



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2026 – 07:55 am BST

PDB ID : 2JCM / pdb_00002jcm
Title : Crystal structure of Human Cytosolic 5'-Nucleotidase II in complex with beryllium trifluoride
Authors : Wallden, K.; Stenmark, P.; Arrowsmith, C.; Berglund, H.; Busam, R.; Collins, R.; Edwards, A.; Ehn, M.; Flodin, S.; Flores, A.; Graslund, S.; Hammarstrom, M.; Hallberg, B.M.; Holmberg Schiavone, L.; Hogbom, M.; Karlberg, T.; Kotenyova, T.; Magnusdottir, A.; Nilsson-Ehle, P.; Nyman, T.; Ogg, D.; Persson, C.; Sagemark, J.; Sundstrom, M.; Uppenberg, J.; Thorsell, A.G.; Van Den Berg, S.; Loppnau, P.; Weigelt, J.; Welin, M.; Nordlund, P.
Deposited on : 2006-12-27
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

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	: 1.8.4, CSD as541be (2020)
Xtriage (Phenix)	: 2.0
EDS	: 3.0
Percentile statistics	: 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	: 9.0.015 (Gargrove)
Density-Fitness	: 1.0.12

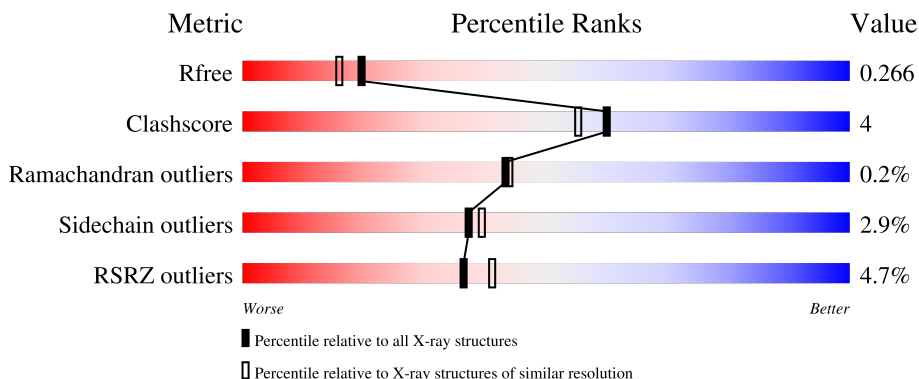
1 Overall quality at a glance

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}	180053	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (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 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	555	

Ideal geometry (proteins) : Engh & Huber (2001)
 Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
 Validation Pipeline (wwPDB-VP) : 2.50

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 4109 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 CYTOSOLIC PURINE 5'-NUCLEOTIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	467	Total	C	N	O	S	0	3	0
			3833	2476	639	697	21			

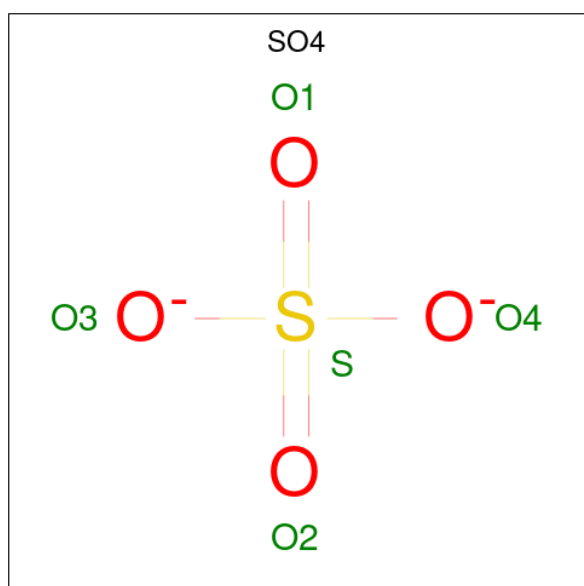
- Molecule 2 is UNKNOWN LIGAND (CCD ID: UNX) (formula: X).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	X	0	0
			1	1		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

