAL:O	1:A:551:GLY:HA3	2.19	0.42
1:A:591:LEU:HG	1:A:595:LEU:HD13	2.01	0.42
1:A:482:ARG:HD3	1:A:608:THR:O	2.19	0.42
1:A:126:ILE:HD11	1:A:176:LEU:CD2	2.47	0.42
1:B:520:LEU:HD22	1:B:579:MET:HE3	2.01	0.42
1:B:335:ASP:HA	1:B:336:PRO:HD3	1.89	0.42
1:A:555:PHE:HA	1:A:558:LEU:HB2	2.01	0.42
1:A:564:GLU:HB3	1:A:565:PRO:CD	2.49	0.42
2:F:410:TYR:CE2	2:F:412:LEU:HD13	2.55	0.42
1:B:270:MET:HB3	1:B:271:TRP:CE3	2.55	0.42
1:A:35:GLU:OE1	1:A:72:PHE:HE1	2.03	0.42
1:A:245:ARG:HG2	1:A:245:ARG:HH21	1.83	0.42
1:B:388:GLN:HB3	1:B:389:PRO:HD2	2.02	0.42
2:F:337:VAL:HA	2:F:387:PHE:HB2	2.02	0.42
1:A:407:ILE:N	1:A:407:ILE:CD1	2.78	0.42
1:A:321:PRO:HD2	1:A:380:GLN:OE1	2.19	0.42
1:B:557:MET:HG2	1:B:573:VAL:HG21	2.02	0.42
1:A:296:ALA:HA	1:A:299:ASP:CB	2.48	0.42
1:A:327:PHE:CE2	1:A:358:ILE:HG13	2.55	0.42
1:B:111:ASP:O	1:B:115:ARG:CB	2.65	0.42
1:B:96:GLN:HG2	1:B:392:LEU:HD13	2.01	0.41
1:A:311:ALA:HA	1:A:373:HIS:NE2	2.34	0.41
1:A:315:PHE:C	1:A:317:SER:N	2.72	0.41
1:A:407:ILE:HA	1:A:407:ILE:HD12	1.57	0.41
1:B:591:LEU:HG	1:B:595:LEU:HD22	2.02	0.41
1:B:418:LEU:O	1:B:423:LEU:O	2.38	0.41
1:A:232:GLU:HB2	1:A:581:VAL:HG11	2.01	0.41
1:B:257:SER:HA	1:B:258:PRO:HD3	1.87	0.41
1:B:281:LEU:HD12	1:B:282:THR:HG23	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:ILE:HG23	1:A:179:LEU:HB3	2.02	0.41
1:A:244:VAL:O	1:A:245:ARG:C	2.57	0.41
2:F:395:ARG:HH11	2:F:395:ARG:HB2	1.86	0.41
1:B:380:GLN:HA	1:B:383:MET:HE2	2.02	0.41
2:F:439:LYS:HG2	2:F:480:ASP:OD1	2.20	0.41
1:B:146:PRO:HB2	1:B:147:GLY:H	1.67	0.41
2:F:387:PHE:CE1	2:F:496:VAL:HB	2.56	0.41
1:A:306:ARG:O	1:A:310:GLU:HB2	2.20	0.41
1:A:468:ILE:HD12	1:A:476:LYS:HG3	2.02	0.41
1:B:499:ASP:N	1:B:500:PRO:CD	2.84	0.41
1:B:51:ASN:HD22	1:B:343:VAL:CG1	2.32	0.41
1:A:145:GLU:HA	1:A:146:PRO:HA	1.81	0.41
1:A:134:ASN:HB2	1:A:137:ASN:O	2.21	0.41
2:F:409:ASN:N	2:F:409:ASN:HD22	2.19	0.41
1:B:545:SER:O	1:B:546:ASN:HB2	2.20	0.41
1:A:582:ARG:HB2	1:A:582:ARG:HE	1.45	0.41
1:B:421:ILE:HA	1:B:421:ILE:HD13	1.90	0.41
1:B:174:LYS:HA	1:B:496:THR:O	2.20	0.41
1:A:212:VAL:O	1:A:213:ASP:C	2.60	0.41
1:A:249:MET:HG2	1:A:256:ILE:HB	2.03	0.41
1:B:53:ASN:ND2	1:B:340:GLN:HE21	2.19	0.40
1:B:535:HIS:CE1	1:B:542:CYS:HA	2.56	0.40
1:B:534:LYS:HA	1:B:534:LYS:HD3	1.96	0.40
1:A:271:TRP:CD2	1:A:503:LEU:HD23	2.56	0.40
1:B:478:TRP:O	1:B:482:ARG:HG3	2.21	0.40
1:A:420:SER:HB2	4:A:1546:NAG:C6	2.51	0.40
1:B:359:LEU:C	1:B:359:LEU:HD13	2.42	0.40
1:B:226:VAL:O	1:B:229:THR:HG22	2.20	0.40
1:B:22:GLU:HG2	1:B:26:LYS:HE3	2.03	0.40
1:B:283:VAL:HA	1:B:284:PRO:HD3	1.90	0.40
1:B:503:LEU:HD23	1:B:504:PHE:H	1.86	0.40
1:A:178:PRO:O	1:A:181:GLU:HB2	2.21	0.40
1:A:85:LEU:C	1:A:87:GLU:H	2.25	0.40
1:A:261:CYS:CB	1:A:488:VAL:HG22	2.36	0.40
1:A:156:LEU:HD21	1:A:281:LEU:HD11	2.02	0.40
2:E:397:ILE:N	2:E:397:ILE:HD12	2.37	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	A	595/597 (100%)	536 (90%)	55 (9%)	4 (1%)	26	63
1	B	595/597 (100%)	529 (89%)	56 (9%)	10 (2%)	11	38
2	E	170/180 (94%)	137 (81%)	24 (14%)	9 (5%)	2	8
2	F	170/180 (94%)	143 (84%)	19 (11%)	8 (5%)	3	11
All	All	1530/1554 (98%)	1345 (88%)	154 (10%)	31 (2%)	9	33

