



wwPDB X-ray Structure Validation Summary Report ⓘ

Nov 3, 2022 – 12:18 pm GMT

PDB ID : 7ZSC
Title : Crystal structure of the heterodimeric human C-P4H-II with truncated alpha subunit (C-P4H-II delta281)
Authors : Lebedev, A.; Koski, M.K.; Wierenga, R.K.; Murthy, A.V.; Sulu, R.
Deposited on : 2022-05-06
Resolution : 3.85 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.31.2
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0267
CCP4 : 7.1.010 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.2

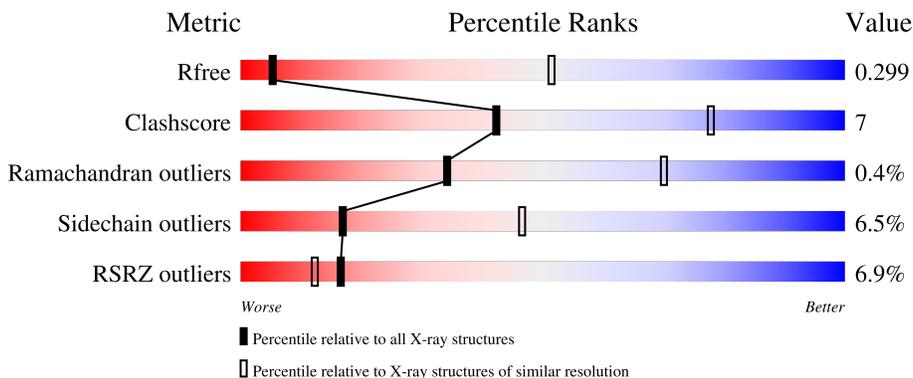
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.85 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}	130704	1048 (4.10-3.62)
Clashscore	141614	1015 (4.08-3.64)
Ramachandran outliers	138981	1069 (4.10-3.62)
Sidechain outliers	138945	1062 (4.10-3.62)
RSRZ outliers	127900	1206 (4.12-3.60)

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	262	 2% (red), 72% (green), 18% (yellow), 10% (grey)
1	B	262	 5% (red), 68% (green), 18% (yellow), 11% (grey)
2	C	492	 9% (red), 79% (green), 13% (yellow), 7% (grey)
2	D	492	 7% (red), 78% (green), 14% (yellow), 7% (grey)

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 11098 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 Prolyl 4-hydroxylase subunit alpha-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	237	Total 1920	C 1213	N 349	O 350	S 8	0	0	0
1	B	233	Total 1897	C 1198	N 347	O 344	S 8	0	1	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	274	MET	-	initiating methionine	UNP O15460
A	275	HIS	-	expression tag	UNP O15460
A	276	HIS	-	expression tag	UNP O15460
A	277	HIS	-	expression tag	UNP O15460
A	278	HIS	-	expression tag	UNP O15460
A	279	HIS	-	expression tag	UNP O15460
A	280	HIS	-	expression tag	UNP O15460
A	281	MET	-	expression tag	UNP O15460
B	274	MET	-	initiating methionine	UNP O15460
B	275	HIS	-	expression tag	UNP O15460
B	276	HIS	-	expression tag	UNP O15460
B	277	HIS	-	expression tag	UNP O15460
B	278	HIS	-	expression tag	UNP O15460
B	279	HIS	-	expression tag	UNP O15460
B	280	HIS	-	expression tag	UNP O15460
B	281	MET	-	expression tag	UNP O15460

