



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 13, 2017 – 02:19 AM EDT

PDB ID : 5CG0
Title : Crystal structure of Spodoptera frugiperda Beta-glycosidase
Authors : Tamaki, F.K.; Souza, D.P.; Souza, V.P.; Farah, S.C.; Marana, S.R.
Deposited on : unknown
Resolution : 2.09 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 : rb-20029824
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) : rb-20029824

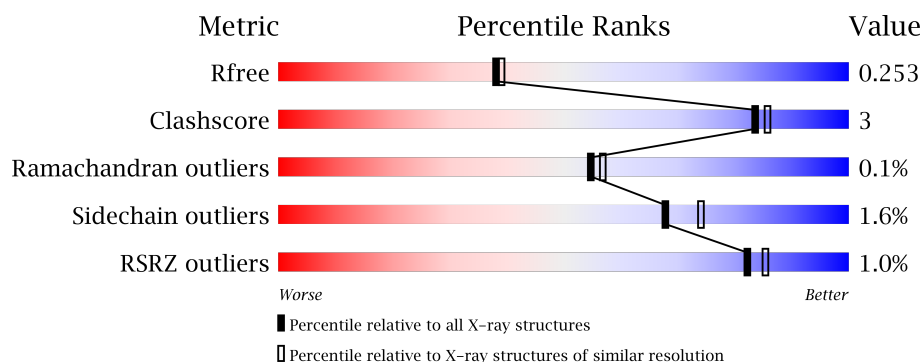
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.09 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	4243 (2.10-2.10)
Clashscore	112137	4788 (2.10-2.10)
Ramachandran outliers	110173	4740 (2.10-2.10)
Sidechain outliers	110143	4741 (2.10-2.10)
RSRZ outliers	101464	4275 (2.10-2.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	489	<div> <div>94%</div> <div>6%</div> </div>
1	B	489	<div> <div>92%</div> <div>8%</div> </div>
1	C	489	<div> <div>92%</div> <div>8%</div> </div>
1	D	489	<div> <div>92%</div> <div>7%</div> </div>
1	E	489	<div> <div>2%</div> <div>89%</div> <div>10%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	489	

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
2	NAG	A	602	X	-	-	-
2	NAG	B	602	X	-	-	-
2	NAG	D	602	X	-	-	-
3	TRS	A	603	-	-	-	X
3	TRS	B	603	-	-	-	X
3	TRS	E	602	-	-	-	X
3	TRS	F	602	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 25131 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.

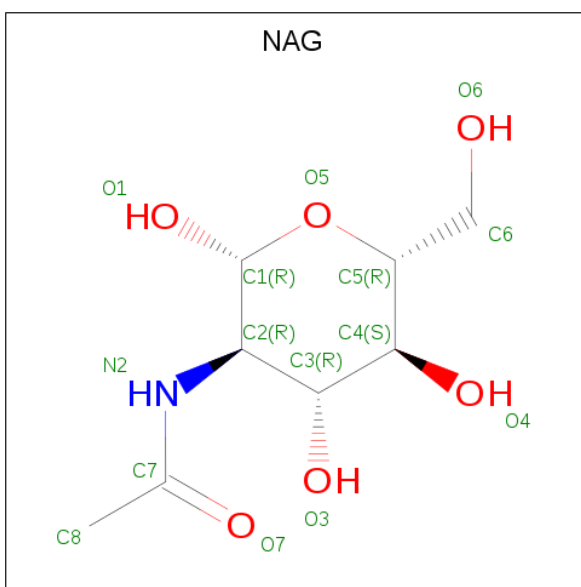
- Molecule 1 is a protein called Beta-glucosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	488	Total	C	N	O	S	0	3	0
			3978	2543	656	761	18			
1	B	488	Total	C	N	O	S	0	1	0
			3966	2537	656	755	18			
1	C	488	Total	C	N	O	S	0	3	0
			3967	2538	653	758	18			
1	D	486	Total	C	N	O	S	0	5	0
			3975	2542	658	757	18			
1	E	485	Total	C	N	O	S	0	0	0
			3934	2519	647	750	18			
1	F	471	Total	C	N	O	S	0	0	0
			3819	2448	633	722	16			

There are 12 discrepancies between the modelled and reference sequences:

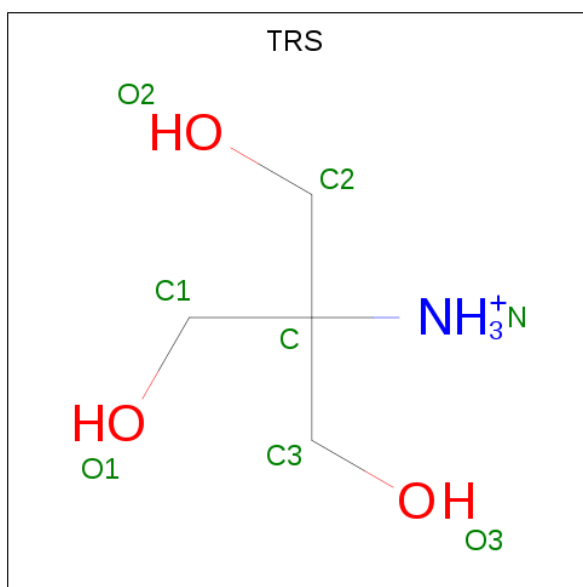
Chain	Residue	Modelled	Actual	Comment	Reference
A	325	PHE	LEU	engineered mutation	UNP O61594
A	326	PHE	ILE	engineered mutation	UNP O61594
B	325	PHE	LEU	engineered mutation	UNP O61594
B	326	PHE	ILE	engineered mutation	UNP O61594
C	325	PHE	LEU	engineered mutation	UNP O61594
C	326	PHE	ILE	engineered mutation	UNP O61594
D	325	PHE	LEU	engineered mutation	UNP O61594
D	326	PHE	ILE	engineered mutation	UNP O61594
E	325	PHE	LEU	engineered mutation	UNP O61594
E	326	PHE	ILE	engineered mutation	UNP O61594
F	325	PHE	LEU	engineered mutation	UNP O61594
F	326	PHE	ILE	engineered mutation	UNP O61594

- Molecule 2 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 3 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: C₄H₁₂NO₃).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	333	Total	O	0	0
			333	333		
4	B	269	Total	O	0	0
			269	269		
4	C	271	Total	O	0	0
			271	271		
4	D	168	Total	O	0	0
			168	168		
4	E	165	Total	O	0	0
			165	165		
4	F	112	Total	O	0	0
			112	112		

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.

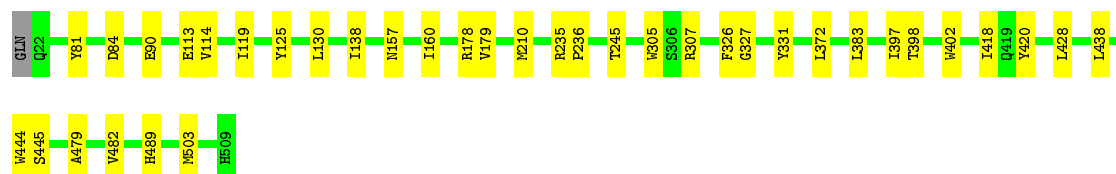
- Molecule 1: Beta-glucosidase

Chain A: 



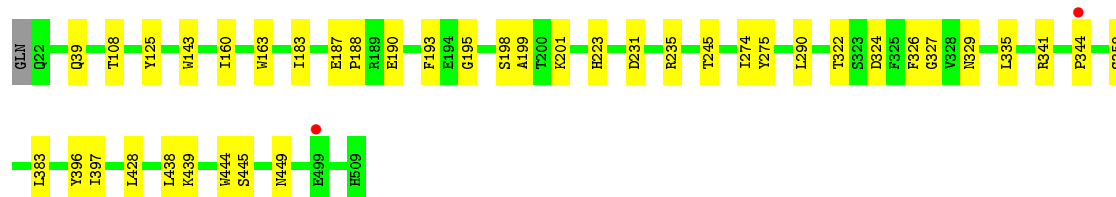
- Molecule 1: Beta-glucosidase

Chain B: 



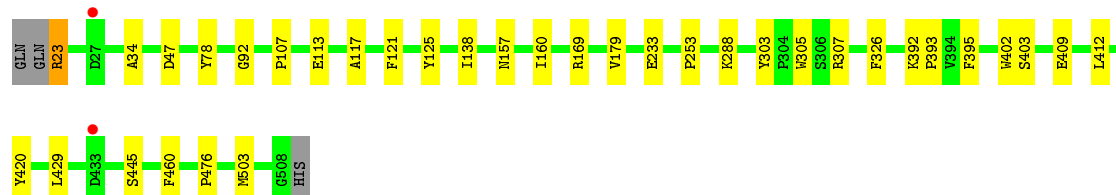
- Molecule 1: Beta-glucosidase

Chain C: 

