



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 07:22 pm GMT

PDB ID : 4MS8
Title : 42F3 TCR pCPB9/H-2Ld Complex
Authors : Birnbaum, M.E.; Adams, J.J.; Garcia, K.C.
Deposited on : 2013-09-18
Resolution : 1.92 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

<http://wwpdb.org/validation/2016/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) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

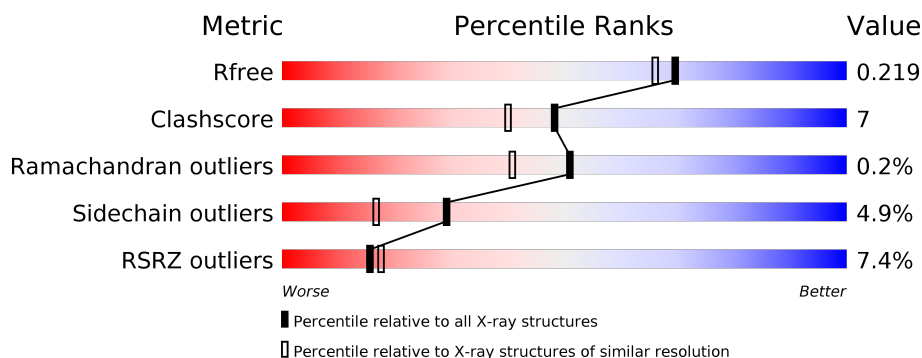
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.92 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}	100719	6276 (1.94-1.90)
Clashscore	112137	7025 (1.94-1.90)
Ramachandran outliers	110173	6947 (1.94-1.90)
Sidechain outliers	110143	6948 (1.94-1.90)
RSRZ outliers	101464	6332 (1.94-1.90)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	C	212	<div> <div>10%</div> <div> <div>81%</div> <div>10%</div> <div>8%</div> </div> </div>
2	D	243	<div> <div>%</div> <div> <div>84%</div> <div>13%</div> <div>••</div> </div> </div>
3	A	180	<div> <div>12%</div> <div> <div>70%</div> <div>18%</div> <div>•</div> <div>10%</div> </div> </div>
4	B	9	<div> <div>56%</div> <div>33%</div> <div>11%</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5275 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 42F3 alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	196	Total	C	N	O	S	0	0	0
			1517	961	249	299	8			

- Molecule 2 is a protein called 42F3 beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	239	Total	C	N	O	S	0	0	0
			1895	1194	329	366	6			

- Molecule 3 is a protein called H-2 class I histocompatibility antigen, L-D alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	162	Total	C	N	O	S	0	0	0
			1335	846	230	252	7			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	EXPRESSION TAG	UNP P01897
A	8	TYR	PHE	ENGINEERED MUTATION	UNP P01897
A	12	THR	VAL	ENGINEERED MUTATION	UNP P01897
A	15	ARG	PRO	ENGINEERED MUTATION	UNP P01897
A	23	THR	ILE	ENGINEERED MUTATION	UNP P01897
A	30	ASP	ASN	ENGINEERED MUTATION	UNP P01897
A	49	VAL	ALA	ENGINEERED MUTATION	UNP P01897
A	131	ARG	LYS	ENGINEERED MUTATION	UNP P01897

- Molecule 4 is a protein called pCPB9.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	B	9	Total	C	N	O	0	1	0
			72	49	9	14			

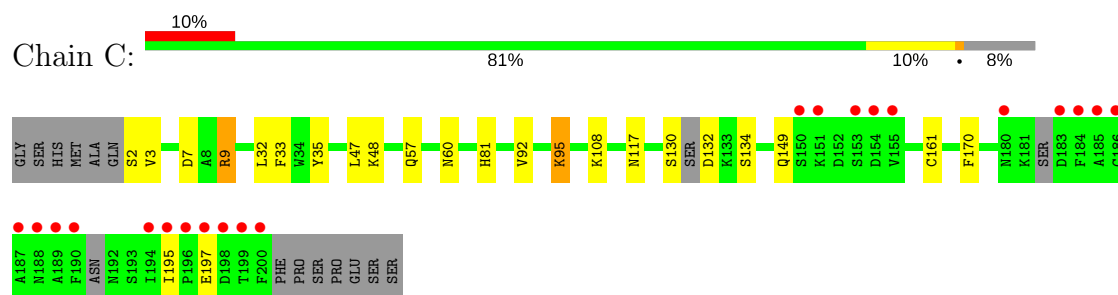
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	164	Total 164	O 164	0	0
5	D	213	Total 213	O 213	0	0
5	A	74	Total 74	O 74	0	0
5	B	5	Total 5	O 5	0	0

