



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:12 PM GMT

PDB ID : 4BTX
Title : Crystal structure of human vascular adhesion protein-1 in complex with pyridazinone inhibitors
Authors : Bligt-Linden, E.; Pihlavisto, M.; Szatmari, I.; Otwinowski, Z.; Smith, D.J.; Lazar, L.; Fulop, F.; Salminen, T.A.
Deposited on : 2013-06-19
Resolution : 2.78 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

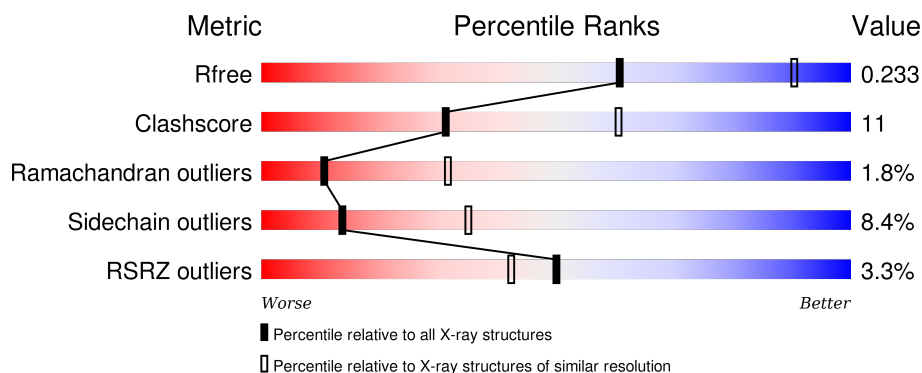
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.78 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	3004 (2.80-2.76)
Clashscore	102246	3480 (2.80-2.76)
Ramachandran outliers	100387	3423 (2.80-2.76)
Sidechain outliers	100360	3425 (2.80-2.76)
RSRZ outliers	91569	3016 (2.80-2.76)

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	737	<div> <div>3%</div> <div>72%</div> <div>20%</div> <div>.</div> <div>.</div> </div>
1	B	737	<div> <div>3%</div> <div>69%</div> <div>23%</div> <div>.</div> <div>.</div> </div>

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	NAG	A	1770	X	-	-	-

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 11686 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 MEMBRANE PRIMARY AMINE OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	709	Total	C	N	O	S	0	0	0
			5592	3589	966	1017	20			
1	B	706	Total	C	N	O	S	0	0	0
			5564	3573	958	1013	20			

- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Cu	0	0
			1	1		
2	A	1	Total	Cu	0	0
			1	1		

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Ca	0	0
			2	2		
3	A	2	Total	Ca	0	0
			2	2		

- Molecule 4 is a polymer of unknown type called SUGAR (3-MER).

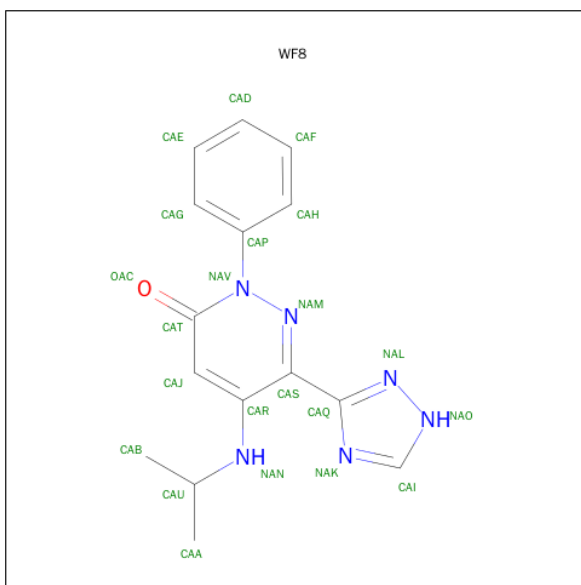
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 5 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 6 is 5-ISOPROPYLAMINO-2-PHENYL-6-(1H-1,2,4-TRIAZOL-5-YL)-3(2H)-PYRIDAZINONE (three-letter code: WF8) (formula: C₁₅H₁₆N₆O).

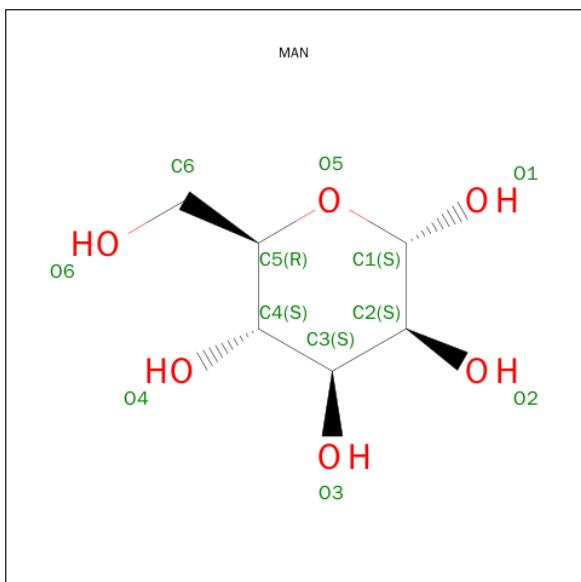


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			22	15	6	1		
6	B	1	Total	C	N	O	0	0
			22	15	6	1		

- Molecule 7 is a polymer of unknown type called SUGAR (3-MER).

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

- Molecule 8 is SUGAR (ALPHA-D-MANNOSE) (three-letter code: MAN) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	B	1	Total	C	O	0	0
			12	6	6		

