



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 07:16 PM GMT

PDB ID : 4RHH  
Title : Crystal structure of the catalytic mutant Xyn52B2-E335G, a GH52 Beta-D-xylosidase from *Geobacillus stearothermophilus* T6  
Authors : Dann, R.; Lansky, S.; Lavid, N.; Zehavi, A.; Belakhov, V.; Baasov, T.; Manjasetty, B.; Belrhali, H.; Shoham, Y.; Shoham, G.  
Deposited on : 2014-10-02  
Resolution : 2.15 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026982
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20026982

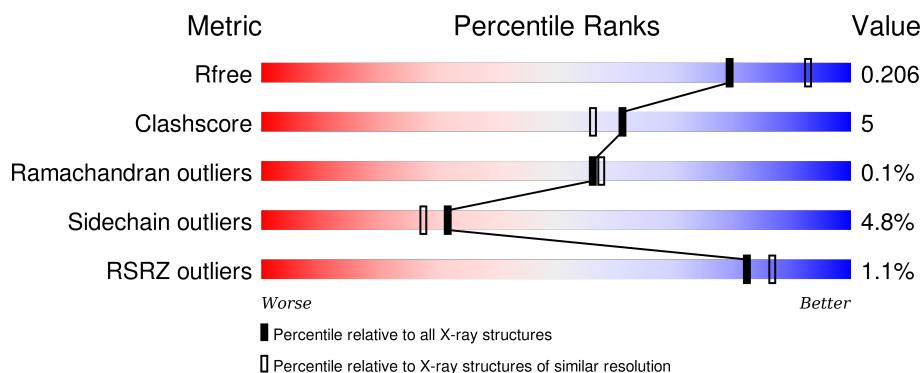
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.15 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	1045 (2.16-2.16)
Clashscore	102246	1152 (2.16-2.16)
Ramachandran outliers	100387	1131 (2.16-2.16)
Sidechain outliers	100360	1131 (2.16-2.16)
RSRZ outliers	91569	1050 (2.16-2.16)

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  $\geq 3$ , 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	705	<div> <div>2%</div> <div>85% 12% ..</div> </div>
1	B	705	<div> <div>2%</div> <div>87% 10% ..</div> </div>
1	C	705	<div> <div>2%</div> <div>86% 11% ..</div> </div>
1	D	705	<div> <div>2%</div> <div>86% 11% ..</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	TRS	C	802	-	-	-	X
3	TRS	D	802	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 24010 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-xylosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	694	Total	C	N	O	S	0	0	0
			5543	3534	942	1041	26			
1	B	694	Total	C	N	O	S	0	2	0
			5565	3546	950	1043	26			
1	C	696	Total	C	N	O	S	0	0	0
			5550	3538	944	1042	26			
1	D	693	Total	C	N	O	S	0	0	0
			5534	3528	941	1039	26			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	335	GLY	GLU	ENGINEERED MUTATION	UNP Q09LZ0
B	335	GLY	GLU	ENGINEERED MUTATION	UNP Q09LZ0
C	335	GLY	GLU	ENGINEERED MUTATION	UNP Q09LZ0
D	335	GLY	GLU	ENGINEERED MUTATION	UNP Q09LZ0

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Ca	0	0
			1	1		
2	A	1	Total	Ca	0	0
			1	1		
2	D	1	Total	Ca	0	0
			1	1		
2	C	2	Total	Ca	0	0
			2	2		

- Molecule 3 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: C<sub>4</sub>H<sub>12</sub>NO<sub>3</sub>).



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

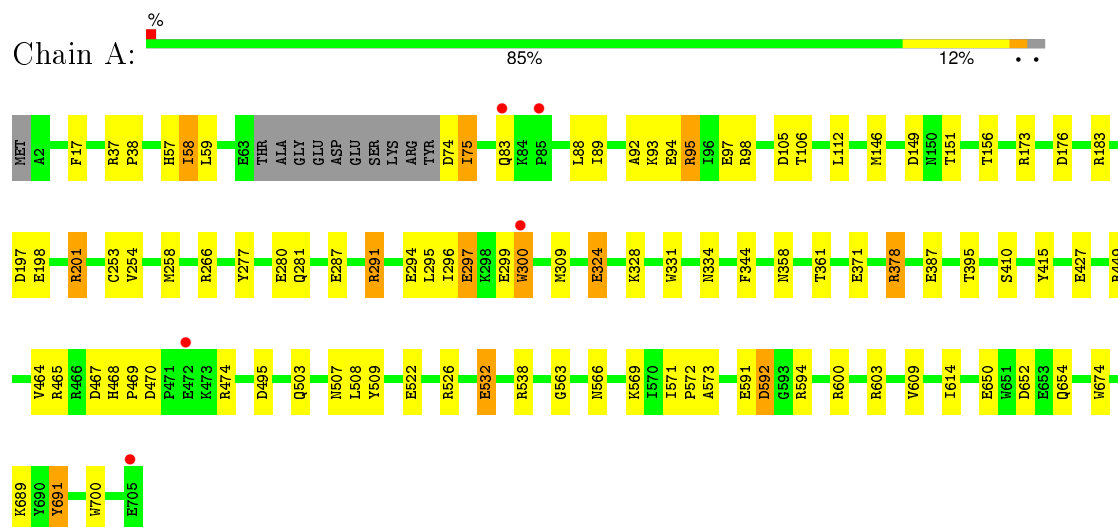
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	434	Total	O	0	0
			434	434		
4	B	439	Total	O	0	0
			439	439		
4	C	454	Total	O	0	0
			454	454		
4	D	430	Total	O	0	0
			430	430		

