



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 01:45 PM UTC

PDB ID : 3AHY / pdb_00003ahy
Title : Crystal structure of beta-glucosidase 2 from fungus *Trichoderma reesei* in complex with Tris
Authors : Jeng, W.-Y.; Liu, C.-I.; Wang, A.H.-J.
Deposited on : 2010-05-06
Resolution : 1.63 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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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

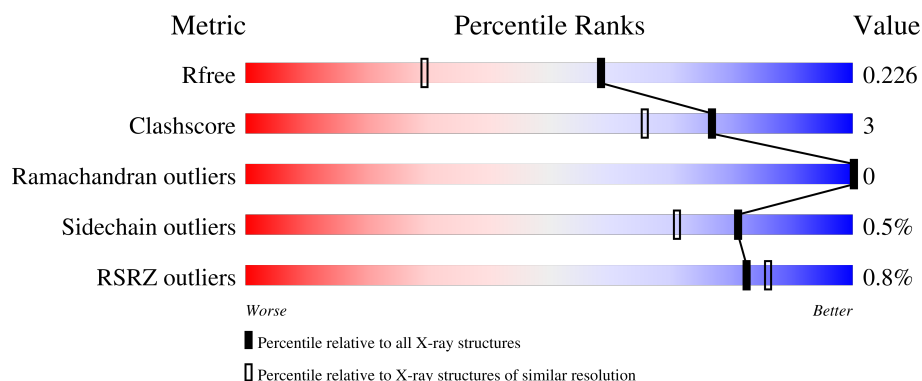
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

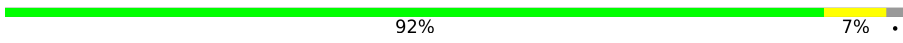
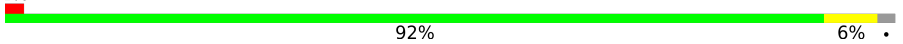
The reported resolution of this entry is 1.63 Å.

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	1141 (1.64-1.64)
Clashscore	190562	1171 (1.64-1.64)
Ramachandran outliers	187476	1151 (1.64-1.64)
Sidechain outliers	187428	1150 (1.64-1.64)
RSRZ outliers	180081	1141 (1.64-1.64)

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	473	 92% 7% .
1	B	473	 89% 9% .
1	C	473	 2% 92% 6% .
1	D	473	 0% 89% 9% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17923 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 Beta-glucosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	465	Total	C	N	O	S	0	0	0
			3694	2365	627	691	11			
1	B	466	Total	C	N	O	S	0	0	0
			3704	2371	630	692	11			
1	C	465	Total	C	N	O	S	0	0	0
			3699	2368	629	691	11			
1	D	465	Total	C	N	O	S	0	0	0
			3694	2365	627	691	11			

There are 28 discrepancies between the modelled and reference sequences:

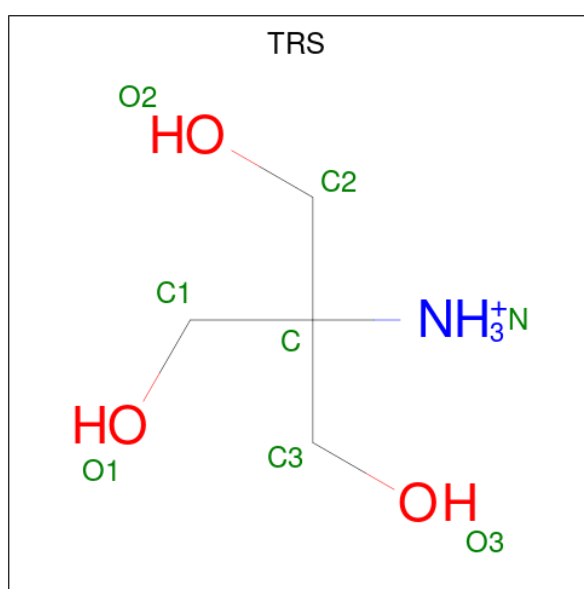
Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	MET	-	expression tag	UNP O93785
A	-5	HIS	-	expression tag	UNP O93785
A	-4	HIS	-	expression tag	UNP O93785
A	-3	HIS	-	expression tag	UNP O93785
A	-2	HIS	-	expression tag	UNP O93785
A	-1	HIS	-	expression tag	UNP O93785
A	0	HIS	-	expression tag	UNP O93785
B	-6	MET	-	expression tag	UNP O93785
B	-5	HIS	-	expression tag	UNP O93785
B	-4	HIS	-	expression tag	UNP O93785
B	-3	HIS	-	expression tag	UNP O93785
B	-2	HIS	-	expression tag	UNP O93785
B	-1	HIS	-	expression tag	UNP O93785
B	0	HIS	-	expression tag	UNP O93785
C	-6	MET	-	expression tag	UNP O93785
C	-5	HIS	-	expression tag	UNP O93785
C	-4	HIS	-	expression tag	UNP O93785
C	-3	HIS	-	expression tag	UNP O93785
C	-2	HIS	-	expression tag	UNP O93785
C	-1	HIS	-	expression tag	UNP O93785
C	0	HIS	-	expression tag	UNP O93785