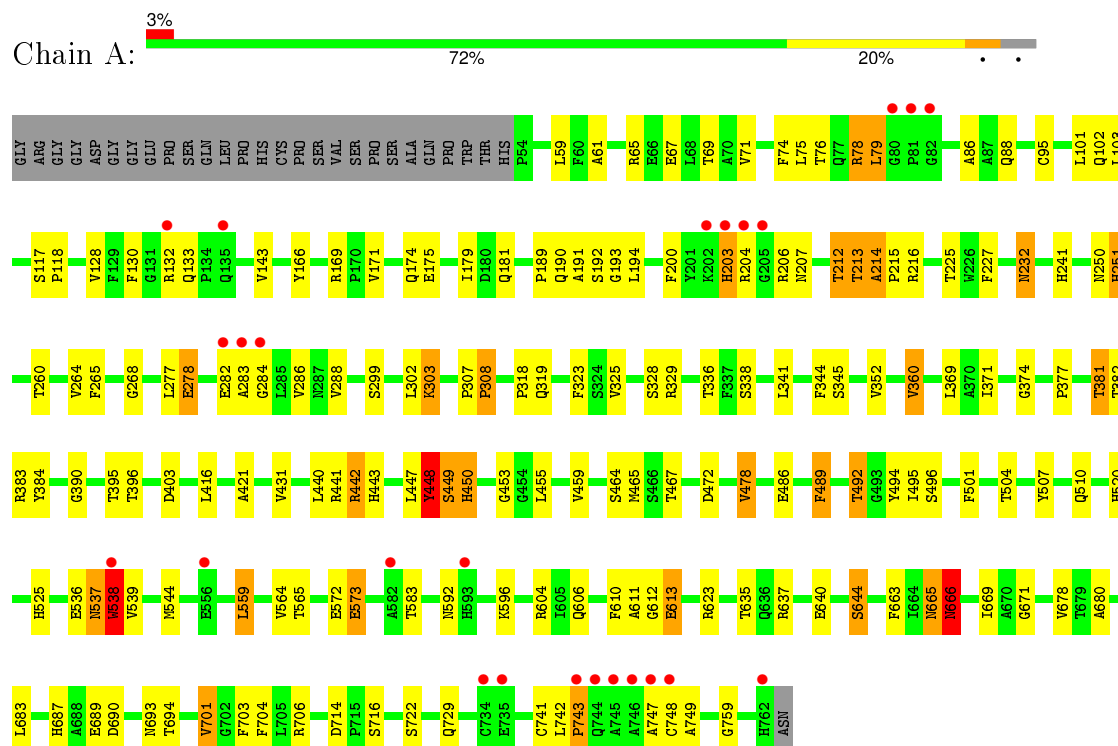
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	159	Total	O	0	0
			159	159		
9	B	133	Total	O	0	0
			133	133		

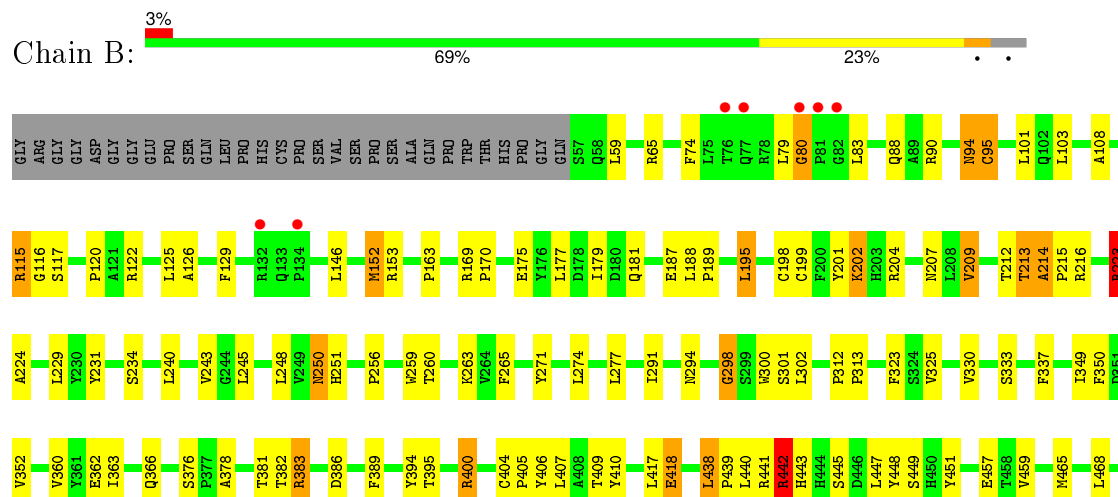
3 Residue-property plots

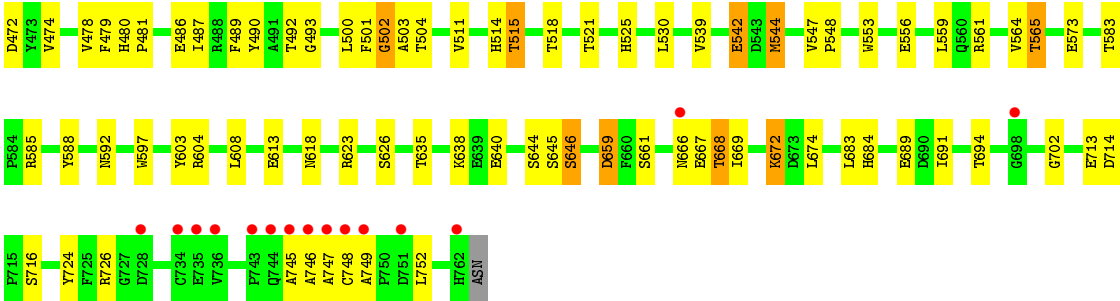
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: MEMBRANE PRIMARY AMINE OXIDASE



• Molecule 1: MEMBRANE PRIMARY AMINE OXIDASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	226.35Å 226.35Å 217.23Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.05 – 2.78 49.01 – 2.78	Depositor EDS
% Data completeness (in resolution range)	80.4 (49.05-2.78) 80.4 (49.01-2.78)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.39 (at 2.77Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.181 , 0.237 0.182 , 0.233	Depositor DCC
R_{free} test set	4125 reflections (6.56%)	DCC
Wilson B-factor (Å ²)	40.2	Xtriage
Anisotropy	0.618	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 49.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	1 of 66236 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11686	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.25% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, NAG, CA, WF8, TPQ, CU, MAN

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	0.73	3/5756 (0.1%)	0.91	10/7848 (0.1%)
1	B	0.69	0/5727	0.86	5/7811 (0.1%)
All	All	0.71	3/11483 (0.0%)	0.88	15/15659 (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	A	0	5
1	B	0	4
All	All	0	9

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	666	ASN	C-N	-9.39	1.12	1.34
1	A	665	ASN	C-N	9.10	1.54	1.34
1	A	207	ASN	CG-OD1	5.10	1.35	1.24

