



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 02:31 AM UTC

PDB ID : 6YQ4 / pdb_00006yq4
Title : Crystal structure of Fusobacterium nucleatum tannase
Authors : Mancheno, J.M.; Anguita, J.; Rodriguez, H.
Deposited on : 2020-04-16
Resolution : 2.40 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 : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

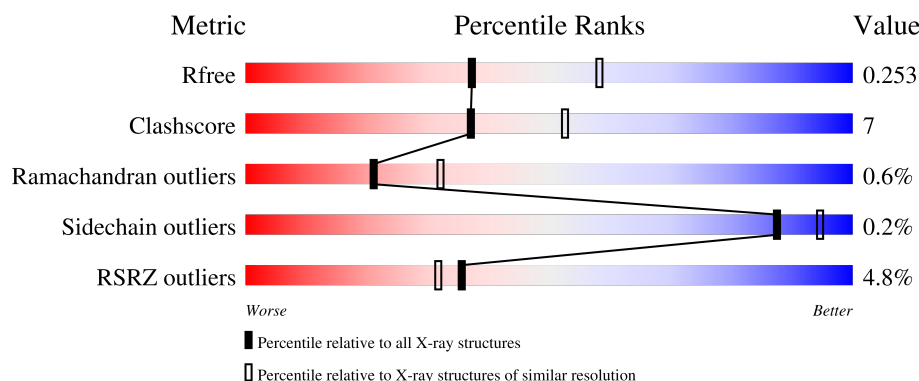
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	A	495	3% 81% 15% .
1	B	495	6% 79% 17% .
1	C	495	5% 82% 14% ..

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 11702 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 Tannase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	478	Total	C	N	O	S	0	1	0
			3752	2377	635	729	11			
1	B	480	Total	C	N	O	S	0	0	0
			3763	2382	638	732	11			
1	C	478	Total	C	N	O	S	0	1	0
			3749	2374	635	729	11			

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).

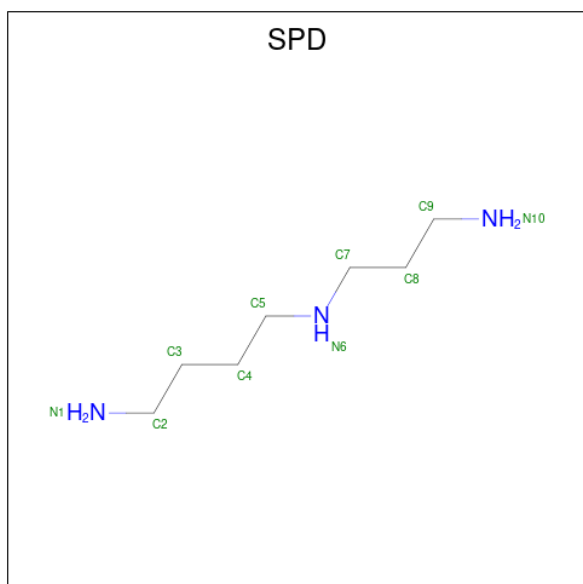


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		

- Molecule 4 is SPERMIDINE (CCD ID: SPD) (formula: C₇H₁₉N₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	N	0	0
			10	7	3		

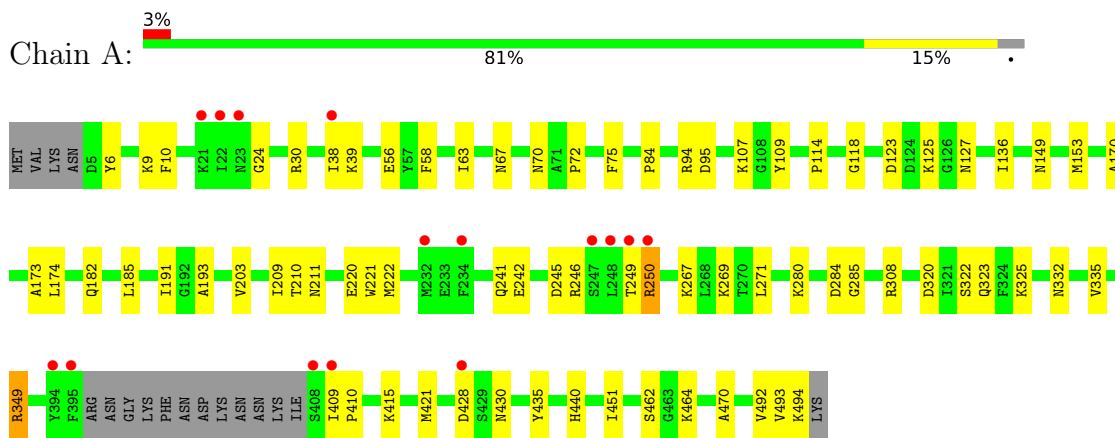
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	180	Total	O	0	0
			180	180		
5	B	107	Total	O	0	0
			107	107		
5	C	120	Total	O	0	0
			120	120		

