



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 07:46 AM UTC

PDB ID : 1VIX / pdb_00001vix
Title : Crystal structure of a putative peptidase T
Authors : Structural GenomiX
Deposited on : 2003-12-01
Resolution : 2.50 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

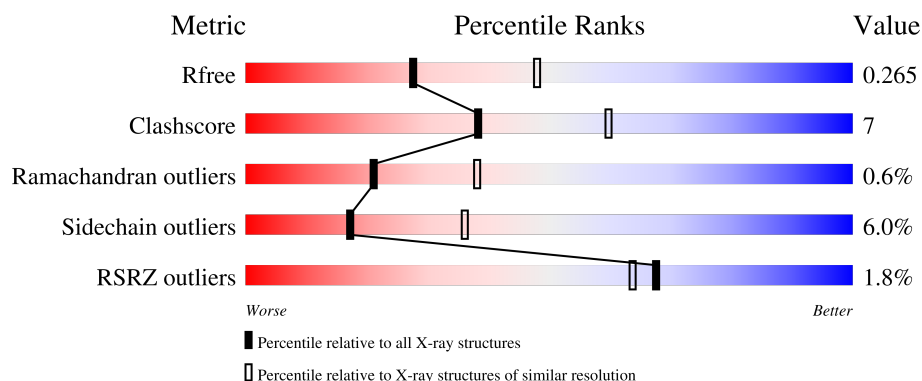
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	420	
1	B	420	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6513 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 Peptidase T.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	411	Total	C	N	O	S	0	2	0
			3167	2001	545	602	19			
1	B	410	Total	C	N	O	S	0	1	0
			3161	1998	542	602	19			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	cloning artifact	UNP P29745
A	0	SER	-	cloning artifact	UNP P29745
A	1	LEU	-	cloning artifact	UNP P29745
A	409	GLU	-	cloning artifact	UNP P29745
A	410	GLY	-	cloning artifact	UNP P29745
A	411	GLY	-	cloning artifact	UNP P29745
A	412	SER	-	cloning artifact	UNP P29745
A	413	HIS	-	cloning artifact	UNP P29745
A	414	HIS	-	cloning artifact	UNP P29745
A	415	HIS	-	cloning artifact	UNP P29745
A	416	HIS	-	cloning artifact	UNP P29745
A	417	HIS	-	cloning artifact	UNP P29745
A	418	HIS	-	cloning artifact	UNP P29745
B	-1	MET	-	cloning artifact	UNP P29745
B	0	SER	-	cloning artifact	UNP P29745
B	1	LEU	-	cloning artifact	UNP P29745
B	409	GLU	-	cloning artifact	UNP P29745
B	410	GLY	-	cloning artifact	UNP P29745
B	411	GLY	-	cloning artifact	UNP P29745
B	412	SER	-	cloning artifact	UNP P29745
B	413	HIS	-	cloning artifact	UNP P29745
B	414	HIS	-	cloning artifact	UNP P29745
B	415	HIS	-	cloning artifact	UNP P29745
B	416	HIS	-	cloning artifact	UNP P29745
B	417	HIS	-	cloning artifact	UNP P29745

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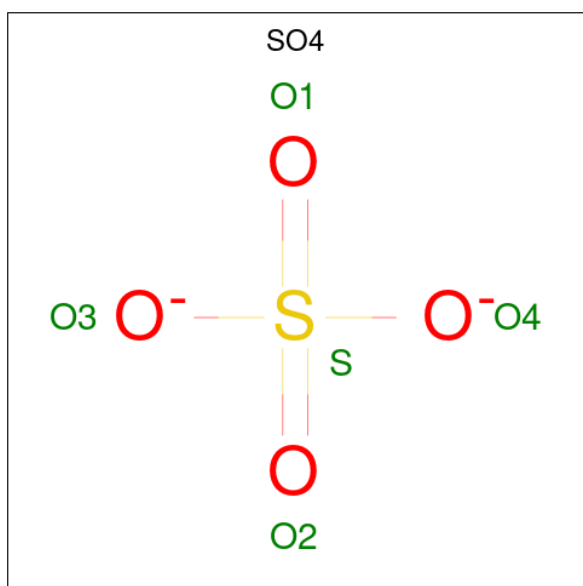
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Chain	Residue	Modelled	Actual	Comment	Reference
B	418	HIS	-	cloning artifact	UNP P29745

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Zn	0	0
			2	2		
2	B	2	Total	Zn	0	0
			2	2		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

