



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:10 PM UTC

PDB ID : 3ZKX / pdb_00003zkx
Title : TERNARY BACE2 XAPERONE COMPLEX
Authors : Kuglstatter, A.; Banner, D.W.; Benz, J.; Bertschinger, J.; Burger, D.; Cuppuleri, S.; Debulpaep, M.; Gast, A.; Grabulovski, D.; Gsell, B.; Hilpert, H.; Huber, W.; Kuszniir, E.; Laeremans, T.; Matile, H.; Rufer, A.; Schlatter, D.; Steyeart, J.; Stihle, M.; Thoma, R.; Weber, M.; Ruf, A.
Deposited on : 2013-01-25
Resolution : 2.37 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

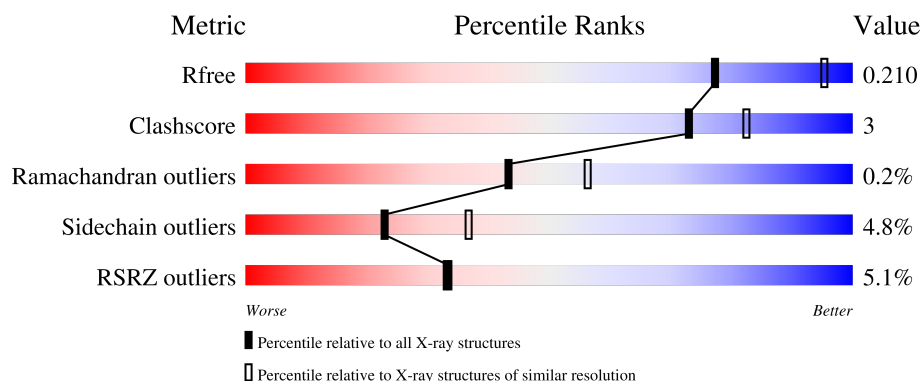
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.37 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	7164 (2.40-2.36)
Clashscore	190562	7722 (2.40-2.36)
Ramachandran outliers	187476	7626 (2.40-2.36)
Sidechain outliers	187428	7627 (2.40-2.36)
RSRZ outliers	180081	7170 (2.40-2.36)

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	386	4% 87% 8% • •
2	B	122	2% 83% 7% • 9%
3	C	122	9% 68% 19% • • 9%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4857 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-SECRETASE 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	370	Total	C	N	O	S	0	0	0
			2866	1848	456	550	12			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	269	ALA	GLU	engineered mutation	UNP Q9Y5Z0

- Molecule 2 is a protein called XA4813.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	111	Total	C	N	O	S	0	0	0
			827	509	150	164	4			

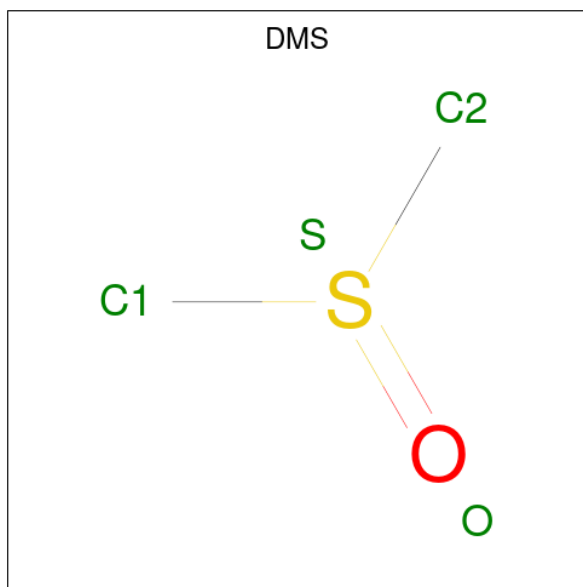
- Molecule 3 is a protein called XA4815.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	111	Total	C	N	O	S	0	0	0
			848	521	165	158	4			

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total	0	0
			1 Cl 1		

- Molecule 5 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C₂H₆OS).



