



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 25, 2020 – 04:38 PM BST

PDB ID : 4FNP
Title : Crystal structure of GH36 alpha-galactosidase AgaA A355E from *Geobacillus stearothermophilus*
Authors : Merceron, R.; Foucault, M.; Haser, R.; Mattes, R.; Watzlawick, H.; Gouet, P.
Deposited on : 2012-06-20
Resolution : 2.80 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.13
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13

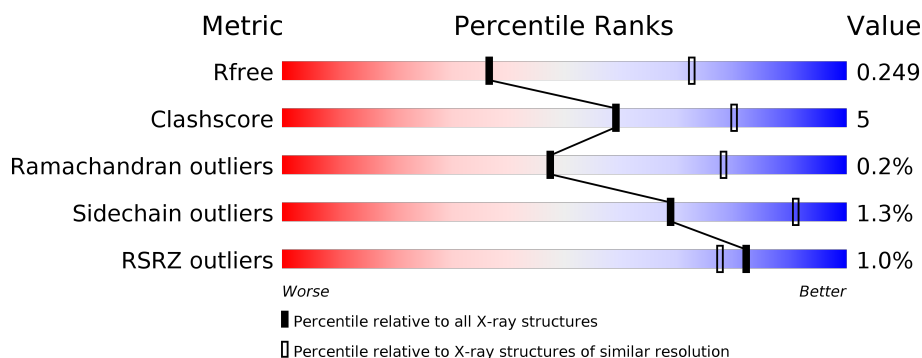
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 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	729	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 10%, orange 10%, orange 86%, yellow 86%, yellow 98%, green 98%, green 100%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 86% 12% </div> </div>
1	B	729	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 10%, orange 10%, orange 85%, yellow 85%, yellow 93%, green 93%, green 100%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 85% 13% </div> </div>
1	C	729	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 10%, orange 10%, orange 85%, yellow 85%, yellow 98%, green 98%, green 100%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 85% 12% </div> </div>
1	D	729	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 10%, orange 10%, orange 85%, yellow 85%, yellow 93%, green 93%, green 100%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 85% 13% </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 23279 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 Alpha-galactosidase AgaA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	718	Total	C	N	O	S	0	0	0
			5787	3673	1020	1072	22			
1	B	718	Total	C	N	O	S	0	0	0
			5787	3673	1020	1072	22			
1	C	718	Total	C	N	O	S	0	0	0
			5787	3673	1020	1072	22			
1	D	718	Total	C	N	O	S	0	0	0
			5787	3673	1020	1072	22			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	355	GLU	ALA	ENGINEERED MUTATION	UNP Q9ALJ4
A	518	LEU	PHE	ENGINEERED MUTATION	UNP Q9ALJ4
A	704	VAL	MET	ENGINEERED MUTATION	UNP Q9ALJ4
B	355	GLU	ALA	ENGINEERED MUTATION	UNP Q9ALJ4
B	518	LEU	PHE	ENGINEERED MUTATION	UNP Q9ALJ4
B	704	VAL	MET	ENGINEERED MUTATION	UNP Q9ALJ4
C	355	GLU	ALA	ENGINEERED MUTATION	UNP Q9ALJ4
C	518	LEU	PHE	ENGINEERED MUTATION	UNP Q9ALJ4
C	704	VAL	MET	ENGINEERED MUTATION	UNP Q9ALJ4
D	355	GLU	ALA	ENGINEERED MUTATION	UNP Q9ALJ4
D	518	LEU	PHE	ENGINEERED MUTATION	UNP Q9ALJ4
D	704	VAL	MET	ENGINEERED MUTATION	UNP Q9ALJ4

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

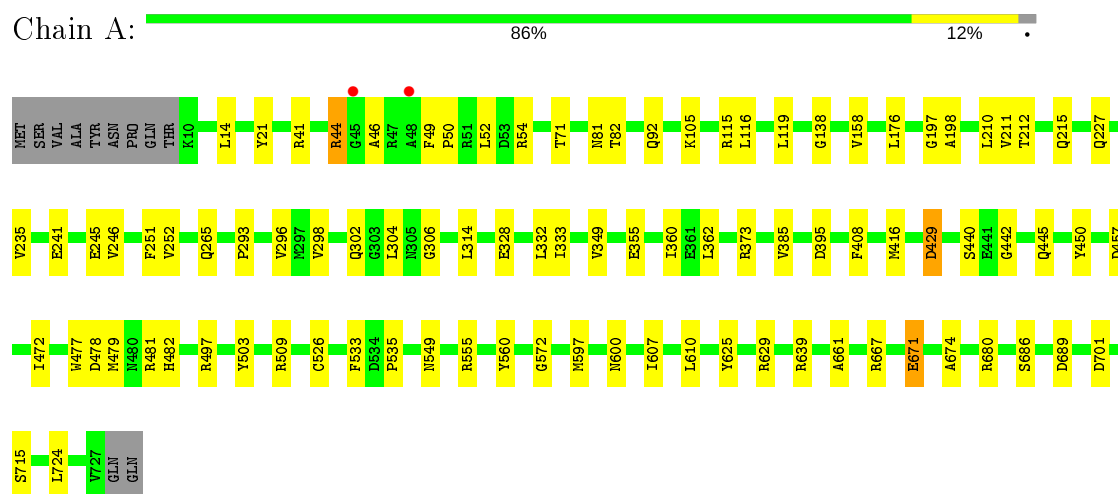
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	22	Total 22	O 22	0	0
3	B	13	Total 13	O 13	0	0
3	C	12	Total 12	O 12	0	0
3	D	24	Total 24	O 24	0	0