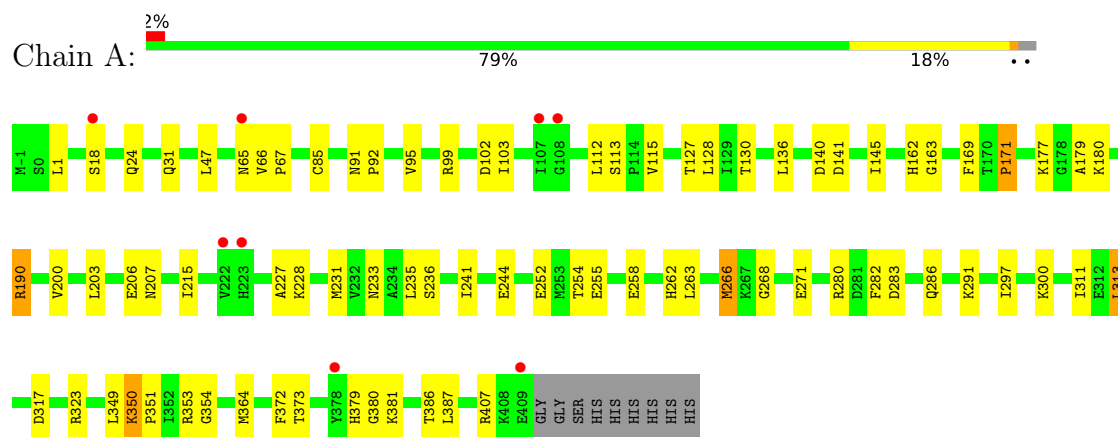
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	65	Total	O	0	0
			65	65		
4	B	56	Total	O	0	0
			56	56		

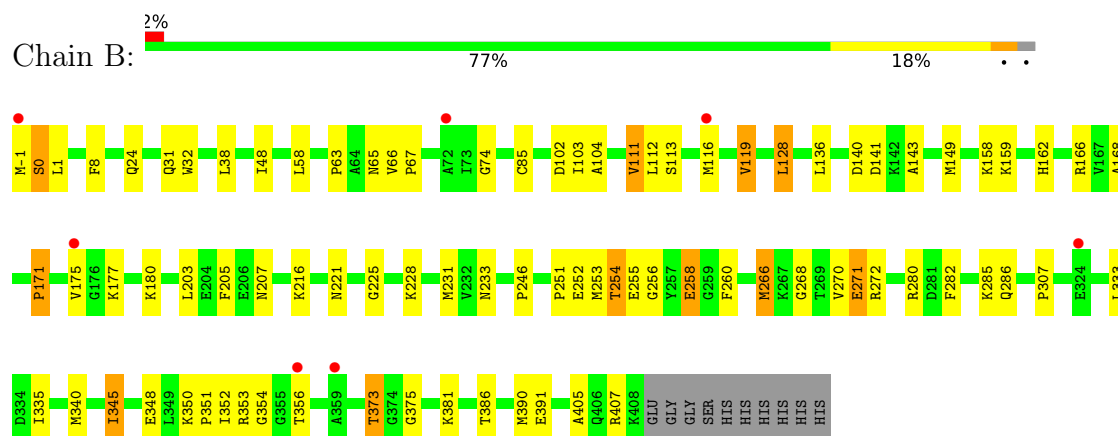
3 Residue-property plots

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: Peptidase T



• Molecule 1: Peptidase T



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	115.27Å 145.79Å 77.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.69 – 2.50 32.69 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (32.69-2.50) 76.3 (32.69-2.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.83 (at 2.51Å)	Xtriage
Refinement program	REFMAC 4.0	Depositor
R, R_{free}	0.256 , 0.303 0.220 , 0.265	Depositor DCC
R_{free} test set	1773 reflections (3.80%)	wwPDB-VP
Wilson B-factor (Å ²)	42.9	Xtriage
Anisotropy	0.179	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 26.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6513	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.36% 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, ZN

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	0.45	0/3240	0.86	1/4386 (0.0%)
1	B	0.44	0/3228	0.88	2/4366 (0.0%)
All	All	0.45	0/6468	0.87	3/8752 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	268	GLY	N-CA-C	5.62	119.89	110.56
1	B	119	VAL	CB-CA-C	-5.25	105.16	112.04
1	B	268	GLY	N-CA-C	5.24	118.36	110.18

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	3167	0	3131	49	0
1	B	3161	0	3140	45	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	30	0	0	0	0
3	B	30	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	65	0	0	1	0
4	B	56	0	0	0	0
All	All	6513	0	6271	87	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 7.

