



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 02:44 AM UTC

PDB ID : 4V1D / pdb_00004v1d
Title : Ternary complex among two human derived single chain antibody fragments and Cn2 toxin from scorpion *Centruroides noxi*.
Authors : Riano-Umbarila, L.; Serrano-Posada, H.; Rojas-Trejo, S.; Rudino-Pinera, E.; Becerril, B.
Deposited on : 2014-09-25
Resolution : 3.10 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
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

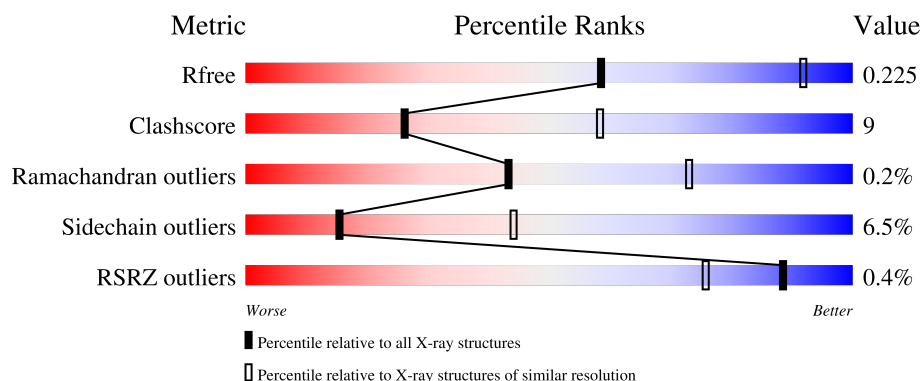
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	117	 85% 15%
2	B	146	 60% 14% 25%
3	C	66	 2% 77% 21%
4	D	124	 76% 21%
5	E	149	 50% 19% 5% 25%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 4046 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 SINGLE CHAIN ANTIBODY FRAGMENT LR, HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	117	Total	C	N	O	S	0	1	0
			894	553	157	179	5			

- Molecule 2 is a protein called SINGLE CHAIN ANTIBODY FRAGMENT LR, LIGHT CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	109	Total	C	N	O	S	0	1	1
			847	530	150	165	2			

- Molecule 3 is a protein called BETA-MAMMAL TOXIN CN2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	66	Total	C	N	O	S	0	0	1
			524	333	89	94	8			

- Molecule 4 is a protein called SINGLE CHAIN ANTIBODY FRAGMENT RU1, HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	124	Total	C	N	O	S	0	0	0
			953	598	170	179	6			

- Molecule 5 is a protein called SINGLE CHAIN ANTIBODY FRAGMENT RU1, LIGHT CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	112	Total	C	N	O	S	0	0	1
			822	506	143	170	3			

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	3	Total 3	O 3	0	0
6	D	1	Total 1	O 1	0	0
6	E	2	Total 2	O 2	0	0

3 Residue-property plots

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.

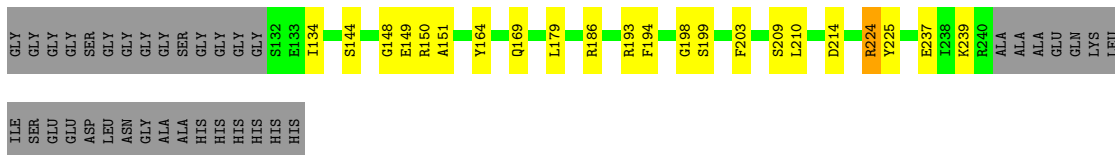
- Molecule 1: SINGLE CHAIN ANTIBODY FRAGMENT LR, HEAVY CHAIN

Chain A: 



- Molecule 2: SINGLE CHAIN ANTIBODY FRAGMENT LR, LIGHT CHAIN

Chain B: 



- Molecule 3: BETA-MAMMAL TOXIN CN2

Chain C: 



- Molecule 4: SINGLE CHAIN ANTIBODY FRAGMENT RU1, HEAVY CHAIN

Chain D: 



- Molecule 5: SINGLE CHAIN ANTIBODY FRAGMENT RU1, LIGHT CHAIN

Chain E: 



S234	L235	L236	L246	T247	V248	L249	G250	ALA	ALA	ALA	ALA	GLU	GLN	LYS	LEU	ILE	SER	GLU	GLU	ASP	LEU	ASN	GLY	ALA	ALA	HIS	HIS	HIS	HIS	HIS
------	------	------	------	------	------	------	------	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	45.58Å 74.64Å 140.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.21 – 3.10 38.21 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (38.21-3.10) 99.8 (38.21-3.10)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.69 (at 3.12Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.185 , 0.229 0.185 , 0.225	Depositor DCC
R_{free} test set	439 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å ²)	44.5	Xtriage
Anisotropy	0.048	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 50.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4046	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.34% 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