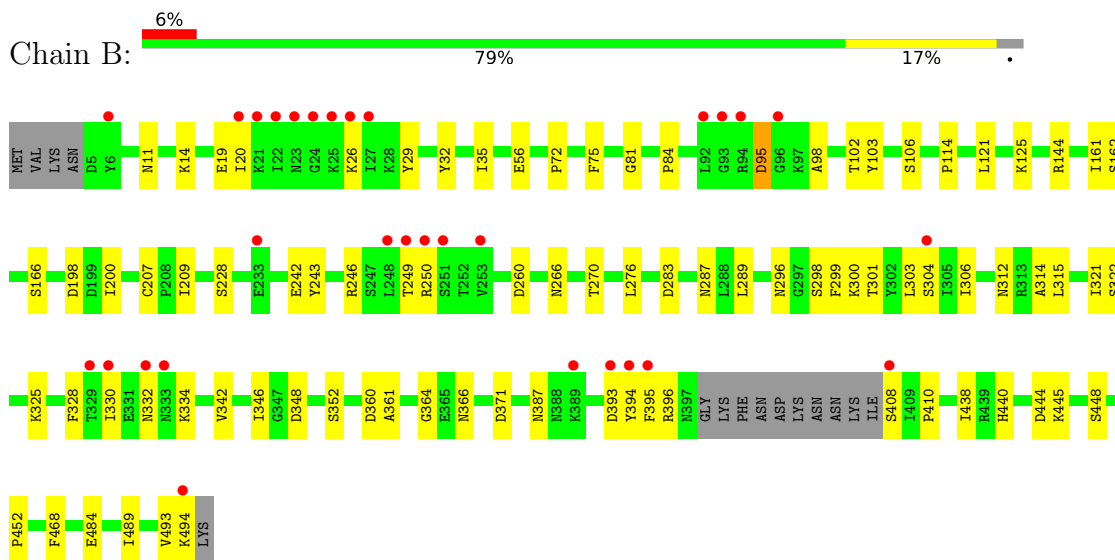
3 Residue-property plots

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.

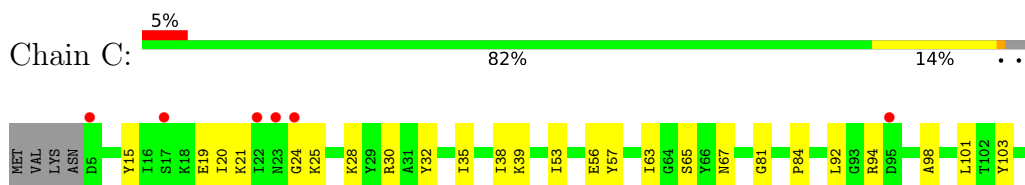
• Molecule 1: Tannase

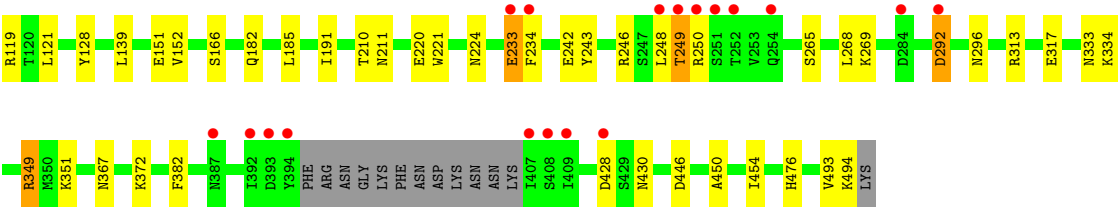


• Molecule 1: Tannase



• Molecule 1: Tannase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	71.84Å 91.07Å 124.07Å 90.00° 101.70° 90.00°	Depositor
Resolution (Å)	47.79 – 2.40 47.79 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.0 (47.79-2.40) 99.0 (47.79-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.85 (at 2.39Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.171 , 0.251 0.173 , 0.253	Depositor DCC
R_{free} test set	3104 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	33.6	Xtriage
Anisotropy	0.532	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 58.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11702	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.06% 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: SPD, GOL, MG