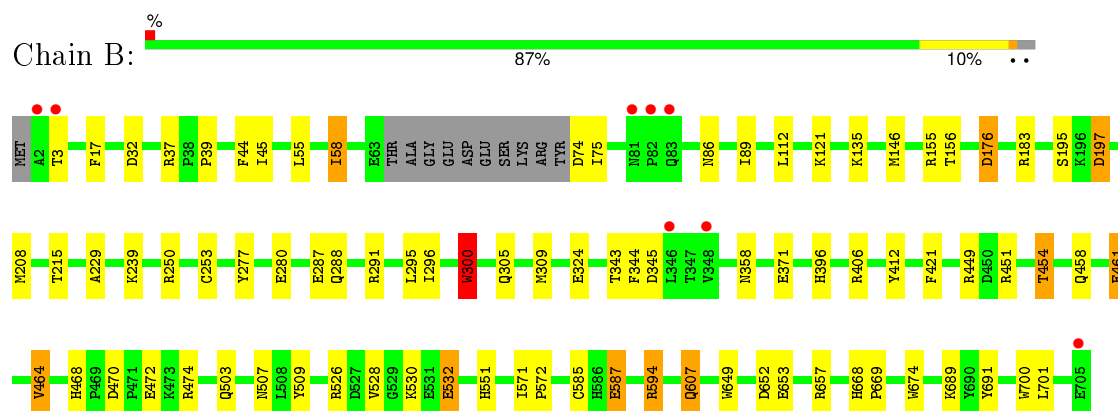
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

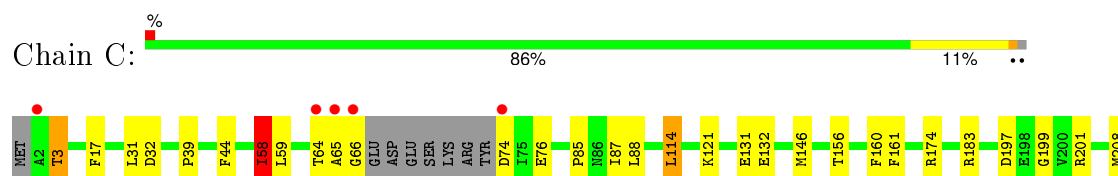
#### • Molecule 1: Beta-xylosidase

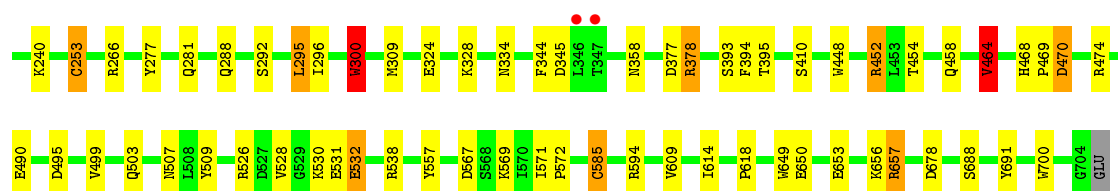


#### • Molecule 1: Beta-xylosidase

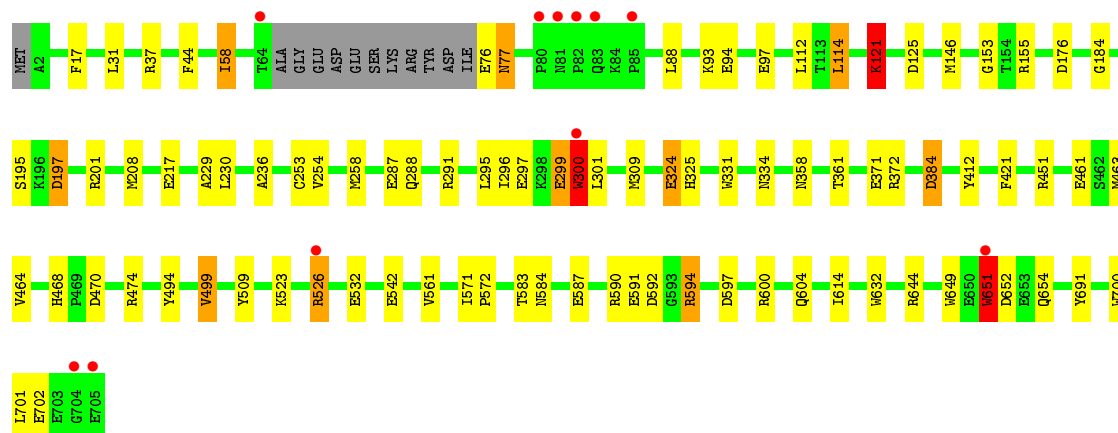
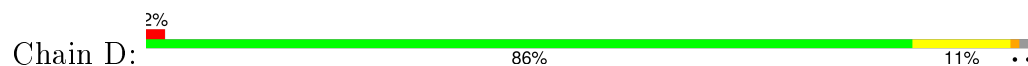


#### • Molecule 1: Beta-xylosidase





• Molecule 1: Beta-xylosidase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	97.45Å 119.08Å 243.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.43 – 2.15 39.40 – 2.15	Depositor EDS
% Data completeness (in resolution range)	98.5 (39.43-2.15) 98.5 (39.40-2.15)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.91 (at 2.16Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.150 , 0.203 0.159 , 0.206	Depositor DCC
$R_{free}$ test set	7592 reflections (5.27%)	DCC
Wilson B-factor (Å <sup>2</sup> )	21.4	Xtriage
Anisotropy	0.042	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 45.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 151616 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	24010	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.08% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