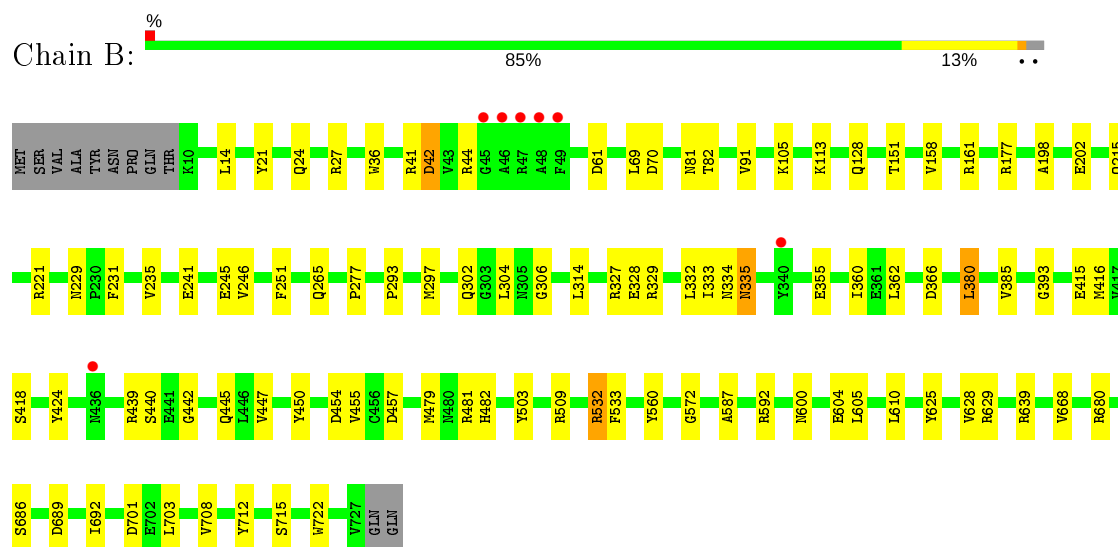
3 Residue-property plots [i](#)

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.