All (87) 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:A:190:ARG:HG3	1:A:190:ARG:HH11	1.11	1.14
1:A:31:GLN:NE2	1:A:171:PRO:O	2.11	0.83
1:A:190:ARG:HG3	1:A:190:ARG:NH1	1.90	0.71
1:B:85:CYS:HB3	1:B:136:LEU:CD2	2.21	0.69
1:B:258:GLU:O	1:B:353:ARG:NH2	2.26	0.68
1:A:85:CYS:HB3	1:A:136:LEU:HG	1.76	0.68
1:B:31:GLN:NE2	1:B:171:PRO:O	2.26	0.68
1:B:102:ASP:OD2	1:B:113:SER:HB2	1.94	0.66
1:B:85:CYS:HB3	1:B:136:LEU:HD21	1.77	0.65
1:B:233:ASN:HA	1:B:270:VAL:HG13	1.79	0.64
1:A:162:HIS:HD2	1:A:163:GLY:O	1.82	0.61
1:B:38:LEU:HD22	1:B:149:MET:HE3	1.84	0.60
1:B:58:LEU:O	1:B:168:ALA:HA	2.02	0.59
1:B:221:ASN:HB3	1:B:271:GLU:HG2	1.85	0.58
1:B:255:GLU:O	1:B:258:GLU:HG3	2.04	0.57
1:A:244:GLU:HB3	1:A:297:ILE:HG12	1.86	0.57
1:A:190:ARG:HH11	1:A:190:ARG:CG	2.03	0.57
1:A:228:LYS:HA	1:B:252:GLU:O	2.05	0.56
1:A:1:LEU:HB3	1:A:387:LEU:HD21	1.88	0.56
1:A:291:LYS:NZ	1:A:317:ASP:OD1	2.32	0.56
1:B:38:LEU:HD22	1:B:149:MET:CE	2.36	0.55
1:B:103:ILE:HD12	1:B:112:LEU:HD23	1.88	0.55
1:A:102:ASP:OD2	1:A:113:SER:HB2	2.07	0.54
1:A:255:GLU:O	1:A:258:GLU:HG2	2.09	0.53
1:B:203:LEU:HB2	1:B:340:MET:HE1	1.90	0.53
1:A:95:VAL:HB	1:A:128:LEU:HB2	1.91	0.53
1:A:65:ASN:HB2	1:A:162:HIS:CE1	2.44	0.52
1:A:136:LEU:CD2	1:A:380:GLY:H	2.22	0.52
1:A:206:GLU:HG2	1:A:323:ARG:HG3	1.93	0.51
1:A:215:ILE:HG12	1:A:313:LEU:HD22	1.92	0.51
1:A:254:THR:O	1:B:228:LYS:HB2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:LEU:CD2	1:A:386:THR:HG22	2.42	0.49
1:A:231:MET:HE2	1:B:251:PRO:HB2	1.95	0.49
1:A:136:LEU:HD21	1:A:380:GLY:H	1.77	0.49
1:A:227:ALA:HB1	1:A:231:MET:HB3	1.94	0.49
1:A:31:GLN:HE22	1:A:171:PRO:C	2.21	0.49
1:A:179:ALA:HB1	1:A:364:MET:HE1	1.95	0.49
1:A:200:VAL:HA	1:A:372:PHE:CD2	2.49	0.48
1:B:1:LEU:HD21	1:B:391:GLU:HB2	1.96	0.48
1:A:65:ASN:HB3	1:A:162:HIS:H	1.78	0.48
1:A:282:PHE:CZ	1:A:351:PRO:HB2	2.49	0.48
1:B:254:THR:HG21	1:B:260:PHE:HD1	1.79	0.47
1:A:231:MET:CE	1:B:251:PRO:HB2	2.44	0.47
1:B:48:ILE:HG12	1:B:63:PRO:HG3	1.95	0.47
1:B:104:ALA:HA	1:B:111:VAL:HG22	1.97	0.47
