



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 07:29 AM EDT

PDB ID : 2HZL / pdb_00002hzl
Title : Crystal structures of a sodium-alpha-keto acid binding subunit from a TRAP transporter in its closed forms
Authors : Gonin, S.; Arnoux, P.; Pierru, B.; Alonso, B.; Sabaty, M.; Pignol, D.
Deposited on : 2006-08-09
Resolution : 1.40 Å(reported)

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

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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

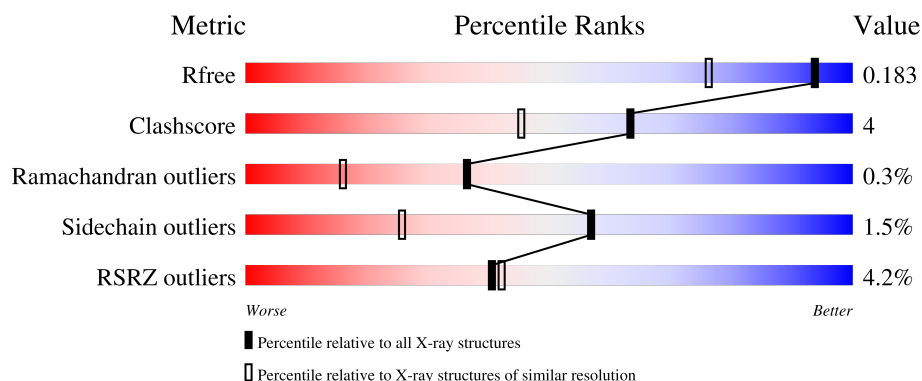
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.40 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	2563 (1.40-1.40)
Clashscore	190562	2660 (1.40-1.40)
Ramachandran outliers	187476	2611 (1.40-1.40)
Sidechain outliers	187428	2610 (1.40-1.40)
RSRZ outliers	180081	2561 (1.40-1.40)

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	365	4% 84% 7% 8%
1	B	365	3% 85% 7% 8%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	PYR	B	603	-	X	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5795 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.

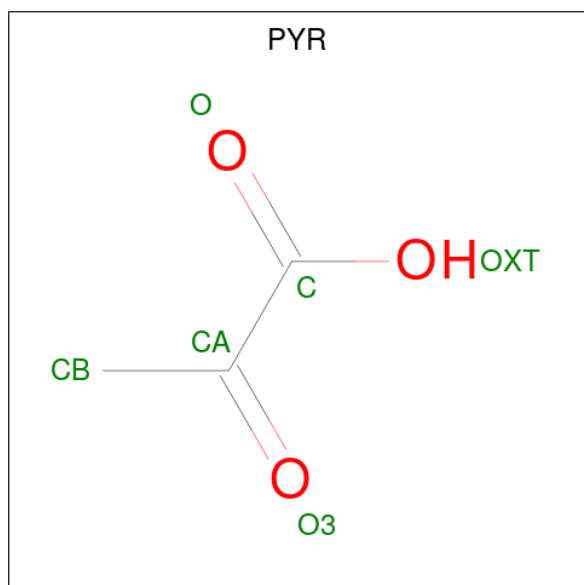
- Molecule 1 is a protein called TRAP-T family sorbitol/mannitol transporter, periplasmic binding protein, SmoM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	334	Total	C	N	O	S	0	4	0
			2632	1701	430	487	14			
1	B	337	Total	C	N	O	S	0	6	0
			2672	1727	437	494	14			

- Molecule 2 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Na	0	0
			1	1		
2	B	1	Total	Na	0	0
			1	1		

- Molecule 3 is PYRUVIC ACID (CCD ID: PYR) (formula: C₃H₄O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total 6	C 3	O 3	0	0
3	B	1	Total 6	C 3	O 3	0	0
3	B	1	Total 6	C 3	O 3	0	0