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

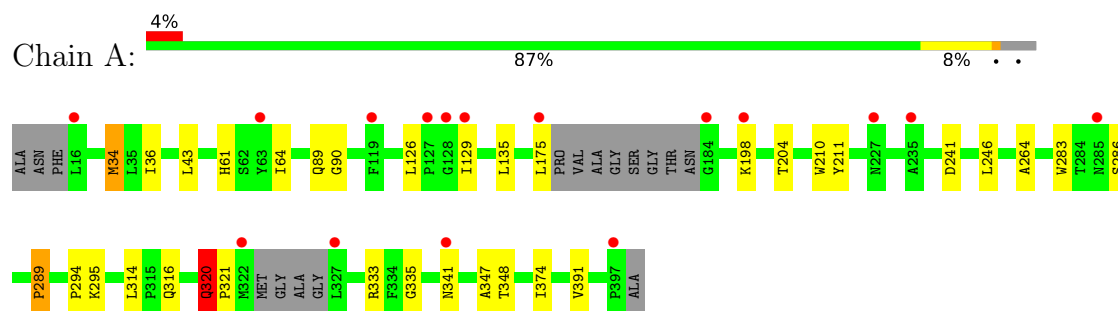
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	202	Total	O	0	0
			202	202		
6	B	53	Total	O	0	0
			53	53		
6	C	56	Total	O	0	0
			56	56		

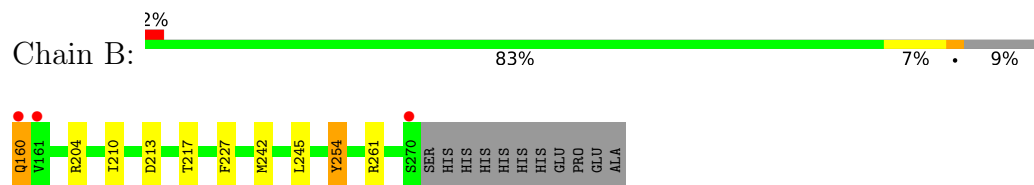
3 Residue-property plots [i](#)

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.

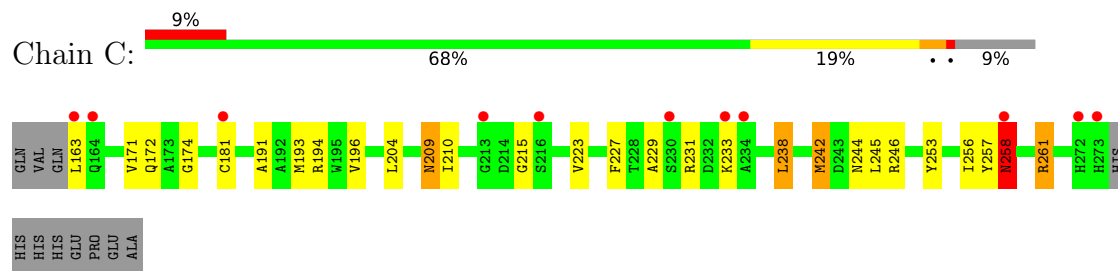
• Molecule 1: BETA-SECRETASE 2



• Molecule 2: XA4813



• Molecule 3: XA4815



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	63.37Å 153.83Å 247.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.70 – 2.37 34.70 – 2.37	Depositor EDS
% Data completeness (in resolution range)	99.6 (34.70-2.37) 99.5 (34.70-2.37)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.15 (at 2.37Å)	Xtriage
Refinement program	BUSTER 2.11.4	Depositor
R, R_{free}	0.185 , 0.208 0.182 , 0.210	Depositor DCC
R_{free} test set	2505 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	56.6	Xtriage
Anisotropy	0.342	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 65.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4857	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.23% 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: DMS, CL