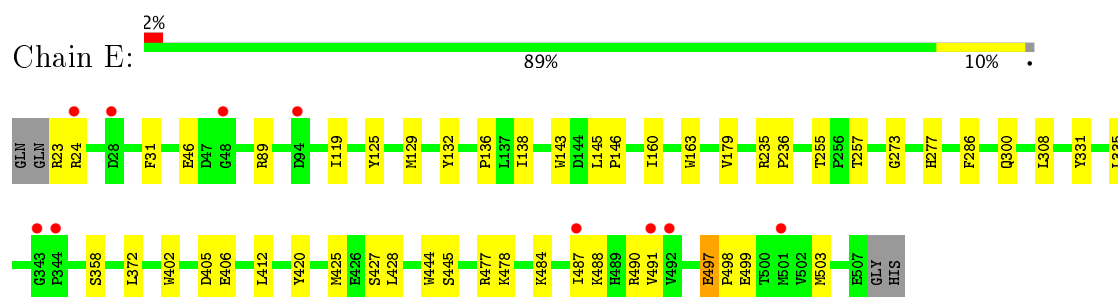


- Molecule 1: Beta-glucosidase

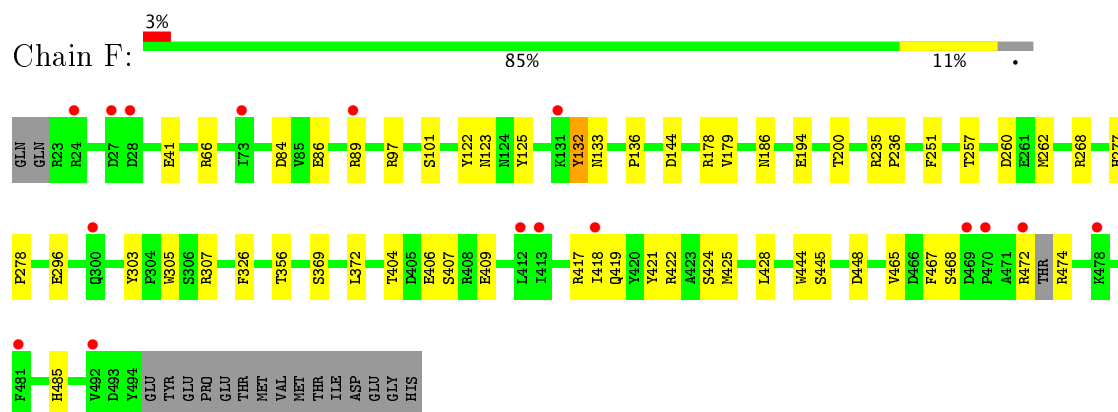
Chain D: 



- Molecule 1: Beta-glucosidase



• Molecule 1: Beta-glucosidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	53.83Å 64.99Å 257.19Å 93.16° 92.04° 112.18°	Depositor
Resolution (Å)	49.75 – 2.09 49.75 – 2.09	Depositor EDS
% Data completeness (in resolution range)	91.0 (49.75-2.09) 88.7 (49.75-2.09)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.28 (at 2.08Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.192 , 0.250 0.198 , 0.253	Depositor DCC
R_{free} test set	8675 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	35.4	Xtriage
Anisotropy	0.144	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 33.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.007 for -h,-k,h+k+l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	25131	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.67% 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 [i](#)

5.1 Standard geometry [i](#)

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.46	0/4099	0.65	0/5575
1	B	0.45	0/4087	0.63	0/5556
1	C	0.44	0/4094	0.63	0/5566
1	D	0.40	0/4098	0.58	0/5571
1	E	0.42	0/4051	0.59	0/5510
1	F	0.39	0/3933	0.58	0/5347
All	All	0.43	0/24362	0.61	0/33125

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	3978	0	3730	13	0
1	B	3966	0	3728	23	0
1	C	3967	0	3730	21	0
1	D	3975	0	3741	17	0
1	E	3934	0	3697	28	0
1	F	3819	0	3593	26	0
2	A	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	28	0	25	0	0
2	C	14	0	13	0	0
2	D	28	0	25	0	0
2	E	14	0	13	0	0
2	F	14	0	13	0	0
3	A	8	0	12	0	0
3	B	8	0	12	0	0
3	C	8	0	12	0	0
3	D	8	0	12	0	0
3	E	8	0	12	0	0
3	F	8	0	12	0	0
4	A	333	0	0	3	0
4	B	269	0	0	1	0
4	C	271	0	0	0	0
4	D	168	0	0	4	0
4	E	165	0	0	3	0
4	F	112	0	0	1	0
All	All	25131	0	22405	128	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 3.