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	0.41	1/3831 (0.0%)	0.58	0/5179
1	B	0.35	0/3839	0.59	1/5190 (0.0%)
1	C	0.36	0/3827	0.56	0/5174
All	All	0.37	1/11497 (0.0%)	0.58	1/15543 (0.0%)

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	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	250	ARG	NE-CZ	5.33	1.39	1.33

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	95	ASP	N-CA-C	-5.49	99.10	110.80

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	393	ASP	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	A	3752	0	3696	59	0
1	B	3763	0	3702	56	0
1	C	3749	0	3698	47	0
2	A	6	0	8	1	0
2	B	6	0	8	1	0
2	C	6	0	8	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
4	B	10	0	19	3	0
5	A	180	0	0	4	0
5	B	107	0	0	5	0
5	C	120	0	0	1	0
All	All	11702	0	11139	160	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 7.

All (160) 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:B:125:LYS:NZ	5:B:601:HOH:O	1.92	1.03
1:A:125:LYS:NZ	5:A:601:HOH:O	1.94	0.97
1:C:246:ARG:NH1	1:C:446:ASP:OD2	2.10	0.84
1:C:243:TYR:O	1:C:246:ARG:NH2	2.14	0.80
1:A:409:ILE:HD12	1:A:410:PRO:HD2	1.62	0.79
1:A:221:TRP:HA	1:A:269:LYS:HG3	1.67	0.76
1:C:210:THR:HG22	1:C:211:ASN:H	1.52	0.75
1:A:38:ILE:HD11	1:A:193:ALA:N	2.04	0.73
1:B:364:GLY:H	4:B:502:SPD:H21	1.54	0.72
1:A:38:ILE:HG13	1:A:191:ILE:O	1.89	0.72
1:A:38:ILE:HD12	1:A:193:ALA:HA	1.73	0.70
1:B:289:LEU:HA	1:B:298:SER:HB2	1.74	0.68
1:B:484:GLU:N	1:B:484:GLU:OE2	2.27	0.68
1:C:221:TRP:HA	1:C:269:LYS:HD2	1.77	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:242:GLU:O	5:A:602:HOH:O	2.12	0.66
1:B:303:LEU:HA	1:B:306:ILE:HD12	1.78	0.66
1:A:38:ILE:CD1	1:A:193:ALA:HA	2.26	0.65
1:A:241:GLN:H	1:A:241:GLN:CD	2.05	0.63
1:A:249:THR:O	1:A:250:ARG:HD3	1.99	0.62
1:B:20:ILE:HG12	1:B:29:TYR:HE2	1.64	0.62
1:C:233:GLU:H	1:C:233:GLU:CD	2.08	0.61
1:B:325:LYS:HE3	1:B:330:ILE:HD12	1.84	0.59
1:C:103:TYR:O	1:C:106:SER:OG	2.14	0.59
1:A:94:ARG:HD3	1:A:245:ASP:OD1	2.02	0.59
1:C:128:TYR:HB2	1:C:372:LYS:HA	1.85	0.59
1:A:493:VAL:O	1:A:494:LYS:HB2	2.03	0.58
1:A:9:LYS:HD3	1:A:149:ASN:OD1	2.04	0.57
1:A:220:GLU:OE1	1:A:349[B]:ARG:NH2	2.27	0.57
1:C:15:TYR:CD1	1:C:30:ARG:HG2	2.40	0.57
1:A:284:ASP:HB3	1:B:121:LEU:HD12	1.86	0.56
1:A:249:THR:C	1:A:250:ARG:HD3	2.31	0.56
1:C:242:GLU:OE1	1:C:249:THR:HG21	2.05	0.55
1:A:38:ILE:CD1	1:A:193:ALA:CA	2.86	0.53
1:B:32:TYR:HB3	1:B:35:ILE:HD11	1.89	0.53
1:A:67:ASN:OD1	1:A:70:ASN:HB2	2.08	0.53
1:C:15:TYR:CE1	1:C:30:ARG:HG2	2.43	0.53
1:C:19:GLU:HG2	1:C:28:LYS:HG2	1.89	0.53
1:A:415:LYS:N	1:A:415:LYS:HD3	2.23	0.53
1:B:11:ASN:CG	1:B:14:LYS:HD3	2.35	0.52
1:B:14:LYS:HD2	1:B:14:LYS:N	2.25	0.52