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-6	MET	-	expression tag	UNP O93785
D	-5	HIS	-	expression tag	UNP O93785
D	-4	HIS	-	expression tag	UNP O93785
D	-3	HIS	-	expression tag	UNP O93785
D	-2	HIS	-	expression tag	UNP O93785
D	-1	HIS	-	expression tag	UNP O93785
D	0	HIS	-	expression tag	UNP O93785

- Molecule 2 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: C₄H₁₂NO₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			8	4	1	3		
2	B	1	Total	C	N	O	0	0
			8	4	1	3		
2	C	1	Total	C	N	O	0	0
			8	4	1	3		
2	D	1	Total	C	N	O	0	0
			8	4	1	3		

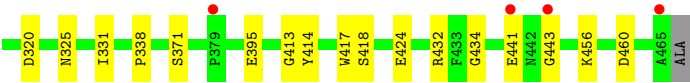
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	752	Total	O	0	0
			752	752		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	804	Total 804	O 804	0	0
3	C	807	Total 807	O 807	0	0
3	D	737	Total 737	O 737	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	92.83Å 103.51Å 94.80Å 90.00° 105.37° 90.00°	Depositor
Resolution (Å)	29.90 – 1.63 29.90 – 1.63	Depositor EDS
% Data completeness (in resolution range)	94.5 (29.90-1.63) 94.4 (29.90-1.63)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.37 (at 1.62Å)	Xtriage
Refinement program	CNS, REFMAC 5.5.0072	Depositor
R, R_{free}	0.164 , 0.218 0.176 , 0.226	Depositor DCC
R_{free} test set	10271 reflections (4.68%)	wwPDB-VP
Wilson B-factor (Å ²)	14.0	Xtriage
Anisotropy	0.120	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 39.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.037 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	17923	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.77% 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: TRS

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	0.37	0/3805	0.92	2/5173 (0.0%)
1	B	0.38	0/3816	0.91	7/5188 (0.1%)
1	C	0.39	0/3811	0.91	5/5181 (0.1%)
1	D	0.38	0/3805	0.90	6/5173 (0.1%)
All	All	0.38	0/15237	0.91	20/20715 (0.1%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	115	ILE	N-CA-C	6.58	117.34	107.80
1	A	18	GLU	N-CA-C	5.84	117.72	111.36
1	C	115	ILE	N-CA-C	5.80	116.21	107.80
1	C	42	ALA	N-CA-C	5.66	118.18	111.33
1	C	444	GLN	N-CA-C	-5.55	104.05	111.87
1	B	434	GLY	N-CA-C	5.54	119.29	112.14
1	B	18	GLU	N-CA-C	5.51	118.47	111.69
1	B	115	ILE	N-CA-C	5.37	115.69	108.17
1	B	405	GLU	N-CA-C	5.35	116.79	111.07
1	D	18	GLU	N-CA-C	5.24	117.07	111.36
1	B	378	LEU	CA-C-N	-5.18	114.38	119.92
1	B	378	LEU	C-N-CA	-5.18	114.38	119.92
1	D	2	LEU	CA-C-N	5.17	125.16	119.89
1	D	2	LEU	C-N-CA	5.17	125.16	119.89
1	D	115	ILE	N-CA-C	5.15	115.27	107.80
1	D	4	LYS	N-CA-C	5.12	117.60	111.71
1	D	434	GLY	N-CA-C	5.06	118.67	112.14
1	B	217	ASP	N-CA-C	5.06	118.99	112.41
1	C	18	GLU	N-CA-C	5.03	117.88	111.69
1	C	289	SER	N-CA-C	5.02	119.04	113.02

There are no chirality outliers.