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



## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CA, 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	0.89	0/5689	0.91	5/7721 (0.1%)
1	B	0.92	4/5711 (0.1%)	0.94	4/7749 (0.1%)
1	C	0.90	2/5696 (0.0%)	0.95	13/7731 (0.2%)
1	D	0.91	4/5680 (0.1%)	0.94	6/7709 (0.1%)
All	All	0.90	10/22776 (0.0%)	0.93	28/30910 (0.1%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	649	TRP	CB-CG	-6.22	1.39	1.50
1	D	651	TRP	CB-CG	5.78	1.60	1.50
1	D	300	TRP	NE1-CE2	-5.67	1.30	1.37
1	C	490	GLU	CD-OE1	5.43	1.31	1.25
1	D	300	TRP	CD2-CE3	-5.42	1.32	1.40
1	B	674	TRP	CB-CG	-5.34	1.40	1.50
1	B	461	GLU	CG-CD	5.26	1.59	1.51
1	B	371	GLU	CG-CD	5.10	1.59	1.51
1	B	300	TRP	CB-CG	-5.03	1.41	1.50
1	C	300	TRP	CD2-CE3	-5.01	1.32	1.40

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	451	ARG	NE-CZ-NH2	-8.84	115.88	120.30
1	D	125	ASP	CB-CG-OD1	5.96	123.66	118.30
1	C	114	LEU	CA-CB-CG	-5.94	101.64	115.30
1	A	495	ASP	CB-CG-OD1	5.84	123.55	118.30
1	A	173	ARG	NE-CZ-NH2	-5.77	117.42	120.30
1	C	201	ARG	NE-CZ-NH1	5.77	123.18	120.30
1	A	258	MET	CG-SD-CE	-5.73	91.04	100.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	201	ARG	NE-CZ-NH2	-5.70	117.45	120.30
1	B	406	ARG	NE-CZ-NH1	5.53	123.06	120.30
1	C	201	ARG	CG-CD-NE	5.48	123.31	111.80
1	C	58	ILE	CA-CB-CG2	5.48	121.85	110.90
1	C	266	ARG	NE-CZ-NH2	-5.47	117.57	120.30
1	D	121	LYS	CD-CE-NZ	5.44	124.21	111.70
1	C	174	ARG	NE-CZ-NH2	-5.33	117.63	120.30
1	C	183	ARG	NE-CZ-NH2	-5.33	117.64	120.30
1	C	594	ARG	CB-CA-C	-5.33	99.75	110.40
1	C	197	ASP	CB-CG-OD1	5.32	123.09	118.30
1	C	378	ARG	NE-CZ-NH1	5.29	122.95	120.30
1	B	58	ILE	CA-CB-CG2	5.24	121.38	110.90
1	C	470	ASP	CB-CG-OD1	5.23	123.01	118.30
1	D	372	ARG	NE-CZ-NH1	5.21	122.91	120.30
1	D	114	LEU	CA-CB-CG	-5.21	103.33	115.30
1	C	464	VAL	CG1-CB-CG2	5.19	119.20	110.90
1	A	105	ASP	CB-CG-OD1	5.11	122.90	118.30
1	B	461	GLU	CA-CB-CG	5.11	124.64	113.40
1	A	173	ARG	NE-CZ-NH1	5.04	122.82	120.30
1	B	250	ARG	NE-CZ-NH2	-5.03	117.78	120.30
1	D	384	ASP	CB-CG-OD2	5.02	122.82	118.30