1:A:428:ASP:HB3	1:A:430:ASN:H	1.75	0.52
1:B:301:THR:HA	1:B:304:SER:HB3	1.90	0.52
1:C:56:GLU:OE2	1:C:63:ILE:HD11	2.10	0.52
1:A:38:ILE:CG1	1:A:191:ILE:O	2.56	0.52
1:A:75:PHE:CZ	1:A:114:PRO:HG3	2.45	0.52
1:C:210:THR:HG22	1:C:211:ASN:N	2.22	0.52
1:C:38:ILE:HG21	1:C:191:ILE:HD12	1.91	0.52
1:B:296:ASN:OD1	1:B:300:LYS:NZ	2.39	0.51
1:A:136:ILE:HG12	1:A:174:LEU:HD13	1.93	0.51
1:B:408:SER:OG	5:B:602:HOH:O	2.15	0.51
1:A:332:ASN:N	5:A:609:HOH:O	2.38	0.50
1:C:292:ASP:HB2	1:C:296:ASN:O	2.11	0.50
1:B:321:ILE:HG23	1:B:328:PHE:HE2	1.76	0.50
1:A:280:LYS:NZ	5:A:611:HOH:O	2.44	0.50
1:C:15:TYR:HB3	1:C:32:TYR:CE2	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:364:GLY:N	4:B:502:SPD:H21	2.27	0.49
1:C:94:ARG:NH1	1:C:248:LEU:HG	2.28	0.49
1:B:161:ILE:HG21	1:B:489:ILE:HD13	1.94	0.49
1:B:322:SER:O	1:B:325:LYS:HG2	2.11	0.49
1:B:342:VAL:O	1:B:346:ILE:HG13	2.13	0.49
1:A:67:ASN:H	1:A:70:ASN:HB3	1.76	0.49
1:A:182:GLN:HA	1:A:185:LEU:HG	1.95	0.49
1:A:209:ILE:HG13	2:A:501:GOL:H12	1.95	0.48
1:C:32:TYR:HB3	1:C:35:ILE:HD11	1.96	0.48
1:C:65:SER:O	1:C:65:SER:OG	2.28	0.48
1:C:166:SER:OG	1:C:476:HIS:NE2	2.33	0.48
1:C:333:ASN:O	1:C:334:LYS:HD2	2.14	0.48
1:C:313:ARG:O	1:C:317:GLU:HG3	2.14	0.48
1:C:220:GLU:HG3	1:C:265:SER:HA	1.96	0.47
1:C:428:ASP:HB3	1:C:430:ASN:H	1.78	0.47
1:A:440:HIS:O	1:A:470:ALA:HA	2.14	0.47
1:C:81:GLY:HA3	1:C:243:TYR:CD2	2.49	0.47
1:B:440:HIS:NE2	1:B:444:ASP:HB3	2.30	0.47
1:C:39:LYS:NZ	5:C:610:HOH:O	2.48	0.47
1:B:242:GLU:CD	1:B:250:ARG:HD3	2.40	0.47
1:A:72:PRO:HB2	1:A:109:TYR:CD2	2.49	0.47
1:C:20:ILE:HD12	1:C:21:LYS:H	1.79	0.47
1:A:56:GLU:CD	1:A:63:ILE:HD11	2.39	0.47
1:B:72:PRO:HG3	1:B:493:VAL:HG11	1.97	0.47
1:A:38:ILE:CD1	1:A:193:ALA:N	2.76	0.46
1:B:56:GLU:OE2	1:B:56:GLU:N	2.43	0.46
1:C:53:ILE:HG12	1:C:110:VAL:HG13	1.97	0.46
1:B:81:GLY:HA3	1:B:243:TYR:CE2	2.50	0.46
1:C:57:TYR:CE1	1:C:67:ASN:HA	2.50	0.46
1:B:266:ASN:O	1:B:270:THR:HG23	2.16	0.46
1:A:308:ARG:HG3	1:A:335:VAL:HG12	1.98	0.46
1:B:387:ASN:OD1	1:B:410:PRO:HB3	2.16	0.46
1:B:75:PHE:CZ	1:B:114:PRO:HG3	2.51	0.46
1:B:98:ALA:HB1	1:B:102:THR:HB	1.98	0.46
1:C:151:GLU:HB3	1:C:152:VAL:HG13	1.98	0.46
1:C:450:ALA:O	1:C:454:ILE:HG12	2.16	0.46
1:A:415:LYS:HD3	1:A:415:LYS:H	1.79	0.46
1:B:371:ASP:HB2	5:B:673:HOH:O	2.15	0.46
1:C:92:LEU:HD23	1:C:98:ALA:HA	1.97	0.46
1:A:435:TYR:CD2	1:A:492:VAL:HG13	2.51	0.45
1:B:228:SER:OG	1:B:348:ASP:OD1	2.34	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:19:GLU:HG2	1:B:26:LYS:HZ3	1.81	0.45
1:B:312:ASN:HA	1:B:315:LEU:HD12	1.98	0.45
1:B:246:ARG:HA	1:B:246:ARG:HD3	1.66	0.45
1:B:276:LEU:HD21	1:B:299:PHE:CG	2.51	0.45
1:C:493:VAL:O	1:C:494:LYS:HB2	2.17	0.45
