



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 27, 2022 – 03:16 PM EDT

PDB ID : 7S4A
Title : MRG15 complex with PALB2 peptide
Authors : Korolev, S.; Deveryshetty, J.
Deposited on : 2021-09-08
Resolution : 2.69 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.28.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.28.1

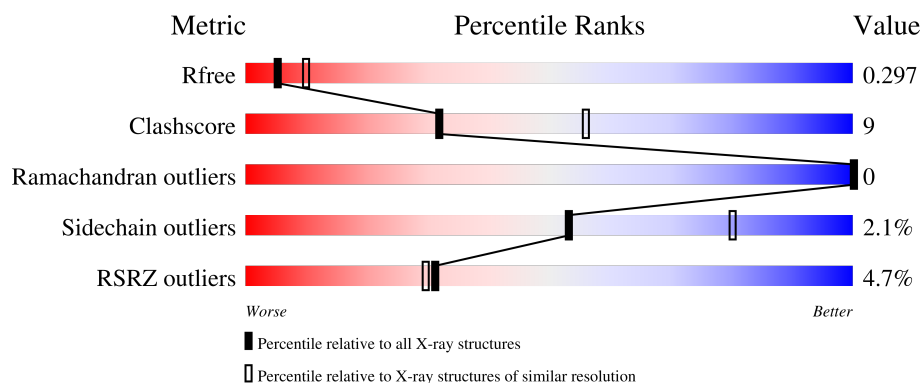
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.69 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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	172	<div> <div>3%</div> <div>70%</div> <div>23%</div> <div>8%</div> </div>
1	C	172	<div> <div>6%</div> <div>74%</div> <div>19%</div> <div>7%</div> </div>
2	B	57	<div> <div>2%</div> <div>47%</div> <div>9%</div> <div>44%</div> </div>
2	D	57	<div> <div>2%</div> <div>53%</div> <div>7%</div> <div>39%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3148 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 Mortality factor 4-like protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	159	Total	C	N	O	S	0	0	0
			1302	856	208	235	3			
1	C	160	Total	C	N	O	S	0	0	0
			1297	853	205	236	3			

- Molecule 2 is a protein called Partner and localizer of BRCA2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	32	Total	C	N	O	S	0	0	0
			259	172	38	48	1			
2	D	35	Total	C	N	O	S	0	0	0
			275	181	41	52	1			

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	574	MET	-	initiating methionine	UNP Q86YC2
B	575	HIS	-	expression tag	UNP Q86YC2
B	576	HIS	-	expression tag	UNP Q86YC2
B	577	HIS	-	expression tag	UNP Q86YC2
B	578	HIS	-	expression tag	UNP Q86YC2
B	579	HIS	-	expression tag	UNP Q86YC2
B	580	HIS	-	expression tag	UNP Q86YC2
B	581	SER	-	expression tag	UNP Q86YC2
B	582	SER	-	expression tag	UNP Q86YC2
B	583	GLY	-	expression tag	UNP Q86YC2
B	584	VAL	-	expression tag	UNP Q86YC2
B	585	ASP	-	expression tag	UNP Q86YC2
B	586	LEU	-	expression tag	UNP Q86YC2
B	587	GLY	-	expression tag	UNP Q86YC2
B	588	THR	-	expression tag	UNP Q86YC2
B	589	GLU	-	expression tag	UNP Q86YC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	590	ASN	-	expression tag	UNP Q86YC2
B	591	LEU	-	expression tag	UNP Q86YC2
B	592	TYR	-	expression tag	UNP Q86YC2
B	593	PHE	-	expression tag	UNP Q86YC2
B	594	GLN	-	expression tag	UNP Q86YC2
B	595	SER	-	expression tag	UNP Q86YC2
B	596	ASN	-	expression tag	UNP Q86YC2
D	574	MET	-	initiating methionine	UNP Q86YC2
D	575	HIS	-	expression tag	UNP Q86YC2
D	576	HIS	-	expression tag	UNP Q86YC2
D	577	HIS	-	expression tag	UNP Q86YC2
D	578	HIS	-	expression tag	UNP Q86YC2
D	579	HIS	-	expression tag	UNP Q86YC2
D	580	HIS	-	expression tag	UNP Q86YC2
D	581	SER	-	expression tag	UNP Q86YC2
D	582	SER	-	expression tag	UNP Q86YC2
D	583	GLY	-	expression tag	UNP Q86YC2
D	584	VAL	-	expression tag	UNP Q86YC2
D	585	ASP	-	expression tag	UNP Q86YC2
D	586	LEU	-	expression tag	UNP Q86YC2
D	587	GLY	-	expression tag	UNP Q86YC2
D	588	THR	-	expression tag	UNP Q86YC2
D	589	GLU	-	expression tag	UNP Q86YC2
D	590	ASN	-	expression tag	UNP Q86YC2
D	591	LEU	-	expression tag	UNP Q86YC2
D	592	TYR	-	expression tag	UNP Q86YC2
D	593	PHE	-	expression tag	UNP Q86YC2
D	594	GLN	-	expression tag	UNP Q86YC2
D	595	SER	-	expression tag	UNP Q86YC2
D	596	ASN	-	expression tag	UNP Q86YC2