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	5543	0	5337	70	0
1	B	5565	0	5361	53	0
1	C	5550	0	5346	55	0
1	D	5534	0	5329	58	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	2	0	0	0	0
2	D	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	16	0	24	5	0
3	B	8	0	12	0	0
3	C	16	0	24	5	0
3	D	16	0	24	2	0
4	A	434	0	0	27	0
4	B	439	0	0	15	0
4	C	454	0	0	12	0
4	D	430	0	0	18	0
All	All	24010	0	21457	240	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 (240) 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:607:GLN:HG3	4:B:1317:HOH:O	1.50	1.12
1:A:468:HIS:HD2	1:A:470:ASP:H	1.16	0.92
1:B:296:ILE:HG21	4:B:1326:HOH:O	1.71	0.89
1:B:253:CYS:SG	4:B:1328:HOH:O	2.30	0.89
1:A:97:GLU:HG3	4:A:928:HOH:O	1.73	0.87
1:D:296:ILE:HG21	4:D:1325:HOH:O	1.78	0.83
1:D:300:TRP:HZ3	1:D:700:TRP:O	1.60	0.83
1:D:253:CYS:SG	4:D:1026:HOH:O	2.35	0.83
1:D:468:HIS:HD2	1:D:470:ASP:H	1.25	0.81
1:C:526:ARG:HH22	1:C:538:ARG:HH12	1.28	0.81
1:D:371:GLU:OE1	3:D:802:TRS:H12	1.80	0.80
1:C:300:TRP:HZ3	1:C:700:TRP:O	1.65	0.80
1:C:526:ARG:HH22	1:C:538:ARG:NH1	1.81	0.79
1:C:253:CYS:SG	4:C:1222:HOH:O	2.40	0.78
1:D:299:GLU:HG2	4:D:1080:HOH:O	1.83	0.78
1:B:45:ILE:HG23	1:B:146:MET:CE	2.12	0.78
1:D:287:GLU:HG2	4:D:1192:HOH:O	1.83	0.77
1:C:526:ARG:NH2	1:C:538:ARG:HH12	1.84	0.75
1:A:573:ALA:HA	4:A:1323:HOH:O	1.87	0.74
1:B:587:GLU:OE2	1:B:594:ARG:NH1	2.21	0.74
1:A:591:GLU:HB3	4:A:926:HOH:O	1.89	0.73
1:B:74:ASP:N	4:B:1289:HOH:O	2.22	0.72
1:B:587:GLU:HG2	4:B:940:HOH:O	1.87	0.72
1:A:291:ARG:NH1	1:A:294:GLU:OE1	2.22	0.72
1:C:585:CYS:SG	4:C:1160:HOH:O	2.49	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:468:HIS:HD2	1:B:470:ASP:H	1.38	0.70
1:D:300:TRP:CZ3	1:D:700:TRP:O	2.45	0.70
1:A:522:GLU:OE1	1:A:594:ARG:NH2	2.25	0.70
1:A:201:ARG:HG3	1:A:201:ARG:HH11	1.58	0.69
1:A:691:TYR:OH	3:A:803:TRS:H12	1.92	0.68
1:A:526:ARG:HH22	1:A:538:ARG:NH1	1.91	0.68
1:A:603:ARG:HD3	4:A:1320:HOH:O	1.94	0.68
1:D:587:GLU:OE2	1:D:594:ARG:NH1	2.27	0.67
1:B:3:THR:HG23	4:B:1201:HOH:O	1.94	0.67
1:A:300:TRP:N	1:A:300:TRP:CD1	2.62	0.66
1:A:266:ARG:HD2	1:A:324:GLU:OE1	1.97	0.65
3:C:803:TRS:H12	4:C:1354:HOH:O	1.94	0.65
3:C:802:TRS:C2	4:C:1118:HOH:O	2.43	0.65
1:B:208:MET:HE1	1:B:229:ALA:HB2	1.78	0.65
1:B:300:TRP:HZ3	1:B:700:TRP:O	1.80	0.65
1:B:454:THR:HG23	4:B:1135:HOH:O	1.96	0.65
1:D:300:TRP:CD1	1:D:300:TRP:N	2.65	0.64
1:C:657:ARG:HD3	4:C:934:HOH:O	1.97	0.64
1:D:184:GLY:HA2	1:D:230:LEU:HD21	1.79	0.64
1:A:297:GLU:CG	4:A:1200:HOH:O	2.47	0.62
1:A:253:CYS:HB3	4:A:1056:HOH:O	2.00	0.62
1:D:526:ARG:HB2	1:D:526:ARG:CZ	2.29	0.61
1:D:523:LYS:HE2	1:D:583:THR:O	2.01	0.61
1:B:464:VAL:HG13	1:B:474:ARG:CZ	2.31	0.61
1:D:384:ASP:OD1	4:D:1016:HOH:O	2.16	0.60
1:C:199:GLY:HA3	4:C:1061:HOH:O	2.01	0.60
1:A:508:LEU:HG	4:A:1323:HOH:O	2.01	0.60
1:A:526:ARG:NH2	1:A:538:ARG:NH1	2.50	0.60
1:B:44:PHE:CD2	1:B:208:MET:HE2	2.37	0.59
1:B:464:VAL:HG13	1:B:474:ARG:NH1	2.18	0.59
1:D:153:GLY:O	1:D:236:ALA:HB1	2.02	0.59
1:C:3:THR:HG22	4:C:1276:HOH:O	2.03	0.58
1:C:44:PHE:CD2	1:C:208:MET:HE2	2.38	0.58
1:B:657:ARG:NH1	4:B:1278:HOH:O	2.37	0.58
1:D:591:GLU:CD	1:D:591:GLU:H	2.06	0.57
1:C:378:ARG:NH1	4:C:1139:HOH:O	2.38	0.57
1:A:689:LYS:HE3	1:B:75:ILE:HG23	1.87	0.57
1:A:464:VAL:HG12	1:A:474:ARG:CZ	2.35	0.57
1:B:208:MET:CE	1:B:229:ALA:HB2	2.35	0.56
1:C:468:HIS:HD2	1:C:470:ASP:H	1.53	0.56
1:D:195:SER:OG	1:D:197:ASP:HB2	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:GLU:CG	4:A:928:HOH:O	2.40	0.56
1:A:526:ARG:HD2	4:A:1220:HOH:O	2.05	0.56
1:D:334:ASN:HB3	4:D:988:HOH:O	2.04	0.56
3:A:802:TRS:H21	4:A:1193:HOH:O	2.06	0.56
1:A:280:GLU:HG3	4:A:1032:HOH:O	2.05	0.56
1:D:58:ILE:CD1	1:D:88:LEU:HD22	2.36	0.56
1:A:300:TRP:HZ3	1:A:700:TRP:O	1.88	0.55
1:B:461:GLU:HG3	4:B:1190:HOH:O	2.05	0.55