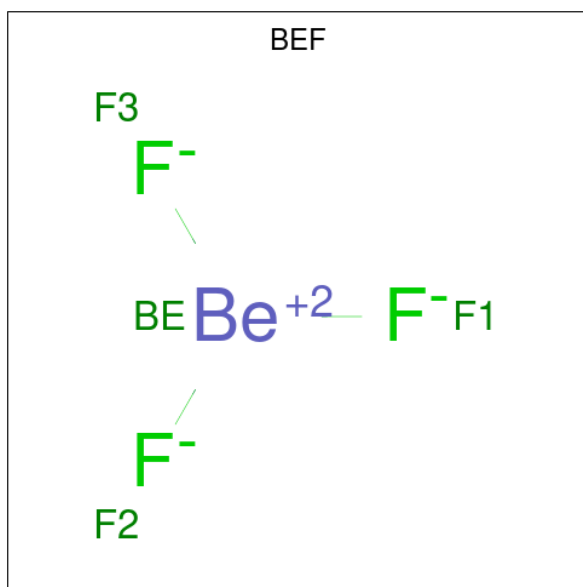
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 5 is BERYLLIUM TRIFLUORIDE ION (CCD ID: BEF) (formula: BeF_3).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	Be	F	0	0
			4	1	3		

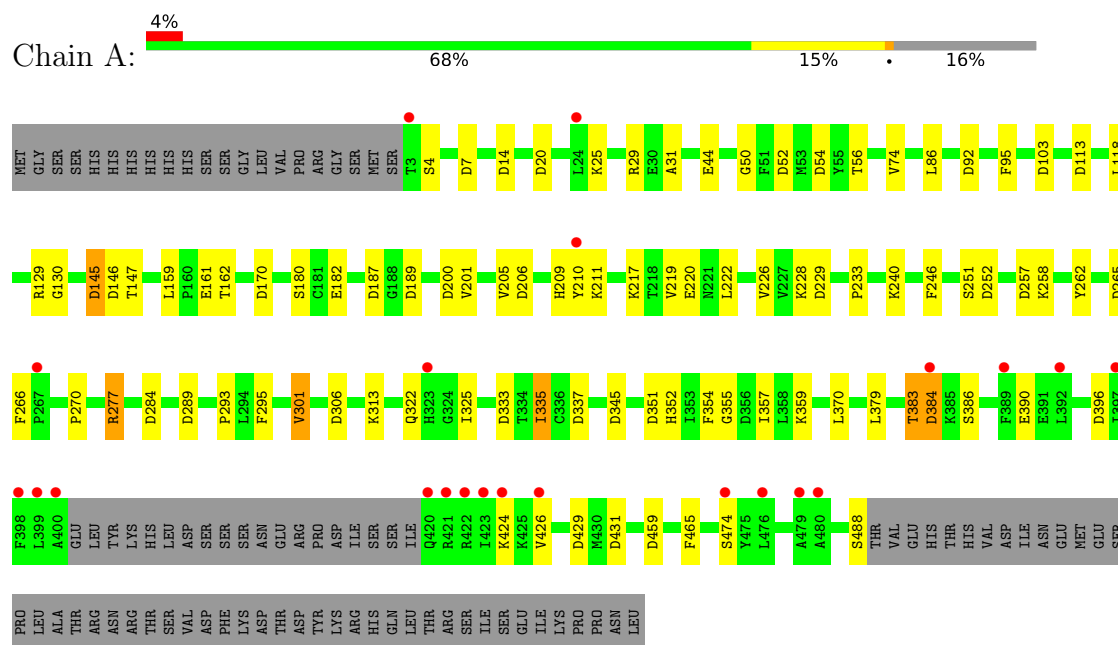
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	260	Total	O	0	0
			260	260		

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: CYTOSOLIC PURINE 5'-NUCLEOTIDASE



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	91.48Å 128.26Å 131.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.67 – 2.15 20.67 – 2.15	Depositor EDS
% Data completeness (in resolution range)	98.1 (20.67-2.15) 98.0 (20.67-2.15)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.17 (at 2.15Å)	Xtriage
Refinement program	REFMAC 5.3.0021	Depositor
R, R_{free}	0.210 , 0.266 0.217 , 0.266	Depositor DCC
R_{free} test set	2089 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	21.5	Xtriage
Anisotropy	0.558	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 39.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.015 for -h,-l,-k	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4109	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.24% 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: MG, UNX, SO4, BEF

