



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 6, 2026 – 06:41 PM UTC

PDB ID : 1RZI / pdb_00001rzi
Title : Crystal structure of human anti-HIV-1 gp120-reactive antibody 47e fab
Authors : Huang, C.C.; Venturi, M.; Majeed, S.; Moore, M.J.; Phogat, S.; Zhang, M.-Y.; Dimitrov, D.S.; Hendrickson, W.A.; Robinson, J.; Sodroski, J.; Wyatt, R.; Choe, H.; Farzan, M.; Kwong, P.D.
Deposited on : 2003-12-24
Resolution : 2.90 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	:	4-5-2 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

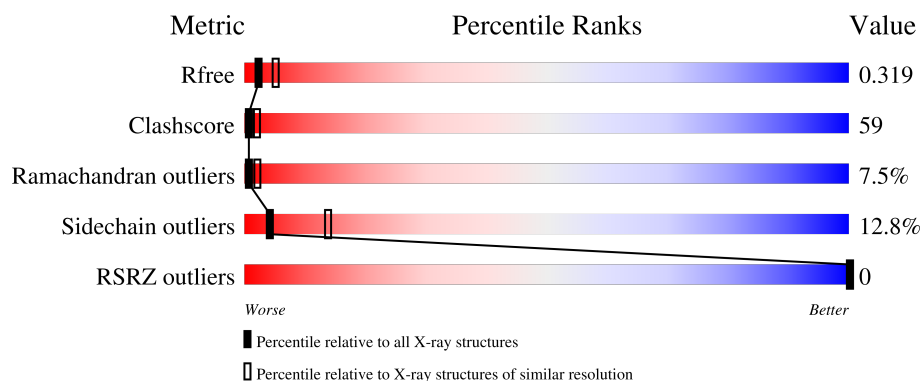
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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.



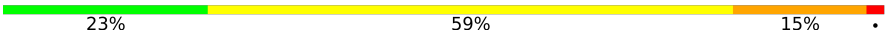
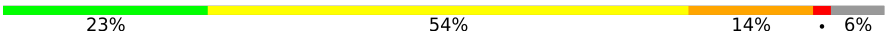
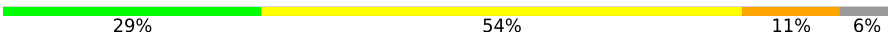
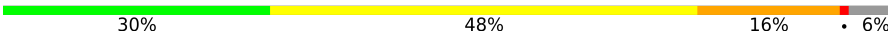
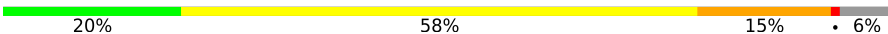
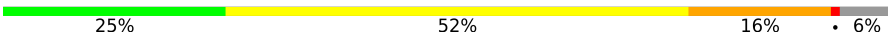
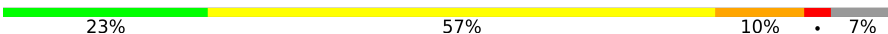
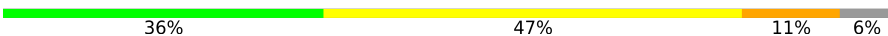
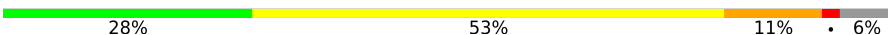
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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 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	212	25% 59% 12% .
1	C	212	26% 51% 22% .
1	E	212	25% 56% 16% .
1	G	212	30% 53% 14% .
1	I	212	29% 56% 15% .

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Mol	Chain	Length	Quality of chain
1	K	212	 26% 57% 13% .
1	M	212	 31% 53% 15% .
1	O	212	 23% 59% 15% .
2	B	230	 23% 54% 14% . 6%
2	D	230	 29% 54% 11% 6%
2	F	230	 30% 48% 16% . 6%
2	H	230	 20% 58% 15% . 6%
2	J	230	 25% 52% 16% . 6%
2	L	230	 23% 57% 10% . 7%
2	N	230	 36% 47% 11% 6%
2	P	230	 28% 53% 11% . 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 25973 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 Fab 47e light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	211	Total	C	N	O	S	0	0	0
			1620	1012	272	331	5			
1	C	212	Total	C	N	O	S	0	0	0
			1628	1016	273	334	5			
1	E	211	Total	C	N	O	S	0	0	0
			1620	1012	272	331	5			
1	G	211	Total	C	N	O	S	0	0	0
			1620	1012	272	331	5			
1	I	211	Total	C	N	O	S	0	0	0
			1620	1012	272	331	5			
1	K	211	Total	C	N	O	S	0	0	0
			1620	1012	272	331	5			
1	M	211	Total	C	N	O	S	0	0	0
			1620	1012	272	331	5			
1	O	211	Total	C	N	O	S	0	0	0
			1620	1012	272	331	5			