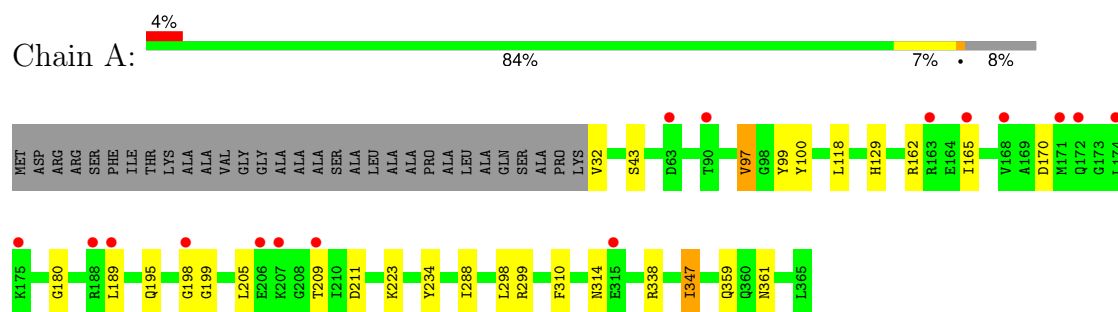
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	222	Total 222	O 222	0	0
4	B	249	Total 249	O 249	0	0

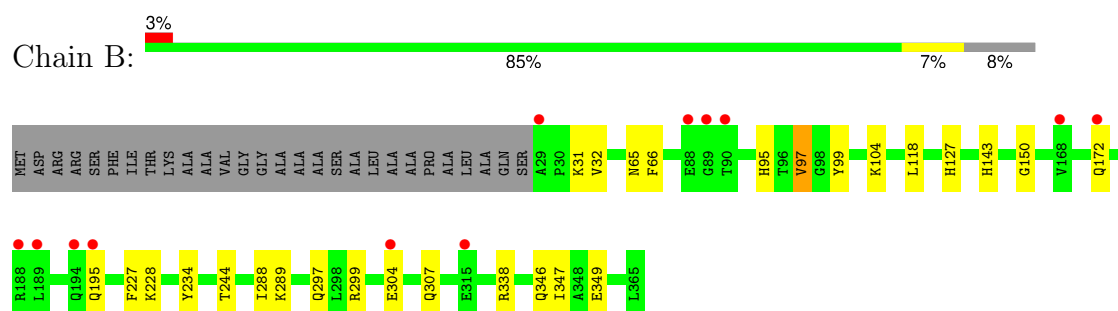
3 Residue-property plots [i](#)

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: TRAP-T family sorbitol/mannitol transporter, periplasmic binding protein, SmoM



- Molecule 1: TRAP-T family sorbitol/mannitol transporter, periplasmic binding protein, SmoM



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	118.02Å 78.06Å 95.14Å 90.00° 124.98° 90.00°	Depositor
Resolution (Å)	39.01 – 1.40 39.01 – 1.40	Depositor EDS
% Data completeness (in resolution range)	100.0 (39.01-1.40) 98.8 (39.01-1.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	79.12 (at 1.40Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.173 , 0.184 0.172 , 0.183	Depositor DCC
R_{free} test set	6908 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	9.4	Xtriage
Anisotropy	0.068	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 38.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5795	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.03% 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: NA, PYR

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	0/2723	0.70	0/3703
1	B	0.45	0/2758	0.68	0/3751
All	All	0.45	0/5481	0.69	0/7454