There are no planarity outliers.

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	3694	0	3518	16	0
1	B	3704	0	3525	25	0
1	C	3699	0	3520	17	0
1	D	3694	0	3518	26	0
2	A	8	0	12	1	0
2	B	8	0	12	1	0
2	C	8	0	12	1	0
2	D	8	0	12	1	0
3	A	752	0	0	4	1
3	B	804	0	0	8	1
3	C	807	0	0	4	0
3	D	737	0	0	4	1
All	All	17923	0	14129	84	2

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 3.

All (84) 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:C:440:TYR:HB2	3:C:1293:HOH:O	1.52	1.06
1:D:217:ASP:OD2	3:D:1206:HOH:O	1.87	0.89
1:B:394:ASN:HD21	1:B:398:ARG:NH1	1.80	0.79
1:B:394:ASN:HD21	1:B:398:ARG:HH12	1.33	0.76
1:D:65:LYS:HE3	1:D:109:ALA:O	1.89	0.72
1:A:320:ASP:OD2	3:A:1234:HOH:O	2.08	0.72
1:C:382:LYS:HA	1:C:382:LYS:HE2	1.72	0.70
1:B:92:ASN:ND2	3:B:1299:HOH:O	2.27	0.68
1:C:298:TYR:CE2	2:C:500:TRS:H22	2.30	0.67
1:B:53:SER:OG	1:B:420:MET:HE1	1.96	0.66
1:D:320:ASP:OD1	3:D:1230:HOH:O	2.13	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:65:LYS:CE	1:D:109:ALA:O	2.44	0.65
1:D:212:LYS:HE3	1:D:218:GLY:O	1.99	0.63
1:A:56:ARG:NH1	3:A:1226:HOH:O	2.30	0.63
1:B:1:MET:HG2	3:B:1314:HOH:O	2.00	0.62
1:C:244:GLU:HG2	3:C:717:HOH:O	2.00	0.61
1:B:394:ASN:ND2	1:B:398:ARG:NH1	2.49	0.60
1:C:248:GLU:HA	1:C:252:ALA:HB3	1.85	0.58
1:D:441:GLU:C	1:D:443:GLY:H	2.11	0.58
1:D:456:LYS:HE2	1:D:460:ASP:OD2	2.05	0.56
1:B:237:PRO:HA	1:B:240:LYS:HE3	1.88	0.56
1:A:166:PRO:HB3	1:A:196:ILE:HG21	1.88	0.55
1:D:155:PRO:O	1:D:158:ARG:NH2	2.35	0.55
1:B:246:ARG:NE	3:B:1275:HOH:O	2.39	0.55
1:B:166:PRO:HB3	1:B:196:ILE:HG21	1.89	0.55
1:D:248:GLU:HA	1:D:252:ALA:HB3	1.88	0.54
1:D:7:GLN:O	1:D:413:GLY:HA2	2.08	0.54
1:A:4:LYS:HG2	3:A:1193:HOH:O	2.09	0.53
1:B:10:PHE:CE1	1:B:435:VAL:HG21	2.44	0.52
1:A:246:ARG:NH2	3:A:1161:HOH:O	2.41	0.52
1:B:246:ARG:CG	3:B:1275:HOH:O	2.57	0.52
1:C:380:LYS:HD2	1:C:380:LYS:C	2.34	0.51
1:D:338:PRO:HB2	3:D:1086:HOH:O	2.11	0.50
1:A:248:GLU:HA	1:A:252:ALA:HB3	1.93	0.49
1:B:1:MET:HE3	3:B:1314:HOH:O	2.13	0.49
1:C:13:ALA:HB2	1:C:119:HIS:CE1	2.47	0.49
1:D:279:THR:OG1	1:D:281:GLU:HG2	2.13	0.49