All (128) 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:274:ILE:HD11	1:C:290:LEU:HD23	1.65	0.78
1:C:231:ASP:OD1	1:C:235:ARG:NH1	2.24	0.70
1:B:114:VAL:CG1	1:B:119:ILE:HD11	2.21	0.70
1:A:133[B]:ASN:ND2	4:A:701:HOH:O	2.29	0.64
1:A:245:THR:HG22	1:A:327:GLY:HA3	1.80	0.64
1:D:47:ASP:HB3	1:D:117:ALA:CB	2.31	0.61
1:F:136:PRO:HG2	1:F:179:VAL:HG22	1.85	0.58
1:D:403:SER:HB3	1:D:460:PHE:CZ	2.40	0.57
1:C:198:SER:OG	1:C:199:ALA:N	2.38	0.56
1:F:97:ARG:NH2	1:F:186:ASN:OD1	2.39	0.56
1:B:114:VAL:HG11	1:B:119:ILE:HD11	1.86	0.56
1:E:499:GLU:N	1:E:499:GLU:OE1	2.38	0.56
1:B:90:GLU:HB3	1:B:503:MET:CG	2.37	0.55
1:B:210:MET:HE2	1:B:307:ARG:HG2	1.87	0.55
1:E:145:LEU:HD12	1:E:146:PRO:HD2	1.89	0.55
1:D:303:TYR:CE1	1:D:307:ARG:HD2	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:160:ILE:HD12	1:C:163:TRP:HE3	1.73	0.54
1:B:418:ILE:HD12	1:B:482:VAL:HG21	1.90	0.54
1:F:86:GLU:O	1:F:89:ARG:HG2	2.08	0.54
1:D:169[B]:ARG:NH1	1:D:233:GLU:OE1	2.38	0.53
1:B:245:THR:HG22	1:B:327:GLY:HA3	1.91	0.53
1:E:412:LEU:HA	1:E:477:ARG:HG2	1.91	0.52
1:F:303:TYR:CE1	1:F:307:ARG:HD2	2.45	0.52
1:C:187:GLU:CD	1:C:329:ASN:HD22	2.13	0.52
1:D:92:GLY:HA2	4:D:820:HOH:O	2.08	0.52
1:B:489:HIS:HE1	4:B:841:HOH:O	1.93	0.51
1:B:428:LEU:HD22	1:B:438:LEU:HD21	1.92	0.51
1:F:136:PRO:CG	1:F:179:VAL:HG22	2.41	0.51
1:A:387:ASN:OD1	1:A:392:LYS:HA	2.11	0.51
1:C:183:ILE:HG21	1:C:245:THR:HG23	1.93	0.51
1:C:396:TYR:CE2	1:C:439:LYS:HB2	2.46	0.50
1:E:89:ARG:HD2	1:E:132:TYR:CD2	2.47	0.50
1:E:31:PHE:CB	1:E:487:ILE:HD12	2.41	0.50
1:F:448:ASP:HB2	1:F:465:VAL:HG22	1.94	0.50
1:D:393:PRO:HG2	1:D:395:PHE:CZ	2.47	0.50
1:C:235:ARG:NH2	1:C:324:ASP:OD2	2.45	0.50
1:E:129:MET:SD	1:E:136:PRO:HB3	2.52	0.49
1:C:335:LEU:O	1:C:358:SER:HA	2.12	0.49
1:B:138:ILE:HD12	1:B:179:VAL:HG21	1.95	0.49
1:F:262:MET:HE1	4:F:809:HOH:O	2.12	0.48
1:F:421:TYR:O	1:F:425:MET:HG2	2.14	0.48
1:D:402:TRP:CE3	1:D:420:TYR:CD2	3.02	0.48
1:F:404:THR:OG1	1:F:417:ARG:NH2	2.39	0.48
1:E:484:LYS:HD2	1:E:503:MET:HE1	1.96	0.48
