



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 03:08 am GMT

PDB ID : 4O4U
Title : Crystal structure of the vaccine antigen Transferrin Binding Protein B (TbpB) mutant Trp-176-Ala from Haemophilus parasuis Hp5
Authors : Calmettes, C.; Yu, R.H.; Schryvers, A.B.; Moraes, T.F.
Deposited on : 2013-12-19
Resolution : 2.64 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

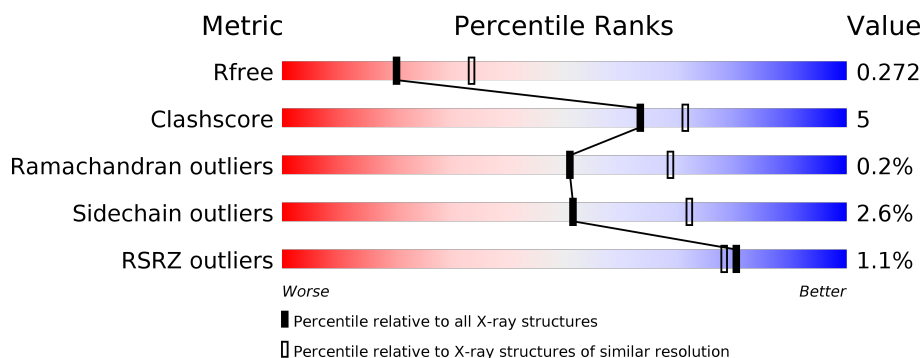
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.64 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}	100719	1044 (2.66-2.62)
Clashscore	112137	1092 (2.66-2.62)
Ramachandran outliers	110173	1077 (2.66-2.62)
Sidechain outliers	110143	1077 (2.66-2.62)
RSRZ outliers	101464	1047 (2.66-2.62)

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	532	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 12%, green 82%, grey 5%);"></div> <div style="display: flex; justify-content: space-between; font-size: 0.8em;"> % 82% 12% • 5% </div> </div>
1	B	532	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 12%, green 81%, grey 6%);"></div> <div style="display: flex; justify-content: space-between; font-size: 0.8em;"> % 81% 12% • 6% </div> </div>
1	C	532	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 2%, orange 1%, yellow 12%, green 81%, grey 6%);"></div> <div style="display: flex; justify-content: space-between; font-size: 0.8em;"> 2% 81% 12% • 6% </div> </div>
1	D	532	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 13%, green 82%, grey 5%);"></div> <div style="display: flex; justify-content: space-between; font-size: 0.8em;"> % 82% 13% 5% </div> </div>

2 Entry composition [i](#)

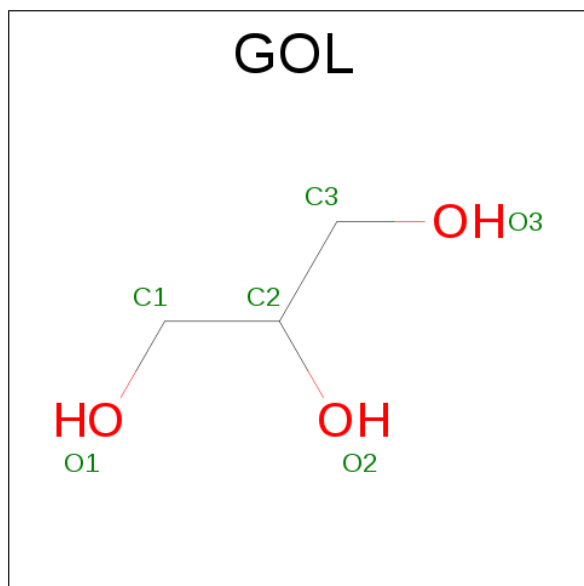
There are 3 unique types of molecules in this entry. The entry contains 15892 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 TbpB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	505	Total	C	N	O	S	0	0	0
			3912	2465	655	782	10			
1	B	500	Total	C	N	O	S	0	0	0
			3871	2441	644	776	10			
1	C	499	Total	C	N	O	S	0	0	0
			3861	2434	643	774	10			
1	D	507	Total	C	N	O	S	0	2	0
			3944	2487	659	788	10			