- Molecule 2 is a protein called Fab 47e heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	216	Total	C	N	O	S	0	0	0
			1576	992	263	314	7			
2	D	216	Total	C	N	O	S	0	0	0
			1576	992	263	314	7			
2	F	217	Total	C	N	O	S	0	0	0
			1586	998	265	316	7			
2	H	216	Total	C	N	O	S	0	0	0
			1576	992	263	314	7			
2	J	216	Total	C	N	O	S	0	0	0
			1576	992	263	314	7			
2	L	215	Total	C	N	O	S	0	0	0
			1572	990	262	313	7			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	N	216	Total	C	N	O	S	0	0	0
			1576	992	263	314	7			
2	P	216	Total	C	N	O	S	0	0	0
			1576	992	263	314	7			

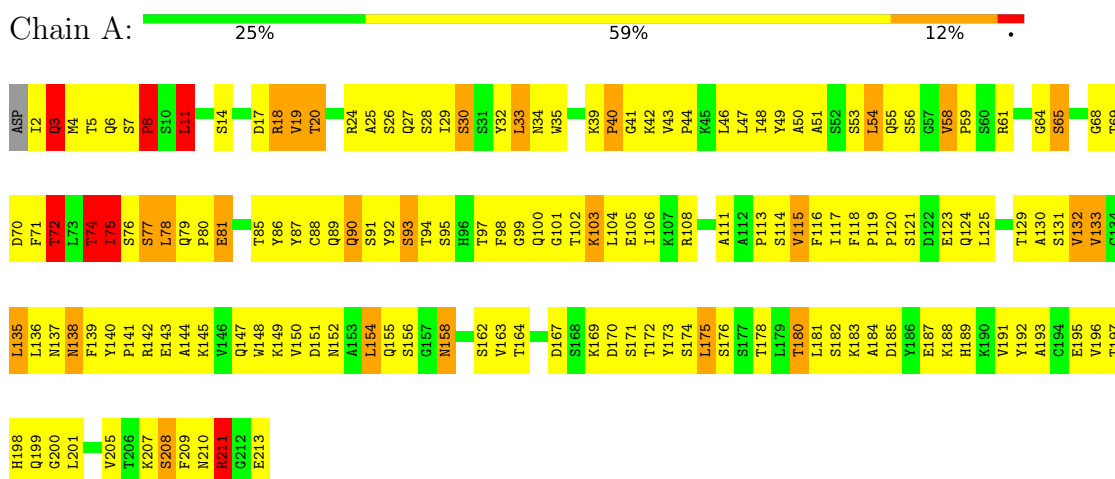
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	34	Total	O	0	0
			34	34		
3	B	29	Total	O	0	0
			29	29		
3	C	25	Total	O	0	0
			25	25		
3	D	27	Total	O	0	0
			27	27		
3	E	26	Total	O	0	0
			26	26		
3	F	18	Total	O	0	0
			18	18		
3	G	23	Total	O	0	0
			23	23		
3	H	31	Total	O	0	0
			31	31		
3	I	21	Total	O	0	0
			21	21		
3	J	17	Total	O	0	0
			17	17		
3	K	21	Total	O	0	0
			21	21		
3	L	20	Total	O	0	0
			20	20		
3	M	23	Total	O	0	0
			23	23		
3	N	31	Total	O	0	0
			31	31		
3	O	22	Total	O	0	0
			22	22		
3	P	23	Total	O	0	0
			23	23		

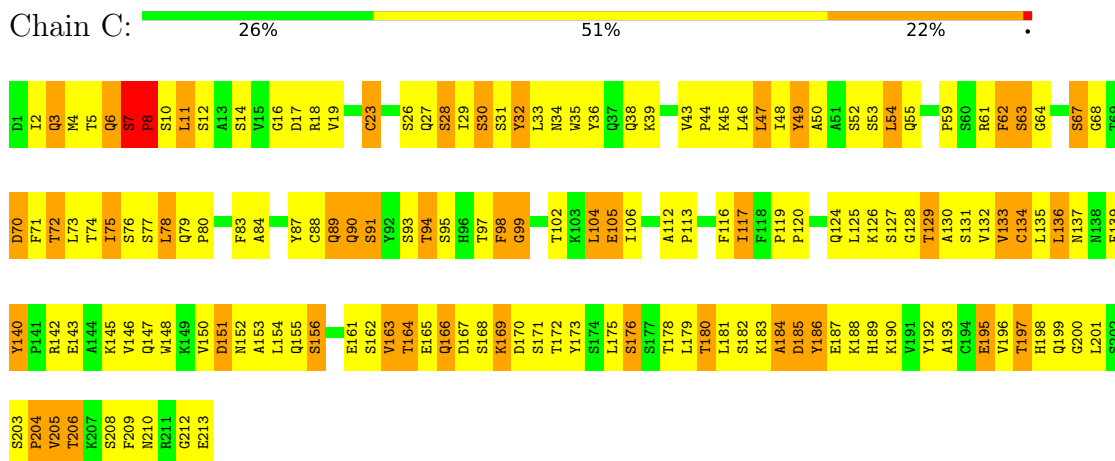
3 Residue-property plots

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.