1:F:194:GLU:HG2	1:F:200:THR:OG1	2.14	0.47
1:F:406:GLU:HG3	1:F:407:SER:N	2.29	0.47
1:B:245:THR:CG2	1:B:398:THR:HG21	2.45	0.47
1:E:138:ILE:HD12	1:E:179:VAL:HG21	1.97	0.47
1:A:447:MET:HE1	1:A:474:ARG:CZ	2.45	0.46
1:D:253:PRO:HD3	4:D:818:HOH:O	2.15	0.46
1:C:195:GLY:HA2	1:C:201:LYS:O	2.16	0.46
1:D:412:LEU:HD12	1:D:476:PRO:O	2.15	0.46
1:E:286:PHE:CE2	1:E:308:LEU:HD11	2.51	0.46
1:F:41:GLU:OE2	1:F:122:TYR:OH	2.27	0.46
1:D:409:GLU:N	4:D:701:HOH:O	2.44	0.46
1:D:78:TYR:HA	1:D:121:PHE:CE1	2.51	0.46
1:E:425:MET:HA	1:E:428:LEU:HD12	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:424:SER:O	1:F:428:LEU:HG	2.16	0.46
1:D:107:PRO:HD2	1:D:113:GLU:HG3	1.98	0.45
1:F:251:PHE:HB2	1:F:268:ARG:HG2	1.98	0.45
1:F:132:TYR:CD1	1:F:132:TYR:N	2.84	0.45
1:E:497:GLU:HG3	1:E:498:PRO:HD3	1.99	0.45
1:E:160:ILE:HA	1:E:163:TRP:CE3	2.52	0.45
1:B:235:ARG:N	1:B:236:PRO:CD	2.80	0.45
1:F:123:ASN:OD1	1:F:178:ARG:NH2	2.41	0.45
1:A:403:SER:HB3	1:A:460:PHE:CZ	2.52	0.45
1:E:23:ARG:HA	4:E:853:HOH:O	2.16	0.45
1:E:405:ASP:OD1	1:E:406:GLU:N	2.49	0.45
1:F:467:PHE:HA	1:F:472:ARG:HH11	1.81	0.45
1:E:402:TRP:CE3	1:E:420:TYR:CD1	3.05	0.44
1:E:273:GLY:O	1:E:277:HIS:HB2	2.18	0.44
1:D:34:ALA:O	1:D:445:SER:OG	2.26	0.43
1:F:194:GLU:CG	1:F:200:THR:OG1	2.66	0.43
1:B:114:VAL:HG12	1:B:119:ILE:HD11	1.99	0.43
1:D:503:MET:HA	4:D:789:HOH:O	2.18	0.43
1:B:418:ILE:CD1	1:B:479:ALA:HA	2.48	0.43
1:C:183:ILE:CG2	1:C:245:THR:HG23	2.47	0.43
1:E:412:LEU:HB2	4:E:813:HOH:O	2.18	0.43
1:D:138:ILE:CD1	1:D:179:VAL:HG21	2.49	0.43
1:F:444:TRP:HA	1:F:445:SER:HA	1.73	0.43
1:B:331:TYR:HB3	1:B:372:LEU:HD13	2.01	0.43
1:B:81:TYR:O	1:B:84:ASP:HB2	2.18	0.43
1:A:257:THR:HG22	4:A:845:HOH:O	2.19	0.43
1:B:245:THR:HG21	1:B:398:THR:HG21	2.00	0.43
1:A:162:ASP:OD1	1:A:225:LYS:NZ	2.52	0.43
1:F:235:ARG:HB3	1:F:236:PRO:HD3	2.01	0.43
1:A:300:GLN:HB3	4:A:977:HOH:O	2.19	0.42
1:A:75:ALA:HB1	1:A:447:MET:HE1	2.01	0.42
1:E:255:THR:OG1	1:E:257:THR:HG22	2.19	0.42
1:E:143:TRP:N	1:E:143:TRP:CD1	2.87	0.42
1:C:108:THR:O	1:C:108:THR:HG22	2.20	0.42