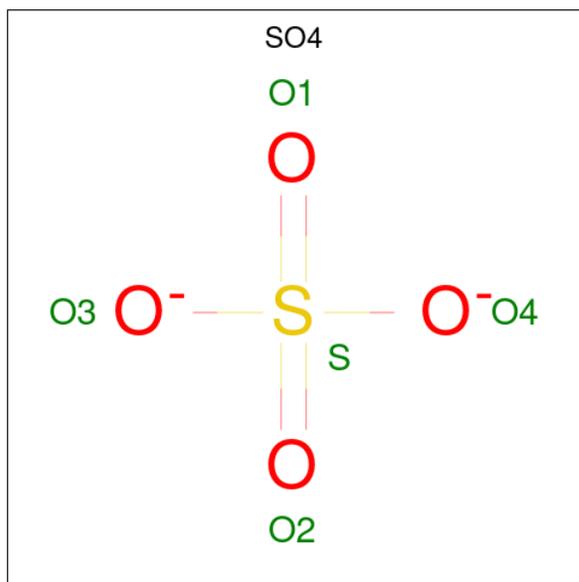
- Molecule 2 is a protein called Protein disulfide-isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	456	Total 3631	C 2329	N 593	O 700	S 9	0	0	0
2	D	458	Total 3640	C 2334	N 595	O 702	S 9	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	17	MET	-	initiating methionine	UNP P07237
D	17	MET	-	initiating methionine	UNP P07237

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).

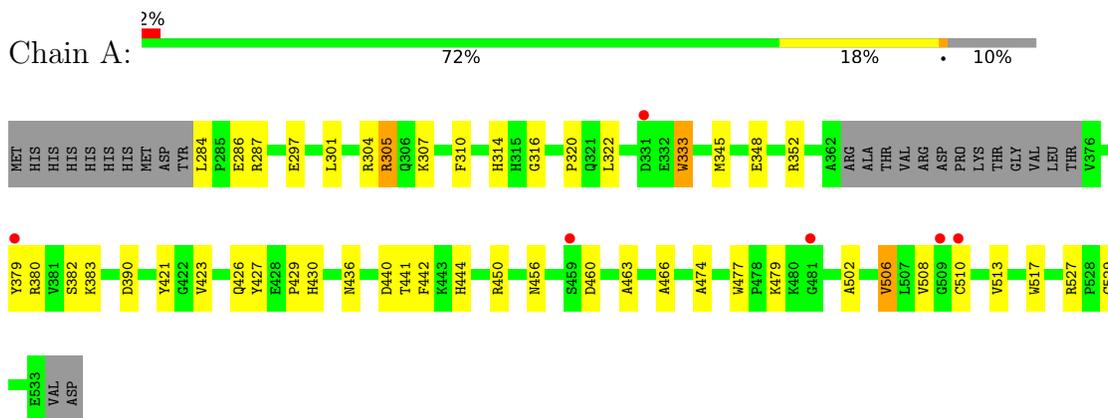


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

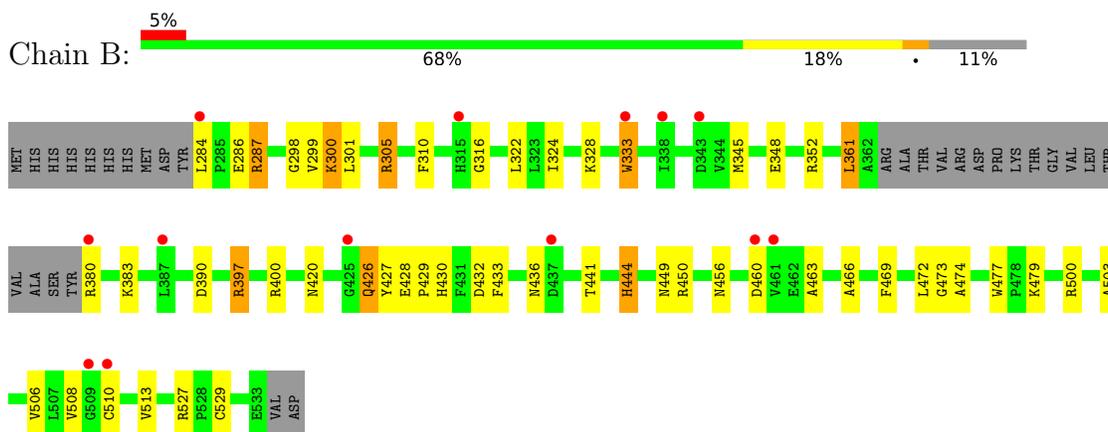
3 Residue-property plots [i](#)

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.