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	C	O	0	0
			6	3	3		

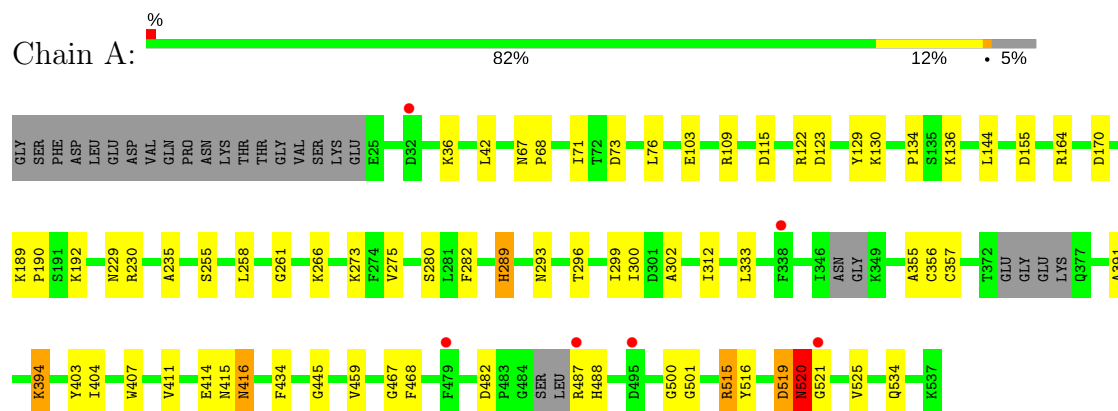
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	79	Total 79	O 79	0	0
3	B	91	Total 91	O 91	0	0
3	C	48	Total 48	O 48	0	0
3	D	80	Total 80	O 80	0	0

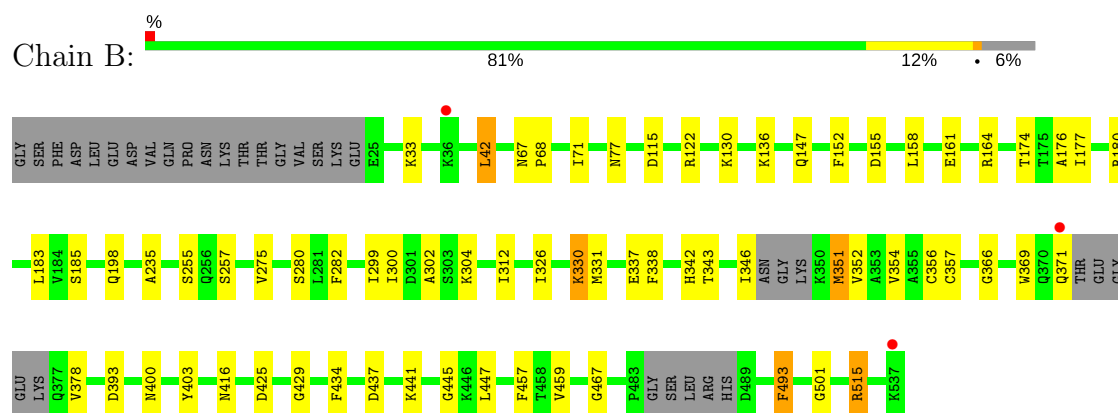
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