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

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	2632	0	2521	25	0
1	B	2672	0	2561	23	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	6	0	0	0	0
3	B	12	0	0	0	0
4	A	222	0	0	3	0
4	B	249	0	0	6	0
All	All	5795	0	5082	42	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 (42) 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:349:GLU:HG3	4:B:747:HOH:O	1.60	1.00
1:A:180:GLY:H	1:A:195:GLN:HE21	1.27	0.83
1:B:297:GLN:HE21	1:B:299:ARG:HH22	1.32	0.76
1:A:359:GLN:NE2	1:B:307:GLN:HE22	1.87	0.73
1:A:198:GLY:HA2	4:A:823:HOH:O	1.89	0.72
1:A:359:GLN:HE22	1:B:307:GLN:HE22	1.37	0.72
1:A:162:ARG:NH1	1:A:211:ASP:HA	2.08	0.68
1:A:199:GLY:N	4:A:823:HOH:O	2.14	0.68
1:B:228:LYS:HD2	4:B:704:HOH:O	1.94	0.67
1:A:43:SER:O	1:A:223:LYS:HE2	1.96	0.66
1:A:97[A]:VAL:HG11	1:A:99:TYR:CE2	2.32	0.64
1:A:180:GLY:H	1:A:195:GLN:NE2	1.95	0.63
1:A:361:ASN:HD21	1:B:346:GLN:HE22	1.45	0.62
1:A:162:ARG:HH12	1:A:211:ASP:HA	1.65	0.61
1:A:97[A]:VAL:HG11	1:A:99:TYR:CZ	2.37	0.60
1:B:227:PHE:CZ	1:B:228:LYS:HE3	2.38	0.58
1:B:97[B]:VAL:HG11	1:B:99:TYR:CE2	2.38	0.58
1:A:100:TYR:HA	1:A:195:GLN:HE22	1.69	0.57
1:B:127:HIS:HE1	1:B:244:THR:OG1	1.87	0.56
1:B:95:HIS:HD2	4:B:610:HOH:O	1.87	0.55
1:B:97[B]:VAL:HG11	1:B:99:TYR:CZ	2.44	0.53
1:A:361:ASN:HD21	1:B:346:GLN:NE2	2.07	0.52
1:A:347[B]:ILE:HD12	1:B:288[B]:ILE:HD13	1.93	0.51
1:B:118:LEU:HD21	1:B:338:ARG:HA	1.92	0.51
1:A:288:ILE:HD11	1:A:298:LEU:HD21	1.92	0.51
1:A:310:PHE:CE1	1:A:314[A]:ASN:ND2	2.77	0.51
1:A:32:VAL:N	4:A:745:HOH:O	2.43	0.51
1:A:347[B]:ILE:CD1	1:B:288[B]:ILE:HD13	2.42	0.49
1:A:162:ARG:NH2	1:A:205:LEU:O	2.45	0.49
1:B:297:GLN:HE21	1:B:299:ARG:NH2	2.05	0.49
1:A:129:HIS:HD2	4:B:805:HOH:O	1.96	0.48
1:B:32[A]:VAL:HG13	1:B:66:PHE:HA	1.94	0.48
1:B:104:LYS:HE3	1:B:143:HIS:CE1	2.48	0.47
1:B:31:LYS:H	1:B:65:ASN:HD22	1.63	0.46
1:B:127:HIS:HD2	4:B:643:HOH:O	1.99	0.46
1:B:95:HIS:HE1	1:B:150:GLY:O	1.99	0.46
1:A:234:TYR:HA	1:A:299:ARG:O	2.16	0.46
1:A:118:LEU:HD21	1:A:338:ARG:HA	2.00	0.43
1:B:234:TYR:HA	1:B:299:ARG:O	2.19	0.43
1:B:127:HIS:CD2	4:B:643:HOH:O	2.70	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:165:ILE:HG22	1:A:170:ASP:HB3	2.01	0.42
1:A:97[A]:VAL:CG1	1:A:99:TYR:CE2	3.03	0.41

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	336/365 (92%)	325 (97%)	9 (3%)	2 (1%)	21	5
1	B	341/365 (93%)	330 (97%)	9 (3%)	2 (1%)	21	5
All	All	677/730 (93%)	655 (97%)	18 (3%)	4 (1%)	36	5

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	347[A]	ILE
1	A	347[B]	ILE
1	B	347[A]	ILE
1	B	347[B]	ILE

5.3.2 Protein sidechains [i](#)

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	268/282 (95%)	264 (98%)	4 (2%)	57	26
1	B	272/282 (96%)	266 (98%)	6 (2%)	45	14
All	All	540/564 (96%)	530 (98%)	10 (2%)	57	18

