



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 6, 2026 – 03:23 PM JST

PDB ID : 9LE9 / pdb_00009le9
Title : Complex structure of THC-4447 Fab bound to SARS-CoV-2 XBB.1.5 RBD
Authors : Wang, X.; Guo, F.
Deposited on : 2025-01-07
Resolution : 3.14 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
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

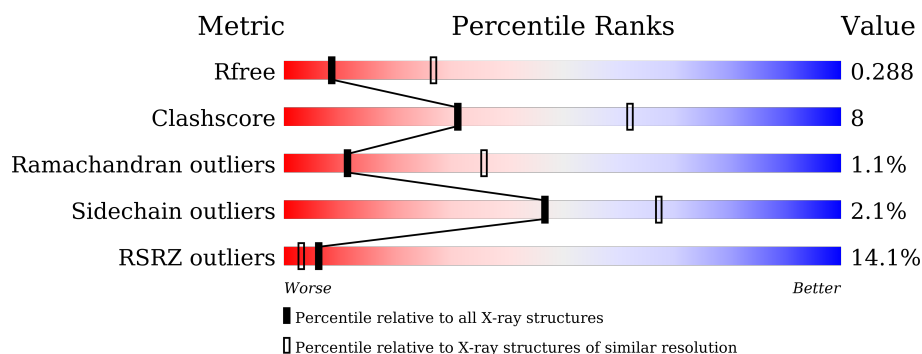
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.14 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	2351 (3.18-3.10)
Clashscore	190562	2452 (3.18-3.10)
Ramachandran outliers	187476	2324 (3.18-3.10)
Sidechain outliers	187428	2324 (3.18-3.10)
RSRZ outliers	180081	2351 (3.18-3.10)

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	B	219	13% 74% 14% 11%
1	E	219	13% 72% 15% 11%
2	A	232	14% 77% 21% •
2	D	232	21% 82% 18%
3	C	218	11% 91% 9%
3	H	218	8% 86% 13% •

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9848 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	E	194	Total	C	N	O	S	0	0	0
			1540	994	258	280	8			
1	B	194	Total	C	N	O	S	0	0	0
			1540	994	258	280	8			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	339	HIS	GLY	variant	UNP P0DTC2
E	346	THR	ARG	variant	UNP P0DTC2
E	368	ILE	LEU	variant	UNP P0DTC2
E	371	PHE	SER	variant	UNP P0DTC2
E	373	PRO	SER	variant	UNP P0DTC2
E	375	PHE	SER	variant	UNP P0DTC2
E	376	ALA	THR	variant	UNP P0DTC2
E	405	ASN	ASP	variant	UNP P0DTC2
E	408	SER	ARG	variant	UNP P0DTC2
E	417	ASN	LYS	variant	UNP P0DTC2
E	440	LYS	ASN	variant	UNP P0DTC2
E	445	PRO	VAL	variant	UNP P0DTC2
E	446	SER	GLY	variant	UNP P0DTC2
E	460	LYS	ASN	variant	UNP P0DTC2
E	477	ASN	SER	variant	UNP P0DTC2
E	478	LYS	THR	variant	UNP P0DTC2
E	484	ALA	GLU	variant	UNP P0DTC2
E	486	PRO	PHE	variant	UNP P0DTC2
E	490	SER	PHE	variant	UNP P0DTC2
E	498	ARG	GLN	variant	UNP P0DTC2
E	501	TYR	ASN	variant	UNP P0DTC2
E	505	HIS	TYR	variant	UNP P0DTC2
B	339	HIS	GLY	variant	UNP P0DTC2
B	346	THR	ARG	variant	UNP P0DTC2
B	368	ILE	LEU	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	371	PHE	SER	variant	UNP P0DTC2
B	373	PRO	SER	variant	UNP P0DTC2
B	375	PHE	SER	variant	UNP P0DTC2
B	376	ALA	THR	variant	UNP P0DTC2
B	405	ASN	ASP	variant	UNP P0DTC2
B	408	SER	ARG	variant	UNP P0DTC2
B	417	ASN	LYS	variant	UNP P0DTC2
B	440	LYS	ASN	variant	UNP P0DTC2
B	445	PRO	VAL	variant	UNP P0DTC2
B	446	SER	GLY	variant	UNP P0DTC2
B	460	LYS	ASN	variant	UNP P0DTC2
B	477	ASN	SER	variant	UNP P0DTC2
B	478	LYS	THR	variant	UNP P0DTC2
B	484	ALA	GLU	variant	UNP P0DTC2
B	486	PRO	PHE	variant	UNP P0DTC2
B	490	SER	PHE	variant	UNP P0DTC2
B	498	ARG	GLN	variant	UNP P0DTC2
B	501	TYR	ASN	variant	UNP P0DTC2
B	505	HIS	TYR	variant	UNP P0DTC2

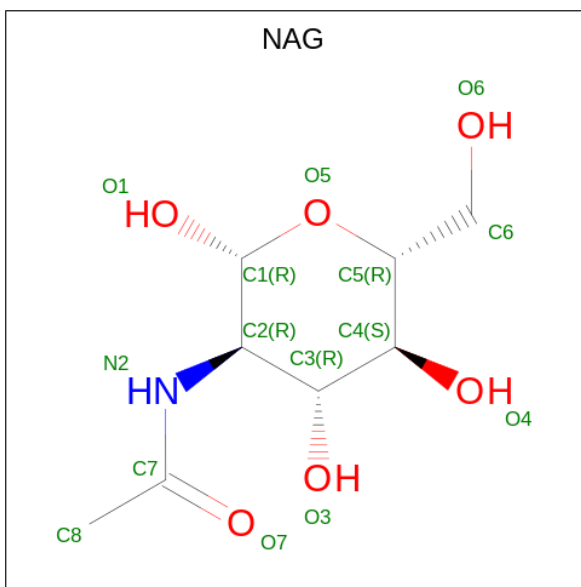
- Molecule 2 is a protein called THC-4447 Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	232	Total	C	N	O	S	0	0	0
			1767	1120	296	342	9			
2	D	232	Total	C	N	O	S	0	0	0
			1767	1120	296	342	9			

- Molecule 3 is a protein called THC-4447 Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	218	Total	C	N	O	S	0	0	0
			1603	999	270	329	5			
3	H	218	Total	C	N	O	S	0	0	0
			1603	999	270	329	5			

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆) (labeled as "Ligand of Interest" by depositor).



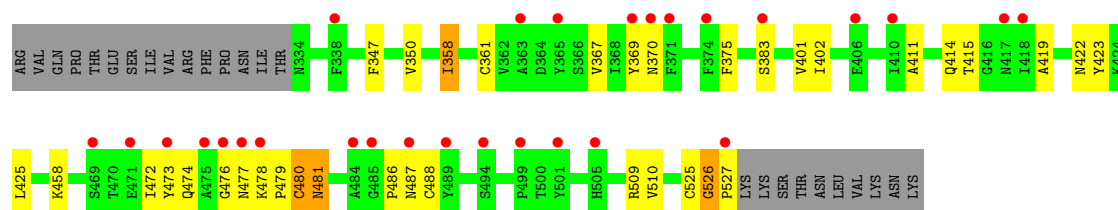
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	E	1	Total	C	N	O	14	0
			14	8	1	5		
4	B	1	Total	C	N	O	14	0
			14	8	1	5		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