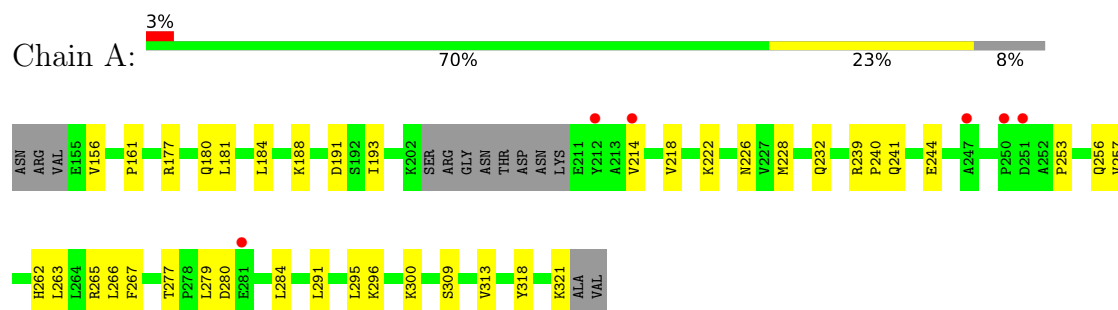
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	4	Total O 4 4	0	0
3	B	2	Total O 2 2	0	0
3	C	8	Total O 8 8	0	0
3	D	1	Total O 1 1	0	0

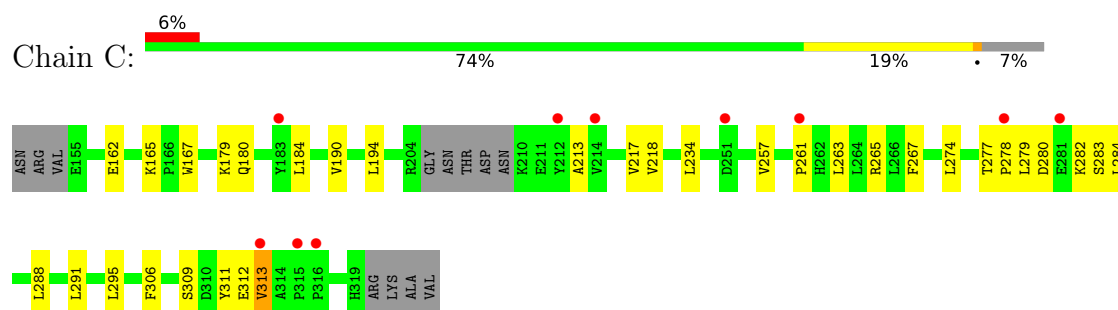
3 Residue-property plots [i](#)

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.

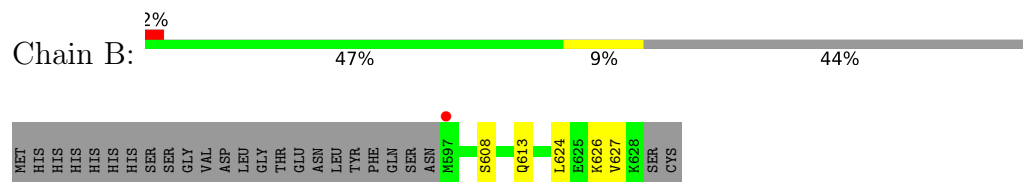
- Molecule 1: Mortality factor 4-like protein 1