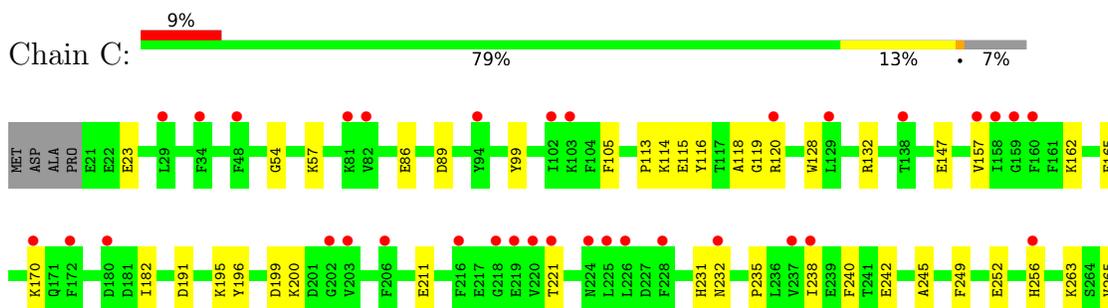
- Molecule 1: Prolyl 4-hydroxylase subunit alpha-2

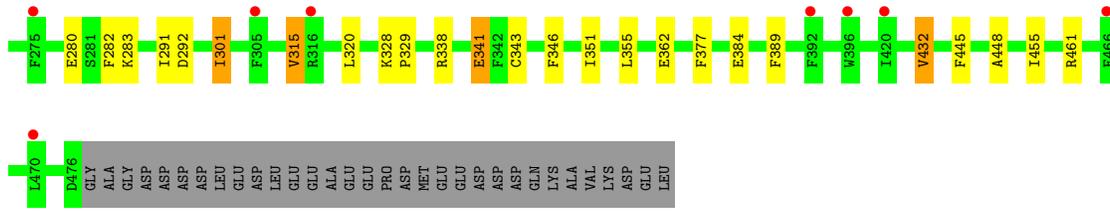


- Molecule 1: Prolyl 4-hydroxylase subunit alpha-2

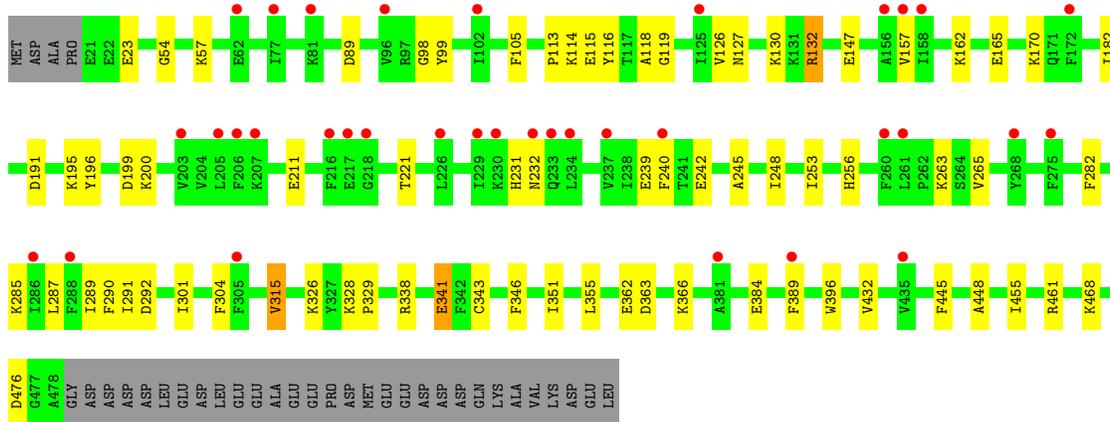
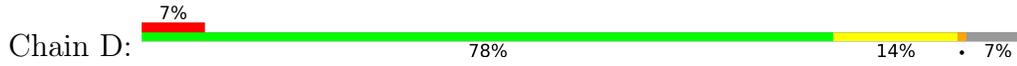