1:C:39:GLN:O	1:C:449:ASN:HB2	2.20	0.42
1:A:114:VAL:HG12	1:A:119:ILE:HD12	2.02	0.42
1:A:39:GLN:O	1:A:449:ASN:HB2	2.18	0.42
1:C:187:GLU:HB3	1:C:190:GLU:OE1	2.20	0.42
1:C:223:HIS:NE2	1:C:322:THR:O	2.48	0.42
1:A:187:GLU:HB3	1:A:190:GLU:OE1	2.20	0.42
1:F:132:TYR:N	1:F:132:TYR:HD1	2.17	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:157:ASN:O	1:B:160:ILE:HG22	2.20	0.42
1:B:383:LEU:HD11	1:B:397:ILE:HD11	2.01	0.42
1:F:257:THR:HG23	1:F:260:ASP:H	1.84	0.41
1:E:488:LYS:NZ	4:E:712:HOH:O	2.53	0.41
1:C:383:LEU:HD21	1:C:397:ILE:HD11	2.02	0.41
1:E:497:GLU:HG3	1:E:498:PRO:CD	2.50	0.41
1:B:130:LEU:HD11	1:B:178:ARG:HB3	2.02	0.41
1:F:418:ILE:O	1:F:422:ARG:HG3	2.20	0.41
1:C:193:PHE:O	1:C:198:SER:HB3	2.20	0.41
1:F:84:ASP:OD1	1:F:474:ARG:NH1	2.53	0.41
1:B:90:GLU:HB3	1:B:503:MET:HG2	2.02	0.41
1:D:23:ARG:HG3	1:D:429:LEU:HD13	2.03	0.41
1:E:31:PHE:CG	1:E:487:ILE:HD12	2.55	0.41
1:E:478:LYS:CG	1:E:498:PRO:HG3	2.50	0.41
1:F:277:HIS:N	1:F:278:PRO:CD	2.84	0.41
1:B:402:TRP:CE3	1:B:420:TYR:CD1	3.09	0.41
1:C:188:PRO:HB2	1:C:275:TYR:CG	2.56	0.41
1:E:235:ARG:N	1:E:236:PRO:CD	2.84	0.41
1:E:24:ARG:HA	1:E:490:ARG:O	2.21	0.41
1:E:331:TYR:HB3	1:E:372:LEU:HD13	2.03	0.41
1:E:335:LEU:O	1:E:358:SER:HA	2.21	0.41
1:D:157:ASN:O	1:D:160:ILE:HG22	2.21	0.40
1:B:418:ILE:HD11	1:B:479:ALA:HA	2.02	0.40
1:C:444:TRP:HA	1:C:445:SER:HA	1.85	0.40
1:A:331:TYR:HB3	1:A:372:LEU:HD13	2.03	0.40
1:C:245:THR:HG22	1:C:327:GLY:HA3	2.02	0.40
1:C:428:LEU:HD22	1:C:438:LEU:HD21	2.02	0.40
1:F:369:SER:HB2	1:F:372:LEU:HB3	2.03	0.40
1:B:444:TRP:HA	1:B:445:SER:HA	1.87	0.40
1:E:444:TRP:HA	1:E:445:SER:HA	1.78	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	489/489 (100%)	471 (96%)	17 (4%)	1 (0%)	51	52
1	B	487/489 (100%)	467 (96%)	20 (4%)	0	100	100
1	C	489/489 (100%)	469 (96%)	18 (4%)	2 (0%)	38	35
1	D	489/489 (100%)	468 (96%)	21 (4%)	0	100	100
1	E	483/489 (99%)	451 (93%)	32 (7%)	0	100	100
1	F	467/489 (96%)	428 (92%)	38 (8%)	1 (0%)	51	52
All	All	2904/2934 (99%)	2754 (95%)	146 (5%)	4 (0%)	55	57