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	97[A]	VAL
1	A	97[B]	VAL
1	A	189	LEU
1	A	209	THR
1	B	97[A]	VAL
1	B	97[B]	VAL
1	B	172	GLN
1	B	195	GLN
1	B	289	LYS
1	B	304	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	124	ASN
1	A	129	HIS
1	A	194	GLN
1	A	195	GLN
1	A	262	GLN
1	A	275	ASN
1	A	307	GLN
1	A	346	GLN
1	A	359	GLN
1	A	361	ASN
1	B	65	ASN
1	B	95	HIS
1	B	127	HIS
1	B	166	ASN
1	B	275	ASN
1	B	297	GLN
1	B	314	ASN

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Mol	Chain	Res	Type
1	B	361	ASN

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 5 ligands modelled in this entry, 2 are 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	$\# Z > 2$
3	PYR	B	603	-	5,5,5	2.97	3 (60%)	3,6,6	1.92	2 (66%)
3	PYR	B	602	2	5,5,5	2.82	3 (60%)	3,6,6	1.89	2 (66%)
3	PYR	A	601	2	5,5,5	2.74	3 (60%)	3,6,6	1.76	2 (66%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PYR	B	603	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PYR	B	602	2	-	0/4/4/4	-
3	PYR	A	601	2	-	0/4/4/4	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	603	PYR	CA-C	-4.38	1.39	1.54
3	B	602	PYR	CA-C	-4.15	1.40	1.54
3	A	601	PYR	CA-C	-4.10	1.40	1.54
3	B	603	PYR	O-C	3.60	1.31	1.22
3	B	602	PYR	O3-CA	3.46	1.31	1.23
3	B	603	PYR	O3-CA	3.45	1.31	1.23
3	B	602	PYR	O-C	3.25	1.30	1.22
3	A	601	PYR	O-C	3.23	1.30	1.22
3	A	601	PYR	O3-CA	3.22	1.30	1.23

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	603	PYR	OXT-C-CA	2.59	120.77	113.59
3	B	602	PYR	OXT-C-O	-2.46	118.05	123.90
3	A	601	PYR	OXT-C-O	-2.22	118.62	123.90
3	B	602	PYR	OXT-C-CA	2.14	119.52	113.59
3	B	603	PYR	OXT-C-O	-2.05	119.02	123.90
3	A	601	PYR	OXT-C-CA	2.03	119.22	113.59

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	603	PYR	OXT-C-CA-CB
3	B	603	PYR	O-C-CA-CB

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

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	334/365 (91%)	0.02	16 (4%) 35 38	4, 10, 21, 27	4 (1%)
1	B	337/365 (92%)	-0.11	12 (3%) 46 48	4, 9, 18, 25	6 (1%)
All	All	671/730 (91%)	-0.05	28 (4%) 40 42	4, 9, 20, 27	10 (1%)

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	172	GLN	3.4
1	A	165	ILE	3.2
1	B	29	ALA	3.1
1	A	189	LEU	2.9
1	A	198	GLY	2.9
1	A	188	ARG	2.7
1	B	194	GLN	2.6
1	A	90	THR	2.5
1	B	304	GLU	2.4
1	A	209	THR	2.4
1	A	206	GLU	2.4
1	B	195	GLN	2.3
1	A	315	GLU	2.3
1	B	315	GLU	2.3
1	A	168	VAL	2.3
1	B	168	VAL	2.3
1	A	174	LEU	2.2
1	A	163	ARG	2.2
1	A	63	ASP	2.2
1	B	88	GLU	2.2
1	B	172	GLN	2.1
1	B	189	LEU	2.1
1	B	188	ARG	2.1
1	B	89	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	207	LYS	2.1
1	A	175	LYS	2.0
1	B	90	THR	2.0
1	A	171	MET	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
3	PYR	B	603	6/6	0.85	0.13	24,25,26,26	0
3	PYR	B	602	6/6	0.98	0.05	8,8,9,10	0
3	PYR	A	601	6/6	0.98	0.05	8,9,9,10	0
2	NA	B	501	1/1	1.00	0.05	4,4,4,4	0
2	NA	A	500	1/1	1.00	0.03	4,4,4,4	0

6.5 Other polymers [i](#)

There are no such residues in this entry.