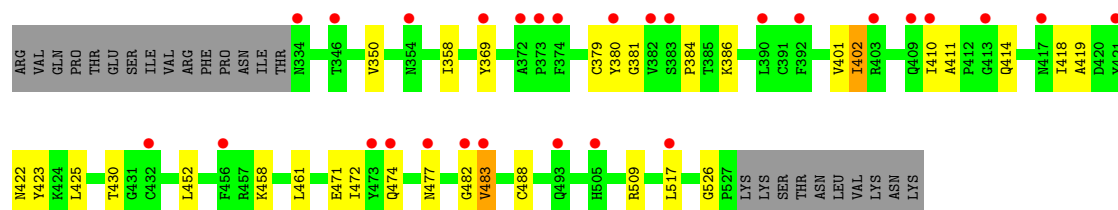
• Molecule 1: Spike protein S1

Chain E: 



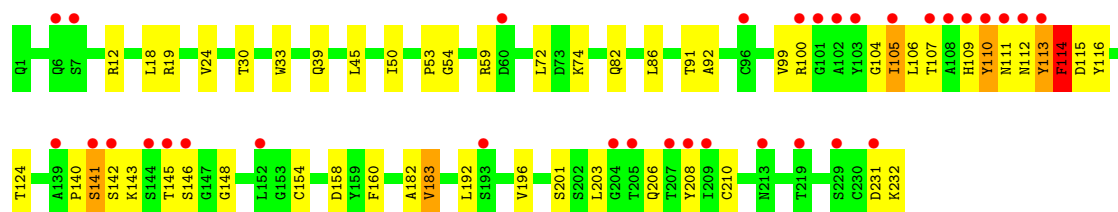
• Molecule 1: Spike protein S1

Chain B: 



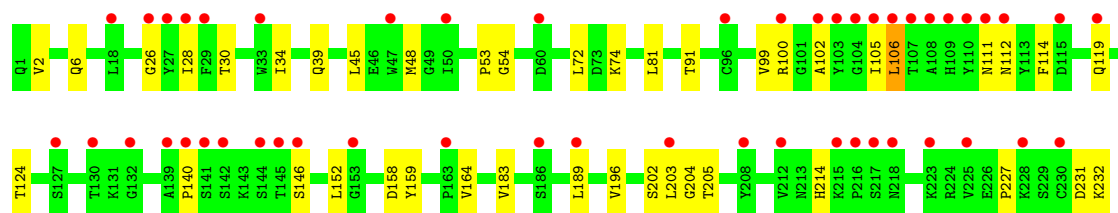
• Molecule 2: THC-4447 Heavy chain

Chain A: 

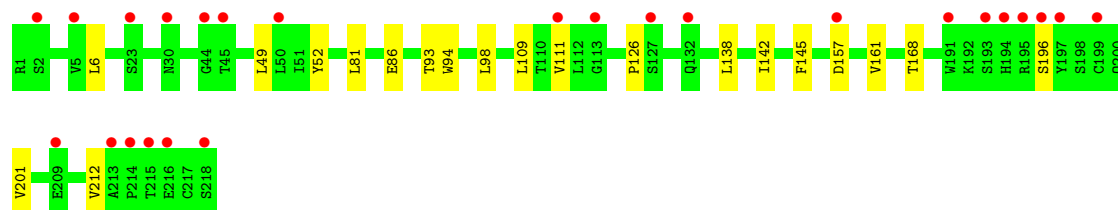
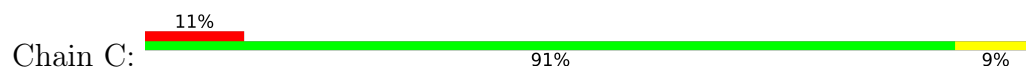


• Molecule 2: THC-4447 Heavy chain

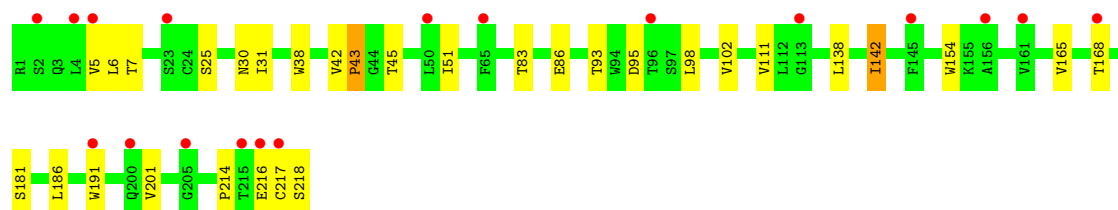
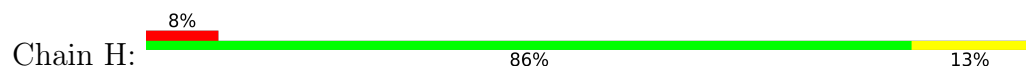
Chain D: 



● Molecule 3: THC-4447 Light Chain



● Molecule 3: THC-4447 Light Chain



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	154.40Å 142.13Å 95.86Å 90.00° 123.98° 90.00°	Depositor
Resolution (Å)	47.79 – 3.14 47.79 – 3.14	Depositor EDS
% Data completeness (in resolution range)	99.7 (47.79-3.14) 99.7 (47.79-3.14)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.30 (at 3.12Å)	Xtriage
Refinement program	PHENIX (1.18.2_3874: ???)	Depositor
R, R_{free}	0.281 , 0.291 0.281 , 0.288	Depositor DCC
R_{free} test set	1992 reflections (6.63%)	wwPDB-VP
Wilson B-factor (Å ²)	76.0	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	9848	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.55% 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

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	B	0.96	0/1589	1.14	0/2166
1	E	0.95	0/1589	1.12	0/2166
2	A	1.01	1/1813 (0.1%)	1.22	5/2470 (0.2%)
2	D	1.01	0/1813	1.16	2/2470 (0.1%)
3	C	1.00	0/1641	1.17	0/2243
3	H	1.03	1/1641 (0.1%)	1.16	2/2243 (0.1%)
All	All	0.99	2/10086 (0.0%)	1.16	9/13758 (0.1%)

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	E	0	1
2	A	0	1
All	All	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	43	PRO	N-CD	-8.54	1.35	1.47
2	A	113	TYR	CA-C	5.95	1.61	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	104	GLY	N-CA-C	-9.21	101.82	115.72
2	D	202	SER	N-CA-C	-8.16	100.37	111.87
2	A	146	SER	N-CA-C	7.24	119.17	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	202	SER	CB-CA-C	5.99	120.75	111.28
3	H	43	PRO	CA-N-CD	5.66	119.92	112.00
2	A	201	SER	N-CA-C	-5.64	105.22	111.36
2	A	114	PHE	N-CA-CB	-5.42	101.32	110.49
3	H	43	PRO	N-CA-CB	-5.33	98.57	103.31
2	A	113	TYR	CA-C-O	5.02	125.44	120.32

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	A	160	PHE	Peptide
1	E	526	GLY	Peptide

5.2 Too-close contacts ⓘ

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	B	1540	0	1465	14	24
1	E	1540	0	1465	20	85
2	A	1767	0	1732	56	65
2	D	1767	0	1730	37	4
3	C	1603	0	1556	28	0
3	H	1603	0	1556	22	0
4	B	14	0	13	0	0
4	E	14	0	13	0	0
All	All	9848	0	9530	157	89

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 8.

