



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 05:46 PM UTC

PDB ID : 5MRJ / pdb_00005mrj
Title : Crystal structure of Endo-1,4-beta-xylanase-like protein from *Acremonium chrysogenum*
Authors : Gabdulkhakov, A.; Tishchenko, S.; Lisov, A.; Leontievsky, A.
Deposited on : 2016-12-23
Resolution : 2.70 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

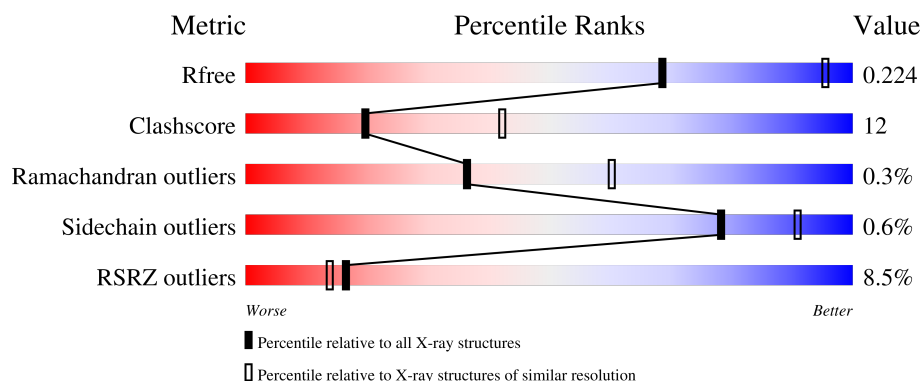
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	396	2% 67% 11% 22%
1	B	396	12% 53% 24% 23%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4817 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-xylanase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	307	Total	C	N	O	S	0	0	0
			2386	1505	427	446	8			
1	B	305	Total	C	N	O	S	0	0	0
			2375	1499	425	443	8			

- Molecule 2 is SULFATE ION (CCD ID: 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	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	27	Total 27	O 27	0	0
3	B	14	Total 14	O 14	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	88.21Å 124.83Å 165.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.81 – 2.70 46.81 – 2.70	Depositor EDS
% Data completeness (in resolution range)	93.3 (46.81-2.70) 94.3 (46.81-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.88 (at 2.65Å)	Xtriage
Refinement program	PHENIX (1.11.1 _2575: ???)	Depositor
R, R_{free}	0.221 , 0.285 0.235 , 0.224	Depositor DCC
R_{free} test set	1264 reflections (4.63%)	wwPDB-VP
Wilson B-factor (Å ²)	37.2	Xtriage
Anisotropy	1.127	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 35.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4817	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.49% 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: 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	0.38	0/2442	0.58	0/3322
1	B	0.44	1/2431 (0.0%)	0.66	0/3307
All	All	0.41	1/4873 (0.0%)	0.62	0/6629

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	375	ASN	C-O	-5.04	1.21	1.23

There are no bond angle outliers.

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	2386	0	2309	24	0
1	B	2375	0	2299	84	0
2	A	10	0	0	1	0
2	B	5	0	0	0	0
3	A	27	0	0	0	0
3	B	14	0	0	0	0
All	All	4817	0	4608	108	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 12.