1:A:532:GLU:CD	1:A:532:GLU:H	2.10	0.55
1:C:58:ILE:HD11	1:C:160:PHE:HZ	1.71	0.55
1:C:300:TRP:CZ3	1:C:700:TRP:O	2.53	0.55
1:D:94:GLU:CD	1:D:94:GLU:H	2.10	0.55
1:A:468:HIS:CD2	1:A:469:PRO:HD2	2.41	0.55
1:C:58:ILE:HD11	1:C:160:PHE:CZ	2.41	0.55
1:D:571:ILE:HB	1:D:572:PRO:HD3	1.88	0.55
1:D:526:ARG:CB	1:D:526:ARG:CZ	2.85	0.55
1:A:468:HIS:CD2	1:A:470:ASP:H	2.08	0.54
1:D:208:MET:HE1	1:D:229:ALA:HB2	1.88	0.54
1:C:532:GLU:H	1:C:532:GLU:CD	2.11	0.54
1:D:464:VAL:HG13	1:D:474:ARG:CZ	2.38	0.54
1:A:201:ARG:HG3	1:A:201:ARG:NH1	2.21	0.53
1:A:112:LEU:HD11	1:A:146:MET:CE	2.39	0.53
1:C:526:ARG:CZ	1:C:526:ARG:HB2	2.38	0.53
1:D:654:GLN:HG3	4:D:1037:HOH:O	2.07	0.53
1:B:300:TRP:N	1:B:300:TRP:CD1	2.74	0.53
1:C:146:MET:CE	1:C:161:PHE:HB2	2.39	0.53
1:C:65:ALA:N	1:C:66:GLY:HA3	2.25	0.52
1:D:526:ARG:HD2	4:D:1106:HOH:O	2.09	0.52
1:D:44:PHE:CD2	1:D:208:MET:HE2	2.44	0.52
1:D:702:GLU:HB3	4:D:997:HOH:O	2.10	0.51
1:C:76:GLU:OE2	1:D:37:ARG:HD3	2.10	0.51
1:B:300:TRP:HH2	1:B:701:LEU:O	1.94	0.51
1:A:149:ASP:OD1	1:A:151:THR:HG23	2.11	0.51
1:A:603:ARG:CD	4:A:1320:HOH:O	2.55	0.51
1:C:528:VAL:HG23	1:C:530:LYS:HG3	1.93	0.51
1:C:334:ASN:HB3	4:C:1002:HOH:O	2.10	0.51
1:A:652:ASP:H	1:A:654:GLN:HE22	1.57	0.51
1:C:44:PHE:CG	1:C:208:MET:HE2	2.46	0.51
1:A:395:THR:HB	1:A:410:SER:HB3	1.93	0.51
1:B:45:ILE:HG23	1:B:146:MET:HE1	1.91	0.50
1:A:465:ARG:HH11	1:A:465:ARG:HG2	1.75	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:495:ASP:OD1	3:C:803:TRS:H21	2.11	0.50
1:B:571:ILE:HB	1:B:572:PRO:HD3	1.94	0.50
1:A:328:LYS:HE2	4:A:1261:HOH:O	2.10	0.50
1:C:146:MET:HE1	1:C:161:PHE:HB2	1.94	0.50
3:C:802:TRS:H21	4:C:1118:HOH:O	2.06	0.49
1:A:464:VAL:O	1:A:467:ASP:HB3	2.12	0.49
1:A:287:GLU:HG3	4:A:1328:HOH:O	2.12	0.49
1:D:253:CYS:CB	4:D:1026:HOH:O	2.59	0.49
1:A:112:LEU:HD11	1:A:146:MET:HE3	1.94	0.49
1:D:464:VAL:HG13	1:D:474:ARG:NH1	2.28	0.48
1:D:208:MET:CE	1:D:229:ALA:HB2	2.43	0.48
1:B:195:SER:OG	1:B:197:ASP:HB2	2.13	0.48
1:A:296:ILE:O	1:A:296:ILE:HG22	2.13	0.48
3:D:802:TRS:H31	4:D:1147:HOH:O	2.12	0.48
1:A:98:ARG:HA	1:A:106:THR:O	2.14	0.48
1:A:57:HIS:HA	1:A:89:ILE:O	2.14	0.48
1:D:201:ARG:NH1	1:D:217:GLU:OE1	2.35	0.48
1:C:58:ILE:HG23	1:C:59:LEU:O	2.14	0.48
1:A:650:GLU:HB2	4:A:1187:HOH:O	2.11	0.48
1:A:334:ASN:HB3	4:A:1003:HOH:O	2.14	0.48
1:D:324:GLU:HG2	4:D:1218:HOH:O	2.14	0.47
1:C:296:ILE:HD12	1:C:309:MET:HE1	1.96	0.47
1:A:503:GLN:O	1:A:507:ASN:HB2	2.14	0.47
1:C:526:ARG:NH2	1:C:538:ARG:NH1	2.52	0.47
1:B:454:THR:HG22	1:B:458:GLN:HE21	1.79	0.47
1:C:618:PRO:O	1:C:657:ARG:NH2	2.48	0.47
1:B:287:GLU:OE2	1:B:291:ARG:HD3	2.15	0.47
1:C:85:PRO:HB2	1:C:87:ILE:HG23	1.97	0.47
1:A:526:ARG:NH2	1:A:538:ARG:HH12	2.11	0.47
1:D:93:LYS:HG2	4:D:1139:HOH:O	2.14	0.47
1:A:674:TRP:CG	1:A:689:LYS:HD2	2.50	0.47
1:B:291:ARG:NE	4:B:1159:HOH:O	2.48	0.47
1:D:499:VAL:HG13	1:D:561:VAL:HG21	1.96	0.47
1:D:604:GLN:HG3	4:D:925:HOH:O	2.14	0.46
1:A:58:ILE:HG23	1:A:59:LEU:O	2.15	0.46
1:B:530:LYS:HB3	1:B:532:GLU:OE2	2.15	0.46
1:C:468:HIS:CD2	1:C:469:PRO:HD2	2.49	0.46
1:C:58:ILE:HD13	1:C:88:LEU:HB3	1.97	0.46
1:B:532:GLU:HB3	4:B:1121:HOH:O	2.13	0.46
1:A:571:ILE:HB	1:A:572:PRO:HD3	1.98	0.46
1:C:32:ASP:HB2	1:C:39:PRO:HG3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:ALA:HB3	1:A:95:ARG:HG3	1.97	0.46
1:B:503:GLN:O	1:B:507:ASN:HB2	2.16	0.46
1:C:300:TRP:N	1:C:300:TRP:CD1	2.80	0.46
1:B:652:ASP:HB3	1:B:653:GLU:OE1	2.16	0.46
1:A:74:ASP:N	4:A:1295:HOH:O	2.49	0.46
1:A:58:ILE:CD1	1:A:88:LEU:HD22	2.45	0.45
1:D:31:LEU:N	1:D:31:LEU:HD12	2.31	0.45
1:B:412:TYR:CE1	1:B:421:PHE:HA	2.51	0.45
1:D:112:LEU:HD11	1:D:146:MET:HE3	1.97	0.45
1:D:300:TRP:HH2	1:D:701:LEU:O	2.00	0.45
1:A:297:GLU:HG3	4:A:1200:HOH:O	2.13	0.45
1:B:112:LEU:HD23	1:B:112:LEU:C	2.36	0.45