All (157) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:113:TYR:CD1	3:C:49:LEU:HD22	1.64	1.31

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:113:TYR:CD1	3:C:49:LEU:CD2	2.16	1.29
2:D:99:VAL:HG13	2:D:111:ASN:OD1	1.08	1.24
2:A:113:TYR:HD1	3:C:49:LEU:CD2	1.51	1.22
2:D:99:VAL:CG1	2:D:111:ASN:OD1	1.92	1.18
2:D:203:LEU:HD12	2:D:204:GLY:N	1.67	1.07
2:A:143:LYS:HE3	2:A:145:THR:OG1	1.61	1.01
2:A:183:VAL:HG22	3:C:168:THR:CG2	1.93	0.99
2:A:100:ARG:HG3	2:A:113:TYR:OH	1.62	0.99
2:A:113:TYR:CE1	2:A:115:ASP:HB2	2.00	0.96
2:D:159:TYR:CE1	2:D:164:VAL:HG23	2.00	0.96
2:A:113:TYR:HD1	3:C:49:LEU:HD22	0.77	0.92
2:A:113:TYR:CD1	3:C:49:LEU:HD21	2.05	0.90
2:D:159:TYR:CE1	2:D:164:VAL:CG2	2.56	0.89
3:H:30:ASN:OD1	3:H:31:ILE:N	2.06	0.88
2:D:183:VAL:HG22	3:H:168:THR:HB	1.57	0.87
2:D:203:LEU:HD12	2:D:204:GLY:H	1.39	0.87
3:H:42:VAL:HG22	3:H:43:PRO:HD2	1.57	0.84
3:C:142:ILE:HG22	3:C:145:PHE:CD1	2.15	0.82
2:D:159:TYR:CD1	2:D:164:VAL:CG2	2.65	0.80
2:D:2:VAL:HA	2:D:26:GLY:HA3	1.65	0.77
2:A:105:ILE:HG22	2:A:105:ILE:O	1.85	0.77
3:C:142:ILE:HG22	3:C:145:PHE:CE1	2.19	0.77
2:A:183:VAL:HG22	3:C:168:THR:HG22	1.66	0.76
1:E:478:LYS:HD3	1:E:487:ASN:HB2	1.66	0.76
1:B:474:GLN:HG3	1:B:488:CYS:SG	2.25	0.76
2:D:159:TYR:CD1	2:D:164:VAL:HG21	2.21	0.75
2:A:183:VAL:HG22	3:C:168:THR:HG23	1.67	0.73
2:A:100:ARG:HG3	2:A:113:TYR:HH	1.52	0.72
2:D:99:VAL:HG13	2:D:111:ASN:CG	2.10	0.72
2:A:113:TYR:CZ	2:A:115:ASP:HB2	2.24	0.71
2:A:154:CYS:CB	2:A:210:CYS:HG	2.04	0.70
2:D:99:VAL:CG1	2:D:111:ASN:CG	2.63	0.70
2:A:113:TYR:CE1	2:A:115:ASP:CB	2.73	0.70
3:H:216:GLU:C	3:H:218:SER:H	2.02	0.68
2:D:203:LEU:CD1	2:D:204:GLY:H	2.06	0.68
2:D:164:VAL:HG22	2:D:214:HIS:HD2	1.57	0.67
2:A:100:ARG:HG3	2:A:113:TYR:CZ	2.30	0.66
3:H:42:VAL:CG2	3:H:43:PRO:HD2	2.26	0.66
2:D:159:TYR:CE1	2:D:164:VAL:HG21	2.31	0.65
2:A:154:CYS:CB	2:A:210:CYS:SG	2.85	0.65
2:A:12:ARG:HE	2:A:18:LEU:HD13	1.62	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:168:THR:HG22	3:H:181:SER:H	1.63	0.64
1:E:480:CYS:O	1:E:481:ASN:HB2	1.98	0.63
2:D:54:GLY:HA2	2:D:74:LYS:HD3	1.81	0.63
2:D:99:VAL:HG22	2:D:114:PHE:CE1	2.35	0.62
1:E:474:GLN:O	1:E:487:ASN:O	2.17	0.62
1:E:367:VAL:O	2:A:105:ILE:HG21	1.99	0.61
3:H:42:VAL:HG22	3:H:43:PRO:CD	2.30	0.61
2:D:203:LEU:CD1	2:D:204:GLY:N	2.54	0.61
2:A:91:THR:HG23	2:A:124:THR:HA	1.82	0.60
2:A:110:TYR:HD1	3:C:94:TRP:HE3	1.49	0.60
2:A:113:TYR:CD1	2:A:113:TYR:O	2.55	0.60
2:A:30:THR:HA	2:A:53:PRO:HB2	1.83	0.59
1:B:411:ALA:HB3	1:B:414:GLN:HG3	1.84	0.59
3:H:6:LEU:HD11	3:H:93:THR:HG22	1.85	0.59
1:E:369:TYR:CZ	2:A:105:ILE:HG13	2.39	0.58
2:A:182:ALA:HB2	2:A:192:LEU:HD23	1.86	0.58
2:A:113:TYR:HB3	3:C:52:TYR:HB2	1.86	0.58
2:D:91:THR:HG23	2:D:124:THR:HA	1.86	0.57
1:E:472:ILE:CG2	1:E:488:CYS:HB3	2.35	0.57
3:C:81:LEU:HD11	3:C:109:LEU:HD21	1.87	0.57
1:E:472:ILE:HG21	1:E:488:CYS:HB3	1.86	0.56
1:E:486:PRO:O	1:E:487:ASN:HB2	2.06	0.56
2:A:54:GLY:H	2:A:74:LYS:HD2	1.72	0.55
3:H:42:VAL:CG2	3:H:43:PRO:CD	2.84	0.55
3:C:142:ILE:HG12	3:C:201:VAL:HG21	1.88	0.55
1:E:525:CYS:HB3	1:E:527:PRO:HD3	1.89	0.54
2:D:158:ASP:HA	2:D:189:LEU:HB3	1.88	0.54
2:A:105:ILE:O	2:A:105:ILE:CG2	2.56	0.54
2:A:140:PRO:O	2:A:141:SER:C	2.50	0.53
2:A:154:CYS:HG	2:A:210:CYS:HG	0.72	0.52
2:A:99:VAL:HA	2:A:112:ASN:HD21	1.74	0.51
2:D:159:TYR:CD1	2:D:164:VAL:HG23	2.38	0.51
2:A:154:CYS:SG	2:A:210:CYS:CB	2.98	0.51
2:A:19:ARG:HG3	2:A:82:GLN:HG2	1.93	0.51
1:B:350:VAL:HG22	1:B:422:ASN:HB3	1.92	0.51
3:C:6:LEU:HD11	3:C:93:THR:HG22	1.92	0.51