All (108) 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:108:PHE:CE2	1:B:359:ILE:HD11	2.00	0.96
1:A:301:LEU:HD13	1:A:310:ALA:HB2	1.62	0.81
1:B:229:LEU:HD11	1:B:240:GLU:HB3	1.63	0.81
1:B:139:MET:HA	1:B:144:ILE:CD1	2.11	0.79
1:B:339:ALA:HB3	1:B:342:ASP:HB2	1.64	0.79
1:B:318:ARG:HH22	1:B:352:GLN:HE21	1.32	0.78
1:B:103:LYS:HD2	1:B:396:LEU:HD12	1.66	0.76
1:A:310:ALA:HB3	1:A:342:ASP:OD1	1.87	0.74
1:B:261:ILE:HG12	1:B:278:MET:HE1	1.72	0.72
1:B:139:MET:HA	1:B:144:ILE:HD11	1.74	0.69
1:B:186:LYS:HA	1:B:233:VAL:HG21	1.74	0.68
1:B:94:GLY:HA2	1:B:99:LYS:HZ3	1.59	0.68
1:B:365:SER:OG	1:B:380:ASP:O	2.12	0.67
1:B:331:THR:O	1:B:360:THR:HG23	1.94	0.67
1:B:251:ARG:HH22	1:B:293:ASP:CG	2.04	0.66
1:B:139:MET:HA	1:B:144:ILE:HD13	1.76	0.66
1:A:276:ARG:O	1:A:281:LYS:NZ	2.28	0.65
1:B:350:CYS:HA	1:B:353:ILE:HG22	1.80	0.64
1:B:108:PHE:CZ	1:B:359:ILE:HD11	2.32	0.64
1:B:251:ARG:NH2	1:B:293:ASP:OD2	2.32	0.63
1:A:310:ALA:HA	1:A:313:VAL:HG13	1.82	0.62
1:A:290:VAL:HG12	1:A:292:ILE:HG13	1.82	0.62
1:B:223:PHE:HE2	1:B:278:MET:HE3	1.65	0.61
1:B:343:TYR:HB3	1:B:392:LEU:HD11	1.83	0.60
1:B:92:GLY:C	1:B:94:GLY:H	2.09	0.60
1:B:225:GLU:HB3	1:B:271:TYR:OH	2.01	0.60
1:A:301:LEU:HD12	1:A:335:ILE:HG12	1.84	0.59
1:B:290:VAL:HG12	1:B:292:ILE:HD12	1.85	0.58
1:B:318:ARG:NH2	1:B:352:GLN:HE21	1.99	0.58
1:B:349:ALA:O	1:B:353:ILE:HG22	2.04	0.58
1:B:395:ALA:O	1:B:396:LEU:HD13	2.03	0.58
1:B:351:LEU:HD21	1:B:396:LEU:HD11	1.86	0.58
1:B:247:PHE:HB3	1:B:291:PRO:HD2	1.86	0.57
1:B:301:LEU:HD12	1:B:335:ILE:HG12	1.86	0.56
1:B:247:PHE:CE1	1:B:292:ILE:HG13	2.41	0.56
1:B:330:VAL:HG13	1:B:359:ILE:HB	1.87	0.56
1:B:334:ASP:HB2	1:B:376:PRO:HB2	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:325:VAL:O	1:B:355:LYS:NZ	2.29	0.55
1:A:188:ILE:HD13	1:A:193:THR:HG22	1.89	0.55
1:A:97:HIS:CE1	1:A:131:GLY:HA3	2.42	0.55
1:A:293:ASP:OD1	1:A:326:LYS:NZ	2.40	0.55
1:B:310:ALA:HB3	1:B:342:ASP:OD2	2.07	0.54
1:B:162:PHE:O	1:B:166:ASN:ND2	2.37	0.54
1:B:130:PHE:O	1:B:168:LYS:HE3	2.07	0.54
1:B:301:LEU:HD13	1:B:310:ALA:HB2	1.90	0.54
1:B:386:LYS:HD3	1:B:387:ALA:N	2.23	0.54
1:B:228:ASN:ND2	1:B:229:LEU:H	2.06	0.53
1:B:386:LYS:HD3	1:B:388:ALA:H	1.73	0.53
1:B:241:ASP:O	1:B:245:ILE:HG13	2.08	0.53
1:B:295:ILE:HG21	1:B:320:LEU:HD22	1.91	0.52
1:B:105:LYS:NZ	1:B:351:LEU:HA	2.24	0.52
1:B:103:LYS:CD	1:B:396:LEU:HD12	2.38	0.51
1:A:142:ASP:HA	1:A:182:PRO:HG3	1.93	0.51
1:B:94:GLY:HA2	1:B:99:LYS:NZ	2.24	0.51
1:B:96:LEU:HB2	1:B:389:TYR:HE1	1.76	0.51