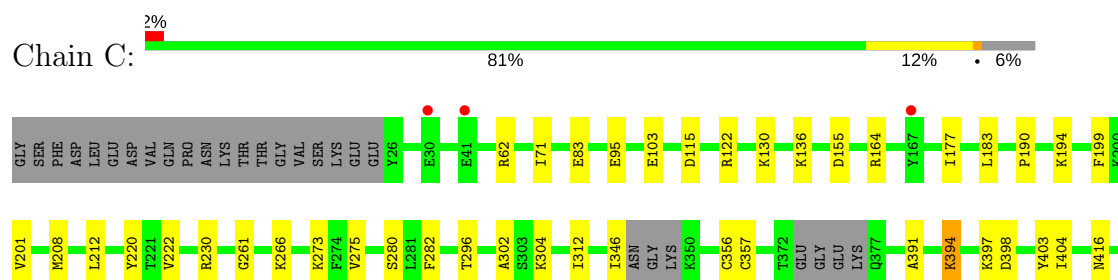
• Molecule 1: TbpB

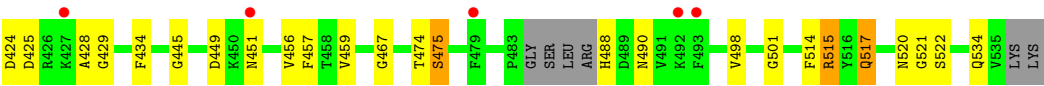


• Molecule 1: TbpB

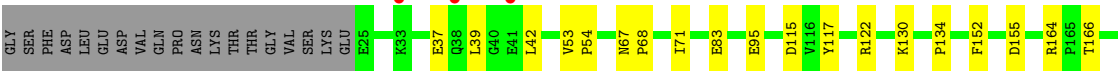
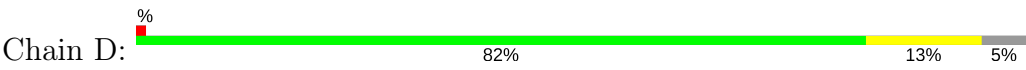


• Molecule 1: TbpB





● Molecule 1: TbpB



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.03Å 135.70Å 103.96Å 90.00° 100.05° 90.00°	Depositor
Resolution (Å)	48.39 – 2.64 48.39 – 2.64	Depositor EDS
% Data completeness (in resolution range)	99.5 (48.39-2.64) 74.3 (48.39-2.64)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 2.65Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.214 , 0.266 0.218 , 0.272	Depositor DCC
R_{free} test set	2393 reflections (5.21%)	DCC
Wilson B-factor (Å ²)	51.6	Xtriage
Anisotropy	0.272	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 23.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	15892	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 37.79 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 4.0646e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL

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	0.24	0/3987	0.42	0/5372
1	B	0.25	0/3945	0.44	0/5317
1	C	0.25	0/3936	0.43	0/5308
1	D	0.25	0/4024	0.44	0/5424
All	All	0.24	0/15892	0.43	0/21421

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3912	0	3795	38	0
1	B	3871	0	3752	41	0
1	C	3861	0	3734	34	0
1	D	3944	0	3828	38	0
2	D	6	0	8	1	0
3	A	79	0	0	2	0
3	B	91	0	0	2	0
3	C	48	0	0	1	0
3	D	80	0	0	1	0
All	All	15892	0	15117	149	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 5.