2:A:50:ILE:HG13	2:A:59:ARG:HB2	1.92	0.51
2:D:6:GLN:H	2:D:119:GLN:HE21	1.58	0.51
3:H:216:GLU:C	3:H:218:SER:N	2.69	0.51
3:C:157:ASP:N	3:C:196:SER:HB3	2.26	0.50
3:H:5:VAL:HG21	3:H:102:VAL:HG11	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:142:ILE:CG1	3:C:201:VAL:HG21	2.41	0.50
2:D:39:GLN:HB2	2:D:45:LEU:HD23	1.93	0.50
1:E:526:GLY:N	1:E:527:PRO:HD3	2.27	0.50
2:D:99:VAL:HG11	2:D:111:ASN:CG	2.37	0.50
3:C:157:ASP:H	3:C:196:SER:HB3	1.75	0.50
3:H:5:VAL:HG23	3:H:6:LEU:HG	1.95	0.49
3:C:142:ILE:CG2	3:C:145:PHE:CD1	2.92	0.49
3:H:142:ILE:HG12	3:H:201:VAL:HG21	1.96	0.48
2:A:50:ILE:HD13	2:A:110:TYR:HB2	1.94	0.48
2:D:231:ASP:O	2:D:232:LYS:C	2.57	0.48
2:A:113:TYR:O	2:A:114:PHE:C	2.56	0.48
3:C:142:ILE:CG2	3:C:145:PHE:CE1	2.94	0.48
2:D:30:THR:HA	2:D:53:PRO:HB2	1.95	0.48
3:H:42:VAL:HG13	3:H:45:THR:OG1	2.14	0.47
3:C:86:GLU:HB2	3:C:111:VAL:HG23	1.96	0.47
1:E:358:ILE:HG23	1:E:361:CYS:SG	2.54	0.47
2:A:39:GLN:HB2	2:A:45:LEU:HD23	1.97	0.47
1:E:411:ALA:HB3	1:E:414:GLN:HG3	1.95	0.47
1:E:350:VAL:HG22	1:E:422:ASN:HB3	1.96	0.47
2:D:99:VAL:HG11	2:D:111:ASN:CB	2.46	0.46
3:H:154:TRP:CD1	3:H:165:VAL:HG23	2.50	0.46
2:D:140:PRO:HD2	2:D:227:PRO:HA	1.96	0.46
2:A:206:GLN:HG2	2:A:208:TYR:H	1.80	0.46
1:E:401:VAL:HG22	1:E:509:ARG:HG2	1.96	0.45
2:D:99:VAL:HG22	2:D:114:PHE:CD1	2.51	0.45
2:A:113:TYR:HE1	2:A:115:ASP:CB	2.27	0.45
1:B:401:VAL:HG22	1:B:509:ARG:HG2	1.99	0.45
3:C:126:PRO:HD3	3:C:138:LEU:HG	1.98	0.44
1:B:402:ILE:HG12	1:B:410:ILE:HD11	1.99	0.44
3:H:86:GLU:HB2	3:H:111:VAL:HG23	1.99	0.44
1:E:370:ASN:HB2	2:A:106:LEU:HD12	1.98	0.44
1:B:381:GLY:HA3	1:B:430:THR:HG23	2.00	0.44
3:H:7:THR:HB	3:H:25:SER:HB3	2.00	0.44
2:A:18:LEU:HB2	2:A:86:LEU:HD11	2.00	0.43
2:A:106:LEU:HD23	2:A:106:LEU:HA	1.91	0.43
3:H:38:TRP:HB2	3:H:51:ILE:HB	2.00	0.43
2:A:113:TYR:HE1	2:A:115:ASP:CG	2.27	0.43
3:H:83:THR:HA	3:H:111:VAL:HG21	2.01	0.43
1:B:482:GLY:O	1:B:483:VAL:HB	2.17	0.43
1:E:526:GLY:H	1:E:527:PRO:HD3	1.83	0.43
2:A:107:THR:HG22	2:A:109:HIS:H	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:152:LEU:HB2	2:D:196:VAL:HG13	2.01	0.43
1:E:419:ALA:HA	1:E:423:TYR:O	2.19	0.42
2:A:113:TYR:CE1	3:C:49:LEU:CD2	2.92	0.42
2:A:182:ALA:HA	2:A:192:LEU:HB3	2.01	0.42
1:E:367:VAL:O	2:A:105:ILE:CG2	2.66	0.42
2:D:106:LEU:HD12	2:D:106:LEU:HA	1.88	0.42
2:A:113:TYR:HD1	2:A:113:TYR:O	2.01	0.42
2:A:183:VAL:CG2	3:C:168:THR:HG23	2.45	0.42
2:D:102:ALA:HB2	2:D:112:ASN:HB2	2.01	0.42
2:D:99:VAL:CG1	2:D:111:ASN:CB	2.98	0.42
3:C:98:LEU:HD23	3:C:98:LEU:HA	1.89	0.41
2:A:113:TYR:CB	3:C:52:TYR:HB2	2.48	0.41
3:H:95:ASP:HB3	3:H:98:LEU:HB2	2.02	0.41
1:E:425:LEU:HD23	1:E:425:LEU:HA	1.87	0.41
2:A:33:TRP:HB2	2:A:99:VAL:HB	2.03	0.41
2:A:113:TYR:O	3:C:49:LEU:HD22	2.21	0.41
3:C:86:GLU:HG3	3:C:109:LEU:O	2.20	0.41
3:H:138:LEU:HD13	3:H:186:LEU:HD12	2.03	0.41
2:A:154:CYS:HG	2:A:210:CYS:CB	2.25	0.41
1:E:402:ILE:HD11	1:E:510:VAL:HG21	2.02	0.41
1:B:369:TYR:O	2:D:105:ILE:HB	2.21	0.41
1:B:419:ALA:HA	1:B:423:TYR:O	2.20	0.41
2:D:164:VAL:HG22	2:D:214:HIS:CD2	2.46	0.41
1:B:472:ILE:HG21	1:B:488:CYS:HB3	2.02	0.41
2:A:109:HIS:HB3	2:A:110:TYR:H	1.66	0.40
3:H:191:TRP:CZ2	3:H:214:PRO:HA	2.56	0.40
2:A:99:VAL:HG13	2:A:112:ASN:CG	2.45	0.40
1:B:461:LEU:HD23	1:B:461:LEU:HA	1.98	0.40
1:B:350:VAL:HG21	1:B:418:ILE:HG23	2.02	0.40
1:B:425:LEU:HD23	1:B:425:LEU:HA	1.91	0.40
1:B:379:CYS:SG	1:B:384:PRO:HG3	2.62	0.40
2:D:203:LEU:CG	2:D:204:GLY:H	2.34	0.40
2:D:48:MET:HE1	2:D:81:LEU:HD21	2.03	0.40