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	537	ASN	C-N-CA	13.35	155.07	121.70
1	A	537	ASN	O-C-N	-10.38	106.09	122.70
1	A	538	TRP	O-C-N	9.91	138.56	122.70
1	A	538	TRP	CA-C-N	-9.76	95.72	117.20
1	A	442	ARG	NE-CZ-NH1	7.47	124.04	120.30
1	B	223	ARG	NE-CZ-NH1	7.27	123.94	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	442	ARG	NE-CZ-NH2	-6.77	116.92	120.30
1	B	223	ARG	NE-CZ-NH2	-6.63	116.99	120.30
1	A	538	TRP	C-N-CA	-6.27	106.02	121.70
1	B	544	MET	CG-SD-CE	-5.28	91.75	100.20
1	A	759	GLY	N-CA-C	5.21	126.12	113.10
1	A	706	ARG	NE-CZ-NH2	-5.19	117.70	120.30
1	B	65	ARG	NE-CZ-NH1	5.15	122.87	120.30
1	A	449	SER	C-N-CA	5.14	134.56	121.70
1	A	537	ASN	CA-C-N	5.14	128.50	117.20

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	203	HIS	Peptide
1	A	213	THR	Peptide
1	A	448	TYR	Peptide
1	A	537	ASN	Mainchain
1	A	538	TRP	Mainchain
1	B	202	LYS	Peptide
1	B	213	THR	Peptide
1	B	298	GLY	Peptide
1	B	94	ASN	Peptide

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	5592	0	5339	118	0
1	B	5564	0	5301	142	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	39	0	34	0	0
5	A	42	0	38	1	0
5	B	56	0	52	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	22	0	16	0	0
6	B	22	0	16	2	0
7	B	39	0	34	2	0
8	B	12	0	12	0	0
9	A	159	0	0	8	0
9	B	133	0	0	6	0
All	All	11686	0	10842	241	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 11.