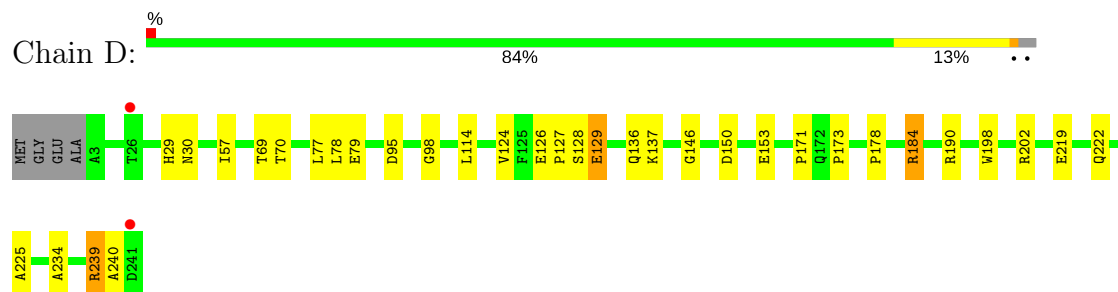
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

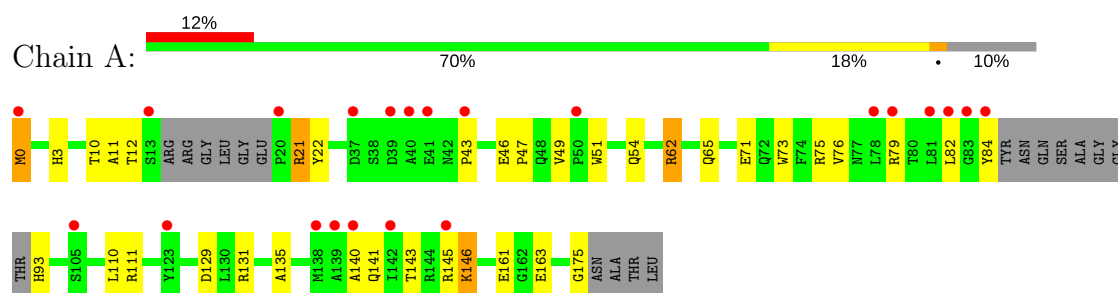
• Molecule 1: 42F3 alpha



• Molecule 2: 42F3 beta



• Molecule 3: H-2 class I histocompatibility antigen, L-D alpha chain



• Molecule 4: pCPB9



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	175.26Å 60.62Å 70.10Å 90.00° 96.06° 90.00°	Depositor
Resolution (Å)	34.85 – 1.92 43.57 – 1.92	Depositor EDS
% Data completeness (in resolution range)	93.1 (34.85-1.92) 93.1 (43.57-1.92)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.32 (at 1.92Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.180 , 0.219 0.179 , 0.219	Depositor DCC
R_{free} test set	2655 reflections (5.10%)	DCC
Wilson B-factor (Å ²)	31.0	Xtriage
Anisotropy	0.134	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 57.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5275	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.86% 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](#)

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	C	0.81	1/1551 (0.1%)	0.81	0/2098
2	D	0.77	0/1947	0.80	1/2655 (0.0%)
3	A	0.64	0/1373	0.69	0/1861
4	B	0.62	0/77	0.56	0/103
All	All	0.75	1/4948 (0.0%)	0.77	1/6717 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	35	TYR	CD1-CE1	5.67	1.47	1.39

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	184	ARG	NE-CZ-NH2	-5.86	117.37	120.30

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	C	1517	0	1439	14	0
2	D	1895	0	1791	26	0
3	A	1335	0	1216	26	0
4	B	72	0	68	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	74	0	0	12	0
5	B	5	0	0	0	0
5	C	164	0	0	9	1
5	D	213	0	0	13	2
All	All	5275	0	4514	64	2

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.

