



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 05:14 AM UTC

PDB ID : 4ZNC / pdb_00004znc
Title : Fc fragment of human IgG in complex with the C domain of staphylococcal protein A mutant - Q9W
Authors : Deis, L.N.; Oas, T.G.
Deposited on : 2015-05-04
Resolution : 2.28 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

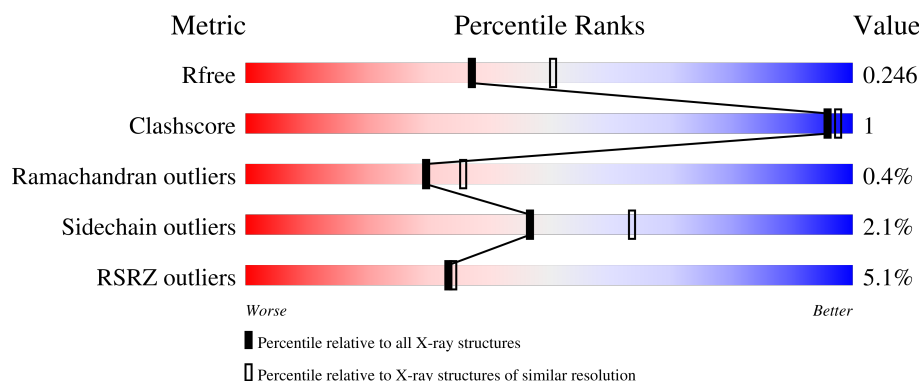
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

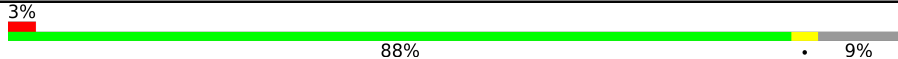
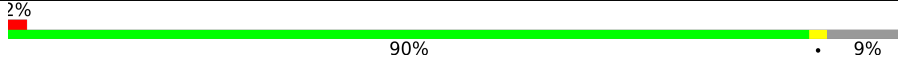
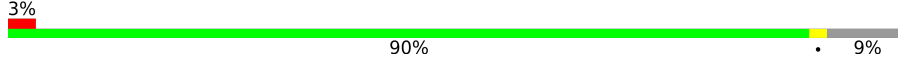
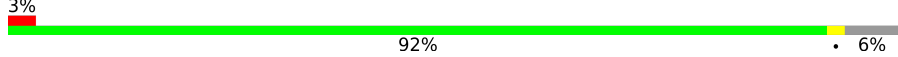
The reported resolution of this entry is 2.28 Å.

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	9078 (2.30-2.26)
Clashscore	190562	9802 (2.30-2.26)
Ramachandran outliers	187476	9690 (2.30-2.26)
Sidechain outliers	187428	9691 (2.30-2.26)
RSRZ outliers	180081	9085 (2.30-2.26)

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	58	
1	B	58	
1	C	58	
2	D	220	
2	E	220	

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Mol	Chain	Length	Quality of chain
2	F	220	10% 80% 7% 11%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12392 atoms, of which 6069 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 Immunoglobulin G-binding protein A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	53	Total	C	H	N	O	0	2	0
			892	286	439	78	89			
1	B	53	Total	C	H	N	O	0	0	0
			852	275	419	73	85			
1	C	53	Total	C	H	N	O	0	0	0
			852	275	419	73	85			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	TRP	GLN	engineered mutation	UNP P38507
B	9	TRP	GLN	engineered mutation	UNP P38507
C	9	TRP	GLN	engineered mutation	UNP P38507

- Molecule 2 is a protein called Ig gamma-3 chain C region.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	D	207	Total	C	H	N	O	S	0	0	0
			3290	1057	1629	278	318	8			
2	E	207	Total	C	H	N	O	S	0	0	0
			3290	1057	1629	278	318	8			
2	F	195	Total	C	H	N	O	S	0	0	0
			3090	993	1534	258	297	8			

There are 39 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	296	PHE	TYR	conflict	UNP P01860
D	435	HIS	ARG	conflict	UNP P01860
D	436	TYR	PHE	conflict	UNP P01860
D	448	GLY	-	expression tag	UNP P01860
D	449	SER	-	expression tag	UNP P01860