1:D:296:ILE:N	1:D:296:ILE:HD13	2.31	0.45
1:C:393:SER:OG	1:C:394:PHE:N	2.50	0.45
1:B:37:ARG:CZ	1:B:37:ARG:HB3	2.45	0.44
1:A:600:ARG:HD2	4:A:1133:HOH:O	2.16	0.44
1:C:377:ASP:C	1:C:377:ASP:OD1	2.56	0.44
1:C:292:SER:O	1:C:295:LEU:HB2	2.16	0.44
1:C:503:GLN:O	1:C:507:ASN:HB2	2.17	0.44
1:B:32:ASP:HB2	1:B:39:PRO:HG3	1.99	0.44
1:B:121:LYS:NZ	1:B:288:GLN:HE22	2.14	0.44
1:A:331:TRP:CD1	1:A:361:THR:HB	2.53	0.44
1:C:395:THR:HB	1:C:410:SER:HB3	1.99	0.44
1:C:44:PHE:HB2	1:C:208:MET:HE3	2.00	0.44
1:B:343:THR:HG23	1:B:396:HIS:CE1	2.53	0.44
1:A:75:ILE:HG23	1:B:689:LYS:HE3	2.00	0.44
1:D:76:GLU:O	1:D:77:ASN:HB3	2.18	0.44
1:C:653:GLU:HA	1:C:656:LYS:HD3	2.00	0.43
1:B:305:GLN:O	1:B:309:MET:HG3	2.18	0.43
1:B:528:VAL:HG23	1:B:530:LYS:HG3	2.00	0.43
1:A:37:ARG:O	1:A:38:PRO:C	2.56	0.43
1:D:253:CYS:HB3	4:D:1026:HOH:O	2.17	0.43
1:C:571:ILE:HB	1:C:572:PRO:HD3	2.01	0.43
1:B:454:THR:CG2	4:B:1135:HOH:O	2.62	0.43
1:A:464:VAL:CG1	1:A:474:ARG:CZ	2.97	0.43
1:B:532:GLU:H	1:B:532:GLU:CD	2.21	0.43
1:C:678:ASP:HB3	1:C:688:SER:HB2	2.00	0.43
1:D:590:ARG:NH2	1:D:592:ASP:OD2	2.52	0.43
1:D:112:LEU:HD11	1:D:146:MET:CE	2.48	0.43
1:D:542:GLU:HG2	4:D:1156:HOH:O	2.17	0.43
1:C:464:VAL:HG13	1:C:474:ARG:CZ	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:591:GLU:O	1:D:600:ARG:NH1	2.52	0.43
1:A:563:GLY:N	4:A:1243:HOH:O	2.52	0.43
1:C:121:LYS:NZ	1:C:288:GLN:HE22	2.16	0.43
1:B:45:ILE:HG23	1:B:146:MET:HE2	1.99	0.43
1:A:532:GLU:CD	1:A:532:GLU:N	2.72	0.43
1:D:644:ARG:HD3	1:D:651:TRP:CE3	2.54	0.43
1:C:448:TRP:CE2	1:C:452:ARG:HD3	2.54	0.42
1:B:668:HIS:HA	1:B:669:PRO:HD3	1.82	0.42
1:C:65:ALA:N	1:C:66:GLY:CA	2.83	0.42
1:A:378:ARG:NH2	1:A:387:GLU:OE1	2.53	0.42
1:D:597:ASP:OD2	1:D:600:ARG:NH2	2.45	0.42
1:A:198:GLU:HB2	4:A:1276:HOH:O	2.20	0.42
1:D:331:TRP:CD1	1:D:361:THR:HB	2.55	0.42
1:D:258:MET:HE2	1:D:325:HIS:CD2	2.55	0.42
1:D:121:LYS:HD3	1:D:288:GLN:NE2	2.34	0.42
1:D:584:ASN:N	4:D:1203:HOH:O	2.35	0.42
1:B:468:HIS:CD2	1:B:470:ASP:H	2.27	0.42
3:A:802:TRS:C2	4:A:1193:HOH:O	2.65	0.42
1:A:415:TYR:CD1	1:A:415:TYR:N	2.86	0.42
1:D:301:LEU:HD13	1:D:309:MET:HE1	2.01	0.42
1:B:55:LEU:HA	1:B:55:LEU:HD23	1.84	0.42
3:C:802:TRS:C3	4:C:1121:HOH:O	2.68	0.41
1:A:253:CYS:CB	4:A:1056:HOH:O	2.63	0.41
1:C:58:ILE:CD1	1:C:88:LEU:HD22	2.50	0.41
1:B:551:HIS:HE1	4:B:1186:HOH:O	2.03	0.41
1:C:31:LEU:HD12	1:C:31:LEU:N	2.34	0.41
1:D:334:ASN:CB	4:D:988:HOH:O	2.65	0.41
1:A:592:ASP:CG	4:A:1281:HOH:O	2.57	0.41
1:A:592:ASP:OD1	1:A:592:ASP:N	2.48	0.41
1:A:371:GLU:OE1	3:A:802:TRS:H22	2.20	0.41
1:D:412:TYR:CE1	1:D:421:PHE:HA	2.56	0.41
1:C:454:THR:O	1:C:458:GLN:HG3	2.20	0.41
1:C:569:LYS:CG	1:C:609:VAL:HG12	2.51	0.41
1:C:74:ASP:HA	4:C:956:HOH:O	2.19	0.41
1:B:607:GLN:CG	4:B:1317:HOH:O	2.31	0.41
1:A:566:ASN:HB2	4:A:1105:HOH:O	2.20	0.41
1:B:44:PHE:CG	1:B:208:MET:HE2	2.56	0.41
1:A:569:LYS:CG	1:A:609:VAL:HG12	2.51	0.41
1:A:371:GLU:OE1	3:A:802:TRS:C2	2.69	0.41
1:C:296:ILE:HD12	1:C:309:MET:CE	2.51	0.41
1:B:135:LYS:HE2	1:B:277:TYR:CD2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:494:TYR:HA	1:D:632:TRP:CH2	2.55	0.41
1:D:463:MET:HE2	1:D:463:MET:HB3	1.92	0.41
1:B:280:GLU:HG3	4:B:1111:HOH:O	2.21	0.40
1:B:176:ASP:OD1	1:B:176:ASP:N	2.54	0.40
1:C:557:TYR:OH	1:C:567:ASP:OD1	2.35	0.40
1:A:93:LYS:HE3	4:A:1262:HOH:O	2.20	0.40
1:C:277:TYR:CZ	1:C:281:GLN:HG3	2.56	0.40
1:A:522:GLU:OE2	4:A:1282:HOH:O	2.21	0.40
1:B:451[B]:ARG:HH11	1:B:451[B]:ARG:HD2	1.71	0.40
1:A:277:TYR:CZ	1:A:281:GLN:HG3	2.57	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	690/705 (98%)	664 (96%)	25 (4%)	1 (0%)	56	57
1	B	692/705 (98%)	675 (98%)	17 (2%)	0	100	100
1	C	692/705 (98%)	675 (98%)	16 (2%)	1 (0%)	56	57
1	D	689/705 (98%)	662 (96%)	25 (4%)	2 (0%)	46	42
All	All	2763/2820 (98%)	2676 (97%)	83 (3%)	4 (0%)	56	57