All (64) 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:130:SER:HG	1:C:132:ASP:N	1.62	0.96
1:C:9:ARG:NH2	5:C:435:HOH:O	2.00	0.93
2:D:136:GLN:O	5:D:374:HOH:O	1.95	0.84
3:A:84:TYR:O	5:A:241:HOH:O	2.00	0.80
3:A:111:ARG:NH1	5:A:237:HOH:O	2.16	0.78
5:C:464:HOH:O	4:B:4[A]:GLU:HG2	1.84	0.78
3:A:146:LYS:NZ	5:A:247:HOH:O	2.25	0.70
2:D:240:ALA:O	5:D:483:HOH:O	2.08	0.70
2:D:219:GLU:OE2	5:D:356:HOH:O	2.09	0.69
2:D:129:GLU:HG3	5:D:479:HOH:O	1.93	0.69
2:D:153:GLU:OE2	5:D:493:HOH:O	2.12	0.67
1:C:108:LYS:NZ	5:C:414:HOH:O	2.27	0.65
2:D:95:ASP:OD2	5:D:477:HOH:O	2.14	0.65
3:A:129:ASP:O	3:A:131:ARG:NH1	2.30	0.64
1:C:7:ASP:HB2	5:C:390:HOH:O	1.96	0.64
2:D:222:GLN:NE2	5:D:449:HOH:O	2.27	0.62
2:D:150:ASP:OD1	2:D:173:PRO:HG2	2.00	0.62
1:C:60:ASN:OD1	5:C:455:HOH:O	2.17	0.60
2:D:225:ALA:O	5:D:395:HOH:O	2.17	0.57
2:D:77:LEU:HG	2:D:79:GLU:HG3	1.86	0.57
2:D:222:GLN:O	5:D:464:HOH:O	2.17	0.57
3:A:12:THR:HA	3:A:93:HIS:O	2.04	0.57
1:C:95:LYS:HE3	5:C:400:HOH:O	2.04	0.57
2:D:146:GLY:O	2:D:184:ARG:HD3	2.05	0.57
3:A:161:GLU:OE1	5:A:257:HOH:O	2.17	0.56
3:A:79:ARG:HA	3:A:82:LEU:HD12	1.86	0.56
3:A:143:THR:HG21	4:B:9:LEU:HD23	1.89	0.54
1:C:57:GLN:HG2	5:C:356:HOH:O	2.08	0.53
3:A:141:GLN:O	5:A:245:HOH:O	2.19	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:69:THR:HG23	2:D:70:THR:HG23	1.92	0.51
1:C:60:ASN:ND2	5:C:459:HOH:O	2.27	0.51
1:C:32:LEU:HD23	1:C:92:VAL:HG12	1.93	0.51
3:A:65:GLN:HG3	5:A:258:HOH:O	2.11	0.50
3:A:46:GLU:HG3	3:A:47:PRO:HD2	1.92	0.50
3:A:145:ARG:HA	5:A:265:HOH:O	2.11	0.49
1:C:33:PHE:CZ	2:D:98:GLY:HA3	2.48	0.48
3:A:0:MET:SD	3:A:3:HIS:NE2	2.87	0.48
3:A:21:ARG:O	3:A:21:ARG:HG3	2.12	0.48
3:A:11:ALA:HA	3:A:22:TYR:HA	1.95	0.48
3:A:62:ARG:NH2	5:A:234:HOH:O	2.46	0.47
3:A:73:TRP:NE1	4:B:7:PHE:O	2.35	0.47
4:B:4[A]:GLU:HG3	4:B:6:GLY:N	2.30	0.47
1:C:195:ILE:H	1:C:195:ILE:HD12	1.80	0.47
2:D:126:GLU:OE1	2:D:239:ARG:NH1	2.47	0.47
2:D:129:GLU:HB2	5:D:463:HOH:O	2.13	0.47
3:A:163:GLU:HG3	5:A:224:HOH:O	2.15	0.46
3:A:62:ARG:CZ	5:A:234:HOH:O	2.62	0.46
3:A:71:GLU:O	3:A:75:ARG:HG3	2.16	0.45
1:C:57:GLN:CG	5:C:356:HOH:O	2.65	0.45
2:D:30:ASN:OD1	3:A:76:VAL:HG21	2.16	0.44
3:A:175:GLY:O	5:A:244:HOH:O	2.21	0.43
3:A:49:VAL:HG21	3:A:51:TRP:CE2	2.53	0.43
3:A:10:THR:OG1	5:A:208:HOH:O	2.21	0.43
3:A:135:ALA:HB1	3:A:140:ALA:CB	2.49	0.43
2:D:178:PRO:HG3	5:D:468:HOH:O	2.19	0.43
2:D:29:HIS:CD2	2:D:95:ASP:HB3	2.52	0.43
2:D:202:ARG:NH1	5:D:465:HOH:O	2.41	0.43
2:D:78:LEU:HD12	2:D:78:LEU:N	2.35	0.41
2:D:171:PRO:HA	5:D:510:HOH:O	2.21	0.41
2:D:124:VAL:HG23	2:D:234:ALA:HB3	2.02	0.41
1:C:170:PHE:CE2	2:D:137:LYS:HE2	2.55	0.41
3:A:135:ALA:HB1	3:A:140:ALA:HB1	2.01	0.41
2:D:127:PRO:HD2	2:D:198:TRP:CZ2	2.55	0.40
1:C:170:PHE:CD2	2:D:137:LYS:HE2	2.56	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:D:504:HOH:O	5:D:507:HOH:O[4_546]	1.83	0.37