1:B:66:VAL:HB	1:B:67:PRO:HD2	1.96	0.47
1:A:241:ILE:HD11	1:A:311:ILE:HD13	1.97	0.47
1:B:140:ASP:HA	1:B:141:ASP:HA	1.72	0.47
1:A:283:ASP:HB3	1:A:286:GLN:HB2	1.97	0.46
1:B:128:LEU:HD22	1:B:386:THR:HG22	1.97	0.46
1:B:216:LYS:HG3	1:B:272:ARG:NH2	2.31	0.46
1:A:128:LEU:HD22	1:A:386:THR:HG22	1.97	0.46
1:A:203:LEU:HB3	1:A:349:LEU:HD23	1.98	0.46
1:A:233:ASN:HB3	1:A:236:SER:HB2	1.97	0.46
1:B:373:THR:HG22	1:B:375:GLY:H	1.81	0.45
1:A:207:ASN:HB2	1:A:354:GLY:O	2.16	0.45
1:B:280:ARG:NH1	3:B:423:SO4:S	2.90	0.45
1:A:262:HIS:CE1	1:B:231:MET:HE1	2.52	0.45
1:A:66:VAL:HB	1:A:67:PRO:HD2	1.99	0.44
1:A:353:ARG:HD3	1:B:225:GLY:O	2.17	0.44
1:B:143:ALA:HA	1:B:390:MET:HE3	2.00	0.44
1:B:352:ILE:HG21	1:B:356:THR:HG23	2.00	0.44
1:A:92:PRO:HA	1:A:130:THR:O	2.17	0.44
1:A:103:ILE:HD13	1:A:112:LEU:HD23	1.98	0.43
1:A:136:LEU:HD21	1:A:379:HIS:HA	1.99	0.43
1:B:282:PHE:CZ	1:B:351:PRO:HB2	2.54	0.43
1:A:145:ILE:HG23	1:A:169:PHE:CE2	2.54	0.42
1:A:350:LYS:HB2	1:A:351:PRO:HD2	2.01	0.42
1:B:251:PRO:HG3	1:B:260:PHE:CE2	2.54	0.42
1:B:340:MET:HB3	1:B:345:ILE:HB	2.01	0.42
1:A:140:ASP:HA	1:A:141:ASP:HA	1.78	0.42
1:B:255:GLU:H	1:B:258:GLU:HG3	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:-1:MET:HG2	1:B:0:SER:H	1.85	0.42
1:A:263:LEU:HD21	1:A:266:MET:HE2	2.00	0.42
1:B:335:ILE:HD11	1:B:407:ARG:HD2	2.02	0.41
1:A:280:ARG:NH1	4:A:427:HOH:O	2.44	0.41
1:B:207:ASN:HB2	1:B:354:GLY:O	2.20	0.41
1:B:8:PHE:CG	1:B:390:MET:HE1	2.56	0.41
1:B:158:LYS:O	1:B:159:LYS:C	2.64	0.41
1:B:32:TRP:HB3	3:B:426:SO4:O1	2.21	0.41
1:B:205:PHE:HB3	1:B:333:LEU:HD21	2.01	0.41
1:A:136:LEU:HD21	1:A:379:HIS:CA	2.51	0.40
1:A:227:ALA:HB1	1:A:231:MET:CB	2.51	0.40
1:B:162:HIS:HA	1:B:405:ALA:HB2	2.03	0.40
1:A:235:LEU:HD22	1:B:266:MET:HG2	2.03	0.40
1:B:74:GLY:HA2	1:B:166:ARG:O	2.22	0.40
1:A:127:THR:O	1:A:386:THR:HA	2.20	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	411/420 (98%)	395 (96%)	15 (4%)	1 (0%)	43	63
1	B	409/420 (97%)	392 (96%)	13 (3%)	4 (1%)	12	24
All	All	820/840 (98%)	787 (96%)	28 (3%)	5 (1%)	21	38