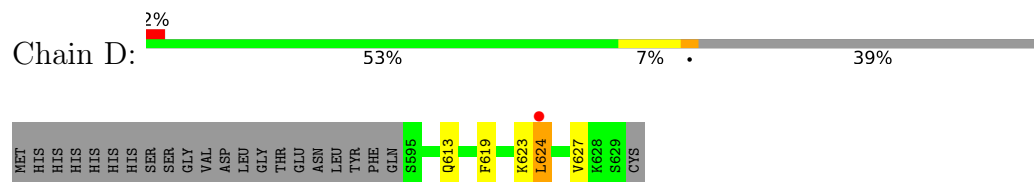
- Molecule 1: Mortality factor 4-like protein 1



- Molecule 2: Partner and localizer of BRCA2



- Molecule 2: Partner and localizer of BRCA2



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2 ₁	Depositor
Cell constants a, b, c, α , β , γ	54.08 Å 60.16 Å 131.85 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.67 – 2.69 29.67 – 2.69	Depositor EDS
% Data completeness (in resolution range)	98.4 (29.67-2.69) 88.6 (29.67-2.69)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.53 (at 2.68 Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, R_{free}	0.227 , 0.297 0.227 , 0.297	Depositor DCC
R_{free} test set	1228 reflections (10.01%)	wwPDB-VP
Wilson B-factor (Å ²)	44.5	Xtriage
Anisotropy	0.306	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 46.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3148	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 23.93 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.2336e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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](#)

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	0.43	0/1337	0.63	0/1817
1	C	0.48	0/1332	0.63	0/1813
2	B	0.52	0/263	0.70	0/352
2	D	0.56	0/279	0.77	0/374
All	All	0.47	0/3211	0.65	0/4356

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 [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	1302	0	1297	31	0
1	C	1297	0	1277	20	0
2	B	259	0	274	6	0
2	D	275	0	283	4	0
3	A	4	0	0	0	0
3	B	2	0	0	1	0
3	C	8	0	0	0	0
3	D	1	0	0	0	0
All	All	3148	0	3131	56	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 9.