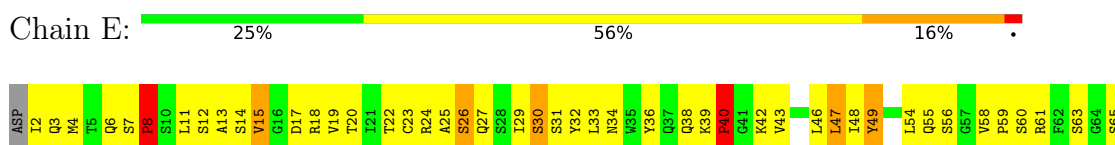
• Molecule 1: Fab 47e light chain

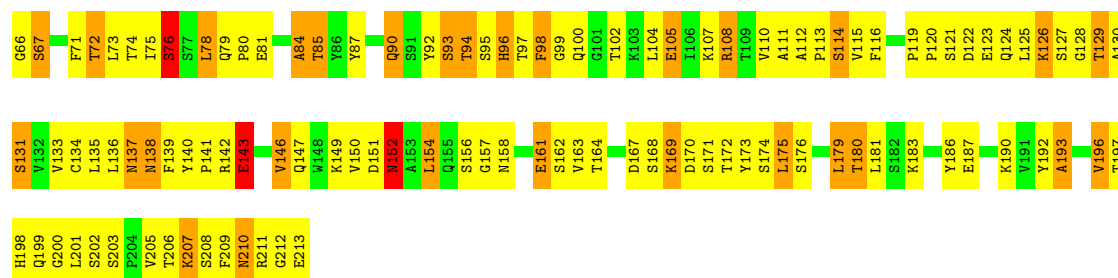


• Molecule 1: Fab 47e light chain



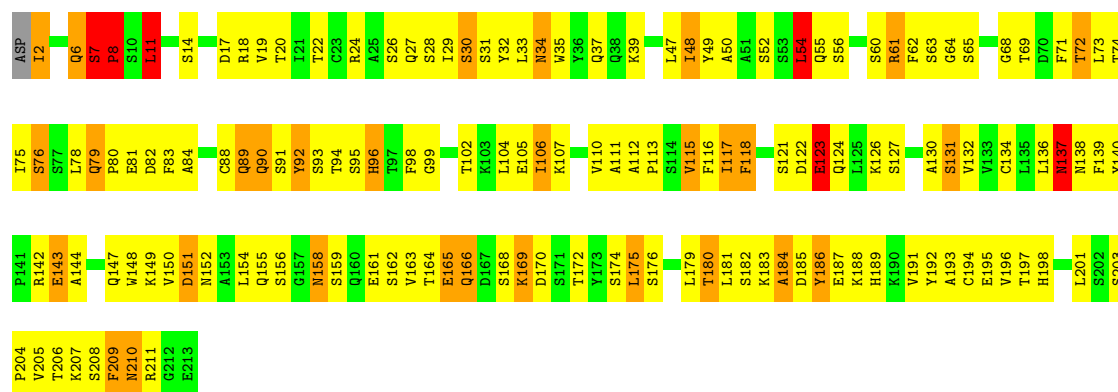
• Molecule 1: Fab 47e light chain





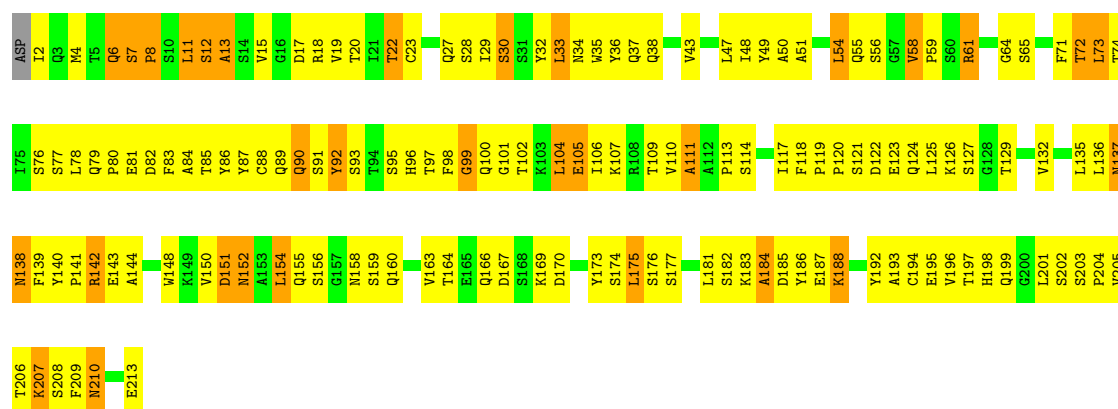
- Molecule 1: Fab 47e light chain

Chain G: 30% 53% 14%



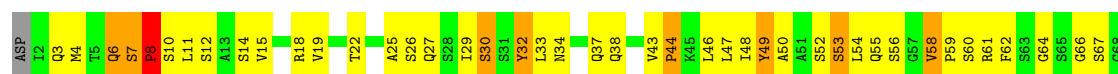
- Molecule 1: Fab 47e light chain

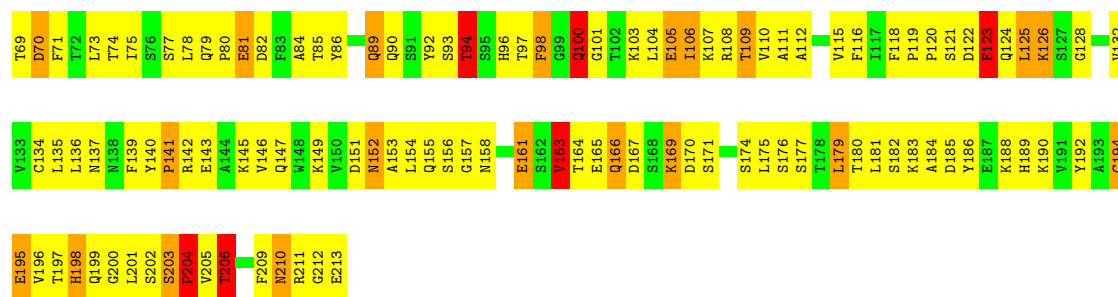
Chain I: 29% 56% 15%



- Molecule 1: Fab 47e light chain

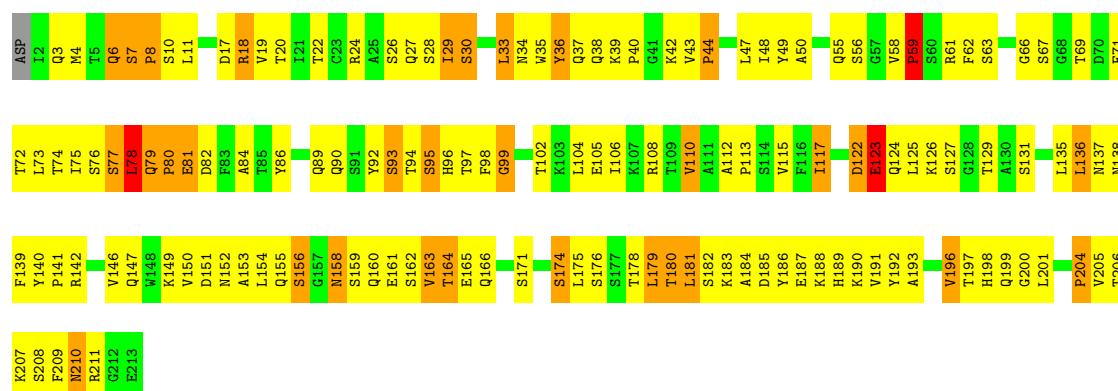
Chain K: 26% 57% 13%





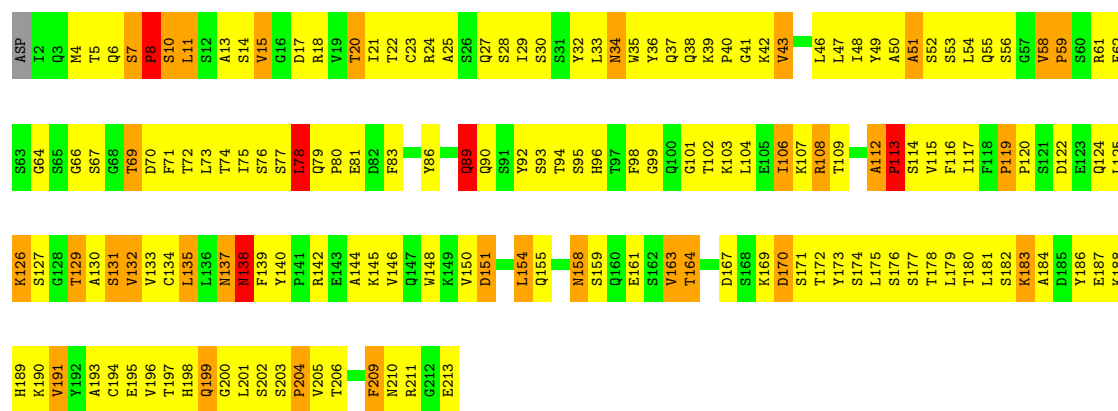
• Molecule 1: Fab 47e light chain

Chain M: 31% 53% 15%



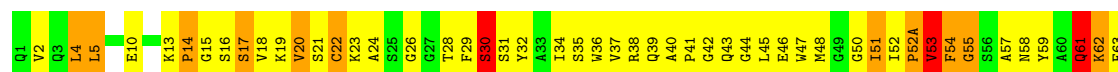
• Molecule 1: Fab 47e light chain

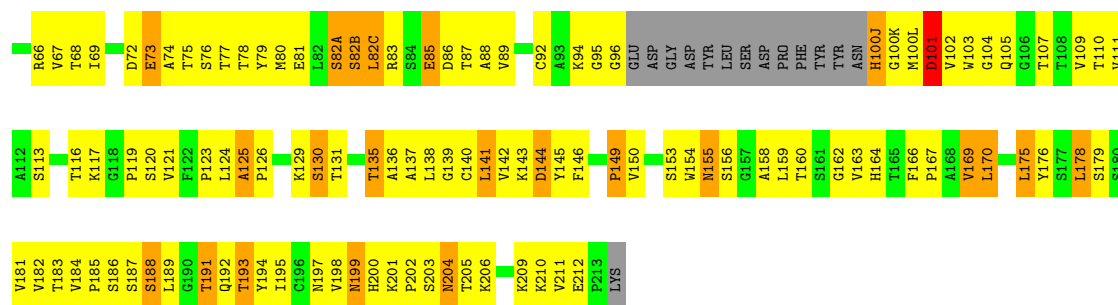
Chain O: 23% 59% 15%



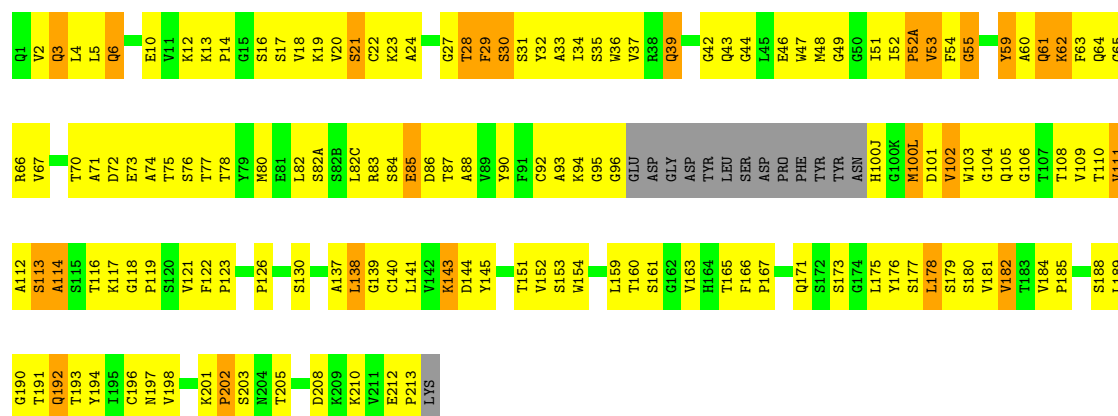
• Molecule 2: Fab 47e heavy chain