1:A:323:GLN:H	1:A:323:GLN:HG3	1.54	0.45
1:A:246:ARG:HA	1:A:250:ARG:HH22	1.81	0.44
1:B:352:SER:H	4:B:502:SPD:H91	1.81	0.44
1:A:285:GLY:HA2	1:B:84:PRO:HG2	1.99	0.44
1:B:209:ILE:HG13	2:B:501:GOL:H2	1.99	0.44
1:B:314:ALA:CB	1:B:321:ILE:HD11	2.47	0.44
1:B:166:SER:HA	1:B:207:CYS:O	2.18	0.44
1:C:84:PRO:N	1:C:118:GLY:HA3	2.33	0.44
1:A:84:PRO:N	1:A:118:GLY:HA3	2.32	0.44
1:A:462:SER:O	1:A:464:LYS:HE2	2.17	0.44
1:B:360:ASP:C	1:B:360:ASP:OD2	2.61	0.44
1:B:283:ASP:OD1	1:B:287:ASN:HB2	2.19	0.43
1:B:325:LYS:HD2	1:B:325:LYS:HA	1.80	0.43
1:A:95:ASP:OD1	1:A:95:ASP:N	2.50	0.43
1:B:162:SER:HB2	1:B:200:ILE:HD12	2.00	0.43
1:B:348:ASP:O	1:B:445:LYS:NZ	2.50	0.43
1:A:123:ASP:OD1	1:A:127:ASN:N	2.51	0.43
1:A:30:ARG:HD2	1:A:58:PHE:CE2	2.53	0.43
1:B:332:ASN:O	1:B:334:LYS:N	2.52	0.43
1:C:248:LEU:O	1:C:249:THR:OG1	2.27	0.43
1:B:81:GLY:HA3	1:B:243:TYR:CD2	2.53	0.43
1:C:182:GLN:HA	1:C:185:LEU:HD12	2.00	0.43
1:B:325:LYS:HE3	1:B:330:ILE:CD1	2.49	0.43
1:C:139:LEU:HD23	1:C:139:LEU:HA	1.81	0.43
1:A:250:ARG:HE	1:A:250:ARG:HB2	1.63	0.42
1:A:107:LYS:HA	1:A:107:LYS:HD3	1.88	0.42
1:A:170:ALA:CB	1:A:210:THR:HG23	2.49	0.42
1:A:6:TYR:O	1:A:39:LYS:HG2	2.19	0.42
1:A:173:ALA:HA	1:A:203:VAL:HG11	2.01	0.42
1:A:322:SER:HA	1:A:325:LYS:HD2	2.02	0.42
1:B:438:ILE:O	1:B:468:PHE:HA	2.20	0.42
1:C:121:LEU:HD12	1:C:121:LEU:HA	1.85	0.42
1:A:222:MET:HE3	1:A:222:MET:HB2	1.91	0.42
1:A:435:TYR:CG	1:A:492:VAL:HG13	2.54	0.42
1:B:103:TYR:O	1:B:106:SER:HB3	2.20	0.42
1:B:394:TYR:HA	5:B:610:HOH:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:234:PHE:HD1	1:C:234:PHE:HA	1.67	0.42
1:A:267:LYS:O	1:A:271:LEU:HD13	2.20	0.41
1:C:101:LEU:H	1:C:101:LEU:HD12	1.84	0.41
1:C:249:THR:HA	1:C:250:ARG:NE	2.34	0.41
1:C:334:LYS:HD2	1:C:334:LYS:HA	1.77	0.41
1:C:349[B]:ARG:NE	1:C:351:LYS:O	2.45	0.41
1:A:494:LYS:HD3	1:A:494:LYS:HA	1.72	0.41
1:C:119:ARG:HD2	1:C:367:ASN:HB3	2.03	0.41
1:C:220:GLU:O	1:C:224:ASN:HB2	2.20	0.41
1:A:170:ALA:HB1	1:A:210:THR:HG23	2.03	0.41
1:A:320:ASP:OD2	1:A:322:SER:HB3	2.21	0.41
1:C:25:LYS:HD2	1:C:25:LYS:HA	1.79	0.41
1:A:174:LEU:HD11	1:A:421:MET:HA	2.03	0.41
1:B:260:ASP:OD2	1:B:260:ASP:N	2.54	0.41
1:B:493:VAL:O	1:B:494:LYS:HD2	2.20	0.41
1:B:448:SER:O	1:B:452:PRO:HD2	2.21	0.41
1:A:10:PHE:CD1	1:A:153:MET:HE2	2.56	0.40
1:A:38:ILE:HD11	1:A:193:ALA:CA	2.51	0.40
1:B:361:ALA:HA	1:B:366:ASN:CG	2.45	0.40
1:B:396:ARG:HG2	5:B:610:HOH:O	2.20	0.40
1:A:211:ASN:HB3	1:A:451:ILE:HD12	2.02	0.40
1:C:268:LEU:HD21	1:C:382:PHE:CD1	2.56	0.40
1:B:144:ARG:HG3	1:B:198:ASP:HB3	2.04	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	475/495 (96%)	455 (96%)	19 (4%)	1 (0%)	43 58
1	B	476/495 (96%)	450 (94%)	23 (5%)	3 (1%)	21 32