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Chain	Residue	Modelled	Actual	Comment	Reference
D	450	LEU	-	expression tag	UNP P01860
D	451	GLU	-	expression tag	UNP P01860
D	452	HIS	-	expression tag	UNP P01860
D	453	HIS	-	expression tag	UNP P01860
D	454	HIS	-	expression tag	UNP P01860
D	455	HIS	-	expression tag	UNP P01860
D	456	HIS	-	expression tag	UNP P01860
D	457	HIS	-	expression tag	UNP P01860
E	296	PHE	TYR	conflict	UNP P01860
E	435	HIS	ARG	conflict	UNP P01860
E	436	TYR	PHE	conflict	UNP P01860
E	448	GLY	-	expression tag	UNP P01860
E	449	SER	-	expression tag	UNP P01860
E	450	LEU	-	expression tag	UNP P01860
E	451	GLU	-	expression tag	UNP P01860
E	452	HIS	-	expression tag	UNP P01860
E	453	HIS	-	expression tag	UNP P01860
E	454	HIS	-	expression tag	UNP P01860
E	455	HIS	-	expression tag	UNP P01860
E	456	HIS	-	expression tag	UNP P01860
E	457	HIS	-	expression tag	UNP P01860
F	296	PHE	TYR	conflict	UNP P01860
F	435	HIS	ARG	conflict	UNP P01860
F	436	TYR	PHE	conflict	UNP P01860
F	448	GLY	-	expression tag	UNP P01860
F	449	SER	-	expression tag	UNP P01860
F	450	LEU	-	expression tag	UNP P01860
F	451	GLU	-	expression tag	UNP P01860
F	452	HIS	-	expression tag	UNP P01860
F	453	HIS	-	expression tag	UNP P01860
F	454	HIS	-	expression tag	UNP P01860
F	455	HIS	-	expression tag	UNP P01860
F	456	HIS	-	expression tag	UNP P01860
F	457	HIS	-	expression tag	UNP P01860

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	24	Total O 24 24	0	0
3	B	19	Total O 19 19	0	0

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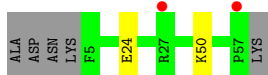
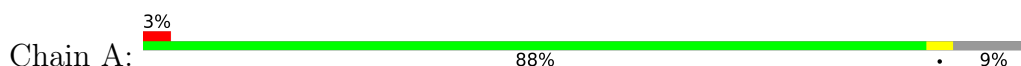
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	1	Total 1	O 1	0	0
3	D	42	Total 42	O 42	0	0
3	E	18	Total 18	O 18	0	0
3	F	22	Total 22	O 22	0	0

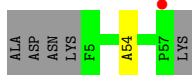
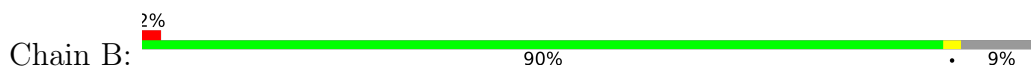
3 Residue-property plots [i](#)

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: Immunoglobulin G-binding protein A



- Molecule 1: Immunoglobulin G-binding protein A



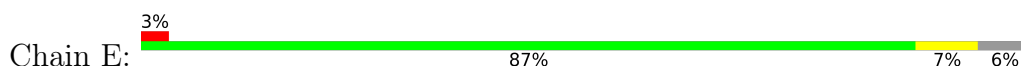
- Molecule 1: Immunoglobulin G-binding protein A



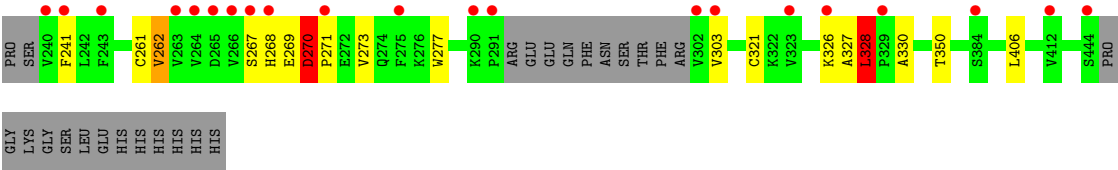
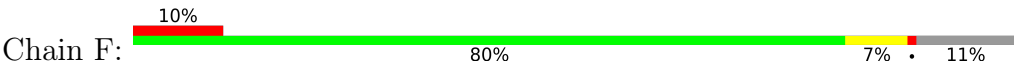
- Molecule 2: Ig gamma-3 chain C region