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	0.32	0/911	0.78	0/1227
2	B	0.44	1/866 (0.1%)	0.92	0/1175
3	C	0.48	1/538 (0.2%)	0.87	1/725 (0.1%)
4	D	0.28	0/975	0.73	0/1316
5	E	0.64	1/841 (0.1%)	0.92	1/1146 (0.1%)
All	All	0.44	3/4131 (0.1%)	0.84	2/5589 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
5	E	0	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	65	CYS	C-N	-8.05	1.22	1.33
5	E	249	LEU	C-N	-7.22	1.23	1.33
2	B	239	LYS	C-N	-6.79	1.23	1.33

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	63	LYS	N-CA-C	5.72	117.73	108.41
5	E	196	SER	CB-CA-C	-5.13	110.25	117.23

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
5	E	194	ARG	Peptide
5	E	199	PRO	Peptide
5	E	233	ASP	Peptide

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	894	0	850	11	0
2	B	847	0	825	13	1
3	C	524	0	492	8	0
4	D	953	0	913	15	1
5	E	822	0	772	30	0
6	A	3	0	0	0	0
6	D	1	0	0	0	0
6	E	2	0	0	0	0
All	All	4046	0	3852	74	1

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 (74) 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:62:ASN:ND2	5:E:234:SER:OG	1.97	0.97
5:E:200:ASP:C	5:E:202:PHE:H	1.91	0.77
1:A:58:ILE:O	3:C:18:LYS:NZ	2.22	0.73
5:E:193:GLN:O	5:E:193:GLN:HG2	1.88	0.72
5:E:200:ASP:O	5:E:202:PHE:N	2.26	0.67
5:E:194:ARG:HD3	5:E:195:PRO:HD2	1.76	0.67
4:D:83:MET:HE2	4:D:86:LEU:HD21	1.75	0.66
5:E:179:LEU:HD22	5:E:224:ALA:HB2	1.79	0.65
5:E:193:GLN:HE21	5:E:198:VAL:HB	1.62	0.64
5:E:198:VAL:O	5:E:200:ASP:N	2.31	0.64
2:B:134:ILE:HD11	2:B:225:TYR:CD2	2.34	0.63
1:A:83:MET:HE2	1:A:86:LEU:HD21	1.81	0.61
5:E:198:VAL:C	5:E:200:ASP:H	2.09	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:2:GLU:HG2	3:C:49:THR:HA	1.84	0.59
4:D:67:ARG:NH1	4:D:90:ASP:OD2	2.35	0.59
4:D:91:THR:HG23	4:D:121:THR:HA	1.84	0.58
5:E:234:SER:HA	5:E:235:LEU:C	2.28	0.57
5:E:201:ARG:NH2	5:E:222:ASP:OD2	2.37	0.57
5:E:198:VAL:N	5:E:199:PRO:HD3	2.19	0.57
1:A:68:PHE:CZ	1:A:83:MET:HE3	2.41	0.55
2:B:151:ALA:HB2	2:B:210:LEU:HD11	1.88	0.55
5:E:199:PRO:HB2	5:E:201:ARG:HD3	1.89	0.55
5:E:188:ILE:HG12	5:E:204:ALA:HB2	1.89	0.54
2:B:186:ARG:NH2	2:B:194:PHE:O	2.39	0.54
1:A:13:GLN:H	1:A:13:GLN:CD	2.15	0.54
5:E:196:SER:O	5:E:198:VAL:N	2.41	0.53
2:B:134:ILE:HD11	2:B:225:TYR:HD2	1.72	0.53
2:B:164:TYR:HD2	2:B:224[B]:ARG:HB2	1.74	0.53
5:E:192:ASN:HB2	5:E:193:GLN:HB3	1.89	0.52
1:A:2:VAL:HG12	1:A:106:THR:HG21	1.90	0.52
4:D:101:ARG:NH2	4:D:108:ASP:OD2	2.42	0.52
3:C:10:THR:OG1	3:C:12:CYS:SG	2.55	0.52
4:D:106:CYS:HB3	4:D:108:ASP:OD1	2.11	0.51
5:E:223:GLU:HB3	5:E:248:VAL:HG23	1.91	0.51
4:D:105:LEU:H	4:D:105:LEU:HD23	1.74	0.51
4:D:22:CYS:HB3	4:D:79:LEU:HB3	1.91	0.51
2:B:144:SER:HA	2:B:237:GLU:O	2.09	0.51
5:E:153:PRO:HD3	5:E:250:GLY:N	2.26	0.51
2:B:169:GLN:HB2	2:B:179:LEU:HD11	1.92	0.51
2:B:193:ARG:NE	2:B:214:ASP:OD2	2.43	0.50
1:A:27:PHE:CE2	1:A:29:PHE:HA	2.46	0.50
4:D:93:VAL:HG22	4:D:119:LEU:HD23	1.93	0.49
5:E:199:PRO:CB	5:E:201:ARG:HD3	2.42	0.49
5:E:200:ASP:O	5:E:201:ARG:HB2	2.13	0.46
2:B:198:GLY:HA3	2:B:203:PHE:HA	1.98	0.46
2:B:164:TYR:HD2	2:B:224[A]:ARG:HG2	1.81	0.46
5:E:192:ASN:N	5:E:192:ASN:OD1	2.49	0.45
5:E:193:GLN:O	5:E:194:ARG:O	2.34	0.45
1:A:34:MET:HB3	1:A:79:LEU:HD22	1.98	0.45
4:D:63:SER:O	4:D:67:ARG:NH2	2.43	0.45
4:D:73:ASP:OD1	4:D:75:SER:OG	2.30	0.45
5:E:201:ARG:O	5:E:215:ILE:HA	2.17	0.45
4:D:97:ALA:HB1	4:D:111:PHE:HB3	1.97	0.45
4:D:20:LEU:HD12	4:D:81:LEU:HD23	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:224[B]:ARG:NH1	3:C:42:TYR:OH	2.49	0.44
3:C:11:GLY:HA2	3:C:60:LEU:HD12	1.99	0.44
4:D:2:VAL:HG12	4:D:113:LEU:HD11	2.00	0.44
5:E:200:ASP:N	5:E:200:ASP:OD1	2.51	0.44
1:A:2:VAL:HG13	1:A:27:PHE:CD1	2.53	0.43
1:A:91:THR:HG23	1:A:114:THR:HA	2.00	0.43
2:B:164:TYR:CD2	2:B:224[A]:ARG:HG2	2.52	0.43
4:D:4:LEU:HG	4:D:113:LEU:HD13	1.99	0.43
5:E:200:ASP:O	5:E:202:PHE:HD2	2.01	0.43
5:E:219:GLN:O	5:E:248:VAL:HG21	2.20	0.42
2:B:149:GLU:O	2:B:209:SER:HA	2.19	0.42
3:C:6:VAL:HB	3:C:58:TRP:HB2	2.02	0.42
1:A:60:TYR:HB2	1:A:65:LYS:HG3	2.01	0.42
1:A:6:GLU:CD	1:A:110:GLY:H	2.27	0.42
3:C:26:LEU:HD11	3:C:30:LYS:NZ	2.35	0.41
5:E:159:ILE:HD11	5:E:246:LEU:HD12	2.03	0.41
4:D:104:LEU:HD12	4:D:104:LEU:HA	1.76	0.41
5:E:196:SER:O	5:E:198:VAL:HG22	2.21	0.40
5:E:144:THR:OG1	5:E:162:SER:HB3	2.21	0.40
5:E:198:VAL:C	5:E:200:ASP:N	2.73	0.40