All (89) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:479:PRO:C	2:A:232:LYS:CE[3_455]	0.68	1.52
1:E:478:LYS:O	2:A:232:LYS:CB[3_455]	0.75	1.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:480:CYS:N	2:A:232:LYS:CE[3_455]	0.79	1.41
1:E:474:GLN:CB	2:A:143:LYS:CE[3_455]	0.92	1.28
1:E:477:ASN:N	2:A:232:LYS:O[3_455]	0.97	1.23
1:E:474:GLN:CG	2:A:143:LYS:CD[3_455]	1.01	1.19
1:E:474:GLN:CB	2:A:143:LYS:CD[3_455]	1.02	1.18
1:E:478:LYS:O	2:A:232:LYS:CG[3_455]	1.03	1.17
1:E:478:LYS:C	2:A:232:LYS:CG[3_455]	1.04	1.16
1:B:477:ASN:OD1	2:D:232:LYS:NZ[3_546]	1.05	1.15
1:E:478:LYS:CA	2:A:232:LYS:N[3_455]	1.08	1.12
1:E:480:CYS:CA	2:A:232:LYS:NZ[3_455]	1.08	1.12
1:E:478:LYS:N	2:A:232:LYS:CA[3_455]	1.12	1.08
1:E:383:SER:CB	1:B:369:TYR:CD1[1_454]	1.14	1.06
1:E:480:CYS:C	2:A:232:LYS:NZ[3_455]	1.14	1.06
1:E:478:LYS:N	2:A:232:LYS:C[3_455]	1.15	1.05
1:B:477:ASN:OD1	2:D:232:LYS:CE[3_546]	1.21	0.99
1:E:473:TYR:CD2	2:A:145:THR:CG2[3_455]	1.25	0.95
1:E:474:GLN:C	2:A:143:LYS:NZ[3_455]	1.26	0.94
1:E:477:ASN:CA	2:A:232:LYS:O[3_455]	1.26	0.94
1:E:369:TYR:CG	1:B:386:LYS:NZ[1_454]	1.27	0.93
1:E:474:GLN:CA	2:A:143:LYS:NZ[3_455]	1.27	0.93
1:E:369:TYR:CG	1:B:386:LYS:CE[1_454]	1.28	0.92
1:E:476:GLY:C	2:A:232:LYS:O[3_455]	1.28	0.92
1:E:477:ASN:O	2:A:231:ASP:O[3_455]	1.28	0.92
1:E:369:TYR:CB	1:B:386:LYS:NZ[1_454]	1.29	0.91
1:E:478:LYS:C	2:A:232:LYS:CB[3_455]	1.31	0.89
1:E:369:TYR:CB	1:B:386:LYS:CE[1_454]	1.32	0.88
1:E:477:ASN:C	2:A:232:LYS:CA[3_455]	1.32	0.88
1:E:480:CYS:N	2:A:232:LYS:CD[3_455]	1.40	0.80
1:E:474:GLN:CD	2:A:143:LYS:CG[3_455]	1.42	0.78
1:E:369:TYR:CD2	1:B:386:LYS:CE[1_454]	1.44	0.76
1:E:473:TYR:CE2	2:A:145:THR:CG2[3_455]	1.45	0.75
1:E:383:SER:OG	1:B:369:TYR:CD1[1_454]	1.46	0.74
1:E:479:PRO:N	2:A:232:LYS:CG[3_455]	1.46	0.74
1:E:383:SER:CB	1:B:369:TYR:CE1[1_454]	1.50	0.70
1:E:474:GLN:CD	2:A:143:LYS:CD[3_455]	1.51	0.69
1:E:479:PRO:C	2:A:232:LYS:CD[3_455]	1.51	0.69
1:E:369:TYR:CA	1:B:386:LYS:NZ[1_454]	1.52	0.68
1:E:479:PRO:O	2:A:232:LYS:CE[3_455]	1.52	0.68
1:E:474:GLN:CB	2:A:143:LYS:NZ[3_455]	1.54	0.66
1:E:369:TYR:CG	1:B:386:LYS:CD[1_454]	1.57	0.63
1:E:480:CYS:N	2:A:232:LYS:NZ[3_455]	1.59	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:477:ASN:N	2:A:232:LYS:C[3_455]	1.62	0.58
1:B:458:LYS:NZ	2:D:146:SER:CB[3_546]	1.63	0.57
1:E:476:GLY:O	2:A:232:LYS:O[3_455]	1.65	0.55
1:E:478:LYS:CA	2:A:232:LYS:CA[3_455]	1.66	0.54
1:E:481:ASN:N	2:A:232:LYS:NZ[3_455]	1.66	0.54
1:E:474:GLN:OE1	2:A:143:LYS:CG[3_455]	1.67	0.53
1:E:477:ASN:O	2:A:231:ASP:C[3_455]	1.70	0.50
1:E:474:GLN:O	2:A:143:LYS:NZ[3_455]	1.73	0.47
1:E:477:ASN:CA	2:A:232:LYS:C[3_455]	1.74	0.46
1:E:369:TYR:CB	1:B:386:LYS:CD[1_454]	1.76	0.44
1:B:477:ASN:CG	2:D:232:LYS:NZ[3_546]	1.78	0.42
1:E:474:GLN:OE1	2:A:143:LYS:CB[3_455]	1.79	0.41
1:E:477:ASN:O	2:A:232:LYS:CA[3_455]	1.79	0.41
1:E:477:ASN:C	2:A:232:LYS:N[3_455]	1.81	0.39
1:E:458:LYS:NZ	2:A:148:GLY:CA[3_455]	1.82	0.38
1:E:474:GLN:OE1	2:A:143:LYS:CD[3_455]	1.83	0.37
1:E:478:LYS:C	2:A:232:LYS:CA[3_455]	1.85	0.35
1:E:480:CYS:CA	2:A:232:LYS:CE[3_455]	1.86	0.34
1:E:477:ASN:C	2:A:232:LYS:O[3_455]	1.87	0.33
1:E:474:GLN:CG	2:A:143:LYS:CG[3_455]	1.88	0.32
1:E:383:SER:N	1:B:369:TYR:CE1[1_454]	1.90	0.30
1:E:478:LYS:O	2:A:232:LYS:CA[3_455]	1.90	0.30
1:E:369:TYR:CD2	1:B:386:LYS:CD[1_454]	1.91	0.29
1:E:477:ASN:O	2:A:232:LYS:N[3_455]	1.92	0.28
1:E:474:GLN:CA	2:A:143:LYS:CE[3_455]	1.93	0.27
1:E:479:PRO:CA	2:A:232:LYS:CG[3_455]	1.93	0.27
1:E:478:LYS:CA	2:A:232:LYS:CB[3_455]	1.94	0.26
1:E:369:TYR:C	1:B:386:LYS:NZ[1_454]	1.97	0.23
1:E:383:SER:CA	1:B:369:TYR:CE1[1_454]	1.97	0.23
1:E:478:LYS:C	2:A:232:LYS:N[3_455]	1.97	0.23
1:E:476:GLY:C	2:A:232:LYS:C[3_455]	1.98	0.22
1:E:480:CYS:CB	2:A:232:LYS:NZ[3_455]	1.98	0.22
1:E:383:SER:OG	1:B:369:TYR:CE1[1_454]	1.99	0.21
1:E:477:ASN:C	2:A:231:ASP:O[3_455]	1.99	0.21
1:E:479:PRO:C	2:A:232:LYS:NZ[3_455]	2.06	0.14
1:E:375:PHE:O	1:B:380:TYR:O[1_454]	2.07	0.13
1:E:477:ASN:C	2:A:231:ASP:C[3_455]	2.07	0.13
1:E:479:PRO:CA	2:A:232:LYS:CE[3_455]	2.08	0.12
1:E:369:TYR:CD1	1:B:386:LYS:NZ[1_454]	2.10	0.10
1:E:478:LYS:O	2:A:232:LYS:CD[3_455]	2.11	0.09
1:E:479:PRO:O	2:A:232:LYS:NZ[3_455]	2.13	0.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:369:TYR:CA	1:B:386:LYS:CE[1_454]	2.15	0.05
1:E:473:TYR:CG	2:A:145:THR:CG2[3_455]	2.15	0.05
1:E:369:TYR:N	1:B:386:LYS:NZ[1_454]	2.16	0.04
1:E:479:PRO:C	2:A:232:LYS:CG[3_455]	2.17	0.03
1:E:478:LYS:N	2:A:232:LYS:CB[3_455]	2.19	0.01