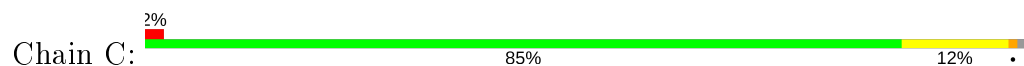
• Molecule 1: Alpha-galactosidase AgaA

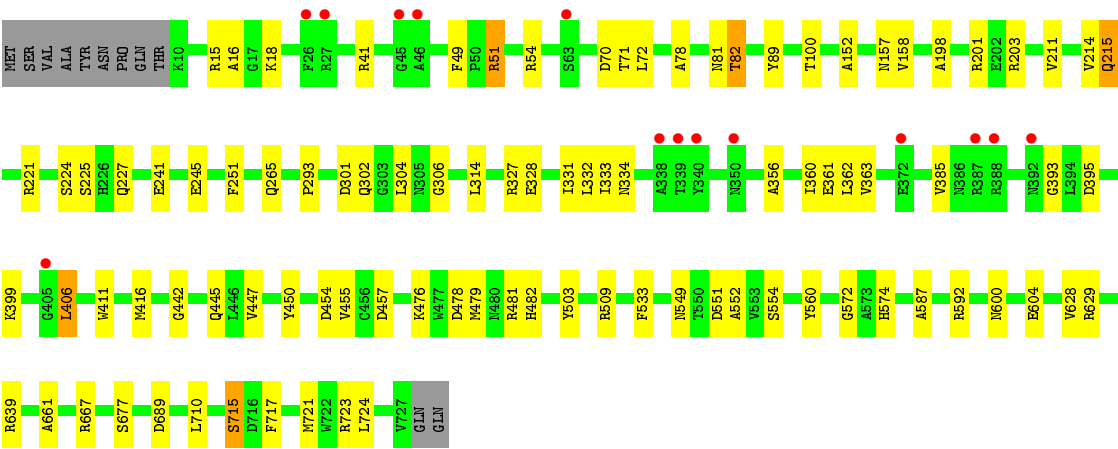


• Molecule 1: Alpha-galactosidase AgaA

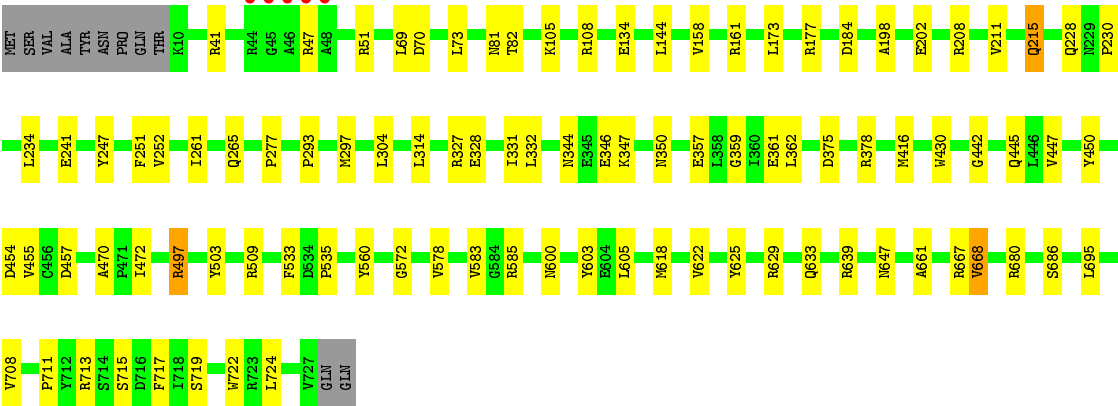
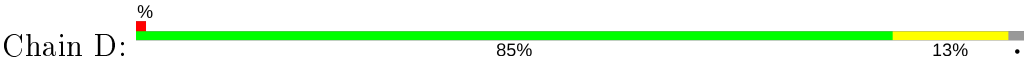


• Molecule 1: Alpha-galactosidase AgaA





● Molecule 1: Alpha-galactosidase AgaA



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	150.50Å 150.50Å 234.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.83 – 2.80 19.83 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.6 (19.83-2.80) 99.6 (19.83-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.27 (at 2.79Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.194 , 0.247 0.196 , 0.249	Depositor DCC
R_{free} test set	3761 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	20.9	Xtriage
Anisotropy	0.105	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 35.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.031 for -h,-k,l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	23279	wwPDB-VP
Average B, all atoms (Å ²)	29.0	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.*

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: SO4

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	0.22	0/5933	0.40	0/8045
1	B	0.22	0/5933	0.40	0/8045
1	C	0.22	0/5933	0.40	0/8045
1	D	0.22	0/5933	0.40	0/8045
All	All	0.22	0/23732	0.40	0/32180