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	143	TRP
1	C	143	TRP
1	F	468	SER
1	C	344	PRO

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	417/415 (100%)	411 (99%)	6 (1%)	71	78
1	B	415/415 (100%)	411 (99%)	4 (1%)	80	85
1	C	417/415 (100%)	414 (99%)	3 (1%)	87	91
1	D	417/415 (100%)	411 (99%)	6 (1%)	71	78
1	E	412/415 (99%)	405 (98%)	7 (2%)	66	72
1	F	398/415 (96%)	385 (97%)	13 (3%)	43	45
All	All	2476/2490 (99%)	2437 (98%)	39 (2%)	68	74

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	125	TYR
1	A	198	SER
1	A	305	TRP
1	A	326	PHE
1	A	358	SER
1	A	474	ARG
1	B	113	GLU
1	B	125	TYR
1	B	305	TRP
1	B	326	PHE
1	C	125	TYR
1	C	326	PHE
1	C	341	ARG
1	D	23	ARG
1	D	125	TYR
1	D	288	LYS
1	D	305	TRP
1	D	326	PHE
1	D	392	LYS
1	E	46	GLU
1	E	119	ILE
1	E	125	TYR
1	E	300	GLN
1	E	427	SER
1	E	491	VAL
1	E	497	GLU
1	F	66	ARG
1	F	101	SER
1	F	125	TYR
1	F	132	TYR
1	F	133	ASN
1	F	144	ASP
1	F	296	GLU
1	F	305	TRP
1	F	326	PHE
1	F	356	THR
1	F	409	GLU
1	F	419	GLN
1	F	485	HIS

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 ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

15 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	NAG	A	601	1,2	14,14,15	0.50	0	15,19,21	0.85	0
2	NAG	A	602	2	14,14,15	0.56	0	15,19,21	1.92	3 (20%)
3	TRS	A	603	-	7,7,7	0.35	0	9,9,9	0.49	0
2	NAG	B	601	1,2	14,14,15	0.55	0	15,19,21	1.32	2 (13%)
2	NAG	B	602	2	14,14,15	0.47	0	15,19,21	1.60	2 (13%)
3	TRS	B	603	-	7,7,7	0.39	0	9,9,9	0.27	0
2	NAG	C	601	1	14,14,15	0.51	0	15,19,21	1.09	2 (13%)
3	TRS	C	602	-	7,7,7	0.30	0	9,9,9	0.59	0
2	NAG	D	601	1,2	14,14,15	0.52	0	15,19,21	1.26	2 (13%)
2	NAG	D	602	2	14,14,15	0.59	0	15,19,21	1.95	5 (33%)
3	TRS	D	603	-	7,7,7	0.29	0	9,9,9	0.36	0
2	NAG	E	601	1	14,14,15	0.49	0	15,19,21	1.29	2 (13%)
3	TRS	E	602	-	7,7,7	0.55	0	9,9,9	5.43	6 (66%)
2	NAG	F	601	1	14,14,15	0.35	0	15,19,21	1.51	1 (6%)
3	TRS	F	602	-	7,7,7	0.35	0	9,9,9	0.55	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	NAG	A	601	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	602	2	1/1/5/7	0/6/23/26	0/1/1/1
3	TRS	A	603	-	-	0/9/9/9	0/0/0/0
2	NAG	B	601	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	602	2	1/1/5/7	0/6/23/26	0/1/1/1
3	TRS	B	603	-	-	0/9/9/9	0/0/0/0
2	NAG	C	601	1	-	0/6/23/26	0/1/1/1
3	TRS	C	602	-	-	0/9/9/9	0/0/0/0
2	NAG	D	601	1,2	-	0/6/23/26	0/1/1/1
2	NAG	D	602	2	1/1/5/7	0/6/23/26	0/1/1/1
3	TRS	D	603	-	-	0/9/9/9	0/0/0/0
2	NAG	E	601	1	-	0/6/23/26	0/1/1/1
3	TRS	E	602	-	-	0/9/9/9	0/0/0/0
2	NAG	F	601	1	-	0/6/23/26	0/1/1/1
3	TRS	F	602	-	-	0/9/9/9	0/0/0/0

There are no bond length outliers.