Chain B: 23% 54% 14% 6%

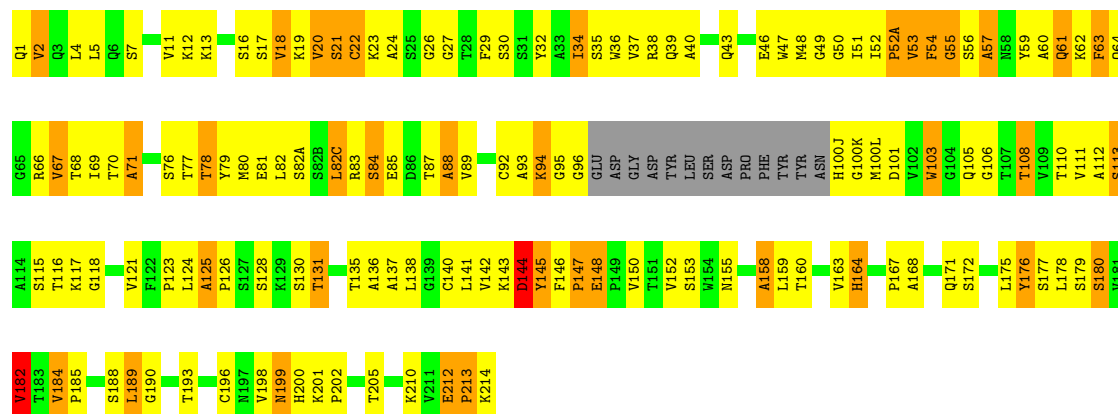




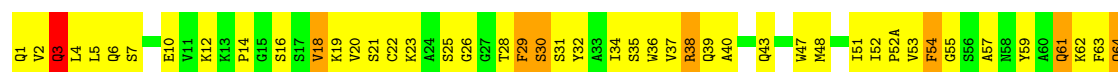
• Molecule 2: Fab 47e heavy chain

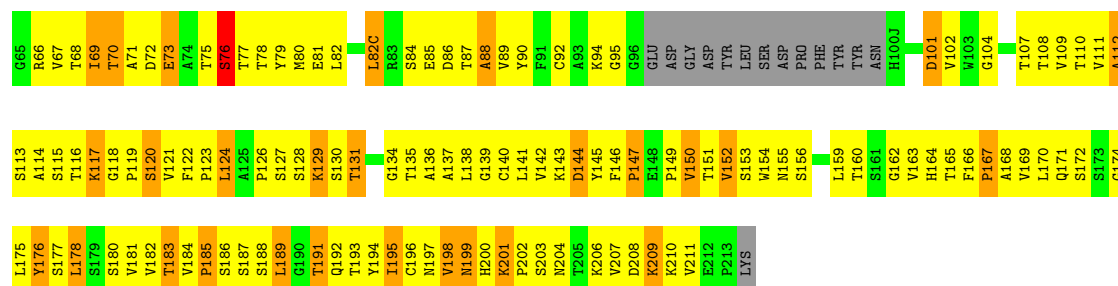


• Molecule 2: Fab 47e heavy chain



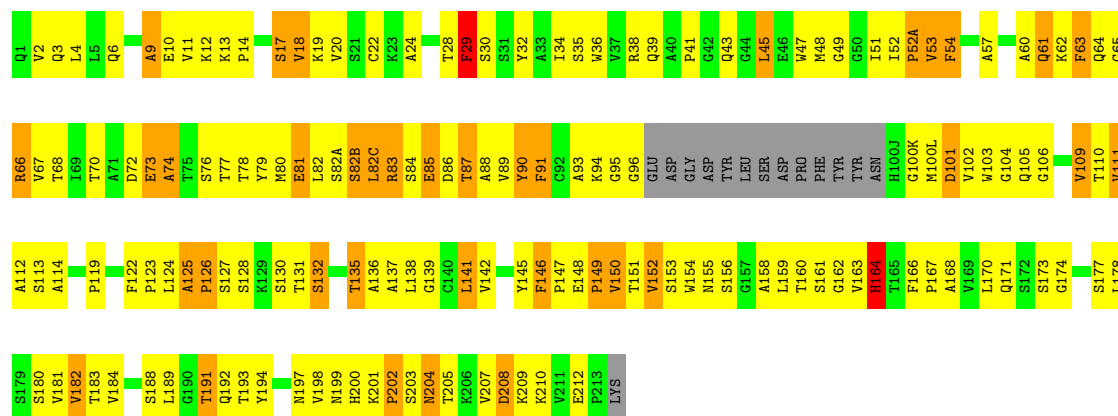
• Molecule 2: Fab 47e heavy chain





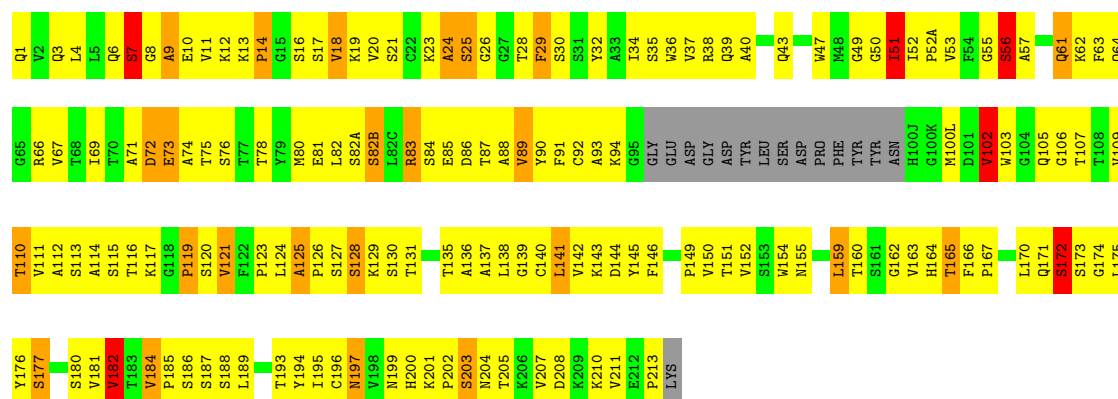
• Molecule 2: Fab 47e heavy chain

Chain J: 25% 52% 16% • 6%



• Molecule 2: Fab 47e heavy chain

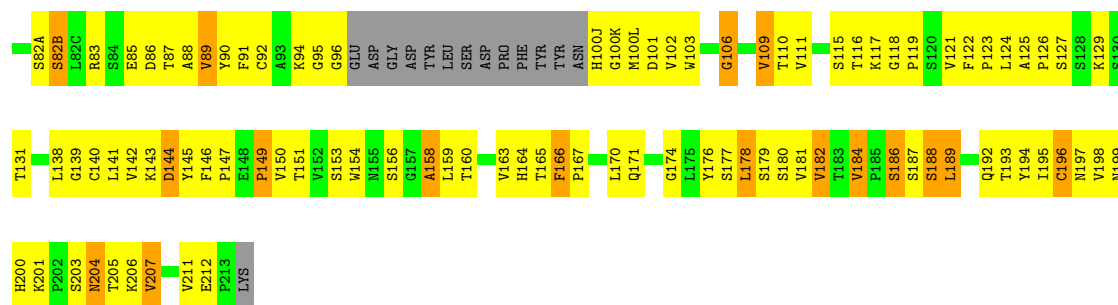
Chain L: 23% 57% 10% • 7%



• Molecule 2: Fab 47e heavy chain

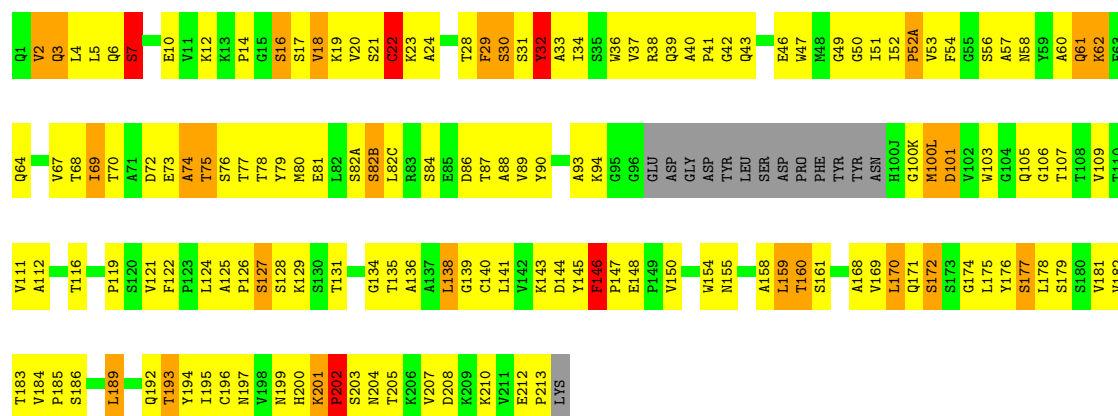
Chain N: 36% 47% 11% 6%





• Molecule 2: Fab 47e heavy chain

Chain P: 28% 53% 11% 6%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	73.43Å 111.64Å 133.31Å 85.48° 90.00° 89.71°	Depositor