1:B:96:LEU:HD11	1:B:108:PHE:CD2	2.46	0.51
1:B:282:VAL:HG21	1:B:295:ILE:HD11	1.94	0.50
1:B:350:CYS:HA	1:B:353:ILE:CG2	2.42	0.49
1:A:330:VAL:HG22	1:A:359:ILE:HG12	1.95	0.49
1:B:248:ARG:HH12	1:B:290:VAL:HA	1.78	0.48
1:B:103:LYS:HD2	1:B:396:LEU:CD1	2.39	0.48
1:B:191:ARG:CZ	1:B:238:LEU:HD23	2.44	0.48
1:B:96:LEU:HB2	1:B:389:TYR:CE1	2.48	0.47
1:B:96:LEU:HD11	1:B:108:PHE:CE2	2.49	0.47
1:B:221:GLU:HA	1:B:263:ASP:CG	2.39	0.47
1:B:266:LEU:N	1:B:267:ASP:OD1	2.47	0.47
1:B:290:VAL:HG12	1:B:292:ILE:CD1	2.43	0.47
1:A:110:THR:OG1	1:A:111:GLU:N	2.48	0.47
1:B:310:ALA:HA	1:B:313:VAL:HG23	1.97	0.47
1:B:353:ILE:HG23	1:B:356:CYS:HB3	1.96	0.46
1:B:186:LYS:HG3	1:B:233:VAL:HG21	1.97	0.46
1:A:135:HIS:NE2	1:A:159:VAL:HB	2.30	0.46
1:A:344:LEU:HD23	1:A:344:LEU:HA	1.76	0.46
1:A:334:ASP:HB2	1:A:376:PRO:HB2	1.98	0.45
1:B:318:ARG:HH22	1:B:352:GLN:NE2	2.08	0.45
1:B:359:ILE:O	1:B:359:ILE:HG13	2.17	0.45
1:A:152:SER:HB2	2:A:401:SO4:O2	2.16	0.45
1:B:228:ASN:CG	1:B:229:LEU:N	2.74	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:314:PRO:O	1:B:318:ARG:HG3	2.17	0.45
1:A:238:LEU:O	1:A:241:ASP:HB2	2.17	0.45
1:A:221:GLU:HG2	1:A:264:TYR:CE1	2.51	0.45
1:B:195:THR:HG22	1:B:245:ILE:HD12	1.97	0.44
1:B:331:THR:C	1:B:360:THR:HG23	2.42	0.44
1:B:194:LEU:HD23	1:B:238:LEU:HD12	1.99	0.44
1:A:139:MET:HE2	1:A:209:TYR:CE2	2.52	0.44
1:B:240:GLU:OE1	1:B:285:TRP:NE1	2.40	0.44
1:B:351:LEU:HA	1:B:351:LEU:HD12	1.79	0.44
1:B:97:HIS:CE1	1:B:107:PHE:HA	2.53	0.43
1:B:128:SER:OG	1:B:129:SER:N	2.51	0.43
1:B:386:LYS:CD	1:B:388:ALA:H	2.32	0.43
1:A:328:ILE:O	1:A:356:CYS:HA	2.19	0.43
1:B:96:LEU:HD21	1:B:392:LEU:HB3	2.01	0.42
1:B:186:LYS:HG3	1:B:233:VAL:CG2	2.50	0.42
1:B:110:THR:O	1:B:133:VAL:HA	2.19	0.42
1:B:261:ILE:CG1	1:B:278:MET:HE1	2.47	0.42
1:B:388:ALA:O	1:B:392:LEU:HD12	2.20	0.42
1:B:136:GLU:HG2	1:B:137:ASN:ND2	2.35	0.42
1:B:232:SER:O	1:B:236:ARG:HD3	2.19	0.42
1:B:330:VAL:HG12	1:B:359:ILE:HG22	2.02	0.42
1:A:173:HIS:HA	1:A:174:THR:HA	1.85	0.42
1:B:243:VAL:O	1:B:247:PHE:HD2	2.03	0.42
1:B:328:ILE:O	1:B:356:CYS:HA	2.20	0.41
1:B:122:LEU:HD23	1:B:122:LEU:C	2.46	0.41
1:A:301:LEU:HD12	1:A:335:ILE:CG1	2.49	0.41
1:B:109:GLY:HA2	1:B:130:PHE:HB3	2.03	0.41
1:B:299:CYS:HB3	1:B:333:LEU:HD12	2.02	0.41
1:A:336:ALA:CB	1:A:374:ASP:HB3	2.51	0.41
1:A:283:ASN:O	1:A:287:SER:OG	2.33	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	305/396 (77%)	294 (96%)	11 (4%)	0	100	100
1	B	303/396 (76%)	279 (92%)	22 (7%)	2 (1%)	18	41
All	All	608/792 (77%)	573 (94%)	33 (5%)	2 (0%)	36	60