All (1) 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)
2:B:148:GLY:O	4:D:17:SER:OG[2_667]	2.18	0.02

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	116/117 (99%)	115 (99%)	1 (1%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	108/146 (74%)	106 (98%)	2 (2%)	0	100	100
3	C	64/66 (97%)	64 (100%)	0	0	100	100
4	D	122/124 (98%)	120 (98%)	2 (2%)	0	100	100
5	E	110/149 (74%)	96 (87%)	13 (12%)	1 (1%)	14	44
All	All	520/602 (86%)	501 (96%)	18 (4%)	1 (0%)	43	73

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
5	E	194	ARG

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	95/94 (101%)	94 (99%)	1 (1%)	65	78
2	B	92/111 (83%)	88 (96%)	4 (4%)	26	57
3	C	54/55 (98%)	53 (98%)	1 (2%)	50	73
4	D	98/98 (100%)	90 (92%)	8 (8%)	10	36
5	E	92/111 (83%)	77 (84%)	15 (16%)	2	11
All	All	431/469 (92%)	402 (93%)	29 (7%)	15	43

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	LEU
2	B	150	ARG
2	B	199	SER
2	B	224[A]	ARG
2	B	224[B]	ARG
3	C	65	CYS
4	D	21	SER

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Mol	Chain	Res	Type
4	D	58	LYS
4	D	89	GLU
4	D	102	ASP
4	D	105	LEU
4	D	113	LEU
4	D	119	LEU
4	D	121	THR
5	E	140	ASN
5	E	144	THR
5	E	158	THR
5	E	167	ASN
5	E	177	ARG
5	E	179	LEU
5	E	186	LEU
5	E	187	LEU
5	E	188	ILE
5	E	193	GLN
5	E	194	ARG
5	E	200	ASP
5	E	218	LEU
5	E	235	LEU
5	E	236	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	GLN
1	A	39	GLN
1	A	82	GLN
2	B	170	GLN
3	C	9	ASN
3	C	62	ASN
4	D	3	ASN
4	D	57	ASN

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 oligosaccharides 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 [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(Å²)	Q<0.9	
1	A	117/117 (100%)	-0.32	0	100	100	21, 43, 68, 93	1 (0%)
2	B	109/146 (74%)	-0.18	0	100	100	22, 46, 62, 73	1 (0%)
3	C	66/66 (100%)	-0.29	1 (1%)	72	52	27, 39, 60, 94	0
4	D	124/124 (100%)	-0.31	0	100	100	32, 47, 65, 87	0
5	E	112/149 (75%)	-0.12	1 (0%)	81	63	31, 43, 71, 82	0
All	All	528/602 (87%)	-0.24	2 (0%)	88	76	21, 45, 68, 94	2 (0%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	66	SER	2.3
5	E	194	ARG	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.