1:B:310:ALA:HB2	1:B:318:ASN:HB2	1.95	0.49
1:B:295:MET:HE2	1:B:297:HIS:HB2	1.96	0.48
1:D:281:GLU:H	1:D:281:GLU:CD	2.22	0.48
1:A:7:GLN:O	1:A:413:GLY:HA2	2.14	0.48
1:C:37:GLN:HB2	1:C:40:LYS:HG3	1.96	0.48
1:B:212:LYS:HB3	3:B:1266:HOH:O	2.13	0.47
1:C:216:GLY:HA2	3:C:1246:HOH:O	2.14	0.47
1:B:167:LEU:C	1:B:167:LEU:HD23	2.39	0.47
1:B:458:LEU:HD12	3:B:1285:HOH:O	2.14	0.47
1:B:7:GLN:O	1:B:413:GLY:HA2	2.15	0.47
1:A:171:ILE:HB	1:A:172:PRO:HD3	1.95	0.47
1:A:310:ALA:HB2	1:A:318:ASN:HB2	1.98	0.46
1:C:232:TRP:CD1	1:C:325:ASN:HA	2.50	0.46
1:C:171:ILE:HB	1:C:172:PRO:HD3	1.98	0.45
1:A:457:PRO:O	1:A:461:GLU:HG3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:298:TYR:CE2	2:A:500:TRS:H22	2.52	0.44
1:C:56:ARG:NH2	1:C:444:GLN:HG3	2.32	0.44
1:C:20:ALA:HB1	1:C:23:GLN:HB3	1.99	0.44
1:C:385:GLU:HG3	3:C:1304:HOH:O	2.15	0.44
1:C:394:ASN:O	1:C:398:ARG:HG3	2.17	0.44
1:A:295:MET:HE2	1:A:297:HIS:HB2	2.00	0.43
1:B:8:TRP:HA	1:B:414:TYR:O	2.18	0.43
1:A:417:TRP:HA	1:A:418:SER:HA	1.81	0.43
1:D:232:TRP:CD1	1:D:325:ASN:HA	2.54	0.43
1:D:298:TYR:CE2	2:D:500:TRS:H22	2.53	0.43
1:D:166:PRO:HB3	1:D:196:ILE:HG21	2.00	0.43
1:D:371:SER:HA	1:D:432:ARG:O	2.18	0.43
1:A:48:VAL:HG13	1:A:49:THR:N	2.34	0.43
1:B:56:ARG:NH1	3:B:1087:HOH:O	2.45	0.43
1:C:166:PRO:HB3	1:C:196:ILE:HG21	2.01	0.43
1:A:232:TRP:CD1	1:A:325:ASN:HA	2.53	0.42
1:B:171:ILE:HB	1:B:172:PRO:HD3	2.01	0.42
1:D:417:TRP:HA	1:D:418:SER:HA	1.78	0.42
1:B:298:TYR:CE2	2:B:500:TRS:H22	2.54	0.42
1:D:395:GLU:HG2	3:D:1084:HOH:O	2.18	0.42
1:A:165:GLU:CD	1:A:296:ASN:HD22	2.28	0.42
1:D:115:ILE:HD12	1:D:157:VAL:HG21	2.01	0.42
1:B:248:GLU:HA	1:B:252:ALA:HB3	2.02	0.42
1:B:232:TRP:CD1	1:B:325:ASN:HA	2.55	0.42
1:B:417:TRP:HA	1:B:418:SER:HA	1.79	0.42
1:D:302:TYR:CE2	1:D:331:ILE:HG12	2.55	0.41
1:D:424:GLU:HG3	1:D:424:GLU:O	2.20	0.41
1:D:441:GLU:C	1:D:443:GLY:N	2.75	0.41
1:D:268:ARG:HG2	1:D:275:LEU:HD23	2.02	0.41
1:C:417:TRP:HA	1:C:418:SER:HA	1.84	0.41
1:D:8:TRP:HA	1:D:414:TYR:O	2.21	0.40
1:D:167:LEU:HD23	1:D:167:LEU:C	2.47	0.40