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	0.88	1/2937 (0.0%)	1.21	11/3996 (0.3%)
2	B	0.81	0/841	1.24	4/1138 (0.4%)
3	C	0.90	1/865 (0.1%)	1.28	4/1166 (0.3%)
All	All	0.87	2/4643 (0.0%)	1.23	19/6300 (0.3%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	242	MET	SD-CE	-6.34	1.63	1.79
1	A	34	MET	SD-CE	-5.37	1.66	1.79

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	227	PHE	CA-CB-CG	7.19	120.99	113.80
3	C	227	PHE	CA-CB-CG	7.01	120.81	113.80
3	C	258	ASN	CA-CB-CG	6.50	119.10	112.60
1	A	286	SER	CA-C-N	6.20	129.91	121.05
1	A	286	SER	C-N-CA	6.20	129.91	121.05
1	A	90	GLY	N-CA-C	6.14	117.79	111.95
1	A	347	ALA	CA-C-N	6.02	128.26	120.44
1	A	347	ALA	C-N-CA	6.02	128.26	120.44
2	B	213	ASP	CA-CB-CG	5.90	118.50	112.60
2	B	254	TYR	CA-C-N	5.90	130.62	122.77
2	B	254	TYR	C-N-CA	5.90	130.62	122.77
1	A	320	GLN	N-CA-C	5.80	118.03	109.62
1	A	204	THR	CB-CA-C	5.75	116.25	109.47
1	A	241	ASP	CA-CB-CG	5.68	118.28	112.60
1	A	211	TYR	N-CA-C	-5.64	101.61	110.36
1	A	289	PRO	CA-C-N	5.37	128.22	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	289	PRO	C-N-CA	5.37	128.22	120.38
3	C	215	GLY	N-CA-C	5.20	119.08	110.97
3	C	209	ASN	CA-CB-CG	5.04	117.64	112.60

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	2866	0	2789	12	0
2	B	827	0	809	5	0
3	C	848	0	809	12	0
4	B	1	0	0	0	0
5	B	4	0	6	0	0
6	A	202	0	0	0	0
6	B	53	0	0	0	0
6	C	56	0	0	0	0
All	All	4857	0	4413	29	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 3.

All (29) 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)
3:C:242:MET:HE2	3:C:245:LEU:HD21	1.75	0.69
3:C:193:MET:HG2	3:C:238:LEU:HD22	1.77	0.66
3:C:193:MET:HE3	3:C:238:LEU:HD13	1.89	0.55
1:A:295:LYS:HG2	1:A:314:LEU:HD22	1.92	0.52
2:B:242:MET:HE2	2:B:245:LEU:HD21	1.93	0.51
1:A:283:TRP:CZ3	1:A:289:PRO:HB3	2.47	0.49
3:C:191:ALA:HB3	3:C:257:TYR:CE1	2.48	0.49
3:C:256:ILE:HG12	3:C:261:ARG:HG3	1.94	0.49
2:B:210:ILE:HG13	2:B:217:THR:HG22	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:191:ALA:HB1	3:C:258:ASN:HB2	1.94	0.49
1:A:36:ILE:HD12	1:A:43:LEU:HD13	1.94	0.48
1:A:283:TRP:CH2	1:A:289:PRO:HB3	2.49	0.48
3:C:238:LEU:HD11	3:C:257:TYR:CD1	2.49	0.47
2:B:254:TYR:CD2	2:B:261:ARG:HD2	2.51	0.46
1:A:316:GLN:HB3	1:A:374:ILE:HD12	1.99	0.45
1:A:126:LEU:O	1:A:129:ILE:HG12	2.16	0.45
2:B:242:MET:HB3	2:B:245:LEU:HD21	2.00	0.44
1:A:61:HIS:HB3	1:A:64:ILE:HG13	2.00	0.43
3:C:210:ILE:HD11	3:C:229:ALA:HB1	2.01	0.43
1:A:246:LEU:HD22	1:A:335:GLY:HA2	2.00	0.42
1:A:34:MET:CE	1:A:135:LEU:HD22	2.49	0.42
3:C:194:ARG:HG3	3:C:209:ASN:OD1	2.20	0.42
3:C:181:CYS:HB3	3:C:238:LEU:HB2	2.02	0.42
1:A:320:GLN:HA	1:A:321:PRO:HD3	1.96	0.42
1:A:264:ALA:HB3	1:A:294:PRO:HG3	2.02	0.41
2:B:160:GLN:HE21	2:B:160:GLN:HB3	1.75	0.41
1:A:34:MET:HE1	1:A:135:LEU:HD22	2.03	0.41
3:C:174:GLY:HA2	3:C:244:ASN:HD22	1.86	0.41
3:C:242:MET:HE1	3:C:253:TYR:CZ	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	364/386 (94%)	356 (98%)	8 (2%)	0	100	100
2	B	109/122 (89%)	107 (98%)	2 (2%)	0	100	100
3	C	109/122 (89%)	105 (96%)	3 (3%)	1 (1%)	14	20
All	All	582/630 (92%)	568 (98%)	13 (2%)	1 (0%)	43	56