All (149) 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:D:227:ARG:NH1	1:D:232:THR:OG1	2.21	0.72
1:D:155:ASP:HB3	1:D:164:ARG:HH22	1.55	0.71
1:B:352:VAL:HG21	1:B:371:GLN:NE2	2.09	0.68
1:A:391:ALA:HB3	1:A:394:LYS:HB2	1.75	0.67
1:D:166:THR:H	2:D:601:GOL:H2	1.59	0.67
1:A:115:ASP:OD2	1:A:122:ARG:NH1	2.29	0.66
1:C:155:ASP:HB3	1:C:164:ARG:HH22	1.60	0.66
1:B:403:TYR:HB2	1:B:434:PHE:HB2	1.77	0.65
1:D:166:THR:HG22	1:D:176:ALA:HA	1.79	0.65
1:C:312:ILE:HG22	1:C:515:ARG:HH21	1.62	0.64
1:B:331:MET:HE3	1:B:342:HIS:HB2	1.79	0.64
1:D:190:PRO:HB3	1:D:194:LYS:HE2	1.80	0.64
1:A:515:ARG:NH1	3:A:655:HOH:O	2.32	0.63
1:A:144:LEU:HD13	1:A:289:HIS:HD2	1.63	0.62
1:C:467:GLY:HA2	1:C:501:GLY:HA2	1.82	0.62
1:A:445:GLY:N	1:A:459:VAL:O	2.32	0.60
1:D:425:ASP:HB3	1:D:429:GLY:HA3	1.83	0.60
1:D:155:ASP:HB2	1:D:280:SER:HA	1.83	0.60
1:B:400:ASN:ND2	1:B:437:ASP:OD1	2.31	0.59
1:B:393:ASP:HB2	1:D:205:THR:HG22	1.84	0.59
1:D:95:GLU:HG2	1:D:117:TYR:HE2	1.67	0.59
1:B:77:ASN:ND2	3:B:690:HOH:O	2.36	0.58
1:A:403:TYR:HB2	1:A:434:PHE:HB2	1.86	0.57
1:C:403:TYR:HB2	1:C:434:PHE:HB2	1.86	0.57
1:B:155:ASP:HB3	1:B:164:ARG:HH22	1.69	0.57
1:D:403:TYR:HB2	1:D:434:PHE:HB2	1.87	0.56
1:C:115:ASP:OD2	1:C:122:ARG:NH1	2.36	0.56
1:C:425:ASP:HB3	1:C:429:GLY:HA3	1.88	0.55
1:D:95:GLU:HG2	1:D:117:TYR:CE2	2.40	0.55
1:A:266:LYS:HB3	1:A:296:THR:HG21	1.88	0.55
1:C:190:PRO:HB3	1:C:194:LYS:HE2	1.88	0.55
1:D:498:VAL:HG22	1:D:514:PHE:HB3	1.90	0.54
1:A:333:LEU:HD22	1:A:355:ALA:HB2	1.89	0.54
1:C:445:GLY:N	1:C:459:VAL:O	2.39	0.54
1:A:71:ILE:HG21	1:A:130:LYS:HE2	1.89	0.53
1:B:147:GLN:NE2	1:B:198:GLN:OE1	2.41	0.53
1:B:177:ILE:HG23	1:B:183:LEU:HD13	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:122:ARG:NH1	3:D:707:HOH:O	2.35	0.53
1:A:468:PHE:CZ	1:A:500:GLY:HA3	2.44	0.52
1:C:273:LYS:HD3	1:C:282:PHE:HE1	1.75	0.52
1:B:330:LYS:H	1:B:330:LYS:HE2	1.74	0.52
1:A:467:GLY:HA2	1:A:501:GLY:HA2	1.91	0.52
1:D:467:GLY:HA2	1:D:501:GLY:HA2	1.92	0.51
1:B:312:ILE:HG22	1:B:515:ARG:HH21	1.74	0.51
1:B:337:GLU:HG3	1:B:338:PHE:H	1.76	0.51
1:C:275:VAL:HG22	1:C:282:PHE:HD1	1.76	0.51
1:A:144:LEU:HD13	1:A:289:HIS:CD2	2.46	0.51
1:B:467:GLY:HA2	1:B:501:GLY:HA2	1.92	0.51
1:D:152:PHE:CE2	1:D:185:SER:HB3	2.46	0.50
1:D:115:ASP:OD2	1:D:122:ARG:NH1	2.40	0.50
1:B:416:ASN:OD1	1:B:416:ASN:N	2.45	0.50