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	1.19	33/3932 (0.8%)	1.13	5/5308 (0.1%)

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	206	ASP	CG-OD1	7.94	1.40	1.25
1	A	429	ASP	CG-OD1	7.92	1.40	1.25
1	A	337	ASP	CG-OD1	7.43	1.39	1.25
1	A	345	ASP	CG-OD1	7.15	1.39	1.25
1	A	187	ASP	CG-OD1	7.14	1.39	1.25
1	A	306	ASP	CG-OD1	7.08	1.38	1.25
1	A	431	ASP	CG-OD1	6.85	1.38	1.25
1	A	103	ASP	CG-OD1	6.51	1.37	1.25
1	A	384	ASP	CG-OD1	6.40	1.37	1.25
1	A	252	ASP	CG-OD1	6.39	1.37	1.25
1	A	54	ASP	CG-OD1	6.21	1.37	1.25
1	A	333	ASP	CG-OD1	6.20	1.37	1.25
1	A	14	ASP	CG-OD1	6.19	1.37	1.25
1	A	284	ASP	CG-OD1	6.16	1.37	1.25
1	A	229	ASP	CG-OD1	6.08	1.36	1.25
1	A	257	ASP	CG-OD1	6.08	1.36	1.25
1	A	145[A]	ASP	CG-OD1	6.06	1.36	1.25
1	A	145[B]	ASP	CG-OD1	6.06	1.36	1.25
1	A	396	ASP	CG-OD1	6.04	1.36	1.25
1	A	20	ASP	CG-OD1	5.90	1.36	1.25
1	A	289	ASP	CG-OD1	5.79	1.36	1.25
1	A	92	ASP	CG-OD1	5.76	1.36	1.25
1	A	265	ASP	CG-OD1	5.74	1.36	1.25
1	A	113	ASP	CG-OD1	5.65	1.36	1.25
1	A	146	ASP	CG-OD1	5.60	1.35	1.25
1	A	200	ASP	CG-OD1	5.59	1.35	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	170	ASP	CG-OD1	5.48	1.35	1.25
1	A	189	ASP	CG-OD1	5.30	1.35	1.25
1	A	226	VAL	CA-CB	5.23	1.61	1.54
1	A	351	ASP	CG-OD1	5.15	1.35	1.25
1	A	52	ASP	CG-OD2	5.14	1.35	1.25
1	A	301	VAL	CA-CB	5.11	1.60	1.54
1	A	56	THR	CA-CB	5.01	1.59	1.53

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	459	ASP	N-CA-C	-6.37	104.26	111.07
1	A	95	PHE	O-C-N	5.57	124.11	120.27
1	A	313	LYS	N-CA-C	-5.50	103.27	110.53
1	A	277	ARG	CB-CA-C	5.14	116.84	108.87
1	A	95	PHE	N-CA-CB	5.05	115.19	110.39

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	3833	0	3772	33	0
2	A	1	0	0	0	0
3	A	1	0	0	0	0
4	A	10	0	0	0	0
5	A	4	0	0	0	0
6	A	260	0	0	5	0
All	All	4109	0	3772	33	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 4.