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	93	ASN
1	B	304	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	249/320 (78%)	248 (100%)	1 (0%)	84	93
1	B	248/320 (78%)	246 (99%)	2 (1%)	73	88
All	All	497/640 (78%)	494 (99%)	3 (1%)	78	91

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	91	THR
1	B	179	SER
1	B	186	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	187	GLN
1	A	375	ASN
1	A	381	ASN

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Mol	Chain	Res	Type
1	B	120	GLN
1	B	137	ASN
1	B	228	ASN
1	B	308	ASN
1	B	352	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 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	A	402	-	4,4,4	0.89	0	6,6,6	1.64	3 (50%)
2	SO4	B	401	-	4,4,4	0.30	0	6,6,6	0.36	0
2	SO4	A	401	-	4,4,4	0.33	0	6,6,6	0.55	0

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	402	SO4	O4-S-O1	2.38	121.99	109.56
2	A	402	SO4	O3-S-O2	2.07	120.39	109.56
2	A	402	SO4	O2-S-O1	-2.02	94.84	109.06

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	SO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	307/396 (77%)	0.27	6 (1%) 65 62	18, 33, 48, 60	0
1	B	305/396 (77%)	1.01	46 (15%) 5 4	23, 41, 58, 77	0
All	All	612/792 (77%)	0.64	52 (8%) 16 14	18, 37, 56, 77	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	316	ALA	5.1
1	B	393	VAL	4.2
1	B	227	GLY	4.1
1	A	90	GLY	3.7
1	B	311	SER	3.5
1	B	277	GLY	3.4
1	B	308	ASN	3.4
1	B	353	ILE	3.3
1	B	349	ALA	3.2
1	B	313	VAL	3.1
1	A	277	GLY	3.1
1	B	189	ASN	3.1
1	B	396	LEU	3.1
1	B	354	GLU	2.9
1	A	94	GLY	2.8
1	B	96	LEU	2.8
1	B	320	LEU	2.8
1	B	276	ARG	2.8
1	B	309	PRO	2.7
1	B	304	PRO	2.7
1	B	310	ALA	2.6
1	A	391	ALA	2.6
1	B	312	GLN	2.5
1	B	224	SER	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	342	ASP	2.5
1	B	323	ALA	2.5
1	B	372	GLN	2.5
1	B	257	ALA	2.5
1	B	271	TYR	2.5
1	B	226	ASP	2.4
1	B	270	ASN	2.4
1	B	334	ASP	2.4
1	B	239	GLY	2.4
1	B	319	ALA	2.4
1	B	395	ALA	2.4
1	A	310	ALA	2.4
1	B	347	MET	2.3
1	B	386	LYS	2.3
1	B	294	GLY	2.3
1	B	373	GLY	2.3
1	B	275	THR	2.2
1	B	274	VAL	2.2
1	B	361	VAL	2.2
1	B	392	LEU	2.2
1	B	344	LEU	2.2
1	B	269	ALA	2.2
1	B	260	TYR	2.2
1	B	388	ALA	2.1
1	B	318	ARG	2.1
1	A	91	THR	2.1
1	B	273	LYS	2.1
1	B	268	ASN	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	SO4	A	401	5/5	0.82	0.14	36,37,51,69	0
2	SO4	A	402	5/5	0.83	0.13	57,57,65,93	0
2	SO4	B	401	5/5	0.84	0.16	39,41,55,59	0

6.5 Other polymers [i](#)

There are no such residues in this entry.