All (241) 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:214:ALA:HB2	1:A:382:THR:HA	1.54	0.89
1:B:492:THR:HG23	1:B:694:THR:O	1.73	0.88
1:B:360:VAL:HG12	1:B:530:LEU:HD23	1.58	0.83
1:B:214:ALA:HB2	1:B:382:THR:HG23	1.60	0.82
1:B:381:THR:HG22	9:B:2037:HOH:O	1.82	0.80
1:B:95:CYS:HB3	1:B:129:PHE:HB2	1.61	0.80
1:A:613:GLU:HG3	9:B:2112:HOH:O	1.82	0.80
1:A:538:TRP:CZ3	1:A:592:ASN:HB2	2.19	0.77
1:A:441:ARG:HA	1:B:492:THR:HG21	1.67	0.77
1:A:666:ASN:O	1:A:666:ASN:ND2	2.15	0.77
1:A:86:ALA:HA	1:A:95:CYS:SG	2.25	0.76
1:B:214:ALA:CB	1:B:215:PRO:CD	2.64	0.75
1:B:214:ALA:HB3	1:B:215:PRO:HD2	1.68	0.75
1:B:478:VAL:HG13	1:B:486:GLU:HB3	1.68	0.75
1:A:492:THR:HG21	1:B:441:ARG:HA	1.65	0.75
1:B:212:THR:HG21	1:B:216:ARG:NH2	2.04	0.73
1:A:680:ALA:HB1	1:A:701:VAL:HG13	1.70	0.72
1:A:623:ARG:HB2	9:A:2136:HOH:O	1.90	0.71
1:B:271:TYR:CE2	1:B:277:LEU:HD13	2.25	0.71
6:B:2000:WF8:NAK	9:B:2036:HOH:O	2.25	0.69
1:B:214:ALA:CB	1:B:215:PRO:HD2	2.21	0.69
1:B:501:PHE:O	1:B:502:GLY:O	2.09	0.69
1:A:538:TRP:HA	9:A:2116:HOH:O	1.92	0.69
1:B:101:LEU:HD23	1:B:103:LEU:HD23	1.75	0.68
1:B:223:ARG:HG3	1:B:223:ARG:HH11	1.59	0.68
1:A:449:SER:HA	1:A:450:HIS:HB2	1.75	0.67
1:B:115:ARG:O	1:B:117:SER:N	2.27	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:214:ALA:HB1	1:B:215:PRO:CD	2.25	0.66
1:B:360:VAL:CG1	1:B:530:LEU:HD23	2.26	0.66
1:B:381:THR:O	1:B:381:THR:HG22	1.98	0.64
1:B:573:GLU:OE2	1:B:666:ASN:HA	1.98	0.64
1:B:352:VAL:HB	1:B:360:VAL:HG22	1.81	0.63
1:A:214:ALA:CB	1:A:382:THR:HA	2.28	0.62
1:A:191:ALA:HA	1:A:278:GLU:CG	2.30	0.62
1:B:250:ASN:C	1:B:250:ASN:HD22	2.03	0.62
1:A:371:ILE:O	1:A:520:HIS:HA	1.99	0.62
1:A:536:GLU:HG2	9:A:2114:HOH:O	1.99	0.62
1:B:349:ILE:HD11	1:B:363:ILE:HB	1.82	0.62
1:A:494:TYR:HA	1:A:495:ILE:HD12	1.82	0.61
1:A:377:PRO:HB3	1:B:553:TRP:CZ3	2.35	0.61
1:B:514:HIS:N	9:B:2094:HOH:O	2.22	0.61
1:A:573:GLU:OE2	1:A:666:ASN:HA	2.00	0.61
1:B:175:GLU:O	1:B:179:ILE:HG13	2.01	0.61
1:A:214:ALA:HB1	1:A:215:PRO:HD2	1.83	0.60
1:A:191:ALA:HA	1:A:278:GLU:HG2	1.84	0.60
1:B:298:GLY:HA2	1:B:301:SER:O	2.01	0.60
1:A:232:ASN:ND2	5:A:1768:NAG:H83	2.18	0.59
1:A:495:ILE:HD11	1:A:690:ASP:O	2.03	0.59
1:B:714:ASP:OD2	1:B:716:SER:OG	2.15	0.59
1:B:337:PHE:CE2	1:B:479:PHE:CE2	2.91	0.58
1:B:251:HIS:HA	1:B:259:TRP:CD1	2.39	0.58
6:B:2000:WF8:HAJ	6:B:2000:WF8:HAA3	1.85	0.58
1:A:689:GLU:O	1:B:714:ASP:HB2	2.03	0.58
1:B:125:LEU:C	1:B:125:LEU:HD23	2.24	0.58
1:A:303:LYS:HG2	1:B:724:TYR:CD2	2.39	0.58
1:A:573:GLU:OE2	1:A:666:ASN:CA	2.52	0.58
1:B:440:LEU:HD22	1:B:481:PRO:HG2	1.86	0.57
1:B:212:THR:HG22	1:B:213:THR:N	2.19	0.57
1:B:438:LEU:HD22	1:B:439:PRO:HD2	1.87	0.57
1:B:214:ALA:CB	1:B:382:THR:HA	2.34	0.57
1:B:349:ILE:CD1	1:B:363:ILE:HB	2.34	0.57
1:A:59:LEU:H	1:A:59:LEU:HD23	1.68	0.57
1:B:386:ASP:HB3	1:B:468:LEU:HD13	1.86	0.57
1:B:224:ALA:HB1	1:B:248:LEU:HD11	1.87	0.56
1:B:638:LYS:NZ	1:B:667:GLU:OE2	2.38	0.56
1:A:179:ILE:HD11	1:A:251:HIS:ND1	2.21	0.56
1:B:556:GLU:HA	1:B:556:GLU:OE1	2.05	0.56
1:B:214:ALA:HB1	1:B:383:ARG:H	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:250:ASN:C	1:B:250:ASN:ND2	2.60	0.55
1:B:478:VAL:CG1	1:B:486:GLU:HB3	2.37	0.55
1:A:544:MET:CE	1:B:683:LEU:HD13	2.36	0.55
1:A:544:MET:HE2	1:B:683:LEU:HD13	1.88	0.54
1:A:714:ASP:HB2	1:B:689:GLU:O	2.07	0.54
1:B:152:MET:HG3	1:B:153:ARG:N	2.22	0.54
1:A:74:PHE:O	1:A:78:ARG:HD3	2.08	0.54
1:A:678:VAL:HG11	1:A:703:PHE:CD2	2.43	0.54
1:A:489:PHE:C	1:A:489:PHE:CD1	2.82	0.54
1:B:195:LEU:HB3	1:B:201:TYR:HB2	1.89	0.53
1:B:108:ALA:HB1	1:B:120:PRO:HG3	1.89	0.53
1:A:381:THR:HG23	9:A:2044:HOH:O	2.07	0.53
1:A:495:ILE:N	1:A:495:ILE:HD12	2.24	0.53
1:B:383:ARG:HD2	1:B:383:ARG:N	2.24	0.53
1:B:381:THR:O	1:B:381:THR:CG2	2.56	0.53
1:B:669:ILE:HG22	1:B:674:LEU:HD22	1.90	0.52
1:A:318:PRO:HB3	1:A:747:ALA:CB	2.40	0.52
1:B:447:LEU:O	1:B:449:SER:N	2.36	0.52
1:A:71:VAL:HG13	1:A:143:VAL:HG11	1.90	0.52
1:B:487:ILE:O	1:B:702:GLY:HA3	2.09	0.52
1:B:564:VAL:HG12	1:B:565:THR:N	2.25	0.51
1:A:328:SER:O	1:A:338:SER:HA	2.10	0.51
1:A:117:SER:HB3	1:A:118:PRO:HD2	1.91	0.51
1:B:163:PRO:HB3	7:B:1765:NAG:H82	1.93	0.51
1:A:440:LEU:HD23	1:A:455:LEU:HD23	1.93	0.51
1:A:225:THR:HB	1:A:227:PHE:CE1	2.46	0.51
1:A:212:THR:OG1	1:A:213:THR:N	2.44	0.51
1:B:214:ALA:CB	1:B:383:ARG:H	2.23	0.51
1:B:271:TYR:CD2	1:B:277:LEU:HD13	2.46	0.51
1:B:378:ALA:O	1:B:382:THR:OG1	2.21	0.51
1:A:572:GLU:OE1	1:A:665:ASN:OD1	2.29	0.51
1:A:396:THR:OG1	1:B:442:ARG:NH2	2.40	0.51
1:A:61:ALA:O	1:A:101:LEU:HD22	2.11	0.50
1:A:88:GLN:HA	1:A:174:GLN:HG2	1.93	0.50
1:B:457:GLU:OE2	1:B:480:HIS:ND1	2.44	0.50
1:A:250:ASN:OD1	1:A:250:ASN:C	2.48	0.50
1:B:511:VAL:HG13	1:B:691:ILE:HD11	1.91	0.50
1:B:502:GLY:O	1:B:503:ALA:HB3	2.11	0.50
1:B:129:PHE:CZ	1:B:169:ARG:HB2	2.47	0.50
1:A:478:VAL:HG13	1:A:486:GLU:HB3	1.93	0.50
1:A:241:HIS:NE2	1:A:299:SER:O	2.43	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:344:PHE:HA	1:A:390:GLY:HA2	1.94	0.50
1:A:510:GLN:HG2	1:B:597:TRP:HB3	1.94	0.49
1:A:88:GLN:HA	1:A:174:GLN:CG	2.42	0.49
1:A:189:PRO:C	1:A:191:ALA:H	2.16	0.49
1:B:492:THR:CG2	1:B:694:THR:O	2.52	0.49
1:B:214:ALA:HB3	1:B:215:PRO:CD	2.35	0.49
1:A:573:GLU:OE2	1:A:666:ASN:N	2.46	0.49
1:B:95:CYS:CB	1:B:129:PHE:HB2	2.37	0.49
1:A:403:ASP:OD1	1:B:442:ARG:HG3	2.13	0.48
1:A:369:LEU:HD12	1:A:384:TYR:O	2.13	0.48
1:A:683:LEU:HD13	1:B:544:MET:HE2	1.95	0.48
1:A:693:ASN:ND2	9:A:2146:HOH:O	2.42	0.48
1:B:366:GLN:HB3	1:B:644:SER:HB3	1.96	0.47