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	5787	0	5622	58	0
1	B	5787	0	5622	59	0
1	C	5787	0	5622	57	0
1	D	5787	0	5622	56	0
2	A	15	0	0	1	0
2	B	15	0	0	0	0
2	C	15	0	0	1	0
2	D	15	0	0	1	0
3	A	22	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	13	0	0	0	0
3	C	12	0	0	1	0
3	D	24	0	0	0	0
All	All	23279	0	22488	213	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 (213) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:328:GLU:HB3	1:D:629:ARG:HD3	1.63	0.81
1:C:331:ILE:HG13	1:C:629:ARG:HH21	1.54	0.72
1:D:331:ILE:HG13	1:D:629:ARG:HH21	1.55	0.71
1:D:572:GLY:HA2	1:D:600:ASN:HB3	1.74	0.70
1:A:210:LEU:HA	1:A:215:GLN:HE22	1.57	0.69
1:A:41:ARG:NH2	1:B:689:ASP:OD1	2.25	0.69
1:C:667:ARG:NH2	1:C:715:SER:O	2.25	0.69
1:B:587:ALA:O	1:B:592:ARG:NH1	2.25	0.69
1:C:41:ARG:NH1	1:D:686:SER:O	2.23	0.68
1:A:41:ARG:NH1	1:B:686:SER:O	2.25	0.68
1:A:115:ARG:NH1	1:A:119:LEU:O	2.27	0.68
1:A:572:GLY:HA2	1:A:600:ASN:HB3	1.77	0.67
1:C:587:ALA:O	1:C:592:ARG:NH1	2.28	0.67
1:A:115:ARG:NH1	1:A:116:LEU:O	2.29	0.66
1:B:572:GLY:HA2	1:B:600:ASN:HB3	1.78	0.66
1:A:481:ARG:NH1	1:A:482:HIS:O	2.30	0.66
1:D:47:ARG:HH21	1:D:51:ARG:HH11	1.43	0.65
1:A:457:ASP:OD1	1:A:509:ARG:NH2	2.27	0.65
1:C:18:LYS:HE2	1:C:157:ASN:HD21	1.63	0.64
1:C:481:ARG:NH1	1:C:482:HIS:O	2.31	0.63
1:C:81:ASN:ND2	1:C:100:THR:O	2.32	0.62
1:B:235:VAL:HG12	1:B:246:VAL:HG13	1.84	0.60
1:B:265:GLN:HB2	1:C:198:ALA:HB2	1.84	0.59
1:C:572:GLY:HA2	1:C:600:ASN:HB3	1.84	0.59
1:A:667:ARG:NH2	1:A:715:SER:O	2.34	0.59
1:A:14:LEU:HB2	1:A:21:TYR:HB3	1.84	0.59
1:C:328:GLU:HB3	1:C:629:ARG:HD3	1.85	0.59
1:B:27:ARG:NH1	1:B:61:ASP:OD2	2.36	0.58
1:A:440:SER:O	1:A:481:ARG:NH2	2.36	0.58
1:B:424:TYR:OH	1:B:439:ARG:NH1	2.35	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:457:ASP:OD1	1:B:509:ARG:NH2	2.38	0.57
1:C:361:GLU:OE1	1:C:629:ARG:NH2	2.37	0.57
1:C:689:ASP:OD1	1:D:41:ARG:NH2	2.36	0.57
1:C:395:ASP:OD1	1:C:399:LYS:NZ	2.38	0.56
1:B:158:VAL:HG21	1:B:314:LEU:HD22	1.87	0.56
1:D:161:ARG:HD3	1:D:297:MET:HE1	1.88	0.56
1:A:478:ASP:OD1	1:A:479:MET:N	2.38	0.56
1:B:625:TYR:CE2	1:B:629:ARG:HG3	2.41	0.56
1:C:385:VAL:HG11	1:C:393:GLY:HA2	1.88	0.56
1:A:333:ILE:HG22	1:A:360:ILE:HG21	1.88	0.55
1:A:686:SER:O	1:B:41:ARG:NH1	2.39	0.55
1:C:333:ILE:HG22	1:C:360:ILE:HG21	1.88	0.55
1:B:415:GLU:OE1	1:B:532:ARG:NH2	2.40	0.54
1:C:356:ALA:HB2	1:C:406:LEU:HD11	1.88	0.54
1:C:628:VAL:HG12	1:C:723:ARG:HH11	1.73	0.54
1:B:302:GLN:HG2	1:B:306:GLY:HA3	1.89	0.54
1:A:44:ARG:HE	1:A:46:ALA:HB3	1.73	0.53