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	475/495 (96%)	447 (94%)	24 (5%)	4 (1%)	16	25
All	All	1426/1485 (96%)	1352 (95%)	66 (5%)	8 (1%)	21	32

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	395	PHE
1	C	233	GLU
1	B	95	ASP
1	B	249	THR
1	C	249	THR
1	C	292	ASP
1	A	24	GLY
1	C	24	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	A	403/418 (96%)	401 (100%)	2 (0%)	81	91
1	B	404/418 (97%)	404 (100%)	0	100	100
1	C	403/418 (96%)	401 (100%)	2 (0%)	81	91
All	All	1210/1254 (96%)	1206 (100%)	4 (0%)	87	93

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	349[A]	ARG
1	A	349[B]	ARG
1	C	349[A]	ARG
1	C	349[B]	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	214	ASN
1	B	40	ASN
1	B	59	ASN
1	B	60	ASN
1	B	224	ASN
1	B	254	GLN
1	B	333	ASN
1	B	430	ASN
1	C	40	ASN
1	C	127	ASN
1	C	256	ASN
1	C	287	ASN
1	C	430	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 3 are monoatomic - leaving 4 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	$\# Z > 2$
2	GOL	A	501	-	5,5,5	0.98	0	5,5,5	1.04	0
4	SPD	B	502	-	9,9,9	0.34	0	8,8,8	0.50	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GOL	C	501	-	5,5,5	0.74	0	5,5,5	1.22	0
2	GOL	B	501	-	5,5,5	0.96	0	5,5,5	1.08	1 (20%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	A	501	-	-	2/4/4/4	-
4	SPD	B	502	-	-	4/7/7/7	-
2	GOL	C	501	-	-	1/4/4/4	-
2	GOL	B	501	-	-	2/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	GOL	C3-C2-C1	-2.08	104.15	111.80