All (31) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	146	PRO
1	B	146	PRO
2	E	402	THR
2	E	404	VAL
2	E	416	PHE
1	A	147	GLY
1	B	87	GLU
1	B	110	GLU
1	B	147	GLY
1	B	339	VAL
1	B	519	THR
2	F	370	SER
2	F	402	THR
2	F	404	VAL
1	B	407	ILE
1	B	424	LEU
2	E	395	ARG
2	F	416	PHE
1	B	213	ASP
1	A	86	GLN
2	E	368	GLY
2	E	369	VAL

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Mol	Chain	Res	Type
2	E	370	SER
2	F	369	VAL
2	F	401	GLN
2	E	401	GLN
2	F	368	GLY
1	A	289	PRO
2	E	482	GLY
1	B	364	VAL
2	F	470	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	A	527/527 (100%)	458 (87%)	69 (13%)	5	15
1	B	527/527 (100%)	473 (90%)	54 (10%)	9	27
2	E	152/158 (96%)	134 (88%)	18 (12%)	6	19
2	F	152/158 (96%)	137 (90%)	15 (10%)	10	29
All	All	1358/1370 (99%)	1202 (88%)	156 (12%)	7	21

All (156) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	THR
1	A	21	ILE
1	A	29	LEU
1	A	31	LYS
1	A	37	GLU
1	A	39	LEU
1	A	44	SER
1	A	45	LEU
1	A	60	GLN
1	A	61	ASN
1	A	73	LEU
1	A	79	LEU

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Mol	Chain	Res	Type
1	A	81	GLN
1	A	85	LEU
1	A	102	GLN
1	A	107	VAL
1	A	114	LYS
1	A	125	THR
1	A	129	THR
1	A	149	ASN
1	A	171	GLU
1	A	176	LEU
1	A	198	ASP
1	A	209	VAL
1	A	221	GLN
1	A	225	ASP
1	A	234	LYS
1	A	240	LEU
1	A	245	ARG
1	A	259	ILE
1	A	280	SER
1	A	287	GLN
1	A	294	THR
1	A	316	VAL
1	A	343	VAL
1	A	365	THR
1	A	381	TYR
1	A	385	TYR
1	A	394	ASN
1	A	401	HIS
1	A	407	ILE
1	A	410	LEU
1	A	418	LEU
1	A	429	GLN
1	A	436	ILE
1	A	439	LEU
1	A	444	LEU
1	A	445	THR
1	A	457	GLU
1	A	458	LYS
1	A	460	ARG
1	A	473	TRP
1	A	476	LYS
1	A	481	LYS