All (56) 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:C:282:LYS:HD3	1:C:282:LYS:C	1.78	1.02
1:C:282:LYS:HD3	1:C:283:SER:N	1.83	0.93
1:A:253:PRO:HD2	1:A:256:GLN:HE21	1.36	0.90
1:A:180:GLN:HE21	1:A:313:VAL:HG12	1.45	0.79
1:A:180:GLN:HE21	1:A:313:VAL:CG1	1.99	0.74
1:A:277:THR:HG22	1:A:279:LEU:H	1.54	0.71
2:B:626:LYS:NZ	3:B:701:HOH:O	2.23	0.70
1:A:180:GLN:NE2	1:A:313:VAL:CG1	2.56	0.69
1:C:194:LEU:HB3	1:C:218:VAL:HG13	1.75	0.67
1:A:180:GLN:NE2	1:A:313:VAL:HG13	2.12	0.64
1:A:267:PHE:HE2	1:A:291:LEU:HD22	1.62	0.64
1:C:282:LYS:C	1:C:282:LYS:CD	2.62	0.63
2:B:627:VAL:HB	2:D:624:LEU:HD11	1.84	0.60
1:A:253:PRO:HD2	1:A:256:GLN:NE2	2.16	0.57
1:C:267:PHE:HE2	1:C:291:LEU:HD22	1.72	0.55
2:B:624:LEU:HD13	2:D:627:VAL:HG21	1.90	0.54
1:C:167:TRP:CE2	1:C:306:PHE:HB3	2.42	0.54
1:A:161:PRO:HB3	1:A:296:LYS:HE2	1.91	0.53
1:A:188:LYS:HG2	1:A:193:ILE:HD11	1.92	0.52
1:C:162:GLU:OE1	1:C:165:LYS:HE2	2.10	0.51
1:C:162:GLU:HA	1:C:165:LYS:HG3	1.93	0.50
1:A:184:LEU:HD22	1:A:257:VAL:HG12	1.94	0.50
1:A:240:PRO:O	1:A:244:GLU:HG3	2.11	0.50
1:C:213:ALA:O	1:C:217:VAL:HG13	2.12	0.49
1:A:222:LYS:HG2	1:A:226:ASN:ND2	2.27	0.49
1:A:161:PRO:HD3	1:A:295:LEU:HD13	1.94	0.48
1:C:184:LEU:HD22	1:C:257:VAL:HG12	1.95	0.48
1:A:181:LEU:HD11	1:A:318:TYR:CE1	2.49	0.48
1:A:262:HIS:CD2	1:A:265:ARG:NH1	2.82	0.47
1:A:240:PRO:HB2	1:A:318:TYR:CE2	2.49	0.47
1:C:261:PRO:O	1:C:265:ARG:HG3	2.15	0.47
2:B:613:GLN:OE1	2:D:613:GLN:HB3	2.14	0.47
1:A:191:ASP:OD2	1:A:256:GLN:HG2	2.15	0.47
1:A:181:LEU:HD11	1:A:318:TYR:CZ	2.51	0.46
1:A:177:ARG:NH2	2:B:627:VAL:HG13	2.31	0.46
1:A:228:MET:HG2	1:A:232:GLN:NE2	2.31	0.46
1:C:190:VAL:HG22	1:C:263:LEU:HB2	1.97	0.46
1:C:278:PRO:C	1:C:279:LEU:HD23	2.36	0.45
2:D:619:PHE:O	2:D:623:LYS:HG3	2.17	0.45
1:C:274:LEU:O	1:C:277:THR:OG1	2.34	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:232:GLN:OE1	2:B:608:SER:HB2	2.17	0.45
1:A:188:LYS:HE2	1:A:193:ILE:HD13	1.99	0.45
1:A:321:LYS:HB3	1:A:321:LYS:HE2	1.85	0.44
1:A:156:VAL:HG11	1:A:284:LEU:HD21	1.99	0.44
1:A:296:LYS:O	1:A:300:LYS:HG2	2.18	0.44
1:A:214:VAL:O	1:A:218:VAL:HG22	2.17	0.43
1:C:280:ASP:OD1	1:C:283:SER:OG	2.28	0.43
1:A:280:ASP:OD2	1:A:280:ASP:N	2.52	0.43
1:A:184:LEU:HD12	1:A:241:GLN:CD	2.39	0.43
1:C:180:GLN:HE21	1:C:313:VAL:HG12	1.83	0.42
1:C:295:LEU:HD23	1:C:295:LEU:HA	1.91	0.42
1:C:311:TYR:O	1:C:312:GLU:HG3	2.20	0.42
1:A:180:GLN:NE2	1:A:313:VAL:HG12	2.19	0.41
1:C:284:LEU:O	1:C:288:LEU:HG	2.21	0.41
1:C:190:VAL:HG13	1:C:263:LEU:HD13	2.03	0.40
1:A:263:LEU:O	1:A:266:LEU:HB3	2.21	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	155/172 (90%)	152 (98%)	3 (2%)	0	100	100
1	C	156/172 (91%)	149 (96%)	7 (4%)	0	100	100
2	B	30/57 (53%)	29 (97%)	1 (3%)	0	100	100
2	D	33/57 (58%)	32 (97%)	1 (3%)	0	100	100
All	All	374/458 (82%)	362 (97%)	12 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	138/151 (91%)	136 (99%)	2 (1%)	67	86
1	C	136/151 (90%)	132 (97%)	4 (3%)	42	71
2	B	31/54 (57%)	31 (100%)	0	100	100
2	D	32/54 (59%)	31 (97%)	1 (3%)	40	69
All	All	337/410 (82%)	330 (98%)	7 (2%)	53	80

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	239	ARG
1	A	309	SER
1	C	179	LYS
1	C	234	LEU
1	C	309	SER
1	C	313	VAL
2	D	624	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	180	GLN
1	A	241	GLN
1	A	256	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 [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	159/172 (92%)	0.18	6 (3%) 40 39	33, 54, 73, 81	0
1	C	160/172 (93%)	0.27	10 (6%) 20 19	32, 48, 74, 94	0
2	B	32/57 (56%)	0.04	1 (3%) 49 49	35, 49, 69, 79	0
2	D	35/57 (61%)	-0.04	1 (2%) 51 52	41, 51, 59, 68	0
All	All	386/458 (84%)	0.18	18 (4%) 31 30	32, 51, 73, 94	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	597	MET	4.4
1	C	281	GLU	3.6
1	C	316	PRO	3.3
1	C	313	VAL	3.0
1	A	212	TYR	2.9
1	C	251	ASP	2.7
1	A	250	PRO	2.6
1	A	247	ALA	2.5
1	A	214	VAL	2.4
1	C	212	TYR	2.3
1	A	251	ASP	2.2
1	C	315	PRO	2.2
1	C	214	VAL	2.2
2	D	624	LEU	2.2
1	C	183	TYR	2.1
1	A	281	GLU	2.1
1	C	278	PRO	2.1
1	C	261	PRO	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 monosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.