1:A:278:GLU:HA	1:A:278:GLU:OE1	2.15	0.47
1:A:492:THR:CG2	1:B:441:ARG:HA	2.41	0.47
1:B:212:THR:HG22	1:B:213:THR:O	2.14	0.47
1:A:559:LEU:O	1:A:559:LEU:HD23	2.14	0.47
1:B:659:ASP:OD1	1:B:659:ASP:C	2.52	0.47
1:B:350:PHE:CE1	1:B:362:GLU:HG3	2.50	0.47
1:A:544:MET:HE3	1:B:521:THR:HG22	1.97	0.47
1:B:122:ARG:O	1:B:146:LEU:HD12	2.14	0.47
1:B:187:GLU:HB3	1:B:274:LEU:HD12	1.96	0.47
1:A:214:ALA:HB1	1:A:215:PRO:CD	2.44	0.47
1:A:102:GLN:HE21	1:A:103:LEU:N	2.13	0.46
1:A:282:GLU:C	1:A:284:GLY:H	2.19	0.46
1:A:403:ASP:HB3	1:A:465:MET:CE	2.46	0.46
1:A:525:HIS:CE1	1:A:644:SER:HG	2.30	0.46
1:A:268:GLY:HA3	1:A:501:PHE:CE2	2.51	0.46
1:A:395:THR:HG22	9:A:2079:HOH:O	2.15	0.46
1:B:746:ALA:HB3	1:B:747:ALA:HA	1.98	0.46
1:B:445:SER:HB3	1:B:451:TYR:CE1	2.50	0.46
1:A:449:SER:CA	1:A:450:HIS:HB2	2.43	0.46
1:A:286:VAL:HG12	1:A:288:VAL:HG23	1.98	0.46
1:B:492:THR:HG22	1:B:493:GLY:H	1.81	0.46
1:A:538:TRP:CH2	1:A:592:ASN:HB2	2.50	0.46
1:B:323:PHE:CD1	1:B:323:PHE:C	2.89	0.46
1:A:453:GLY:HA3	1:B:302:LEU:HD13	1.98	0.46
1:B:79:LEU:O	1:B:80:GLY:O	2.34	0.46
1:A:212:THR:OG1	1:A:216:ARG:NH2	2.49	0.45
1:B:501:PHE:HB3	9:B:2046:HOH:O	2.16	0.45
1:B:256:PRO:HA	1:B:259:TRP:CE2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:67:GLU:O	1:A:71:VAL:HG23	2.15	0.45
1:A:472:ASP:OD1	1:B:443:HIS:ND1	2.35	0.45
1:B:542:GLU:HA	1:B:565:THR:O	2.16	0.45
1:B:745:ALA:HA	1:B:746:ALA:C	2.36	0.45
1:A:539:VAL:CG1	1:A:669:ILE:HD12	2.46	0.45
1:B:438:LEU:CD2	1:B:439:PRO:HD2	2.46	0.45
1:A:307:PRO:HA	1:A:308:PRO:HD2	1.82	0.45
1:A:395:THR:CG2	9:A:2079:HOH:O	2.64	0.45
1:A:352:VAL:HB	1:A:360:VAL:HG23	1.99	0.45
1:B:683:LEU:HD12	1:B:684:HIS:N	2.32	0.45
1:B:163:PRO:HB3	7:B:1765:NAG:C8	2.47	0.45
1:A:539:VAL:HG11	1:A:669:ILE:HD12	1.98	0.45
1:A:525:HIS:ND1	1:A:644:SER:OG	2.39	0.44
1:B:400:ARG:HD3	1:B:406:TYR:O	2.17	0.44
1:B:417:LEU:O	1:B:418:GLU:C	2.56	0.44
1:A:69:THR:HG23	9:A:2012:HOH:O	2.16	0.44
1:B:623:ARG:NH1	1:B:661:SER:OG	2.50	0.44
1:B:214:ALA:HB2	1:B:382:THR:HA	1.98	0.44
1:B:445:SER:CB	1:B:451:TYR:CE1	3.01	0.44
1:B:474:VAL:HB	1:B:490:TYR:HB2	1.99	0.44
1:B:465:MET:HG2	1:B:474:VAL:HG22	1.99	0.44
1:A:606:GLN:HG2	1:A:704:PHE:HB2	2.00	0.43
1:A:374:GLY:HA2	1:A:507:TYR:HB3	2.00	0.43
1:A:194:LEU:CD2	1:A:277:LEU:HD21	2.48	0.43
1:A:687:HIS:CD2	1:A:687:HIS:N	2.86	0.43
1:B:214:ALA:HB1	1:B:383:ARG:N	2.34	0.43
1:A:613:GLU:CG	9:B:2112:HOH:O	2.55	0.43
1:A:128:VAL:HG11	1:A:130:PHE:CZ	2.54	0.43
1:B:188:LEU:N	1:B:189:PRO:CD	2.81	0.43
1:A:742:LEU:N	1:A:743:PRO:HD2	2.34	0.43
1:B:108:ALA:CB	1:B:120:PRO:HG3	2.49	0.43
1:B:492:THR:HG22	1:B:493:GLY:N	2.34	0.43
1:A:494:TYR:CA	1:A:495:ILE:HD12	2.48	0.43
1:A:663:PHE:N	1:A:663:PHE:CD1	2.87	0.43
1:A:323:PHE:CE2	1:A:431:VAL:HG12	2.54	0.43
1:B:212:THR:HG21	1:B:216:ARG:HH21	1.81	0.42
1:B:298:GLY:HA2	1:B:301:SER:H	1.84	0.42
1:A:374:GLY:O	1:B:561:ARG:NH2	2.45	0.42
1:A:544:MET:HE2	1:B:683:LEU:HD22	2.02	0.42
1:B:451:TYR:HA	1:B:726:ARG:HA	2.01	0.42
1:A:171:VAL:HG13	1:A:175:GLU:HB3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:214:ALA:HB3	1:B:382:THR:HA	2.02	0.42
1:A:492:THR:CG2	1:A:694:THR:O	2.67	0.42
1:B:94:ASN:HA	1:B:129:PHE:O	2.19	0.42
1:A:193:GLY:N	1:A:278:GLU:OE2	2.53	0.42
1:A:381:THR:O	1:A:381:THR:HG23	2.19	0.42
1:B:407:LEU:HD21	1:B:752:LEU:HD23	2.02	0.42
1:A:538:TRP:CZ2	1:A:671:GLY:HA3	2.55	0.42
1:B:645:SER:OG	1:B:646:SER:N	2.53	0.42
1:A:447:LEU:O	1:A:448:TYR:HB2	2.19	0.42
1:B:294:ASN:HA	1:B:300:TRP:CZ3	2.55	0.42
1:B:214:ALA:HB2	1:B:382:THR:CG2	2.40	0.42
1:A:610:PHE:O	1:A:611:ALA:C	2.57	0.42
1:B:101:LEU:CD2	1:B:103:LEU:HD23	2.46	0.41
1:B:125:LEU:HD23	1:B:126:ALA:N	2.35	0.41
1:B:312:PRO:CB	1:B:313:PRO:CD	2.98	0.41
1:A:189:PRO:O	1:A:191:ALA:N	2.52	0.41
1:A:525:HIS:CE1	1:A:644:SER:OG	2.74	0.41
1:A:166:TYR:O	1:A:169:ARG:HD3	2.20	0.41
1:B:214:ALA:CB	1:B:383:ARG:N	2.83	0.41
1:B:746:ALA:N	1:B:747:ALA:HB2	2.35	0.41
1:A:416:LEU:HD13	1:A:421:ALA:C	2.41	0.41
1:B:240:LEU:O	1:B:515:THR:HG21	2.20	0.41
1:A:264:VAL:HG12	1:A:265:PHE:N	2.35	0.41
1:A:544:MET:CE	1:B:521:THR:HG22	2.50	0.41
1:B:451:TYR:N	1:B:451:TYR:CD1	2.89	0.41
1:B:245:LEU:HD12	1:B:265:PHE:O	2.21	0.41
1:B:502:GLY:O	1:B:503:ALA:CB	2.69	0.41
1:A:396:THR:HG21	1:A:467:THR:HB	2.03	0.41
1:B:231:TYR:HE2	1:B:291:ILE:HG22	1.86	0.41
1:B:389:PHE:HB3	1:B:394:TYR:CE2	2.56	0.41
1:A:538:TRP:HZ3	1:A:592:ASN:HB2	1.79	0.41
1:B:525:HIS:CE1	1:B:644:SER:HG	2.38	0.41
1:B:209:VAL:CG2	5:B:1769:NAG:H82	2.50	0.41
1:A:79:LEU:HA	1:A:79:LEU:HD12	1.95	0.41
1:A:117:SER:HB3	1:A:118:PRO:CD	2.50	0.41
1:B:525:HIS:CE1	1:B:644:SER:OG	2.74	0.41
1:B:668:THR:HG22	1:B:672:LYS:HD2	2.01	0.41
1:B:325:VAL:HG22	1:B:330:VAL:HG22	2.03	0.41
1:B:547:VAL:HA	1:B:548:PRO:HD2	1.98	0.41
1:A:683:LEU:HD13	1:B:544:MET:CE	2.50	0.40
1:A:200:PHE:N	1:A:200:PHE:CD1	2.88	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:404:CYS:HB2	1:B:410:TYR:OH	2.21	0.40
1:A:251:HIS:N	1:A:251:HIS:CD2	2.90	0.40
1:B:588:TYR:HB3	1:B:604:ARG:HA	2.03	0.40
1:B:500:LEU:HD12	1:B:504:THR:HG21	2.02	0.40
1:A:564:VAL:HG12	1:A:565:THR:N	2.36	0.40
1:B:209:VAL:HG23	5:B:1769:NAG:H82	2.03	0.40
1:B:74:PHE:HE2	1:B:152:MET:HB3	1.86	0.40
1:A:443:HIS:ND1	1:B:472:ASP:OD1	2.52	0.40
1:A:75:LEU:O	1:A:76:THR:C	2.60	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	706/737 (96%)	634 (90%)	58 (8%)	14 (2%)	9	28
1	B	703/737 (95%)	640 (91%)	51 (7%)	12 (2%)	11	33
All	All	1409/1474 (96%)	1274 (90%)	109 (8%)	26 (2%)	11	31