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	77	ASN
1	D	614	ILE
1	C	614	ILE
1	A	614	ILE

### 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	584/593 (98%)	556 (95%)	28 (5%)	31	27
1	B	586/593 (99%)	556 (95%)	30 (5%)	29	24
1	C	584/593 (98%)	556 (95%)	28 (5%)	31	27
1	D	583/593 (98%)	558 (96%)	25 (4%)	35	32
All	All	2337/2372 (98%)	2226 (95%)	111 (5%)	31	28

All (111) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	PHE
1	A	58	ILE
1	A	75	ILE
1	A	83	GLN
1	A	94	GLU
1	A	95	ARG
1	A	156	THR
1	A	176	ASP
1	A	183	ARG
1	A	197	ASP
1	A	201	ARG
1	A	254	VAL
1	A	291	ARG
1	A	295	LEU
1	A	297	GLU
1	A	299	GLU
1	A	300	TRP
1	A	309	MET
1	A	324	GLU
1	A	344	PHE
1	A	358	ASN
1	A	378	ARG
1	A	427	GLU
1	A	449	ARG

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Mol	Chain	Res	Type
1	A	509	TYR
1	A	532	GLU
1	A	592	ASP
1	A	691	TYR
1	B	17	PHE
1	B	58	ILE
1	B	86	ASN
1	B	89	ILE
1	B	155	ARG
1	B	156	THR
1	B	176	ASP
1	B	183	ARG
1	B	197	ASP
1	B	215	THR
1	B	239	LYS
1	B	295	LEU
1	B	300	TRP
1	B	324	GLU
1	B	344	PHE
1	B	345	ASP
1	B	358	ASN
1	B	449	ARG
1	B	454	THR
1	B	464	VAL
1	B	472	GLU
1	B	509	TYR
1	B	526	ARG
1	B	532	GLU
1	B	585	CYS
1	B	587	GLU
1	B	594	ARG
1	B	607	GLN
1	B	649	TRP
1	B	691	TYR
1	C	3	THR
1	C	17	PHE
1	C	58	ILE
1	C	64	THR
1	C	114	LEU
1	C	131	GLU
1	C	132	GLU
1	C	156	THR

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Mol	Chain	Res	Type
1	C	240	LYS
1	C	253	CYS
1	C	295	LEU
1	C	300	TRP
1	C	324	GLU
1	C	328	LYS
1	C	344	PHE
1	C	345	ASP
1	C	358	ASN
1	C	452	ARG
1	C	464	VAL
1	C	499	VAL
1	C	509	TYR
1	C	531	GLU
1	C	532	GLU
1	C	585	CYS
1	C	649	TRP
1	C	650	GLU
1	C	657	ARG
1	C	691	TYR
1	D	17	PHE
1	D	58	ILE
1	D	97	GLU
1	D	114	LEU
1	D	121	LYS
1	D	155	ARG
1	D	176	ASP
1	D	197	ASP
1	D	254	VAL
1	D	291	ARG
1	D	295	LEU
1	D	297	GLU
1	D	299	GLU
1	D	300	TRP
1	D	324	GLU
1	D	358	ASN
1	D	461	GLU
1	D	499	VAL
1	D	509	TYR
1	D	526	ARG
1	D	532	GLU
1	D	594	ARG

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Mol	Chain	Res	Type
1	D	651	TRP
1	D	652	ASP
1	D	691	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	288	GLN
1	A	468	HIS
1	A	654	GLN
1	B	288	GLN
1	B	458	GLN
1	B	468	HIS
1	B	551	HIS
1	C	288	GLN
1	C	458	GLN
1	C	468	HIS
1	C	551	HIS
1	D	86	ASN
1	D	288	GLN
1	D	468	HIS

### 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](#)

Of 12 ligands modelled in this entry, 5 are monoatomic - leaving 7 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
3	TRS	A	802	-	7,7,7	2.04	3 (42%)	9,9,9	2.03	5 (55%)
3	TRS	A	803	-	7,7,7	2.33	3 (42%)	9,9,9	2.50	5 (55%)
3	TRS	B	802	-	7,7,7	1.43	1 (14%)	9,9,9	2.43	4 (44%)
3	TRS	C	802	-	7,7,7	1.08	1 (14%)	9,9,9	2.68	4 (44%)
3	TRS	C	803	-	7,7,7	1.29	1 (14%)	9,9,9	2.98	5 (55%)
3	TRS	D	802	-	7,7,7	0.97	0	9,9,9	1.43	1 (11%)
3	TRS	D	803	-	7,7,7	1.30	1 (14%)	9,9,9	2.49	3 (33%)