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	B	192/219 (88%)	187 (97%)	3 (2%)	2 (1%)	12	38
1	E	192/219 (88%)	180 (94%)	10 (5%)	2 (1%)	12	38
2	A	230/232 (99%)	214 (93%)	7 (3%)	9 (4%)	2	12
2	D	230/232 (99%)	218 (95%)	12 (5%)	0	100	100
3	C	216/218 (99%)	205 (95%)	11 (5%)	0	100	100
3	H	216/218 (99%)	212 (98%)	3 (1%)	1 (0%)	24	54
All	All	1276/1338 (95%)	1216 (95%)	46 (4%)	14 (1%)	11	36

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	110	TYR
2	A	141	SER
2	A	142	SER
1	E	481	ASN
2	A	114	PHE
1	B	483	VAL
3	H	217	CYS
1	E	480	CYS
2	A	158	ASP
2	A	105	ILE

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Mol	Chain	Res	Type
2	A	116	TYR
2	A	111	ASN
2	A	92	ALA
1	B	526	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	B	167/192 (87%)	162 (97%)	5 (3%)	36	61
1	E	167/192 (87%)	164 (98%)	3 (2%)	51	69
2	A	198/198 (100%)	192 (97%)	6 (3%)	36	61
2	D	198/198 (100%)	192 (97%)	6 (3%)	36	61
3	C	179/179 (100%)	177 (99%)	2 (1%)	65	75
3	H	179/179 (100%)	178 (99%)	1 (1%)	78	80
All	All	1088/1138 (96%)	1065 (98%)	23 (2%)	47	67

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	347	PHE
1	E	358	ILE
1	E	415	THR
2	A	24	VAL
2	A	72	LEU
2	A	114	PHE
2	A	183	VAL
2	A	196	VAL
2	A	203	LEU
1	B	358	ILE
1	B	402	ILE
1	B	452	LEU
1	B	471	GLU
1	B	517	LEU

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Mol	Chain	Res	Type
3	C	161	VAL
3	C	212	VAL
2	D	28	ILE
2	D	34	ILE
2	D	72	LEU
2	D	100	ARG
2	D	106	LEU
2	D	205	THR
3	H	142	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	185	GLN
2	A	213	ASN
1	B	339	HIS
3	C	40	GLN
3	C	41	GLN
3	C	203	HIS
2	D	1	GLN
2	D	119	GLN
2	D	214	HIS
3	H	34	ASN
3	H	41	GLN
3	H	200	GLN

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	B	601	1	14,14,15	0.22	0	17,19,21	0.59	1 (5%)
4	NAG	E	601	1	14,14,15	0.22	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	B	601	1	-	2/6/23/26	0/1/1/1
4	NAG	E	601	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	601	NAG	C1-O5-C5	2.03	114.94	112.19

There are no chirality outliers.

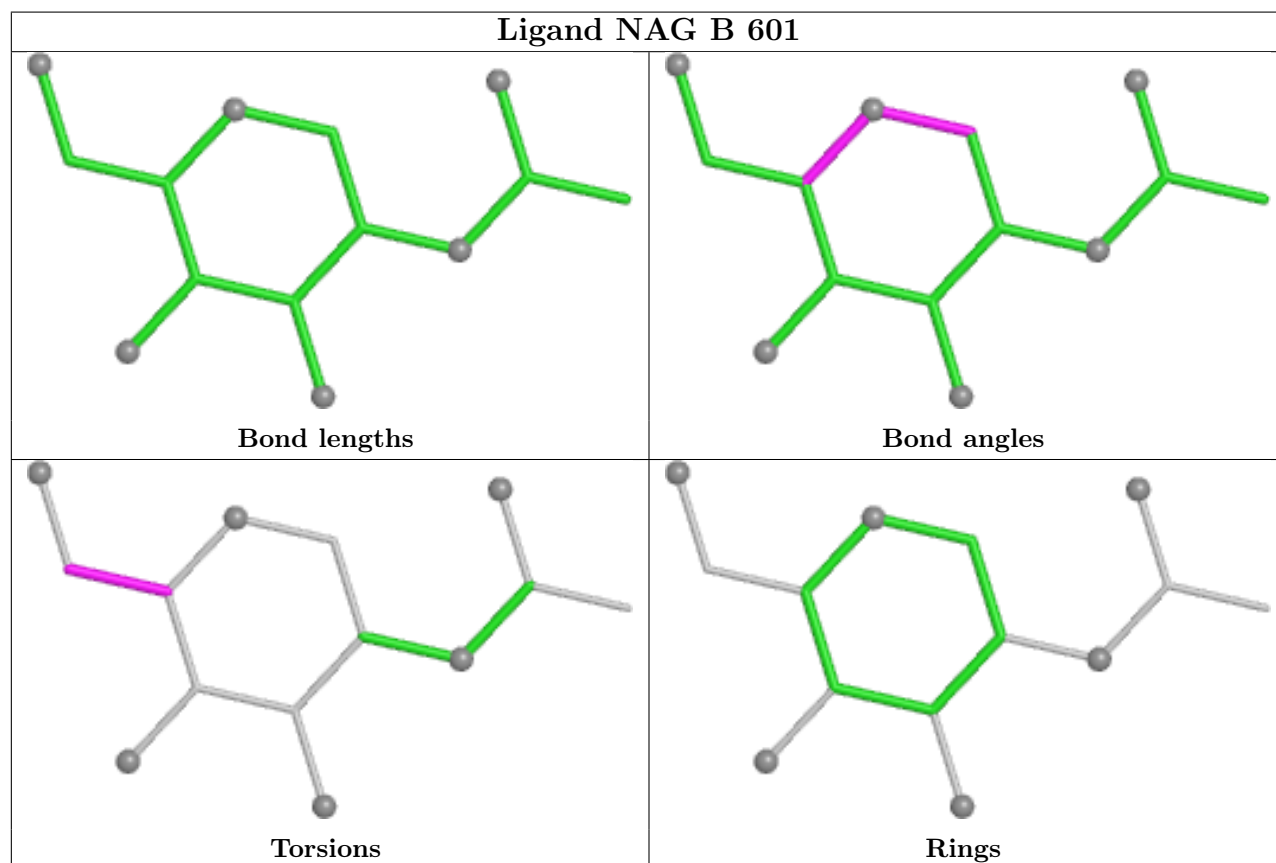
All (4) torsion outliers are listed below:

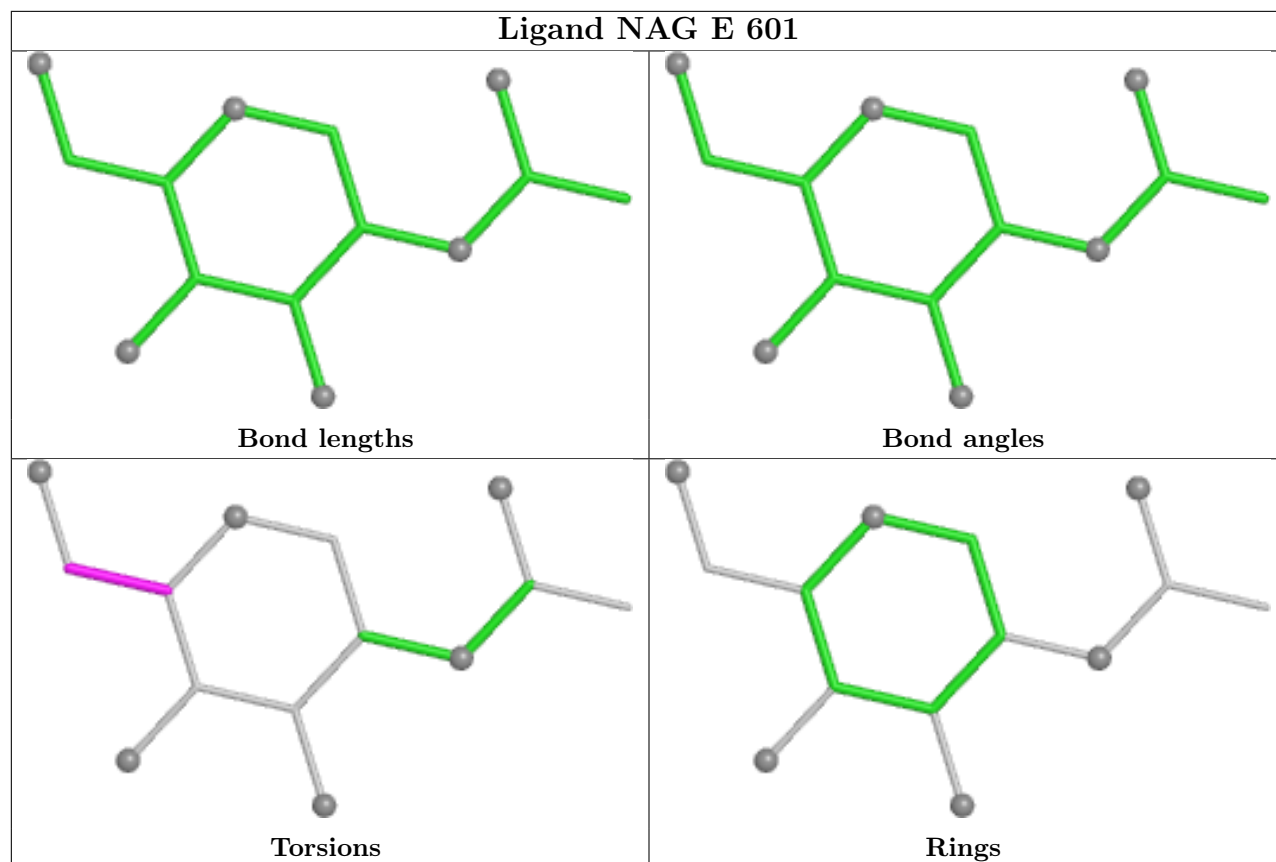
Mol	Chain	Res	Type	Atoms
4	B	601	NAG	O5-C5-C6-O6
4	E	601	NAG	O5-C5-C6-O6
4	B	601	NAG	C4-C5-C6-O6
4	E	601	NAG	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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	B	194/219 (88%)	1.14	28 (14%) 6 3	46, 67, 97, 121	0
1	E	194/219 (88%)	1.07	28 (14%) 6 3	47, 64, 109, 144	0
2	A	232/232 (100%)	1.29	33 (14%) 6 3	30, 74, 116, 147	0
2	D	232/232 (100%)	1.45	49 (21%) 2 1	53, 79, 122, 151	0
3	C	218/218 (100%)	1.05	25 (11%) 9 5	52, 71, 105, 154	0
3	H	218/218 (100%)	1.04	18 (8%) 17 9	57, 80, 104, 159	0
All	All	1288/1338 (96%)	1.18	181 (14%) 6 3	30, 74, 110, 159	0

All (181) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	113	TYR	7.9
2	A	111	ASN	7.5
2	D	109	HIS	7.5
2	D	111	ASN	5.7
2	D	104	GLY	5.4
2	A	208	TYR	5.3
2	D	103	TYR	5.0
2	D	102	ALA	5.0
2	D	146	SER	4.9
1	B	473	TYR	4.5
2	A	112	ASN	4.2
2	A	229	SER	4.1
3	C	44	GLY	4.1
1	B	374	PHE	4.0
1	E	475	ALA	3.9
2	D	105	ILE	3.9
1	E	371	PHE	3.8
2	A	101	GLY	3.8
2	A	108	ALA	3.8

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Mol	Chain	Res	Type	RSRZ
2	A	231	ASP	3.8
1	B	369	TYR	3.7
1	E	476	GLY	3.7
3	H	96	THR	3.7
3	C	199	CYS	3.7
2	D	29	PHE	3.6
1	E	473	TYR	3.6
2	A	145	THR	3.6
1	E	370	ASN	3.5
1	E	487	ASN	3.4
2	D	106	LEU	3.4
3	H	5	VAL	3.4
2	D	203	LEU	3.4
1	B	380	TYR	3.4
2	A	102	ALA	3.3
2	A	141	SER	3.3
1	B	483	VAL	3.3
2	D	142	SER	3.3
3	C	157	ASP	3.3
1	B	477	ASN	3.3
3	C	197	TYR	3.2
3	H	216	GLU	3.2
1	E	383	SER	3.2
2	A	60	ASP	3.2
3	C	193	SER	3.1
1	B	456	PHE	3.1
1	E	410	ILE	3.1
2	D	208	TYR	3.0
2	A	7	SER	3.0
3	H	191	TRP	3.0
1	E	527	PRO	3.0
2	D	144	SER	3.0
3	C	215	THR	2.9
2	D	140	PRO	2.9
2	D	107	THR	2.9
3	H	2	SER	2.9
3	H	65	PHE	2.9
1	B	410	ILE	2.9
2	D	217	SER	2.8
2	D	108	ALA	2.8
3	C	191	TRP	2.8
1	B	373	PRO	2.8

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Mol	Chain	Res	Type	RSRZ
2	A	144	SER	2.8
2	D	27	TYR	2.8
1	E	489	TYR	2.7
1	B	372	ALA	2.7
2	A	139	ALA	2.7
2	A	142	SER	2.7
2	A	107	THR	2.7
3	C	213	ALA	2.7
2	D	127	SER	2.7
3	C	196	SER	2.7
3	C	214	PRO	2.6
3	C	216	GLU	2.6
2	D	47	TRP	2.6
1	B	493	GLN	2.6
3	H	50	LEU	2.6
1	B	505	HIS	2.6
2	A	109	HIS	2.6
2	D	145	THR	2.6
3	H	217	CYS	2.6
2	A	204	GLY	2.6
3	C	194	HIS	2.5
2	D	223	LYS	2.5
3	C	5	VAL	2.5
3	C	30	ASN	2.5
1	B	346	THR	2.5
2	A	205	THR	2.5
1	B	413	GLY	2.5
1	B	474	GLN	2.5
1	B	354	ASN	2.5
2	A	209	ILE	2.5
3	H	4	LEU	2.5
1	E	363	ALA	2.5
2	D	141	SER	2.5
3	H	145	PHE	2.5
2	A	207	THR	2.5
2	D	18	LEU	2.5
2	D	163	PRO	2.5
3	C	2	SER	2.4
1	B	482	GLY	2.4
2	D	28	ILE	2.4
2	A	110	TYR	2.4
2	D	132	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
2	D	228	LYS	2.4
3	C	23	SER	2.4
1	B	432	CYS	2.4
2	A	152	LEU	2.4
2	D	216	PRO	2.4
3	H	168	THR	2.4
1	E	471	GLU	2.4
2	D	153	GLY	2.4
2	A	103	TYR	2.3
3	C	195	ARG	2.3
3	H	161	VAL	2.3
2	D	215	LYS	2.3
1	E	477	ASN	2.3
2	D	100	ARG	2.3
3	C	113	GLY	2.3
1	E	505	HIS	2.3
1	E	369	TYR	2.3
2	D	130	THR	2.3
3	C	127	SER	2.3
2	D	110	TYR	2.3
2	D	119	GLN	2.3
1	E	484	ALA	2.3
2	A	146	SER	2.3
3	C	218	SER	2.3
1	E	417	ASN	2.3
3	H	200	GLN	2.3
2	A	100	ARG	2.2
1	B	382	VAL	2.2
2	D	96	CYS	2.2
2	D	225	VAL	2.2
2	D	33	TRP	2.2
1	E	501	TYR	2.2
2	D	112	ASN	2.2
1	E	418	ILE	2.2
3	H	113	GLY	2.2
1	E	338	PHE	2.2
2	A	219	THR	2.2
2	A	213	ASN	2.2
3	C	209	GLU	2.2
2	D	26	GLY	2.2
1	E	469	SER	2.2
3	H	23	SER	2.2