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:C:455:HOH:O	5:D:446:HOH:O[4_546]	1.98	0.22

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	C	188/212 (89%)	184 (98%)	4 (2%)	0	100	100
2	D	237/243 (98%)	233 (98%)	4 (2%)	0	100	100
3	A	156/180 (87%)	149 (96%)	6 (4%)	1 (1%)	28	16
4	B	8/9 (89%)	8 (100%)	0	0	100	100
All	All	589/644 (92%)	574 (98%)	14 (2%)	1 (0%)	51	41

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	43	PRO

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	C	169/186 (91%)	157 (93%)	12 (7%)	17	7
2	D	206/208 (99%)	200 (97%)	6 (3%)	48	37
3	A	131/148 (88%)	125 (95%)	6 (5%)	31	19

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	B	7/6 (117%)	6 (86%)	1 (14%)	4	1
All	All	513/548 (94%)	488 (95%)	25 (5%)	29	17

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	2	SER
1	C	3	VAL
1	C	9	ARG
1	C	47	LEU
1	C	48	LYS
1	C	81	HIS
1	C	95	LYS
1	C	117	ASN
1	C	134	SER
1	C	149	GLN
1	C	161	CYS
1	C	197	GLU
2	D	57	ILE
2	D	114	LEU
2	D	128	SER
2	D	129	GLU
2	D	190	ARG
2	D	239	ARG
3	A	0	MET
3	A	21	ARG
3	A	54	GLN
3	A	62	ARG
3	A	110	LEU
3	A	146	LYS
4	B	9	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	124	GLN

5.3.3 RNA ⓘ

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 carbohydrates 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	C	196/212 (92%)	0.20	21 (10%) 7 8	23, 40, 73, 96	0
2	D	239/243 (98%)	-0.06	2 (0%) 86 88	24, 36, 63, 88	0
3	A	162/180 (90%)	0.70	22 (13%) 3 4	31, 59, 94, 117	0
4	B	9/9 (100%)	0.22	0 100 100	42, 47, 58, 61	0
All	All	606/644 (94%)	0.23	45 (7%) 15 17	23, 42, 80, 117	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	A	81	LEU	7.6
3	A	142	ILE	5.3
1	C	196	PRO	5.1
1	C	194	ILE	4.9
1	C	200	PHE	4.8
3	A	138	MET	4.3
1	C	189	ALA	4.1
1	C	184	PHE	3.8
1	C	186	CYS	3.8
3	A	40	ALA	3.7
3	A	82	LEU	3.7
1	C	151	LYS	3.6
1	C	187	ALA	3.4
2	D	26	THR	3.3
3	A	83	GLY	3.2
1	C	183	ASP	3.2
3	A	43	PRO	3.2
1	C	198	ASP	3.2
3	A	140	ALA	3.1
3	A	37	ASP	3.0
2	D	241	ASP	3.0

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Mol	Chain	Res	Type	RSRZ
3	A	39	ASP	3.0
1	C	180	ASN	2.9
1	C	197	GLU	2.9
3	A	123	TYR	2.9
1	C	155	VAL	2.8
1	C	188	ASN	2.8
1	C	150	SER	2.7
1	C	199	THR	2.6
3	A	84	TYR	2.6
3	A	0	MET	2.5
1	C	190	PHE	2.5
3	A	41	GLU	2.5
3	A	20	PRO	2.4
3	A	13	SER	2.4
3	A	50	PRO	2.4
1	C	185	ALA	2.3
3	A	145	ARG	2.3
1	C	195	ILE	2.2
1	C	153	SER	2.2
3	A	105	SER	2.2
3	A	139	ALA	2.2
3	A	78	LEU	2.2
3	A	79	ARG	2.1
1	C	154	ASP	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 carbohydrates 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.