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	171	PRO
1	B	246	PRO
1	B	171	PRO

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Mol	Chain	Res	Type
1	B	307	PRO
1	B	256	GLY

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	337/347 (97%)	319 (95%)	18 (5%)	20	42
1	B	338/347 (97%)	316 (94%)	22 (6%)	15	32
All	All	675/694 (97%)	635 (94%)	40 (6%)	17	37

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	SER
1	A	24	GLN
1	A	47	LEU
1	A	91	ASN
1	A	99	ARG
1	A	115	VAL
1	A	177	LYS
1	A	180	LYS
1	A	190	ARG
1	A	252	GLU
1	A	266	MET
1	A	271	GLU
1	A	300	LYS
1	A	313	LEU
1	A	350	LYS
1	A	373	THR
1	A	381	LYS
1	A	407	ARG
1	B	0	SER
1	B	24	GLN
1	B	65	ASN

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Mol	Chain	Res	Type
1	B	111	VAL
1	B	116	MET
1	B	119	VAL
1	B	128	LEU
1	B	175	VAL
1	B	177	LYS
1	B	180	LYS
1	B	253	MET
1	B	254	THR
1	B	258	GLU
1	B	266	MET
1	B	271	GLU
1	B	285	LYS
1	B	286	GLN
1	B	345	ILE
1	B	348	GLU
1	B	350	LYS
1	B	373	THR
1	B	381	LYS

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

Mol	Chain	Res	Type
1	A	31	GLN
1	A	41	GLN
1	A	121	HIS
1	A	162	HIS
1	A	221	ASN
1	A	286	GLN
1	A	329	HIS
1	A	379	HIS
1	B	24	GLN
1	B	41	GLN
1	B	121	HIS
1	B	126	GLN
1	B	223	HIS
1	B	286	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 4 are monoatomic - leaving 12 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	SO4	A	422	-	4,4,4	0.23	0	6,6,6	0.13	0
3	SO4	B	423	-	4,4,4	0.22	0	6,6,6	0.09	0
3	SO4	A	424	-	4,4,4	0.23	0	6,6,6	0.13	0
3	SO4	B	422	-	4,4,4	0.25	0	6,6,6	0.08	0
3	SO4	A	423	-	4,4,4	0.24	0	6,6,6	0.10	0
3	SO4	B	421	-	4,4,4	0.25	0	6,6,6	0.12	0
3	SO4	B	425	-	4,4,4	0.26	0	6,6,6	0.11	0
3	SO4	B	426	-	4,4,4	0.22	0	6,6,6	0.19	0
3	SO4	A	426	-	4,4,4	0.21	0	6,6,6	0.08	0
3	SO4	A	421	-	4,4,4	0.22	0	6,6,6	0.14	0
3	SO4	A	425	-	4,4,4	0.26	0	6,6,6	0.09	0
3	SO4	B	424	-	4,4,4	0.24	0	6,6,6	0.05	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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	423	SO4	1	0
3	B	426	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 [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	411/420 (97%)	0.21	8 (1%) 66 62	27, 46, 60, 85	2 (0%)
1	B	410/420 (97%)	0.21	7 (1%) 69 65	29, 47, 62, 82	1 (0%)
All	All	821/840 (97%)	0.21	15 (1%) 67 64	27, 47, 61, 85	3 (0%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	409	GLU	5.5
1	B	116	MET	2.9
1	B	359	ALA	2.6
1	A	223[A]	HIS	2.6
1	A	108	GLY	2.6
1	B	324	GLU	2.4
1	A	107	ILE	2.4
1	A	378	TYR	2.4
1	A	222	VAL	2.4
1	B	356	THR	2.3
1	A	65	ASN	2.2
1	A	18	SER	2.2
1	B	72	ALA	2.1
1	B	-1	MET	2.0
1	B	175	VAL	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 oligosaccharides in this entry.

6.4 Ligands

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
3	SO4	B	421	5/5	0.75	0.14	103,103,103,103	0
3	SO4	B	426	5/5	0.75	0.21	101,101,101,101	0
3	SO4	A	422	5/5	0.80	0.17	86,87,87,87	0
3	SO4	B	424	5/5	0.81	0.12	124,125,125,125	0
3	SO4	B	423	5/5	0.81	0.17	108,108,109,109	0
3	SO4	A	424	5/5	0.84	0.10	102,102,103,103	0
3	SO4	A	425	5/5	0.84	0.18	89,90,90,90	0
3	SO4	A	423	5/5	0.84	0.09	87,87,87,87	0
3	SO4	B	422	5/5	0.87	0.09	92,92,92,93	0
3	SO4	B	425	5/5	0.88	0.16	85,85,85,85	0
3	SO4	A	426	5/5	0.88	0.15	95,95,95,95	0
2	ZN	B	420	1/1	0.90	0.10	115,115,115,115	0
3	SO4	A	421	5/5	0.93	0.12	76,76,76,77	0
2	ZN	A	420	1/1	0.93	0.15	97,97,97,97	0
2	ZN	B	419	1/1	0.98	0.04	59,59,59,59	0
2	ZN	A	419	1/1	0.99	0.06	56,56,56,56	0

6.5 Other polymers

There are no such residues in this entry.