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Mol	Chain	Res	Type
1	A	488	VAL
1	A	514	ARG
1	A	529	LEU
1	A	546	ASN
1	A	549	GLU
1	A	557	MET
1	A	559	ARG
1	A	560	LEU
1	A	563	SER
1	A	570	LEU
1	A	582	ARG
1	A	585	LEU
1	A	597	ASP
1	A	600	LYS
1	A	602	SER
1	B	21	ILE
1	B	29	LEU
1	B	60	GLN
1	B	62	MET
1	B	89	GLN
1	B	95	LEU
1	B	102	GLN
1	B	114	LYS
1	B	136	ASP
1	B	151	ILE
1	B	156	LEU
1	B	227	GLU
1	B	229	THR
1	B	233	ILE
1	B	244	VAL
1	B	245	ARG
1	B	259	ILE
1	B	273	ARG
1	B	276	THR
1	B	278	LEU
1	B	287	GLN
1	B	317	SER
1	B	335	ASP
1	B	341	LYS
1	B	357	ARG
1	B	359	LEU
1	B	381	TYR

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Mol	Chain	Res	Type
1	B	385	TYR
1	B	388	GLN
1	B	391	LEU
1	B	401	HIS
1	B	418	LEU
1	B	424	LEU
1	B	429	GLN
1	B	444	LEU
1	B	455	MET
1	B	475	LYS
1	B	488	VAL
1	B	491	VAL
1	B	503	LEU
1	B	511	SER
1	B	526	GLN
1	B	529	LEU
1	B	545	SER
1	B	557	MET
1	B	560	LEU
1	B	568	LEU
1	B	570	LEU
1	B	582	ARG
1	B	585	LEU
1	B	589	GLU
1	B	593	THR
1	B	595	LEU
1	B	600	LYS
2	E	323	CYS
2	E	329	PHE
2	E	330	ASN
2	E	337	VAL
2	E	348	CYS
2	E	356	TYR
2	E	369	VAL
2	E	395	ARG
2	E	416	PHE
2	E	417	MET
2	E	423	TRP
2	E	431	THR
2	E	433	THR
2	E	467	CYS
2	E	470	PRO

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Mol	Chain	Res	Type
2	E	472	LEU
2	E	481	TYR
2	E	485	THR
2	F	329	PHE
2	F	337	VAL
2	F	341	GLU
2	F	348	CYS
2	F	356	TYR
2	F	369	VAL
2	F	373	LYS
2	F	402	THR
2	F	416	PHE
2	F	417	MET
2	F	425	THR
2	F	426	ARG
2	F	431	THR
2	F	481	TYR
2	F	485	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	33	ASN
1	A	49	ASN
1	A	101	GLN
1	A	137	ASN
1	A	287	GLN
1	A	338	ASN
1	A	340	GLN
1	A	472	GLN
1	A	493	HIS
1	A	505	HIS
1	A	526	GLN
1	A	580	ASN
1	A	586	ASN
1	A	599	ASN
1	B	24	GLN
1	B	33	ASN
1	B	51	ASN
1	B	53	ASN
1	B	101	GLN

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Mol	Chain	Res	Type
1	B	300	GLN
1	B	373	HIS
1	B	388	GLN
1	B	493	HIS
1	B	505	HIS
1	B	524	GLN
1	B	526	GLN
1	B	535	HIS
1	B	556	ASN
1	B	586	ASN
1	B	599	ASN
1	B	601	ASN
2	E	473	ASN
2	F	473	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