Resolution (Å)	20.00 – 2.90 20.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	81.8 (20.00-2.90) 81.2 (20.00-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.97 (at 2.71Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.233 , 0.321 0.232 , 0.319	Depositor DCC
R_{free} test set	7997 reflections (9.48%)	wwPDB-VP
Wilson B-factor (Å ²)	56.1	Xtriage
Anisotropy	0.401	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 41.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.299 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	25973	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.02% 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	0.81	4/1654 (0.2%)	1.93	30/2244 (1.3%)
1	C	0.70	2/1662 (0.1%)	1.41	25/2255 (1.1%)
1	E	0.68	2/1654 (0.1%)	1.42	29/2244 (1.3%)
1	G	0.63	1/1654 (0.1%)	1.39	20/2244 (0.9%)
1	I	0.76	3/1654 (0.2%)	1.20	16/2244 (0.7%)
1	K	0.70	2/1654 (0.1%)	1.38	23/2244 (1.0%)
1	M	0.73	3/1654 (0.2%)	1.37	19/2244 (0.8%)
1	O	0.75	2/1654 (0.1%)	1.45	29/2244 (1.3%)
2	B	0.57	0/1611	1.06	5/2192 (0.2%)
2	D	0.59	0/1611	1.04	5/2192 (0.2%)
2	F	0.54	0/1621	1.09	13/2203 (0.6%)
2	H	0.52	0/1611	1.05	4/2192 (0.2%)
2	J	0.53	0/1611	1.04	10/2192 (0.5%)
2	L	0.54	0/1607	1.09	13/2187 (0.6%)
2	N	0.58	0/1611	1.09	6/2192 (0.3%)
2	P	0.59	0/1611	1.12	14/2192 (0.6%)
All	All	0.65	19/26134 (0.1%)	1.28	261/35505 (0.7%)

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
1	C	0	1
1	K	0	1
All	All	0	2

The worst 5 of 19 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	7	SER	CA-C	12.63	1.67	1.52
1	K	10	SER	N-CA	10.46	1.58	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	M	6	GLN	CA-C	9.62	1.65	1.53
1	C	7	SER	C-O	-8.79	1.12	1.24
1	I	7	SER	C-N	8.01	1.49	1.34

The worst 5 of 261 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	7	SER	CA-C-N	44.53	175.50	119.84
1	A	7	SER	C-N-CA	44.53	175.50	119.84
1	G	7	SER	CA-C-N	21.96	179.71	127.00
1	G	7	SER	C-N-CA	21.96	179.71	127.00