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	602	TRS	C1-C-N	-9.40	87.75	107.73
3	E	602	TRS	C3-C-N	-8.79	89.04	107.73
3	E	602	TRS	C2-C-N	-8.36	89.94	107.73
2	D	601	NAG	O5-C1-C2	-2.52	107.97	111.47
2	A	602	NAG	O7-C7-C8	-2.30	117.87	122.06
2	C	601	NAG	C1-C2-N2	-2.21	106.71	110.49
2	D	602	NAG	O7-C7-C8	-2.00	118.42	122.06
2	D	601	NAG	C2-N2-C7	2.08	125.98	122.94
2	B	601	NAG	C2-N2-C7	2.19	126.13	122.94
2	D	602	NAG	C4-C3-C2	2.33	114.44	111.02
2	B	601	NAG	C8-C7-N2	2.40	120.43	116.11
2	E	601	NAG	C8-C7-N2	2.45	120.54	116.11
2	C	601	NAG	C1-O5-C5	2.46	115.56	112.17
2	E	601	NAG	C2-N2-C7	2.52	126.62	122.94
2	D	602	NAG	C3-C4-C5	2.53	114.68	110.22
3	E	602	TRS	C2-C-C1	2.70	118.72	111.06
2	B	602	NAG	C1-O5-C5	2.90	116.16	112.17
2	D	602	NAG	C1-O5-C5	2.95	116.23	112.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	602	NAG	C1-O5-C5	3.24	116.64	112.17
3	E	602	TRS	C3-C-C1	3.27	120.34	111.06
3	E	602	TRS	C3-C-C2	3.45	120.83	111.06
2	F	601	NAG	C1-O5-C5	4.37	118.19	112.17
2	B	602	NAG	O5-C1-C2	4.93	118.33	111.47
2	A	602	NAG	O5-C1-C2	5.01	118.44	111.47
2	D	602	NAG	O5-C1-C2	5.16	118.66	111.47

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	602	NAG	C1
2	B	602	NAG	C1
2	D	602	NAG	C1

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	488/489 (99%)	-0.40	0 100 100	20, 31, 49, 71	0
1	B	488/489 (99%)	-0.26	0 100 100	22, 35, 54, 76	0
1	C	488/489 (99%)	-0.35	2 (0%) 92 93	22, 34, 54, 75	0
1	D	486/489 (99%)	-0.22	2 (0%) 92 93	31, 46, 65, 83	0
1	E	485/489 (99%)	0.05	10 (2%) 64 68	30, 48, 71, 95	0
1	F	471/489 (96%)	0.19	16 (3%) 46 53	30, 53, 79, 101	0
All	All	2906/2934 (99%)	-0.17	30 (1%) 82 85	20, 40, 69, 101	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	131	LYS	4.2
1	F	413	ILE	3.5
1	F	89	ARG	3.3
1	C	344	PRO	3.1
1	E	491	VAL	2.9
1	E	28	ASP	2.9
1	E	501	MET	2.8
1	F	28	ASP	2.7
1	F	469	ASP	2.7
1	D	27	ASP	2.6
1	F	73	ILE	2.6
1	F	492	VAL	2.6
1	E	24	ARG	2.6
1	D	433	ASP	2.5
1	F	418	ILE	2.5
1	E	94	ASP	2.4
1	E	492	VAL	2.4
1	C	499	GLU	2.4
1	F	472	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
1	F	481	PHE	2.3
1	F	24	ARG	2.2
1	F	27	ASP	2.2
1	F	470	PRO	2.2
1	F	478	LYS	2.1
1	F	412	LEU	2.1
1	E	48	GLY	2.1
1	F	300	GLN	2.1
1	E	343	GLY	2.1
1	E	344	PRO	2.1
1	E	487	ILE	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(Å ²)	Q<0.9
3	TRS	B	603	8/8	0.91	0.17	2.81	50,56,59,62	0
3	TRS	F	602	8/8	0.80	0.18	2.69	82,85,88,91	0
3	TRS	E	602	8/8	0.91	0.16	2.45	74,78,79,79	0
3	TRS	A	603	8/8	0.87	0.15	2.10	58,64,66,69	0
3	TRS	D	603	8/8	0.92	0.15	1.63	68,69,71,71	0
3	TRS	C	602	8/8	0.90	0.14	1.05	56,59,60,60	0
2	NAG	E	601	14/15	0.85	0.20	-	64,69,76,79	0
2	NAG	B	601	14/15	0.84	0.21	-	62,66,75,88	0
2	NAG	D	602	14/15	0.74	0.30	-	88,93,98,99	0
2	NAG	A	602	14/15	0.80	0.25	-	60,76,84,84	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	NAG	C	601	14/15	0.79	0.17	-	64,69,73,74	0
2	NAG	D	601	14/15	0.85	0.17	-	66,73,78,87	0
2	NAG	B	602	14/15	0.80	0.33	-	94,99,104,105	0
2	NAG	A	601	14/15	0.95	0.10	-	50,60,62,71	0
2	NAG	F	601	14/15	0.88	0.19	-	73,79,81,82	0

6.5 Other polymers [i](#)

There are no such residues in this entry.