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	GOL	C1-C2-C3-O3
2	B	501	GOL	O1-C1-C2-C3
2	B	501	GOL	O1-C1-C2-O2
4	B	502	SPD	C8-C7-N6-C5
4	B	502	SPD	C4-C5-N6-C7
4	B	502	SPD	N6-C7-C8-C9
2	A	501	GOL	O2-C2-C3-O3
4	B	502	SPD	N1-C2-C3-C4
2	C	501	GOL	O2-C2-C3-O3

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	GOL	1	0
4	B	502	SPD	3	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	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	478/495 (96%)	-0.03	15 (3%)	51	47	16, 37, 68, 119	1 (0%)
1	B	480/495 (96%)	0.27	30 (6%)	26	22	22, 45, 85, 131	0
1	C	478/495 (96%)	0.20	24 (5%)	34	30	21, 42, 81, 118	1 (0%)
All	All	1436/1485 (96%)	0.15	69 (4%)	35	32	16, 41, 80, 131	2 (0%)

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	248	LEU	5.5
1	B	249	THR	4.2
1	C	22	ILE	4.2
1	B	494	LYS	3.9
1	A	250	ARG	3.8
1	C	407	ILE	3.8
1	A	22	ILE	3.7
1	B	395	PHE	3.7
1	C	394	TYR	3.6
1	B	94	ARG	3.4
1	C	249	THR	3.4
1	B	22	ILE	3.3
1	B	329	THR	3.2
1	B	394	TYR	3.2
1	C	95	ASP	3.2
1	A	21	LYS	3.0
1	A	408	SER	3.0
1	C	428	ASP	2.9
1	B	248	LEU	2.9
1	C	409	ILE	2.8
1	B	250	ARG	2.8
1	B	393	ASP	2.7
1	B	24	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	233	GLU	2.7
1	A	234	PHE	2.7
1	B	27	ILE	2.7
1	A	23	ASN	2.6
1	C	392	ILE	2.6
1	B	92	LEU	2.6
1	C	5	ASP	2.6
1	B	233	GLU	2.6
1	B	23	ASN	2.6
1	B	253	VAL	2.6
1	A	232	MET	2.5
1	A	249	THR	2.5
1	B	408	SER	2.4
1	A	394	TYR	2.4
1	C	250	ARG	2.4
1	B	26	LYS	2.4
1	C	252	THR	2.4
1	A	395	PHE	2.4
1	A	409	ILE	2.4
1	C	393	ASP	2.4
1	B	21	LYS	2.3
1	B	330	ILE	2.3
1	C	23	ASN	2.3
1	B	93	GLY	2.3
1	B	96	GLY	2.3
1	A	38	ILE	2.3
1	C	251	SER	2.3
1	A	247	SER	2.3
1	C	254	GLN	2.3
1	A	248	LEU	2.2
1	B	25	LYS	2.2
1	B	389	LYS	2.2
1	C	234	PHE	2.2
1	C	408	SER	2.2
1	B	20	ILE	2.2
1	B	251	SER	2.2
1	C	17	SER	2.1
1	C	292	ASP	2.1
1	A	428	ASP	2.1
1	B	304	SER	2.1
1	C	387	ASN	2.1
1	C	24	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	332	ASN	2.0
1	B	333	ASN	2.0
1	B	6	TYR	2.0
1	C	284	ASP	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 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	SPD	B	502	10/10	0.85	0.18	40,61,64,65	0
2	GOL	A	501	6/6	0.92	0.12	33,36,36,39	0
3	MG	C	502	1/1	0.94	0.07	48,48,48,48	0
2	GOL	B	501	6/6	0.94	0.10	32,41,41,45	0
2	GOL	C	501	6/6	0.95	0.09	20,29,35,39	0
3	MG	A	502	1/1	0.97	0.06	41,41,41,41	0
3	MG	B	503	1/1	0.99	0.03	32,32,32,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.