1	M	7	SER	C-N-CD	-21.67	72.93	120.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	140	TYR	Sidechain
1	K	49	TYR	Sidechain

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	1620	0	1575	202	0
1	C	1628	0	1582	179	0
1	E	1620	0	1575	173	0
1	G	1620	0	1575	211	0
1	I	1620	0	1575	191	0
1	K	1620	0	1575	215	0
1	M	1620	0	1575	173	0
1	O	1620	0	1575	175	0
2	B	1576	0	1554	213	0
2	D	1576	0	1554	178	0
2	F	1586	0	1567	219	0
2	H	1576	0	1554	259	0
2	J	1576	0	1554	215	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	L	1572	0	1551	216	0
2	N	1576	0	1554	145	0
2	P	1576	0	1554	176	0
3	A	34	0	0	4	0
3	B	29	0	0	3	0
3	C	25	0	0	6	0
3	D	27	0	0	0	0
3	E	26	0	0	2	0
3	F	18	0	0	2	0
3	G	23	0	0	3	0
3	H	31	0	0	2	0
3	I	21	0	0	2	0
3	J	17	0	0	2	0
3	K	21	0	0	3	0
3	L	20	0	0	3	0
3	M	23	0	0	4	0
3	N	31	0	0	2	0
3	O	22	0	0	1	0
3	P	23	0	0	0	0
All	All	25973	0	25049	2996	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 59.