All (2) 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)
3:A:1233:HOH:O	3:B:1070:HOH:O[1_556]	2.10	0.10
3:D:811:HOH:O	3:D:1201:HOH:O[2_445]	2.13	0.07

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	463/473 (98%)	449 (97%)	14 (3%)	0	100	100
1	B	464/473 (98%)	449 (97%)	15 (3%)	0	100	100
1	C	463/473 (98%)	445 (96%)	18 (4%)	0	100	100
1	D	463/473 (98%)	445 (96%)	18 (4%)	0	100	100
All	All	1853/1892 (98%)	1788 (96%)	65 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	382/389 (98%)	379 (99%)	3 (1%)	73	58
1	B	383/389 (98%)	380 (99%)	3 (1%)	73	58
1	C	383/389 (98%)	383 (100%)	0	100	100
1	D	382/389 (98%)	380 (100%)	2 (0%)	81	71
All	All	1530/1556 (98%)	1522 (100%)	8 (0%)	81	71

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	240	LYS
1	A	283	ARG

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Mol	Chain	Res	Type
1	A	382	LYS
1	B	48	VAL
1	B	269	LYS
1	B	380	LYS
1	D	37	GLN
1	D	281	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	GLN
1	A	23	GLN
1	A	55	ASN
1	A	159	ASN
1	A	219	GLN
1	A	329	ASN
1	B	7	GLN
1	B	37	GLN
1	B	92	ASN
1	B	128	GLN
1	B	145	ASN
1	B	184	GLN
1	B	270	GLN
1	B	296	ASN
1	B	327	GLN
1	B	394	ASN
1	C	7	GLN
1	C	296	ASN
1	C	394	ASN
1	C	444	GLN
1	D	7	GLN
1	D	270	GLN
1	D	327	GLN

5.3.3 RNA ⓘ

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](#)

4 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	$\# Z > 2$
2	TRS	C	500	-	7,7,7	0.31	0	9,9,9	0.62	0
2	TRS	B	500	-	7,7,7	0.31	0	9,9,9	0.59	0
2	TRS	A	500	-	7,7,7	0.29	0	9,9,9	0.37	0
2	TRS	D	500	-	7,7,7	0.26	0	9,9,9	0.34	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	TRS	C	500	-	-	2/9/9/9	-
2	TRS	B	500	-	-	0/9/9/9	-
2	TRS	A	500	-	-	0/9/9/9	-
2	TRS	D	500	-	-	0/9/9/9	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	500	TRS	C1-C-C3-O3
2	C	500	TRS	N-C-C3-O3

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	500	TRS	1	0
2	B	500	TRS	1	0
2	A	500	TRS	1	0
2	D	500	TRS	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	465/473 (98%)	-0.13	0 100 100	6, 12, 19, 27	0
1	B	466/473 (98%)	-0.19	0 100 100	5, 11, 18, 28	0
1	C	465/473 (98%)	-0.22	8 (1%) 69 74	5, 10, 20, 34	0
1	D	465/473 (98%)	-0.08	6 (1%) 75 80	6, 12, 22, 36	0
All	All	1861/1892 (98%)	-0.15	14 (0%) 82 86	5, 11, 20, 36	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	443	GLY	5.1
1	D	443	GLY	2.8
1	D	216	GLY	2.7
1	C	217	ASP	2.6
1	C	216	GLY	2.5
1	D	379	PRO	2.4
1	D	441	GLU	2.4
1	D	273	ASP	2.3
1	C	464	ALA	2.3
1	D	465	ALA	2.2
1	C	440	TYR	2.1
1	C	48	VAL	2.1
1	C	441	GLU	2.1
1	C	214	ALA	2.1

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
2	TRS	B	500	8/8	0.84	0.11	18,18,19,21	0
2	TRS	D	500	8/8	0.86	0.10	19,19,20,21	0
2	TRS	C	500	8/8	0.91	0.08	15,16,17,18	0
2	TRS	A	500	8/8	0.92	0.07	18,20,21,22	0

6.5 Other polymers [i](#)

There are no such residues in this entry.