- Molecule 2: Protein disulfide-isomerase





• Molecule 2: Protein disulfide-isomerase



4 Data and refinement statistics i

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	252.88Å 252.88Å 89.40Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.69 – 3.85 46.69 – 3.85	Depositor EDS
% Data completeness (in resolution range)	99.7 (46.69-3.85) 99.7 (46.69-3.85)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.34 (at 3.88Å)	Xtrriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.244 , 0.278 0.263 , 0.299	Depositor DCC
R_{free} test set	1992 reflections (9.89%)	wwPDB-VP
Wilson B-factor (Å ²)	220.6	Xtrriage
Anisotropy	0.426	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	(Not available) , (Not available)	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	0.058 for h,-h-k,-l	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11098	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.41% 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

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	0.48	0/1970	0.67	1/2665 (0.0%)
1	B	0.51	0/1950	0.68	0/2637
2	C	0.37	0/3715	0.55	0/5016
2	D	0.39	0/3724	0.56	0/5028
All	All	0.42	0/11359	0.60	1/15346 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	314	HIS	C-N-CA	5.08	134.40	121.70

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	1920	0	1853	32	0
1	B	1897	0	1832	42	0
2	C	3631	0	3549	41	0
2	D	3640	0	3557	50	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	11098	0	10791	152	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.

The worst 5 of 152 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:469:PHE:HB3	1:B:472:LEU:HD13	1.36	1.07
2:C:263:LYS:HZ2	2:C:292:ASP:HB2	1.37	0.89
1:B:361:LEU:HD21	1:B:420:ASN:HB3	1.55	0.89
1:A:333:TRP:NE1	1:A:474:ALA:HB2	1.91	0.84
2:D:263:LYS:HZ2	2:D:292:ASP:HB2	1.43	0.83

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	233/262 (89%)	221 (95%)	11 (5%)	1 (0%)	34 70
1	B	230/262 (88%)	217 (94%)	12 (5%)	1 (0%)	34 70
2	C	454/492 (92%)	445 (98%)	7 (2%)	2 (0%)	34 70
2	D	456/492 (93%)	441 (97%)	13 (3%)	2 (0%)	34 70
All	All	1373/1508 (91%)	1324 (96%)	43 (3%)	6 (0%)	34 70

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	298	GLY
2	C	119	GLY

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Mol	Chain	Res	Type
2	D	119	GLY
2	C	448	ALA
2	D	448	ALA

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.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	202/226 (89%)	187 (93%)	15 (7%)	13	43
1	B	200/226 (88%)	181 (90%)	19 (10%)	8	33
2	C	388/419 (93%)	366 (94%)	22 (6%)	20	50
2	D	388/419 (93%)	367 (95%)	21 (5%)	22	51
All	All	1178/1290 (91%)	1101 (94%)	77 (6%)	17	46

5 of 77 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	89	ASP
2	D	341	GLU
2	D	132	ARG
2	D	242	GLU
2	D	389	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
2	D	231	HIS
2	D	232	ASN
2	D	367	GLN
1	B	314	HIS
2	C	231	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	$\# Z > 2$
3	SO4	A	601	-	4,4,4	0.38	0	6,6,6	0.32	0
3	SO4	B	601	-	4,4,4	0.56	0	6,6,6	0.37	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

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	237/262 (90%)	0.43	6 (2%) 57 47	10, 50, 81, 105	0
1	B	233/262 (88%)	0.36	13 (5%) 24 20	3, 44, 75, 95	0
2	C	456/492 (92%)	0.28	42 (9%) 9 7	57, 89, 115, 132	0
2	D	458/492 (93%)	0.26	35 (7%) 13 10	52, 86, 117, 158	0
All	All	1384/1508 (91%)	0.31	96 (6%) 16 12	3, 78, 113, 158	0

The worst 5 of 96 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	216	PHE	11.0
2	D	156	ALA	10.1
2	C	218	GLY	9.6
2	D	233	GLN	8.1
2	C	225	LEU	7.8

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 monosaccharides 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(\AA^2)	Q<0.9
3	SO4	A	601	5/5	0.84	0.69	23,26,27,28	0
3	SO4	B	601	5/5	0.95	0.60	17,21,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.