1:A:671:GLU:HB3	1:A:674:ALA:HB2	1.90	0.53
1:C:245:GLU:OE1	1:D:680:ARG:NH2	2.30	0.53
1:A:235:VAL:HG12	1:A:246:VAL:HG13	1.90	0.53
1:A:597:MET:HE3	1:A:625:TYR:HB2	1.90	0.53
1:C:211:VAL:N	1:C:215:GLN:OE1	2.29	0.53
1:C:363:VAL:HG23	1:C:406:LEU:HD22	1.90	0.53
1:B:177:ARG:HG3	1:B:277:PRO:HG3	1.91	0.52
1:B:708:VAL:HG21	1:B:722:TRP:HZ3	1.73	0.52
1:A:245:GLU:OE1	1:B:680:ARG:NH2	2.38	0.52
1:C:158:VAL:HG21	1:C:314:LEU:HD22	1.91	0.52
1:B:334:ASN:HB3	1:B:604:GLU:HG2	1.90	0.52
1:C:16:ALA:HB2	1:C:152:ALA:HB1	1.91	0.52
1:A:158:VAL:HG21	1:A:314:LEU:HD22	1.91	0.52
1:A:395:ASP:OD1	1:A:395:ASP:N	2.39	0.52
1:D:457:ASP:OD1	1:D:509:ARG:NH2	2.34	0.52
1:B:355:GLU:HB3	1:B:360:ILE:HD12	1.92	0.52
1:B:42:ASP:HB3	1:B:44:ARG:HE	1.76	0.51
1:C:677:SER:HB2	1:C:710:LEU:HB2	1.93	0.51
1:A:560:TYR:CZ	1:A:639:ARG:HD2	2.45	0.51
1:A:686:SER:HB2	1:A:701:ASP:HB3	1.92	0.51
1:B:161:ARG:HD3	1:B:297:MET:HE1	1.93	0.51
1:D:144:LEU:HD12	1:D:173:LEU:HD13	1.91	0.51
1:B:416:MET:HE1	1:B:447:VAL:HG22	1.92	0.50
1:C:241:GLU:HA	1:C:304:LEU:HB2	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:661:ALA:HB3	1:A:724:LEU:HB2	1.94	0.50
1:B:241:GLU:HA	1:B:304:LEU:HB2	1.94	0.50
1:B:14:LEU:HB2	1:B:21:TYR:HB3	1.93	0.50
1:C:454:ASP:OD1	1:C:455:VAL:N	2.44	0.50
1:A:302:GLN:HB2	1:A:306:GLY:HA3	1.93	0.49
1:A:429:ASP:N	1:A:429:ASP:OD1	2.41	0.49
1:A:625:TYR:CE2	1:A:629:ARG:HG3	2.47	0.49
1:C:551:ASP:HB3	1:C:554:SER:HB2	1.94	0.49
1:C:450:TYR:HB2	1:C:503:TYR:CD1	2.47	0.49
1:C:49:PHE:CZ	1:D:711:PRO:HG3	2.48	0.49
1:D:252:VAL:HA	1:D:535:PRO:HB2	1.94	0.49
1:A:265:GLN:HB2	1:D:198:ALA:HB2	1.95	0.49
1:A:328:GLU:HB3	1:A:629:ARG:HD3	1.94	0.48
1:C:15:ARG:NH2	1:C:41:ARG:O	2.36	0.48
1:D:344:ASN:HD21	1:D:347:LYS:HG3	1.78	0.48
1:B:198:ALA:HB2	1:C:265:GLN:HB2	1.95	0.48
1:D:416:MET:HE1	1:D:447:VAL:HG22	1.96	0.48
1:D:375:ASP:OD2	1:D:378:ARG:NH2	2.44	0.48
1:B:625:TYR:HA	1:B:628:VAL:HG22	1.96	0.48
1:B:335:ASN:ND2	1:B:355:GLU:OE2	2.47	0.47
1:A:385:VAL:HG23	3:A:914:HOH:O	2.14	0.47
1:D:215:GLN:NE2	1:D:261:ILE:O	2.47	0.47
1:A:198:ALA:HB2	1:D:265:GLN:HB2	1.97	0.47
1:B:332:LEU:HB3	1:B:362:LEU:HD23	1.96	0.47
1:C:81:ASN:O	1:C:82:THR:OG1	2.23	0.47
1:C:552:ALA:HB2	1:C:592:ARG:HG2	1.97	0.46
1:A:54:ARG:HG3	1:A:71:THR:HG21	1.95	0.46
1:B:385:VAL:HG11	1:B:393:GLY:HA2	1.96	0.46
1:A:549:ASN:ND2	2:C:801:SO4:O2	2.48	0.46
1:D:105:LYS:HD2	2:D:802:SO4:O1	2.16	0.46
1:A:450:TYR:HB2	1:A:503:TYR:CD1	2.50	0.46
1:A:680:ARG:NH2	1:B:245:GLU:OE1	2.34	0.46
1:C:332:LEU:HB3	1:C:362:LEU:HD23	1.97	0.46
1:B:36:TRP:CH2	1:B:297:MET:HB3	2.51	0.46