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	203	HIS
1	A	450	HIS
1	A	504	THR
1	A	666	ASN
1	B	116	GLY
1	B	214	ALA
1	B	502	GLY
1	A	190	GLN
1	A	214	ALA

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Mol	Chain	Res	Type
1	A	232	ASN
1	A	283	ALA
1	B	80	GLY
1	B	83	LEU
1	B	333	SER
1	B	418	GLU
1	B	448	TYR
1	B	618	ASN
1	B	199	CYS
1	A	448	TYR
1	A	749	ALA
1	A	206	ARG
1	B	202	LYS
1	A	743	PRO
1	A	308	PRO
1	A	612	GLY
1	B	749	ALA

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	588/610 (96%)	541 (92%)	47 (8%)	15	37
1	B	584/610 (96%)	532 (91%)	52 (9%)	12	32
All	All	1172/1220 (96%)	1073 (92%)	99 (8%)	14	35

All (99) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	ARG
1	A	78	ARG
1	A	79	LEU
1	A	132	ARG
1	A	133	GLN
1	A	181	GLN

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Mol	Chain	Res	Type
1	A	192	SER
1	A	204	ARG
1	A	212	THR
1	A	251	HIS
1	A	260	THR
1	A	278	GLU
1	A	302	LEU
1	A	303	LYS
1	A	319	GLN
1	A	325	VAL
1	A	329	ARG
1	A	336	THR
1	A	341	LEU
1	A	345	SER
1	A	360	VAL
1	A	381	THR
1	A	383	ARG
1	A	442	ARG
1	A	459	VAL
1	A	464	SER
1	A	478	VAL
1	A	489	PHE
1	A	492	THR
1	A	496	SER
1	A	559	LEU
1	A	573	GLU
1	A	583	THR
1	A	596	LYS
1	A	604	ARG
1	A	613	GLU
1	A	635	THR
1	A	637	ARG
1	A	640	GLU
1	A	644	SER
1	A	666	ASN
1	A	701	VAL
1	A	716	SER
1	A	722	SER
1	A	729	GLN
1	A	741	CYS
1	A	748	CYS
1	B	59	LEU

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Mol	Chain	Res	Type
1	B	88	GLN
1	B	90	ARG
1	B	95	CYS
1	B	115	ARG
1	B	152	MET
1	B	170	PRO
1	B	177	LEU
1	B	181	GLN
1	B	195	LEU
1	B	198	CYS
1	B	204	ARG
1	B	207	ASN
1	B	209	VAL
1	B	223	ARG
1	B	229	LEU
1	B	234	SER
1	B	243	VAL
1	B	250	ASN
1	B	260	THR
1	B	263	LYS
1	B	376	SER
1	B	383	ARG
1	B	395	THR
1	B	400	ARG
1	B	405	PRO
1	B	409	THR
1	B	438	LEU
1	B	442	ARG
1	B	459	VAL
1	B	489	PHE
1	B	515	THR
1	B	518	THR
1	B	539	VAL
1	B	542	GLU
1	B	559	LEU
1	B	565	THR
1	B	583	THR
1	B	585	ARG
1	B	592	ASN
1	B	603	TYR
1	B	608	LEU
1	B	613	GLU