1:C:520:ASN:HA	1:C:522:SER:H	1.75	0.50
1:A:312:ILE:HG22	1:A:515:ARG:HH21	1.76	0.50
1:B:71:ILE:HG21	1:B:130:LYS:HE2	1.94	0.50
1:B:326:ILE:HG21	1:B:351:MET:HG2	1.93	0.50
1:B:180:ARG:NE	3:B:611:HOH:O	2.44	0.49
1:D:411:VAL:HG11	1:D:491:VAL:HG21	1.94	0.49
1:C:404:ILE:HG12	1:C:534:GLN:HG3	1.95	0.49
1:B:152:PHE:CE2	1:B:185:SER:HB3	2.48	0.49
1:A:155:ASP:HB2	1:A:280:SER:HA	1.95	0.49
1:C:520:ASN:HA	1:C:522:SER:N	2.29	0.48
1:C:391:ALA:HB3	1:C:394:LYS:HB2	1.95	0.48
1:D:404:ILE:HG12	1:D:534:GLN:HG3	1.96	0.48
1:D:71:ILE:HG21	1:D:130:LYS:HE2	1.96	0.47
1:D:275:VAL:HG22	1:D:282:PHE:HD1	1.79	0.47
1:C:428:ALA:HB1	1:C:451:ASN:HA	1.97	0.47
1:C:520:ASN:N	1:C:521:GLY:HA3	2.28	0.47
1:A:404:ILE:HG12	1:A:534:GLN:HG3	1.97	0.47
1:D:356:CYS:HA	1:D:357:CYS:HA	1.69	0.47
1:D:239:ASP:O	1:D:249:ASN:ND2	2.48	0.47
1:D:360:LEU:O	1:D:364:LYS:NZ	2.47	0.47
1:D:424:ASP:OD1	1:D:424:ASP:N	2.44	0.47
1:B:425:ASP:HB3	1:B:429:GLY:HA3	1.97	0.46
1:B:136:LYS:HE3	1:B:302:ALA:HB3	1.96	0.46
1:C:122:ARG:NH1	3:C:607:HOH:O	2.34	0.46
1:C:266:LYS:HB3	1:C:296:THR:HG21	1.97	0.46
1:A:356:CYS:HA	1:A:357:CYS:HA	1.74	0.46
1:B:115:ASP:OD2	1:B:122:ARG:NH1	2.38	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:447:LEU:HB2	1:B:457:PHE:HB2	1.97	0.46
1:B:445:GLY:N	1:B:459:VAL:O	2.47	0.46
1:D:155:ASP:O	1:D:164:ARG:NH1	2.48	0.46
1:D:266:LYS:HB3	1:D:296:THR:HG21	1.97	0.46
1:C:424:ASP:OD1	1:C:424:ASP:N	2.41	0.46
1:B:352:VAL:HG21	1:B:371:GLN:CD	2.36	0.45
1:C:201:VAL:HG22	1:C:208:MET:HG3	1.98	0.45
1:A:273:LYS:HD2	1:A:282:PHE:HE1	1.82	0.45
1:A:275:VAL:HG22	1:A:282:PHE:HD1	1.81	0.45
1:B:42:LEU:HD12	1:B:257:SER:OG	2.16	0.45
1:C:416:ASN:N	1:C:416:ASN:OD1	2.50	0.45
1:C:449:ASP:HB2	1:C:456:VAL:HG11	1.97	0.45
1:B:155:ASP:HB2	1:B:280:SER:HA	1.99	0.45
1:B:337:GLU:HG3	1:B:338:PHE:N	2.32	0.45
1:A:487:ARG:HG2	1:A:488:HIS:H	1.81	0.45
1:B:369:TRP:HB2	1:B:378:VAL:HG13	1.98	0.45
1:B:299:ILE:HG13	1:B:300:ILE:HG13	2.00	0.44
1:B:457:PHE:HE1	1:B:493:PHE:HE2	1.64	0.44
1:D:273:LYS:HD2	1:D:282:PHE:HE1	1.82	0.44
1:D:266:LYS:HD3	1:D:296:THR:HG21	2.00	0.44
1:B:158:LEU:HD21	1:B:180:ARG:HD3	1.99	0.44
1:A:134:PRO:HB2	1:A:229:ASN:ND2	2.32	0.44
1:C:177:ILE:HG23	1:C:183:LEU:HD13	1.98	0.44
1:A:73:ASP:OD1	1:A:130:LYS:NZ	2.51	0.43
1:B:275:VAL:HG22	1:B:282:PHE:HD1	1.84	0.43