6 carbohydrates 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	$\# Z > 2$
3	NAG	A	1090	1,3	14,14,15	0.46	0	15,19,21	2.14	4 (26%)
3	NAG	A	1091	3	14,14,15	0.31	0	15,19,21	2.59	1 (6%)
3	BMA	A	1092	3	11,11,12	0.60	0	14,15,17	1.19	1 (7%)
3	NAG	B	1090	1,3	14,14,15	0.52	0	15,19,21	0.66	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	B	1091	3	14,14,15	0.47	0	15,19,21	1.10	1 (6%)
3	BMA	B	1092	3	11,11,12	0.61	0	14,15,17	1.43	3 (21%)

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
3	NAG	A	1090	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	1091	3	-	0/6/23/26	0/1/1/1
3	BMA	A	1092	3	-	0/2/19/22	0/1/1/1
3	NAG	B	1090	1,3	-	0/6/23/26	0/1/1/1
3	NAG	B	1091	3	-	0/6/23/26	0/1/1/1
3	BMA	B	1092	3	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1091	NAG	C3-C4-C5	-2.40	106.01	110.20
3	A	1090	NAG	O7-C7-C8	-2.31	117.82	122.06
3	A	1090	NAG	C3-C4-C5	2.22	114.06	110.20
3	A	1090	NAG	C2-N2-C7	2.31	126.00	123.04
3	B	1092	BMA	O5-C1-C2	2.31	114.61	110.86
3	B	1092	BMA	C1-O5-C5	2.77	115.77	112.25
3	A	1092	BMA	C1-O5-C5	3.26	116.39	112.25
3	B	1092	BMA	C1-C2-C3	3.36	113.52	109.54
3	A	1090	NAG	C1-O5-C5	6.25	120.18	112.25
3	A	1091	NAG	C1-O5-C5	9.51	124.32	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1092	BMA	1	0