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Mol	Chain	Res	Type
1	B	626	SER
1	B	635	THR
1	B	640	GLU
1	B	646	SER
1	B	659	ASP
1	B	668	THR
1	B	672	LYS
1	B	713	GLU
1	B	748	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
1	TPQ	A	471	1	13,14,15	1.31	3 (23%)	15,19,21	1.57	3 (20%)
1	TPQ	B	471	1	13,14,15	1.31	2 (15%)	15,19,21	1.69	4 (26%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	TPQ	A	471	1	-	0/4/22/24	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	TPQ	B	471	1	-	0/4/22/24	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	471	TPQ	C3-C2	-2.11	1.38	1.44
1	B	471	TPQ	C3-C2	-2.07	1.38	1.44
1	B	471	TPQ	C6-C1	2.05	1.40	1.34
1	A	471	TPQ	C6-C1	2.24	1.40	1.34
1	A	471	TPQ	C3-C4	2.46	1.39	1.35

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	471	TPQ	C1-C6-C5	-3.03	121.12	122.97
1	A	471	TPQ	O2-C2-C1	2.25	122.82	120.85
1	A	471	TPQ	O5-C5-C4	2.45	123.08	119.16
1	B	471	TPQ	C3-C2-C1	2.51	120.18	118.30
1	B	471	TPQ	O5-C5-C4	2.58	123.30	119.16
1	B	471	TPQ	C6-C1-C2	2.66	120.32	118.44
1	A	471	TPQ	C6-C1-C2	3.15	120.66	118.44

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates

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
4	NAG	A	1765	1,4	14,14,15	0.61	0	15,19,21	1.30	3 (20%)
4	NAG	A	1766	4	14,14,15	0.50	0	15,19,21	1.47	3 (20%)
4	BMA	A	1767	4	11,11,12	0.60	0	14,15,17	1.04	1 (7%)
7	NAG	B	1765	1,7	14,14,15	0.70	0	15,19,21	1.30	2 (13%)
7	NAG	B	1766	7	14,14,15	0.62	0	15,19,21	1.10	1 (6%)
7	MAN	B	1767	7	11,11,12	0.63	0	14,15,17	1.98	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
4	NAG	A	1765	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	1766	4	-	0/6/23/26	0/1/1/1
4	BMA	A	1767	4	-	0/2/19/22	0/1/1/1
7	NAG	B	1765	1,7	-	0/6/23/26	0/1/1/1
7	NAG	B	1766	7	-	0/6/23/26	0/1/1/1
7	MAN	B	1767	7	-	0/2/19/22	1/1/1/1

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	1767	MAN	C1-C2-C3	-4.46	104.27	109.54
4	A	1766	NAG	C4-C3-C2	-2.55	107.26	111.23
4	A	1766	NAG	C3-C4-C5	-2.46	105.92	110.20
7	B	1765	NAG	O7-C7-C8	-2.26	117.91	122.06
7	B	1766	NAG	O4-C4-C3	-2.25	105.28	110.34
7	B	1765	NAG	O4-C4-C3	-2.23	105.31	110.34
4	A	1766	NAG	O7-C7-C8	-2.23	117.98	122.06
4	A	1765	NAG	O4-C4-C3	-2.13	105.53	110.34
4	A	1767	BMA	C1-C2-C3	2.10	112.02	109.54
4	A	1765	NAG	C1-O5-C5	2.29	115.15	112.25
4	A	1765	NAG	O3-C3-C2	2.82	114.70	109.11
7	B	1767	MAN	O2-C2-C1	2.98	115.17	109.21
7	B	1767	MAN	C1-O5-C5	3.86	117.14	112.25

There are no chirality outliers.

There are no torsion outliers.

All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	B	1767	MAN	C1-C2-C3-C4-C5-O5

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	B	1765	NAG	2	0

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 6 are monoatomic - leaving 10 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	$\# Z > 2$
5	NAG	A	1768	1	14,14,15	0.87	0	15,19,21	1.91	6 (40%)
5	NAG	A	1770	1	14,14,15	0.78	0	15,19,21	3.75	7 (46%)
5	NAG	A	1772	-	14,14,15	0.46	0	15,19,21	1.10	1 (6%)
6	WF8	A	2000	-	19,24,24	2.71	4 (21%)	16,33,33	3.11	3 (18%)
8	MAN	B	1768	-	12,12,12	0.56	0	17,17,17	1.31	3 (17%)
5	NAG	B	1769	1	14,14,15	0.57	0	15,19,21	1.45	3 (20%)
5	NAG	B	1771	1	14,14,15	0.83	0	15,19,21	1.91	5 (33%)
5	NAG	B	1772	1	14,14,15	0.85	1 (7%)	15,19,21	1.70	2 (13%)
5	NAG	B	1773	1	14,14,15	0.58	0	15,19,21	1.05	1 (6%)
6	WF8	B	2000	-	19,24,24	2.42	3 (15%)	16,33,33	2.74	6 (37%)

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
5	NAG	A	1768	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	A	1770	1	1/1/5/7	0/6/23/26	0/1/1/1
5	NAG	A	1772	-	-	0/6/23/26	0/1/1/1
6	WF8	A	2000	-	-	0/9/12/12	0/3/3/3
8	MAN	B	1768	-	-	0/2/22/22	0/1/1/1
5	NAG	B	1769	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1771	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1772	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1773	1	-	0/6/23/26	0/1/1/1
6	WF8	B	2000	-	-	0/9/12/12	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	2000	WF8	CAS-CAQ	-8.24	1.34	1.49
6	A	2000	WF8	NAO-NAL	-7.50	1.21	1.37
6	A	2000	WF8	CAS-CAQ	-6.68	1.37	1.49
6	B	2000	WF8	CAP-NAV	-4.83	1.34	1.44
6	A	2000	WF8	CAP-NAV	-4.41	1.34	1.44
6	B	2000	WF8	NAO-NAL	-2.91	1.31	1.37
6	A	2000	WF8	CAS-NAM	2.16	1.35	1.33
5	B	1772	NAG	C1-C2	2.20	1.55	1.52