1:B:69:LEU:HD12	1:B:81:ASN:O	2.16	0.46
1:C:334:ASN:HB3	1:C:604:GLU:HG2	1.97	0.46
1:D:73:LEU:HD22	1:D:184:ASP:HB3	1.98	0.46
1:D:603:TYR:CZ	1:D:618:MET:HG3	2.51	0.46
1:D:560:TYR:CZ	1:D:639:ARG:HD2	2.51	0.46
1:D:359:GLY:HA3	1:D:622:VAL:HG11	1.97	0.45
1:C:203:ARG:NH2	3:C:901:HOH:O	2.48	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:442:GLY:H	1:C:481:ARG:HE	1.65	0.45
1:D:108:ARG:NH1	1:D:134:GLU:OE1	2.49	0.45
1:A:241:GLU:HA	1:A:304:LEU:HB2	1.99	0.45
1:A:355:GLU:HG3	1:A:607:ILE:HD13	1.98	0.45
1:B:128:GLN:N	1:B:128:GLN:OE1	2.45	0.45
1:B:479:MET:SD	1:B:532:ARG:NH2	2.90	0.45
1:D:647:ASN:OD1	1:D:668:VAL:HG13	2.16	0.45
1:A:212:THR:O	1:D:228:GLN:NE2	2.47	0.45
1:B:481:ARG:NH1	1:B:482:HIS:O	2.50	0.45
1:D:158:VAL:HG21	1:D:314:LEU:HD22	1.98	0.45
1:D:708:VAL:HG11	1:D:722:TRP:HZ3	1.81	0.45
1:B:328:GLU:HB3	1:B:629:ARG:CD	2.47	0.45
1:D:332:LEU:HB3	1:D:362:LEU:HD23	1.99	0.45
1:D:450:TYR:HB2	1:D:503:TYR:CD1	2.52	0.45
1:B:442:GLY:O	1:B:445:GLN:HG2	2.17	0.44
1:C:18:LYS:NZ	1:C:301:ASP:OD1	2.34	0.44
1:D:470:ALA:O	1:D:472:ILE:N	2.51	0.44
1:C:560:TYR:CZ	1:C:639:ARG:HD2	2.53	0.44
1:D:661:ALA:HB3	1:D:724:LEU:HB2	1.98	0.44
1:B:560:TYR:CZ	1:B:639:ARG:HD2	2.52	0.44
1:D:625:TYR:OH	1:D:633:GLN:NE2	2.43	0.44
1:B:440:SER:O	1:B:481:ARG:NH2	2.50	0.44
1:C:302:GLN:HB2	1:C:306:GLY:HA3	2.00	0.44
1:D:605:LEU:HD21	1:D:618:MET:HE1	1.99	0.44
1:A:332:LEU:HB3	1:A:362:LEU:HD23	1.99	0.44
1:B:327:ARG:HA	1:B:327:ARG:HD2	1.85	0.44
1:C:221:ARG:HD3	1:C:482:HIS:CE1	2.53	0.44
1:C:251:PHE:CE1	1:C:293:PRO:HB2	2.53	0.44
1:D:442:GLY:O	1:D:445:GLN:HG2	2.17	0.44
1:A:408:PHE:HD2	1:A:472:ILE:HG12	1.83	0.43
1:D:234:LEU:HB2	1:D:247:TYR:HB2	2.00	0.43
1:B:380:LEU:HB2	1:B:418:SER:OG	2.19	0.43
1:B:70:ASP:HB2	1:B:82:THR:O	2.18	0.43
1:C:225:SER:HB2	1:C:227:GLN:OE1	2.19	0.43
1:C:51:ARG:HH12	1:C:72:LEU:HD13	1.83	0.43
1:A:49:PHE:HA	1:A:50:PRO:HD3	1.72	0.43
2:A:801:SO4:O1	1:C:549:ASN:ND2	2.52	0.43
1:D:454:ASP:OD1	1:D:455:VAL:N	2.51	0.43
1:C:54:ARG:HG3	1:C:71:THR:HG21	2.00	0.43
1:A:477:TRP:O	1:A:526:CYS:N	2.52	0.43
1:B:450:TYR:HB2	1:B:503:TYR:CD1	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:686:SER:HB2	1:B:701:ASP:HB3	2.00	0.43
1:C:411:TRP:HD1	1:C:476:LYS:HD3	1.84	0.43
1:D:241:GLU:HA	1:D:304:LEU:HB2	2.00	0.43
1:D:713:ARG:NH2	1:D:719:SER:O	2.52	0.43
1:A:607:ILE:HA	1:A:610:LEU:HG	2.01	0.43
1:B:701:ASP:OD1	1:B:701:ASP:N	2.51	0.43
1:A:92:GLN:HB2	1:A:176:LEU:HD11	2.01	0.42
1:D:230:PRO:HB2	1:D:251:PHE:HB3	2.01	0.42
1:C:478:ASP:OD1	1:C:479:MET:N	2.52	0.42