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Mol	Chain	Res	Type	RSRZ
2	D	189	LEU	2.2
3	C	50	LEU	2.2
1	E	485	GLY	2.2
1	E	494	SER	2.1
3	C	111	VAL	2.1
1	B	517	LEU	2.1
2	A	6	GLN	2.1
2	D	139	ALA	2.1
1	B	383	SER	2.1
1	E	365	TYR	2.1
1	B	421	TYR	2.1
3	H	215	THR	2.1
1	E	406	GLU	2.1
2	A	193	SER	2.1
2	D	186	SER	2.1
2	D	212	VAL	2.1
1	B	334	ASN	2.1
1	B	417	ASN	2.1
2	D	218	ASN	2.1
1	B	390	LEU	2.1
2	A	105	ILE	2.1
2	D	60	ASP	2.1
3	C	45	THR	2.1
1	E	374	PHE	2.1
1	E	478	LYS	2.0
2	D	115	ASP	2.0
2	A	96	CYS	2.0
3	H	156	ALA	2.0
3	C	132	GLN	2.0
3	H	205	GLY	2.0
1	E	499	PRO	2.0
1	B	392	PHE	2.0
2	D	230	CYS	2.0
1	B	409	GLN	2.0
1	B	403	ARG	2.0
2	D	50	ILE	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 [i](#)

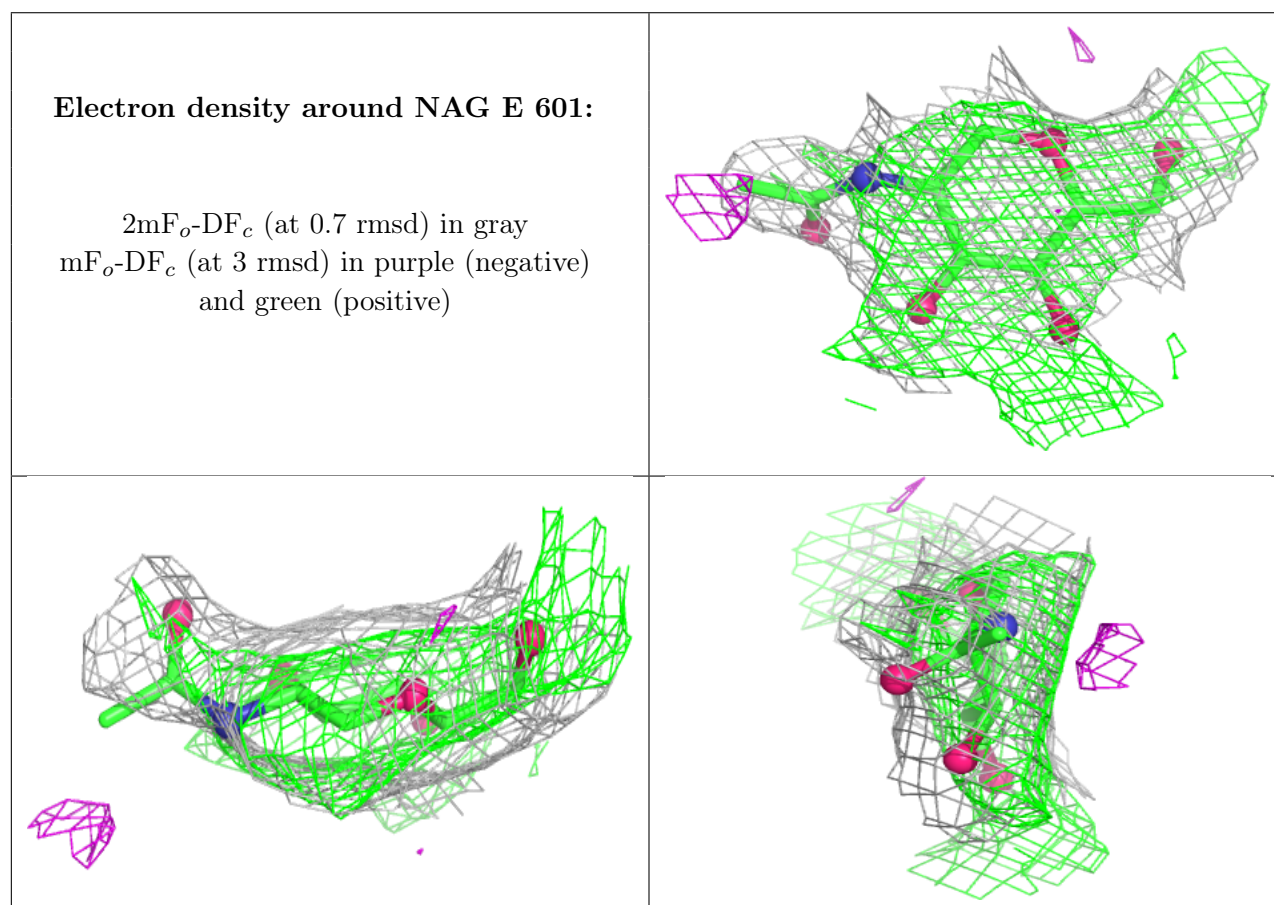
There are no oligosaccharides in this entry.

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.

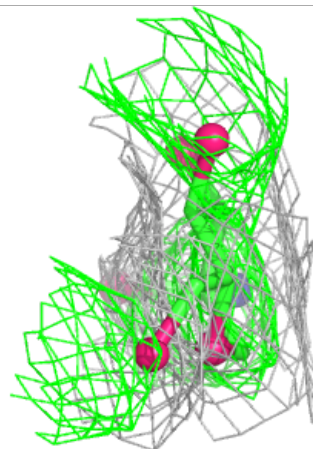
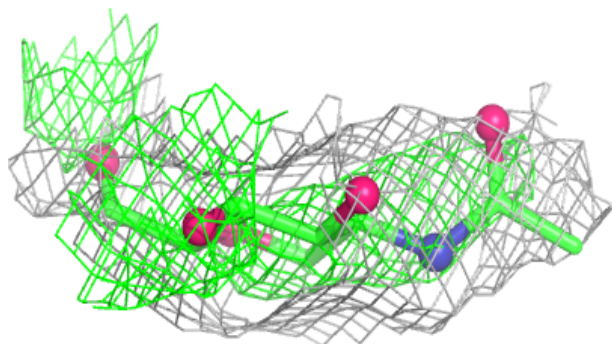
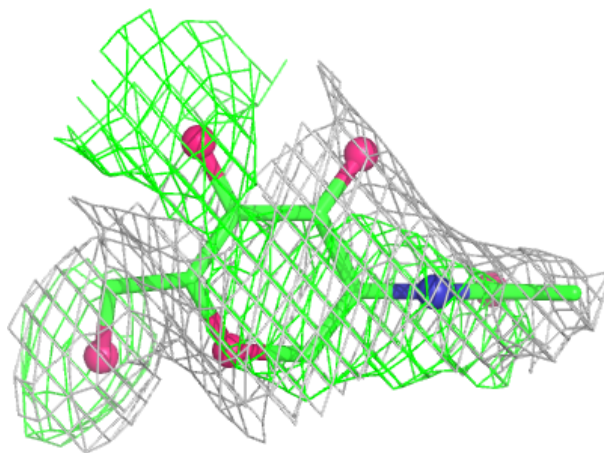
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	NAG	E	601	14/15	-	-	65,66,67,67	14
4	NAG	B	601	14/15	-	-	69,69,71,73	14

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



Electron density around NAG B 601:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.