All (33) 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:219:VAL:HG11	1:A:258:LYS:HD2	1.54	0.89
1:A:386:SER:O	1:A:390:GLU:HG3	1.89	0.72
1:A:210:TYR:CD2	1:A:211:LYS:HG3	2.31	0.65
1:A:352:HIS:CD2	1:A:355:GLY:H	2.16	0.62
1:A:322:GLN:O	1:A:325:ILE:HG22	2.02	0.59
1:A:182:GLU:HG3	6:A:2089:HOH:O	2.02	0.58
1:A:25:LYS:O	1:A:29:ARG:HG2	2.05	0.57
1:A:352:HIS:HD2	1:A:354:PHE:H	1.53	0.55
1:A:233:PRO:HB2	1:A:270:PRO:HA	1.89	0.55
1:A:180:SER:HB3	6:A:2123:HOH:O	2.10	0.52
1:A:357:ILE:HG13	1:A:370:LEU:HD13	1.92	0.52
1:A:209:HIS:HD2	6:A:2137:HOH:O	1.94	0.50
1:A:31:ALA:HB1	1:A:379:LEU:HD22	1.95	0.49
1:A:74:VAL:HG13	1:A:86:LEU:HB3	1.93	0.49
1:A:217:LYS:HA	1:A:220:GLU:HG2	1.96	0.48
1:A:293:PRO:CB	1:A:359:LYS:HG3	2.45	0.47
1:A:383:THR:CG2	1:A:384:ASP:N	2.78	0.47
1:A:293:PRO:HB3	1:A:359:LYS:HG3	1.99	0.45
1:A:222:LEU:HD13	1:A:262:TYR:HB2	1.98	0.45
1:A:209:HIS:HE1	6:A:2107:HOH:O	1.99	0.45
1:A:295:PHE:HZ	1:A:335:ILE:HD13	1.80	0.45
1:A:145[A]:ASP:O	1:A:147:THR:HG23	2.16	0.44
1:A:201:VAL:O	1:A:205:VAL:HG23	2.18	0.44
1:A:4:SER:O	1:A:7:ASP:HB2	2.19	0.43
1:A:159:LEU:HD23	1:A:159:LEU:HA	1.91	0.43
1:A:424:LYS:NZ	1:A:424:LYS:HB3	2.33	0.43
1:A:219:VAL:CG1	1:A:258:LYS:HD2	2.36	0.42
1:A:228:LYS:HG2	1:A:266:PHE:CZ	2.54	0.42
1:A:295:PHE:HZ	1:A:335:ILE:CD1	2.32	0.42
1:A:210:TYR:HB2	6:A:2132:HOH:O	2.20	0.41
1:A:465:PHE:CD1	1:A:465:PHE:C	2.99	0.41
1:A:161:GLU:O	1:A:162:THR:C	2.64	0.41
1:A:50:GLY:HA3	1:A:246:PHE:CZ	2.56	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	466/555 (84%)	454 (97%)	11 (2%)	1 (0%)	43	44

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	130	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	416/497 (84%)	403 (97%)	13 (3%)	35	37

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	GLU
1	A	118	LEU
1	A	129	ARG
1	A	240	LYS
1	A	251[A]	SER
1	A	251[B]	SER
1	A	277	ARG
1	A	301	VAL
1	A	335	ILE
1	A	383	THR
1	A	426	VAL
1	A	474	SER
1	A	488	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	HIS
1	A	174	ASN
1	A	209	HIS
1	A	221	ASN
1	A	352	HIS
1	A	364	GLN
1	A	380	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 1 is unknown and 1 is monoatomic - leaving 3 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	$\# Z > 2$
4	SO4	A	1491	-	4,4,4	0.37	0	6,6,6	0.79	0
5	BEF	A	1493	1	0,3,3	-	-	-		
4	SO4	A	1492	-	4,4,4	0.17	0	6,6,6	0.20	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.

No monomer is involved in short contacts.

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 [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	467/555 (84%)	-0.08	22 (4%)	36 41	2, 10, 34, 56	3 (0%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	398	PHE	4.8
1	A	399	LEU	4.6
1	A	424	LYS	3.8
1	A	3	THR	3.8
1	A	397	ILE	3.5
1	A	323	HIS	3.3
1	A	389	PHE	3.3
1	A	210	TYR	3.2
1	A	420	GLN	3.0
1	A	392	LEU	2.9
1	A	400	ALA	2.7
1	A	474	SER	2.6
1	A	422	ARG	2.5
1	A	24	LEU	2.4
1	A	479	ALA	2.3
1	A	480	ALA	2.3
1	A	423	ILE	2.3
1	A	267	PRO	2.2
1	A	421	ARG	2.2
1	A	426	VAL	2.2
1	A	384	ASP	2.2
1	A	476	LEU	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 [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	UNX	A	1489	1/1	0.73	0.30	42,42,42,42	0
4	SO4	A	1492	5/5	0.92	0.11	64,64,65,66	0
4	SO4	A	1491	5/5	0.96	0.10	24,29,33,33	0
3	MG	A	1490	1/1	0.97	0.08	10,10,10,10	0
5	BEF	A	1493	4/4	0.99	0.10	2,2,3,6	0

6.5 Other polymers [i](#)

There are no such residues in this entry.