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	258	ASN

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	311/319 (98%)	302 (97%)	9 (3%)	37	57
2	B	89/99 (90%)	87 (98%)	2 (2%)	45	65
3	C	84/94 (89%)	72 (86%)	12 (14%)	3	4
All	All	484/512 (94%)	461 (95%)	23 (5%)	23	37

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	89	GLN
1	A	175	LEU
1	A	198	LYS
1	A	210	TRP
1	A	320	GLN
1	A	333	ARG
1	A	341	ASN
1	A	348	THR
1	A	391	VAL
2	B	160	GLN
2	B	204	ARG
3	C	163	LEU
3	C	171	VAL
3	C	172	GLN
3	C	196	VAL
3	C	204	LEU
3	C	223	VAL
3	C	231	ARG
3	C	233	LYS
3	C	238	LEU

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Mol	Chain	Res	Type
3	C	246	ARG
3	C	258	ASN
3	C	261	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	123	ASN
1	A	341	ASN
2	B	160	GLN
2	B	241	GLN
3	C	244	ASN
3	C	266	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	$\# Z > 2$
5	DMS	B	1272	-	3,3,3	0.38	0	3,3,3	0.68	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 [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	370/386 (95%)	0.12	16 (4%) 40 39	44, 62, 93, 133	0
2	B	111/122 (90%)	0.03	3 (2%) 56 56	50, 64, 89, 108	0
3	C	111/122 (90%)	0.41	11 (9%) 13 11	44, 64, 96, 126	0
All	All	592/630 (93%)	0.16	30 (5%) 33 33	44, 63, 93, 133	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	175	LEU	6.8
2	B	270	SER	5.1
3	C	273	HIS	3.8
1	A	285	ASN	3.5
3	C	164	GLN	3.4
1	A	16	LEU	3.3
1	A	397	PRO	3.3
3	C	163	LEU	3.0
1	A	235	ALA	2.8
1	A	341	ASN	2.7
1	A	129	ILE	2.7
1	A	322	MET	2.6
2	B	160	GLN	2.6
3	C	258	ASN	2.5
3	C	216	SER	2.5
1	A	119	PHE	2.4
1	A	227	ASN	2.4
1	A	128	GLY	2.4
1	A	63	TYR	2.3
3	C	234	ALA	2.3
3	C	181	CYS	2.3
2	B	161	VAL	2.3
3	C	230	SER	2.3

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Mol	Chain	Res	Type	RSRZ
3	C	233	LYS	2.2
3	C	272	HIS	2.1
1	A	184	GLY	2.1
1	A	198	LYS	2.1
1	A	327	LEU	2.1
1	A	127	PRO	2.1
3	C	213	GLY	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
5	DMS	B	1272	4/4	0.94	0.22	65,71,72,72	0
4	CL	B	1271	1/1	0.96	0.27	82,82,82,82	0

6.5 Other polymers [i](#)

There are no such residues in this entry.