1:D:47:ARG:HA	1:D:47:ARG:HD3	1.90	0.42
1:A:211:VAL:HG12	1:D:208:ARG:HG3	2.01	0.42
1:A:555:ARG:HA	1:A:555:ARG:HD3	1.95	0.42
1:B:251:PHE:CE1	1:B:293:PRO:HB2	2.55	0.42
1:A:158:VAL:HG22	1:A:298:VAL:HG22	2.00	0.42
1:C:661:ALA:HB3	1:C:724:LEU:HB2	2.01	0.42
1:D:327:ARG:NH1	1:D:328:GLU:H	2.18	0.42
1:A:52:LEU:HD21	1:B:712:TYR:CD1	2.55	0.42
1:B:69:LEU:HD12	1:B:81:ASN:HB3	2.01	0.42
1:B:105:LYS:HE2	1:B:105:LYS:HB2	1.77	0.42
1:B:454:ASP:OD1	1:B:455:VAL:N	2.51	0.42
1:C:442:GLY:O	1:C:445:GLN:HG2	2.19	0.42
1:C:70:ASP:HB2	1:C:82:THR:O	2.20	0.42
1:D:177:ARG:HG3	1:D:277:PRO:HG3	2.01	0.42
1:D:81:ASN:O	1:D:82:THR:OG1	2.25	0.42
1:A:252:VAL:HA	1:A:535:PRO:HB2	2.01	0.42
1:C:78:ALA:HB2	1:C:89:TYR:CE2	2.55	0.42
1:D:497:ARG:NE	1:D:497:ARG:HA	2.34	0.42
1:D:625:TYR:CE2	1:D:629:ARG:HG3	2.55	0.42
1:D:69:LEU:HD12	1:D:81:ASN:O	2.20	0.42
1:A:105:LYS:HD2	1:A:138:GLY:HA3	2.02	0.41
1:B:202:GLU:OE1	1:B:202:GLU:N	2.52	0.41
1:C:457:ASP:OD1	1:C:509:ARG:NH2	2.34	0.41
1:D:361:GLU:OE1	1:D:629:ARG:NH2	2.53	0.41
1:D:70:ASP:HB2	1:D:82:THR:O	2.20	0.41
1:A:442:GLY:O	1:A:445:GLN:HG2	2.20	0.41
1:D:211:VAL:N	1:D:215:GLN:OE1	2.33	0.41
1:A:197:GLY:O	1:A:227:GLN:HA	2.20	0.41
1:D:202:GLU:N	1:D:202:GLU:OE1	2.53	0.41
1:A:158:VAL:HG13	1:A:296:VAL:HG13	2.01	0.41
1:B:229:ASN:O	1:B:231:PHE:N	2.51	0.41
1:B:605:LEU:HD11	1:B:610:LEU:HD11	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:329:ARG:HG3	1:B:600:ASN:HD21	1.85	0.41
1:D:346:GLU:O	1:D:350:ASN:ND2	2.54	0.41
1:D:583:VAL:HG23	1:D:585:ARG:H	1.86	0.41
1:A:81:ASN:O	1:A:82:THR:OG1	2.28	0.41
1:B:221:ARG:HD3	1:B:482:HIS:ND1	2.35	0.41
1:B:333:ILE:HB	1:B:360:ILE:HD13	2.02	0.41
1:C:331:ILE:HG13	1:C:629:ARG:NH2	2.30	0.41
1:A:328:GLU:HB3	1:A:629:ARG:CD	2.50	0.41
1:C:416:MET:HE1	1:C:447:VAL:HG22	2.03	0.41
1:D:251:PHE:CE1	1:D:293:PRO:HB2	2.55	0.41
1:B:113:LYS:HD2	1:B:151:THR:HG21	2.03	0.41
1:C:332:LEU:HD11	1:C:574:HIS:CE1	2.57	0.40
1:B:692:ILE:HD11	1:B:703:LEU:HD21	2.02	0.40
1:C:224:SER:O	1:C:225:SER:OG	2.35	0.40
1:D:668:VAL:HG23	1:D:717:PHE:CD2	2.57	0.40
1:A:251:PHE:CE1	1:A:293:PRO:HB2	2.57	0.40
1:A:416:MET:HG3	1:A:445:GLN:NE2	2.36	0.40
1:A:689:ASP:OD1	1:B:41:ARG:NH2	2.55	0.40
1:C:327:ARG:HD2	1:C:327:ARG:HA	1.73	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	716/729 (98%)	683 (95%)	31 (4%)	2 (0%)	41	72
1	B	716/729 (98%)	688 (96%)	27 (4%)	1 (0%)	51	81
1	C	716/729 (98%)	690 (96%)	24 (3%)	2 (0%)	41	72
1	D	716/729 (98%)	686 (96%)	29 (4%)	1 (0%)	51	81
All	All	2864/2916 (98%)	2747 (96%)	111 (4%)	6 (0%)	47	78