1:A:76:LEU:HD22	1:A:109:ARG:HG2	2.00	0.43
1:A:235:ALA:HB3	1:A:255:SER:HB3	2.01	0.43
1:A:36:LYS:HE3	1:A:36:LYS:HB2	1.75	0.43
1:A:411:VAL:HG13	1:A:525:VAL:HG22	2.01	0.43
1:C:155:ASP:HB2	1:C:280:SER:HA	2.00	0.43
1:B:235:ALA:HB3	1:B:255:SER:HB3	2.00	0.43
1:C:230:ARG:HA	1:C:261:GLY:HA2	2.01	0.43
1:A:134:PRO:HB2	1:A:229:ASN:HD21	1.84	0.42
1:D:83:GLU:H	1:D:83:GLU:CD	2.22	0.42
1:C:356:CYS:HA	1:C:357:CYS:HA	1.71	0.42
1:B:33:LYS:H	1:B:33:LYS:HG2	1.60	0.42
1:C:71:ILE:HG21	1:C:130:LYS:HE2	2.00	0.42
1:B:354:VAL:O	1:B:366:GLY:HA3	2.18	0.42
1:C:498:VAL:HG22	1:C:514:PHE:HB3	2.01	0.42
1:D:432:THR:HG22	1:D:447:LEU:HG	2.02	0.42
1:D:134:PRO:HG3	1:D:316:THR:HG23	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:174:THR:HG22	1:B:176:ALA:H	1.85	0.42
1:A:155:ASP:O	1:A:164:ARG:NH1	2.51	0.42
1:A:230:ARG:HA	1:A:261:GLY:HA2	2.02	0.42
1:A:170:ASP:O	1:A:192:LYS:HG3	2.20	0.42
1:A:519:ASP:N	1:A:519:ASP:OD1	2.53	0.42
1:C:474:THR:O	1:C:475:SER:HB2	2.20	0.42
1:D:349:LYS:N	1:D:370:GLN:HE21	2.18	0.42
1:B:343:THR:HG22	1:B:352:VAL:HG22	2.01	0.41
1:C:136:LYS:HE3	1:C:302:ALA:HB3	2.01	0.41
1:D:253:VAL:HG21	1:D:278:ASP:OD2	2.20	0.41
1:A:299:ILE:HG13	1:A:300:ILE:HG13	2.02	0.41
1:A:520:ASN:HA	1:A:521:GLY:HA2	1.70	0.41
1:B:67:ASN:HA	1:B:68:PRO:HD3	1.85	0.41
1:D:53:VAL:HA	1:D:54:PRO:HD3	1.93	0.41
1:C:212:LEU:HB2	1:C:220:TYR:HB2	2.02	0.41
1:A:414:GLU:HB3	1:A:415:ASN:H	1.55	0.41
1:D:189:LYS:HA	1:D:190:PRO:HD3	1.93	0.41
1:B:356:CYS:HA	1:B:357:CYS:HA	1.74	0.41
1:C:456:VAL:HG23	1:C:457:PHE:CD2	2.56	0.41
1:A:136:LYS:HE2	1:A:302:ALA:HB3	2.03	0.41
1:A:416:ASN:HB2	3:A:673:HOH:O	2.21	0.41
1:B:441:LYS:HA	1:B:441:LYS:HD3	1.91	0.41
1:A:67:ASN:HA	1:A:68:PRO:HD3	1.89	0.41
1:D:456:VAL:HG23	1:D:457:PHE:CD2	2.56	0.41
1:A:293:ASN:HB3	1:C:397:LYS:HD3	2.03	0.40
1:D:67:ASN:HA	1:D:68:PRO:HD3	1.87	0.40
1:D:173:TYR:CD1	1:D:189:LYS:HE2	2.56	0.40
1:A:189:LYS:HA	1:A:190:PRO:HD3	1.86	0.40
1:B:161:GLU:N	1:B:161:GLU:OE1	2.47	0.40
1:C:199:PHE:HE1	1:C:222:VAL:HG21	1.86	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	497/532 (93%)	476 (96%)	20 (4%)	1 (0%)	51	69
1	B	492/532 (92%)	478 (97%)	14 (3%)	0	100	100
1	C	491/532 (92%)	471 (96%)	18 (4%)	2 (0%)	38	54
1	D	503/532 (94%)	487 (97%)	16 (3%)	0	100	100
All	All	1983/2128 (93%)	1912 (96%)	68 (3%)	3 (0%)	51	69