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
3	TRS	A	802	-	-	0/9/9/9	0/0/0/0
3	TRS	A	803	-	-	0/9/9/9	0/0/0/0
3	TRS	B	802	-	-	0/9/9/9	0/0/0/0
3	TRS	C	802	-	-	0/9/9/9	0/0/0/0
3	TRS	C	803	-	-	0/9/9/9	0/0/0/0
3	TRS	D	802	-	-	0/9/9/9	0/0/0/0
3	TRS	D	803	-	-	0/9/9/9	0/0/0/0

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	803	TRS	C1-C	-4.07	1.47	1.53
3	B	802	TRS	C-N	-2.89	1.46	1.50
3	A	803	TRS	C2-C	-2.68	1.49	1.53
3	C	803	TRS	C1-C	-2.62	1.49	1.53
3	C	802	TRS	C1-C	2.02	1.56	1.53
3	D	803	TRS	C1-C	2.22	1.56	1.53
3	A	802	TRS	C2-C	2.53	1.56	1.53
3	A	802	TRS	C1-C	2.57	1.56	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	802	TRS	C3-C	3.41	1.57	1.53
3	A	803	TRS	C-N	3.57	1.55	1.50

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	803	TRS	O2-C2-C	-5.82	97.10	110.92
3	C	802	TRS	C3-C-N	-5.36	98.77	107.88
3	A	803	TRS	C2-C-C1	-5.10	99.89	110.65
3	C	803	TRS	O1-C1-C	-4.67	99.82	110.92
3	C	803	TRS	C2-C-C1	-3.50	103.28	110.65
3	A	802	TRS	C1-C-N	-2.97	102.84	107.88
3	C	802	TRS	C2-C-C1	-2.60	105.17	110.65
3	B	802	TRS	C2-C-C1	-2.52	105.33	110.65
3	D	803	TRS	C1-C-N	-2.35	103.89	107.88
3	B	802	TRS	C3-C-C2	-2.25	105.90	110.65
3	A	802	TRS	C2-C-N	-2.10	104.32	107.88
3	A	803	TRS	C2-C-N	2.10	111.44	107.88
3	A	803	TRS	C3-C-C2	2.11	115.10	110.65
3	C	802	TRS	C2-C-N	2.26	111.72	107.88
3	A	802	TRS	O3-C3-C	2.28	116.34	110.92
3	A	803	TRS	O3-C3-C	2.43	116.68	110.92
3	A	803	TRS	C3-C-N	2.48	112.09	107.88
3	A	802	TRS	O2-C2-C	2.58	117.05	110.92
3	A	802	TRS	C3-C-C2	2.82	116.59	110.65
3	D	803	TRS	C2-C-C1	2.98	116.92	110.65
3	C	803	TRS	C3-C-C2	3.08	117.14	110.65
3	B	802	TRS	C2-C-N	3.18	113.29	107.88
3	D	802	TRS	C2-C-C1	3.37	117.74	110.65
3	C	803	TRS	C2-C-N	3.76	114.27	107.88
3	C	803	TRS	O2-C2-C	3.98	120.38	110.92
3	B	802	TRS	C3-C-C1	4.69	120.53	110.65
3	C	802	TRS	C3-C-C1	4.74	120.63	110.65

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	802	TRS	4	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	803	TRS	1	0
3	C	802	TRS	3	0
3	C	803	TRS	2	0
3	D	802	TRS	2	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	694/705 (98%)	-0.33	5 (0%) 89 91	10, 21, 41, 73	0
1	B	694/705 (98%)	-0.33	8 (1%) 81 85	10, 19, 43, 79	0
1	C	696/705 (98%)	-0.39	7 (1%) 84 88	11, 20, 37, 87	0
1	D	693/705 (98%)	-0.40	11 (1%) 74 81	12, 20, 43, 79	0
All	All	2777/2820 (98%)	-0.36	31 (1%) 82 86	10, 20, 41, 87	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	82	PRO	4.8
1	C	2	ALA	4.6
1	C	65	ALA	4.5
1	A	705	GLU	4.4
1	D	83	GLN	4.2
1	C	66	GLY	4.2
1	D	64	THR	4.1
1	D	81	ASN	4.0
1	C	64	THR	3.9
1	B	2	ALA	3.4
1	B	83	GLN	3.4
1	D	85	PRO	3.1
1	D	651	TRP	3.0
1	A	83	GLN	3.0
1	B	705	GLU	2.9
1	A	300	TRP	2.9
1	B	81	ASN	2.8
1	D	704	GLY	2.7
1	D	80	PRO	2.6
1	D	705	GLU	2.6
1	B	346	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	348	VAL	2.4
1	B	82	PRO	2.4
1	A	85	PRO	2.3
1	D	526	ARG	2.2
1	C	74	ASP	2.2
1	C	347	THR	2.1
1	B	3	THR	2.1
1	D	300	TRP	2.1
1	C	346	LEU	2.0
1	A	472	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	TRS	D	802	8/8	0.80	0.24	5.06	54,65,70,76	0
3	TRS	C	802	8/8	0.78	0.20	4.05	28,31,46,55	0
3	TRS	A	803	8/8	0.90	0.21	1.79	21,25,27,34	0
3	TRS	C	803	8/8	0.90	0.20	1.31	20,25,28,30	0
3	TRS	A	802	8/8	0.75	0.16	1.28	33,36,41,50	0
3	TRS	B	802	8/8	0.96	0.18	0.74	18,23,26,26	0
3	TRS	D	803	8/8	0.94	0.14	0.42	18,20,21,23	0
2	CA	C	801	1/1	0.99	0.06	-0.98	19,19,19,19	0
2	CA	A	801	1/1	0.99	0.04	-2.16	22,22,22,22	0
2	CA	B	801	1/1	0.99	0.06	-2.58	13,13,13,13	0
2	CA	D	801	1/1	0.99	0.04	-3.01	17,17,17,17	0
2	CA	C	804	1/1	0.95	0.06	-	53,53,53,53	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.