The worst 5 of 2996 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:A:11:LEU:CD2	1:A:104:LEU:HD22	1.56	1.32
1:A:11:LEU:CD2	1:A:104:LEU:CD2	2.20	1.18
1:E:199:GLN:HA	1:I:199:GLN:HE22	1.12	1.15
2:H:150:VAL:HG22	2:H:151:THR:H	1.11	1.14
1:O:113:PRO:HB3	1:O:139:PHE:HB3	1.29	1.13

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	209/212 (99%)	170 (81%)	29 (14%)	10 (5%)	2	7
1	C	210/212 (99%)	171 (81%)	24 (11%)	15 (7%)	1	2
1	E	209/212 (99%)	168 (80%)	32 (15%)	9 (4%)	2	8
1	G	209/212 (99%)	167 (80%)	28 (13%)	14 (7%)	1	3
1	I	209/212 (99%)	160 (77%)	39 (19%)	10 (5%)	2	7
1	K	209/212 (99%)	160 (77%)	37 (18%)	12 (6%)	1	4
1	M	209/212 (99%)	166 (79%)	32 (15%)	11 (5%)	1	5
1	O	209/212 (99%)	164 (78%)	34 (16%)	11 (5%)	1	5
2	B	212/230 (92%)	153 (72%)	40 (19%)	19 (9%)	0	1
2	D	212/230 (92%)	152 (72%)	44 (21%)	16 (8%)	1	2
2	F	213/230 (93%)	153 (72%)	42 (20%)	18 (8%)	0	1
2	H	212/230 (92%)	146 (69%)	42 (20%)	24 (11%)	0	1
2	J	212/230 (92%)	151 (71%)	40 (19%)	21 (10%)	0	1
2	L	211/230 (92%)	152 (72%)	39 (18%)	20 (10%)	0	1
2	N	212/230 (92%)	164 (77%)	30 (14%)	18 (8%)	0	1
2	P	212/230 (92%)	152 (72%)	37 (18%)	23 (11%)	0	1
All	All	3369/3536 (95%)	2549 (76%)	569 (17%)	251 (8%)	1	2

5 of 251 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	11	LEU
1	A	211	ARG
2	B	22	CYS
2	B	30	SER
2	B	61	GLN

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	187/188 (100%)	156 (83%)	31 (17%)	2	7
1	C	188/188 (100%)	157 (84%)	31 (16%)	2	8
1	E	187/188 (100%)	158 (84%)	29 (16%)	2	9
1	G	187/188 (100%)	163 (87%)	24 (13%)	4	14
1	I	187/188 (100%)	166 (89%)	21 (11%)	6	19
1	K	187/188 (100%)	166 (89%)	21 (11%)	6	19
1	M	187/188 (100%)	162 (87%)	25 (13%)	4	12
1	O	187/188 (100%)	160 (86%)	27 (14%)	3	11
2	B	176/189 (93%)	150 (85%)	26 (15%)	3	10
2	D	176/189 (93%)	159 (90%)	17 (10%)	8	25
2	F	177/189 (94%)	156 (88%)	21 (12%)	5	16
2	H	176/189 (93%)	158 (90%)	18 (10%)	7	23
2	J	176/189 (93%)	155 (88%)	21 (12%)	5	16
2	L	176/189 (93%)	158 (90%)	18 (10%)	7	23
2	N	176/189 (93%)	157 (89%)	19 (11%)	6	21
2	P	176/189 (93%)	152 (86%)	24 (14%)	3	12
All	All	2906/3016 (96%)	2533 (87%)	373 (13%)	4	14

5 of 373 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	J	132	SER
1	M	77	SER
2	J	208	ASP
2	L	7	SER
1	M	180	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 98 such sidechains are listed below:

Mol	Chain	Res	Type
1	I	199	GLN
2	L	39	GLN
2	J	39	GLN
1	K	89	GLN
1	M	37	GLN

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	211/212 (99%)	-1.61	0 100 100	27, 49, 72, 79	0
1	C	212/212 (100%)	-1.66	0 100 100	24, 45, 68, 75	0
1	E	211/212 (99%)	-1.66	0 100 100	30, 49, 65, 76	0
1	G	211/212 (99%)	-1.62	0 100 100	33, 53, 86, 100	0
1	I	211/212 (99%)	-1.63	0 100 100	31, 50, 71, 82	0
1	K	211/212 (99%)	-1.66	0 100 100	36, 52, 66, 76	0
1	M	211/212 (99%)	-1.65	0 100 100	24, 45, 59, 75	0
1	O	211/212 (99%)	-1.65	0 100 100	25, 44, 61, 67	0
2	B	216/230 (93%)	-1.57	0 100 100	32, 58, 81, 89	0
2	D	216/230 (93%)	-1.60	0 100 100	27, 58, 87, 99	0
2	F	217/230 (94%)	-1.54	0 100 100	28, 64, 93, 105	0
2	H	216/230 (93%)	-1.49	0 100 100	39, 71, 96, 99	0
2	J	216/230 (93%)	-1.53	0 100 100	39, 65, 91, 96	0
2	L	215/230 (93%)	-1.52	0 100 100	37, 63, 82, 100	0
2	N	216/230 (93%)	-1.63	0 100 100	29, 54, 75, 87	0
2	P	216/230 (93%)	-1.58	0 100 100	28, 52, 86, 92	0
All	All	3417/3536 (96%)	-1.60	0 100 100	24, 53, 84, 105	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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.