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	373	ARG
1	B	715	SER
1	C	82	THR
1	D	715	SER
1	A	44	ARG
1	C	715	SER

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	609/619 (98%)	604 (99%)	5 (1%)	81	94
1	B	609/619 (98%)	599 (98%)	10 (2%)	62	88
1	C	609/619 (98%)	601 (99%)	8 (1%)	69	91
1	D	609/619 (98%)	600 (98%)	9 (2%)	65	89
All	All	2436/2476 (98%)	2404 (99%)	32 (1%)	69	91

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	349	VAL
1	A	429	ASP
1	A	497	ARG
1	A	533	PHE
1	A	671	GLU
1	B	24	GLN
1	B	42	ASP
1	B	91	VAL
1	B	215	GLN
1	B	335	ASN
1	B	366	ASP
1	B	380	LEU
1	B	532	ARG
1	B	533	PHE

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Mol	Chain	Res	Type
1	B	668	VAL
1	C	51	ARG
1	C	201	ARG
1	C	214	VAL
1	C	215	GLN
1	C	406	LEU
1	C	533	PHE
1	C	717	PHE
1	C	721	MET
1	D	215	GLN
1	D	357	GLU
1	D	430	TRP
1	D	497	ARG
1	D	533	PHE
1	D	578	VAL
1	D	667	ARG
1	D	668	VAL
1	D	695	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	215	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

12 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	SO4	C	803	-	4,4,4	0.15	0	6,6,6	0.09	0
2	SO4	B	803	-	4,4,4	0.13	0	6,6,6	0.12	0
2	SO4	C	801	-	4,4,4	0.19	0	6,6,6	0.13	0
2	SO4	D	803	-	4,4,4	0.14	0	6,6,6	0.07	0
2	SO4	A	802	-	4,4,4	0.13	0	6,6,6	0.14	0
2	SO4	C	802	-	4,4,4	0.23	0	6,6,6	0.10	0
2	SO4	A	801	-	4,4,4	0.18	0	6,6,6	0.16	0
2	SO4	D	802	-	4,4,4	0.16	0	6,6,6	0.13	0
2	SO4	B	801	-	4,4,4	0.22	0	6,6,6	0.14	0
2	SO4	A	803	-	4,4,4	0.15	0	6,6,6	0.13	0
2	SO4	B	802	-	4,4,4	0.17	0	6,6,6	0.10	0
2	SO4	D	801	-	4,4,4	0.17	0	6,6,6	0.12	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	801	SO4	1	0
2	A	801	SO4	1	0
2	D	802	SO4	1	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	718/729 (98%)	-0.47	2 (0%) 94 93	15, 25, 40, 58	0
1	B	718/729 (98%)	-0.37	7 (0%) 82 77	17, 28, 48, 93	0
1	C	718/729 (98%)	-0.25	14 (1%) 66 59	20, 32, 58, 79	0
1	D	718/729 (98%)	-0.45	5 (0%) 87 84	16, 24, 39, 77	0
All	All	2872/2916 (98%)	-0.39	28 (0%) 82 77	15, 27, 47, 93	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	48	ALA	6.3
1	B	49	PHE	6.1
1	D	45	GLY	5.7
1	B	47	ARG	4.8
1	B	46	ALA	4.8
1	C	388	ARG	4.0
1	B	45	GLY	4.0
1	C	63	SER	3.8
1	D	46	ALA	3.7
1	C	372	GLU	3.4
1	C	340	TYR	3.4
1	A	48	ALA	3.3
1	C	46	ALA	3.3
1	C	338	ALA	3.1
1	C	27	ARG	3.0
1	D	48	ALA	2.9
1	C	339	THR	2.5
1	D	47	ARG	2.4
1	C	350	ASN	2.4
1	C	26	PHE	2.4
1	B	340	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	392	ASN	2.2
1	A	45	GLY	2.1
1	C	405	GLY	2.1
1	B	436	ASN	2.1
1	C	387	ARG	2.0
1	D	44	ARG	2.0
1	C	45	GLY	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 monosaccharides 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(Å ²)	Q<0.9
2	SO4	C	803	5/5	0.93	0.25	49,52,53,53	0
2	SO4	B	803	5/5	0.97	0.21	31,32,42,48	0
2	SO4	D	803	5/5	0.97	0.17	21,25,44,44	0
2	SO4	D	802	5/5	0.97	0.26	14,15,22,29	0
2	SO4	C	802	5/5	0.98	0.18	16,19,30,39	0
2	SO4	A	803	5/5	0.98	0.18	14,22,30,39	0
2	SO4	B	802	5/5	0.98	0.24	18,21,26,40	0
2	SO4	A	802	5/5	0.99	0.21	18,20,20,25	0
2	SO4	B	801	5/5	0.99	0.31	20,27,30,34	0
2	SO4	C	801	5/5	0.99	0.28	20,22,23,27	0
2	SO4	A	801	5/5	0.99	0.28	20,22,31,45	0
2	SO4	D	801	5/5	0.99	0.26	19,19,23,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.