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	520	ASN
1	C	517	GLN
1	C	475	SER

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	417/440 (95%)	403 (97%)	14 (3%)	42	62
1	B	413/440 (94%)	406 (98%)	7 (2%)	66	82
1	C	412/440 (94%)	400 (97%)	12 (3%)	48	68
1	D	421/440 (96%)	411 (98%)	10 (2%)	54	74
All	All	1663/1760 (94%)	1620 (97%)	43 (3%)	51	72

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	LEU
1	A	103	GLU
1	A	123	ASP
1	A	129	TYR
1	A	258	LEU

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Mol	Chain	Res	Type
1	A	289	HIS
1	A	394	LYS
1	A	407	TRP
1	A	416	ASN
1	A	482	ASP
1	A	515	ARG
1	A	516	TYR
1	A	519	ASP
1	A	520	ASN
1	B	42	LEU
1	B	304	LYS
1	B	330	LYS
1	B	346	ILE
1	B	351	MET
1	B	493	PHE
1	B	515	ARG
1	C	62	ARG
1	C	83	GLU
1	C	95	GLU
1	C	103	GLU
1	C	304	LYS
1	C	346	ILE
1	C	394	LYS
1	C	398	ASP
1	C	488	HIS
1	C	490	ASN
1	C	515	ARG
1	C	517	GLN
1	D	37	GLU
1	D	39	LEU
1	D	42	LEU
1	D	213	SER
1	D	304	LYS
1	D	372	THR
1	D	407	TRP
1	D	486	LEU
1	D	490	ASN
1	D	515	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	415	ASN
1	D	400	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	$\# Z > 2$
2	GOL	D	601	-	5,5,5	0.32	0	5,5,5	0.22	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
2	GOL	D	601	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

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
2	D	601	GOL	1	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	505/532 (94%)	-0.21	6 (1%) 79 77	13, 30, 66, 91	0
1	B	500/532 (93%)	-0.24	3 (0%) 89 88	7, 22, 60, 106	0
1	C	499/532 (93%)	-0.02	8 (1%) 72 69	10, 39, 80, 103	0
1	D	507/532 (95%)	-0.15	5 (0%) 82 80	9, 29, 65, 97	0
All	All	2011/2128 (94%)	-0.16	22 (1%) 80 78	7, 30, 68, 106	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	479	PHE	4.4
1	C	492	LYS	4.3
1	C	493	PHE	3.8
1	B	537	LYS	3.7
1	C	427	LYS	3.5
1	C	30	GLU	3.3
1	D	537	LYS	3.2
1	A	521	GLY	3.1
1	B	36	LYS	2.9
1	D	38	GLN	2.6
1	A	487	ARG	2.5
1	A	32	ASP	2.5
1	D	33	LYS	2.4
1	A	495	ASP	2.3
1	A	479	PHE	2.3
1	A	338	PHE	2.2
1	C	167	TYR	2.1
1	B	371	GLN	2.1
1	C	41	GLU	2.1
1	D	346	ILE	2.1
1	C	451	ASN	2.0
1	D	41	GLU	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](#)

There are no carbohydrates 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. 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(\AA^2)	Q<0.9
2	GOL	D	601	6/6	0.91	0.20	-0.03	37,41,45,48	0

6.5 Other polymers [i](#)

There are no such residues in this entry.