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	2000	WF8	NAK-CAQ-NAL	-10.46	106.49	114.88
6	B	2000	WF8	NAK-CAQ-NAL	-7.51	108.85	114.88
6	A	2000	WF8	CAR-CAS-NAM	-4.35	115.87	121.12
5	A	1768	NAG	C3-C4-C5	-3.76	103.64	110.20
5	A	1770	NAG	O7-C7-C8	-3.63	115.40	122.06
5	A	1770	NAG	C4-C3-C2	-3.38	105.98	111.23
5	A	1772	NAG	C2-N2-C7	-2.79	119.45	123.04
6	B	2000	WF8	CAA-CAU-CAB	-2.37	106.23	111.70
6	B	2000	WF8	CAQ-CAS-NAM	-2.20	114.52	117.47
5	A	1768	NAG	O7-C7-N2	-2.16	117.46	121.86
5	B	1771	NAG	O7-C7-C8	-2.08	118.24	122.06
5	A	1768	NAG	C3-C2-N2	-2.01	105.75	110.56
6	B	2000	WF8	CAH-CAP-CAG	-2.01	118.09	121.21
5	B	1769	NAG	O4-C4-C5	2.02	114.59	109.24
8	B	1768	MAN	O5-C5-C6	2.06	111.57	106.36
8	B	1768	MAN	C4-C3-C2	2.16	114.83	110.79
5	B	1771	NAG	C6-C5-C4	2.20	118.43	113.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	B	1768	MAN	C3-C4-C5	2.20	114.04	110.20
5	A	1768	NAG	C8-C7-N2	2.30	120.51	116.11
5	B	1769	NAG	O5-C5-C6	2.34	112.41	107.35
5	B	1771	NAG	O6-C6-C5	2.36	119.12	111.33
5	A	1770	NAG	O3-C3-C2	2.57	114.20	109.11
5	B	1773	NAG	O5-C5-C6	2.63	113.04	107.35
5	B	1771	NAG	O5-C5-C6	2.64	113.06	107.35
5	A	1768	NAG	O4-C4-C5	2.76	116.55	109.24
5	B	1769	NAG	C1-O5-C5	2.92	115.96	112.25
5	A	1768	NAG	O5-C5-C6	3.00	113.84	107.35
6	B	2000	WF8	CAA-CAU-NAN	3.14	115.70	109.67
6	A	2000	WF8	CAR-NAN-CAU	3.24	128.90	124.86
5	A	1770	NAG	C8-C7-N2	3.29	122.40	116.11
5	B	1772	NAG	C1-O5-C5	3.55	116.75	112.25
5	A	1770	NAG	C3-C2-N2	3.56	119.10	110.56
5	A	1770	NAG	O5-C5-C6	3.85	115.69	107.35
5	B	1772	NAG	C2-N2-C7	4.41	128.71	123.04
6	B	2000	WF8	CAR-NAN-CAU	4.64	130.65	124.86
5	B	1771	NAG	C1-O5-C5	4.70	118.21	112.25
5	A	1770	NAG	C1-O5-C5	11.49	126.83	112.25

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	A	1770	NAG	C1

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1768	NAG	1	0
5	B	1769	NAG	2	0
6	B	2000	WF8	2	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	708/737 (96%)	-0.19	25 (3%)	48 40	20, 37, 62, 116	0
1	B	705/737 (95%)	-0.16	22 (3%)	52 45	21, 37, 66, 168	0
All	All	1413/1474 (95%)	-0.17	47 (3%)	50 42	20, 37, 64, 168	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	745	ALA	10.6
1	B	746	ALA	9.8
1	B	744	GLN	8.0
1	B	748	CYS	6.1
1	A	203	HIS	5.3
1	B	762	HIS	5.1
1	B	743	PRO	4.7
1	A	204	ARG	4.1
1	B	735	GLU	3.9
1	B	747	ALA	3.9
1	A	745	ALA	3.5
1	B	81	PRO	3.5
1	A	744	GLN	3.5
1	A	762	HIS	3.3
1	B	82	GLY	3.2
1	A	202	LYS	3.1
1	A	735	GLU	2.9
1	A	284	GLY	2.9
1	B	134	PRO	2.8
1	B	749	ALA	2.7
1	A	747	ALA	2.7
1	A	556	GLU	2.7
1	A	282	GLU	2.6
1	A	283	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	76	THR	2.5
1	B	698	GLY	2.5
1	A	205	GLY	2.5
1	A	80	GLY	2.4
1	B	736	VAL	2.4
1	A	82	GLY	2.4
1	A	81	PRO	2.4
1	A	582	ALA	2.3
1	B	80	GLY	2.3
1	A	734	CYS	2.3
1	B	132	ARG	2.3
1	B	734	CYS	2.2
1	B	751	ASP	2.2
1	A	132	ARG	2.1
1	A	748	CYS	2.1
1	A	743	PRO	2.1
1	A	538	TRP	2.1
1	B	77	GLN	2.1
1	A	593	HIS	2.1
1	B	728	ASP	2.1
1	B	666	ASN	2.1
1	A	746	ALA	2.0
1	A	135	GLN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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
1	TPQ	A	471	14/15	0.97	0.18	-	23,29,34,39	0
1	TPQ	B	471	14/15	0.97	0.23	-	28,30,34,35	0

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
4	NAG	A	1766	14/15	0.90	0.24	1.10	50,68,86,100	0
4	NAG	A	1765	14/15	0.96	0.15	-0.06	42,49,53,59	0
7	NAG	B	1765	14/15	0.97	0.12	-0.73	35,37,40,52	0
4	BMA	A	1767	11/12	0.76	0.34	-	86,101,105,106	0
7	NAG	B	1766	14/15	0.92	0.23	-	52,69,79,92	0
7	MAN	B	1767	11/12	0.71	0.27	-	82,98,106,106	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
5	NAG	B	1769	14/15	0.85	0.18	1.89	54,64,74,76	0
6	WF8	A	2000	22/22	0.96	0.15	-0.06	36,49,55,57	0
5	NAG	A	1768	14/15	0.87	0.16	-0.14	46,50,53,54	0
6	WF8	B	2000	22/22	0.98	0.11	-1.04	27,28,34,36	0
3	CA	A	1763	1/1	0.98	0.06	-1.84	33,33,33,33	0
3	CA	B	1764	1/1	0.98	0.04	-2.68	40,40,40,40	0
3	CA	B	1763	1/1	1.00	0.05	-2.85	30,30,30,30	0
3	CA	A	1764	1/1	0.95	0.06	-3.60	46,46,46,46	0
5	NAG	A	1772	14/15	0.62	0.35	-	97,126,130,139	0
2	CU	B	1762	1/1	1.00	0.16	-	27,27,27,27	0
5	NAG	B	1772	14/15	0.62	0.55	-	92,104,116,117	0
2	CU	A	1762	1/1	0.99	0.15	-	34,34,34,34	0
5	NAG	B	1773	14/15	0.60	0.42	-	81,93,100,100	0
5	NAG	B	1771	14/15	0.71	0.27	-	66,92,104,107	0
8	MAN	B	1768	12/12	0.65	0.45	-	108,123,131,133	0
5	NAG	A	1770	14/15	0.66	0.51	-	72,94,105,107	0

6.5 Other polymers [i](#)

There are no such residues in this entry.