- Molecule 2: Ig gamma-3 chain C region



- Molecule 2: Ig gamma-3 chain C region



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	137.95Å 87.21Å 103.25Å 90.00° 91.06° 90.00°	Depositor
Resolution (Å)	34.82 – 2.28 34.82 – 2.28	Depositor EDS
% Data completeness (in resolution range)	94.5 (34.82-2.28) 94.2 (34.82-2.28)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.83 (at 2.27Å)	Xtriage
Refinement program	PHENIX dev_1664	Depositor
R, R_{free}	0.196 , 0.242 0.204 , 0.246	Depositor DCC
R_{free} test set	2002 reflections (3.61%)	wwPDB-VP
Wilson B-factor (Å ²)	51.9	Xtriage
Anisotropy	0.266	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 32.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.019 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12392	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.96% 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	A	0.91	0/462	1.02	0/623
1	B	1.00	0/442	1.02	0/597
1	C	0.84	0/442	0.92	0/597
2	D	0.88	0/1707	0.94	0/2322
2	E	0.94	1/1707 (0.1%)	0.99	5/2322 (0.2%)
2	F	0.89	2/1598 (0.1%)	1.02	5/2175 (0.2%)
All	All	0.91	3/6358 (0.0%)	0.99	10/8636 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	270	ASP	C-N	9.17	1.41	1.33
2	E	423	PHE	CA-C	-5.22	1.46	1.52
2	F	271	PRO	N-CD	5.01	1.54	1.47

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	342	GLN	CA-C-N	-6.89	112.86	119.76
2	E	342	GLN	C-N-CA	-6.89	112.86	119.76
2	F	270	ASP	CA-C-N	-6.66	113.35	120.94
2	F	270	ASP	C-N-CA	-6.66	113.35	120.94
2	E	240	VAL	N-CA-C	6.54	118.02	108.53
2	F	303	VAL	N-CA-C	5.80	117.03	108.45
2	E	239	SER	N-CA-C	5.46	117.69	109.23
2	F	328	LEU	CA-C-N	-5.17	113.38	119.84
2	F	328	LEU	C-N-CA	-5.17	113.38	119.84
2	E	252	MET	CG-SD-CE	-5.08	89.72	100.90

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	453	439	436	0	0
1	B	433	419	419	1	0
1	C	433	419	419	0	0
2	D	1661	1629	1629	1	0
2	E	1661	1629	1629	6	0
2	F	1556	1534	1533	6	4
3	A	24	0	0	0	0
3	B	19	0	0	0	0
3	C	1	0	0	0	0
3	D	42	0	0	0	0
3	E	18	0	0	0	0
3	F	22	0	0	0	0
All	All	6323	6069	6065	13	4

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 1.

All (13) 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:54:ALA:HB2	2:E:389:ASN:HB2	1.89	0.55
2:E:250:THR:HG1	2:E:313:TRP:CD1	2.29	0.51
2:F:241:PHE:O	2:F:262:VAL:N	2.44	0.47
2:E:343:PRO:HA	2:E:373:TYR:O	2.17	0.45
2:F:327:ALA:O	2:F:328:LEU:HB2	2.16	0.45
2:F:406:LEU:C	2:F:406:LEU:HD12	2.44	0.43
2:D:330:ALA:HB1	2:D:331:PRO:HD2	2.00	0.43
2:E:290:LYS:HB3	2:E:291:PRO:HD2	1.99	0.43
2:E:406:LEU:C	2:E:406:LEU:HD12	2.45	0.41
2:E:241:PHE:CD1	2:E:241:PHE:N	2.89	0.41
2:F:270:ASP:CG	2:F:326:LYS:HG2	2.46	0.41
2:F:328:LEU:HB3	2:F:330:ALA:O	2.20	0.41
2:F:261:CYS:HB2	2:F:277:TRP:CH2	2.56	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:326:LYS:O	2:F:326:LYS:HE3[2_659]	0.97	0.63
2:F:326:LYS:HG3	2:F:326:LYS:HE2[2_659]	1.22	0.38
2:F:326:LYS:O	2:F:326:LYS:CE[2_659]	1.83	0.37
2:F:326:LYS:C	2:F:326:LYS:HE3[2_659]	1.59	0.01