5.6 Ligand geometry

Of 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	NAG	A	1053	1	14,14,15	0.47	0	15,19,21	1.39	1 (6%)
4	NAG	A	1322	1	14,14,15	0.54	0	15,19,21	1.93	1 (6%)
4	NAG	A	1546	1	14,14,15	0.71	0	15,19,21	1.83	3 (20%)
4	NAG	B	1322	1	14,14,15	0.51	0	15,19,21	1.18	1 (6%)
4	NAG	E	1330	2	14,14,15	0.49	0	15,19,21	1.43	2 (13%)
4	NAG	F	1330	2	14,14,15	0.49	0	15,19,21	2.33	3 (20%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	1053	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1322	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1546	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1322	1	-	0/6/23/26	0/1/1/1
4	NAG	E	1330	2	-	0/6/23/26	0/1/1/1
4	NAG	F	1330	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	1330	NAG	C2-N2-C7	-4.47	117.29	123.04
4	F	1330	NAG	C4-C3-C2	-4.22	104.67	111.23
4	A	1546	NAG	C4-C3-C2	-3.86	105.23	111.23
4	A	1546	NAG	C3-C4-C5	-3.52	104.06	110.20
4	E	1330	NAG	O7-C7-C8	-2.25	117.94	122.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1322	NAG	C1-O5-C5	2.37	115.26	112.25
4	E	1330	NAG	C1-O5-C5	3.63	116.86	112.25
4	A	1546	NAG	C1-O5-C5	3.71	116.95	112.25
4	A	1053	NAG	C1-O5-C5	4.58	118.07	112.25
4	F	1330	NAG	C1-O5-C5	6.17	120.08	112.25
4	A	1322	NAG	C1-O5-C5	6.60	120.62	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1546	NAG	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	597/597 (100%)	0.14	17 (2%)	56 50	56, 82, 123, 147	0
1	B	597/597 (100%)	0.39	34 (5%)	27 21	56, 82, 123, 148	0
2	E	174/180 (96%)	0.47	16 (9%)	11 7	75, 96, 142, 153	0
2	F	174/180 (96%)	0.44	14 (8%)	15 10	76, 96, 142, 152	0
All	All	1542/1554 (99%)	0.31	81 (5%)	30 23	56, 88, 131, 153	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	615	ASP	7.6
1	B	136	ASP	7.3
1	B	140	GLU	5.8
2	E	501	PHE	5.5
2	F	354	VAL	5.1
1	B	193	ALA	4.6
2	E	354	VAL	4.4
1	B	339	VAL	4.1
1	B	429	GLN	4.1
1	A	339	VAL	4.1
1	B	131	LYS	4.0
1	B	132	VAL	4.0
1	A	106	SER	4.0
2	E	352	TYR	3.9
1	B	141	CYS	3.8
1	B	168	TRP	3.8
2	F	350	ALA	3.7
2	F	364	PHE	3.7
2	F	443	LEU	3.4
1	B	137	ASN	3.4
1	A	615	ASP	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	108	LEU	3.3
1	A	139	GLN	3.3
1	B	143	LEU	3.2
1	B	82	MET	3.2
1	B	139	GLN	3.2
1	B	89	GLN	3.1
2	E	364	PHE	3.0
2	F	478	LEU	3.0
1	B	135	PRO	2.9
2	E	499	LEU	2.9
2	E	500	SER	2.9
1	B	337	GLY	2.9
2	E	357	ASN	2.8
2	E	372	THR	2.8
1	B	171	GLU	2.8
1	B	194	ASN	2.7
2	F	470	PRO	2.7
2	F	344	LYS	2.7
1	B	59	VAL	2.7
1	A	196	TYR	2.6
1	A	138	PRO	2.6
1	A	415	PRO	2.6
1	B	491	VAL	2.6
2	E	382	VAL	2.6
2	F	355	LEU	2.5
2	E	360	PHE	2.5
2	E	421	LEU	2.5
2	F	357	ASN	2.5
1	B	88	ILE	2.4
1	A	133	CYS	2.4
1	B	23	GLU	2.4
2	F	356	TYR	2.4
1	B	22	GLU	2.4
2	E	355	LEU	2.3
1	A	453	THR	2.3
2	E	410	TYR	2.3
1	A	105	SER	2.3
1	B	614	ALA	2.3
1	B	399	GLY	2.3
1	A	86	GLN	2.3
2	E	350	ALA	2.3
1	B	338	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
2	E	369	VAL	2.2
2	E	351	ASP	2.2
1	B	492	PRO	2.2
1	A	58	ASN	2.2
1	B	19	SER	2.2
1	B	85	LEU	2.2
2	F	421	LEU	2.2
1	A	57	GLU	2.2
1	B	34	HIS	2.2
1	B	129	THR	2.1
1	A	88	ILE	2.1
2	F	442	TYR	2.1
2	F	365	LYS	2.1
1	A	104	GLY	2.1
2	F	358	SER	2.1
1	B	498	CYS	2.0
1	A	141	CYS	2.0
1	B	511	SER	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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	NAG	A	1090	14/15	0.89	0.19	-	79,86,89,93	0
3	NAG	B	1090	14/15	0.93	0.19	-	93,99,100,102	0
3	BMA	A	1092	11/12	0.49	0.32	-	79,80,81,81	0
3	BMA	B	1092	11/12	0.75	0.14	-	83,84,84,85	0
3	NAG	A	1091	14/15	0.90	0.14	-	70,75,77,79	0
3	NAG	B	1091	14/15	0.90	0.17	-	87,90,91,91	0

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
4	NAG	A	1053	14/15	0.76	0.27	0.61	88,89,90,90	0
6	CL	B	1902	1/1	0.87	0.27	0.41	115,115,115,115	0
4	NAG	E	1330	14/15	0.52	0.28	0.21	108,109,112,113	0
4	NAG	F	1330	14/15	0.65	0.28	-0.03	106,107,108,108	0
6	CL	A	902	1/1	0.94	0.14	-3.09	74,74,74,74	0
5	ZN	B	901	1/1	0.98	0.13	-	84,84,84,84	0
4	NAG	A	1546	14/15	0.79	0.25	-	72,74,77,78	0
4	NAG	B	1322	14/15	0.82	0.15	-	72,74,76,77	0
4	NAG	A	1322	14/15	0.86	0.12	-	75,77,78,78	0
5	ZN	A	901	1/1	0.98	0.12	-	74,74,74,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.