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	53/58 (91%)	50 (94%)	3 (6%)	0	100	100
1	B	51/58 (88%)	50 (98%)	1 (2%)	0	100	100
1	C	51/58 (88%)	50 (98%)	1 (2%)	0	100	100
2	D	205/220 (93%)	201 (98%)	4 (2%)	0	100	100
2	E	205/220 (93%)	201 (98%)	3 (2%)	1 (0%)	24	29
2	F	191/220 (87%)	181 (95%)	8 (4%)	2 (1%)	12	13
All	All	756/834 (91%)	733 (97%)	20 (3%)	3 (0%)	30	36

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	269	GLU
2	F	328	LEU
2	E	329	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	A	49/51 (96%)	47 (96%)	2 (4%)	27	38
1	B	47/51 (92%)	47 (100%)	0	100	100
1	C	47/51 (92%)	46 (98%)	1 (2%)	47	63
2	D	195/206 (95%)	192 (98%)	3 (2%)	57	72
2	E	195/206 (95%)	193 (99%)	2 (1%)	68	81
2	F	183/206 (89%)	176 (96%)	7 (4%)	29	42
All	All	716/771 (93%)	701 (98%)	15 (2%)	47	63

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	GLU
1	A	50	LYS
1	C	55	GLN
2	D	356	GLU
2	D	359	THR
2	D	382	GLU
2	E	250	THR
2	E	359	THR
2	F	262	VAL
2	F	267	SER
2	F	268	HIS
2	F	270	ASP
2	F	273	VAL
2	F	321	CYS
2	F	350	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	6	ASN
2	E	295	GLN
2	E	342	GLN
2	E	347	GLN
2	F	421	ASN

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 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

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	53/58 (91%)	0.03	2 (3%) 44 46	29, 56, 77, 91	2 (3%)
1	B	53/58 (91%)	0.04	1 (1%) 66 68	46, 56, 78, 85	0
1	C	53/58 (91%)	0.42	2 (3%) 44 46	54, 76, 88, 96	0
2	D	207/220 (94%)	0.15	6 (2%) 53 55	44, 60, 91, 102	0
2	E	207/220 (94%)	0.38	7 (3%) 48 50	45, 67, 99, 109	0
2	F	195/220 (88%)	0.64	21 (10%) 11 12	46, 71, 104, 112	0
All	All	768/834 (92%)	0.34	39 (5%) 33 34	29, 65, 99, 112	2 (0%)

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	266	VAL	5.4
2	F	302	VAL	4.8
2	F	268	HIS	4.7
2	E	444	SER	4.3
2	F	291	PRO	4.1
2	D	268	HIS	4.0
2	F	326	LYS	4.0
2	F	271	PRO	3.7
2	E	238	PRO	3.7
2	D	444	SER	3.7
2	F	267	SER	3.7
2	F	264	VAL	3.6
2	F	240	VAL	3.5
1	A	57	PRO	3.2
2	F	444	SER	3.1
2	D	238	PRO	3.1
2	D	425	CYS	2.7
1	B	57	PRO	2.7
2	F	263	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
2	F	412	VAL	2.5
2	F	303	VAL	2.5
2	F	290	LYS	2.4
2	D	272	GLU	2.4
2	F	329	PRO	2.3
1	A	27[A]	ARG	2.3
1	C	57	PRO	2.3
2	F	275	PHE	2.2
2	F	265	ASP	2.2
2	F	384	SER	2.2
2	F	241	PHE	2.2
2	E	271	PRO	2.2
2	E	268	HIS	2.2
2	F	323	VAL	2.1
2	F	243	PHE	2.1
1	C	44	ILE	2.1
2	E	292	ARG	2.1
2	D	273	VAL	2.1
2	E	296	PHE	2.1
2	E	326	LYS	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.