



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:32 PM UTC

PDB ID : 4NEE / pdb\_00004nee  
Title : crystal structure of AP-2 alpha/simga2 complex bound to HIV-1 Nef  
Authors : Hurley, J.H.; Bonifacino, J.S.; Ren, X.; Park, S.Y.  
Deposited on : 2013-10-29  
Resolution : 2.88 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](mailto: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>.

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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

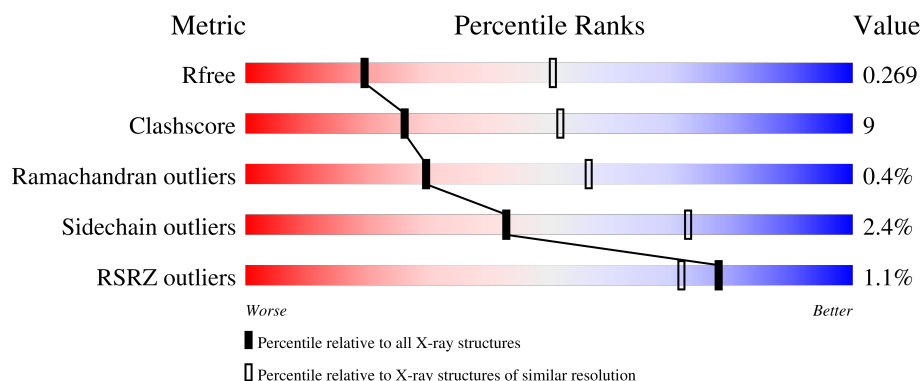
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.88 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	3557 (2.90-2.86)
Clashscore	190562	3801 (2.90-2.86)
Ramachandran outliers	187476	3699 (2.90-2.86)
Sidechain outliers	187428	3702 (2.90-2.86)
RSRZ outliers	180081	3558 (2.90-2.86)

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  $\geq 3$ , 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	D	142	<div> <div>81%</div> <div>17%</div> <div>.</div> </div>
1	F	142	<div> <div>65%</div> <div>33%</div> <div>..</div> </div>
1	I	142	<div> <div>62%</div> <div>36%</div> <div>.</div> </div>
1	L	142	<div> <div>68%</div> <div>30%</div> <div>.</div> </div>
2	A	398	<div> <div>75%</div> <div>21%</div> <div>..</div> </div>

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Mol	Chain	Length	Quality of chain
2	B	398	<div><div></div><div>76%20%<div><div></div><div></div><div></div></div></div></div>
2	G	398	<div>%<div><div></div><div>78%18%<div><div></div><div></div><div></div></div></div></div></div>
2	J	398	<div>%<div><div></div><div>74%23%<div><div></div><div></div><div></div></div></div></div></div>
3	C	155	<div>%<div><div></div><div>65%23%<div><div></div><div></div><div></div></div></div><div>•10%</div></div></div>
3	E	155	<div>%<div><div></div><div>65%23%<div><div></div><div></div><div></div></div></div><div>•10%</div></div></div>
3	H	155	<div>2%<div><div></div><div>66%25%<div><div></div><div></div><div></div></div></div><div>8%</div></div></div>
3	K	155	<div>%<div><div></div><div>72%19%<div><div></div><div></div><div></div></div></div><div>•7%</div></div></div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 21530 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 AP-2 complex subunit sigma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	142	Total	C	N	O	S	0	0	0
			1199	778	200	214	7			
1	L	142	Total	C	N	O	S	0	0	0
			1199	778	200	214	7			
1	F	142	Total	C	N	O	S	0	0	0
			1199	778	200	214	7			
1	I	142	Total	C	N	O	S	0	0	0
			1199	778	200	214	7			

- Molecule 2 is a protein called AP-2 complex subunit alpha-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	388	Total	C	N	O	S	0	0	0
			3047	1931	536	564	16			
2	B	388	Total	C	N	O	S	0	0	0
			3047	1931	536	564	16			
2	J	388	Total	C	N	O	S	0	0	0
			3047	1931	536	564	16			
2	A	385	Total	C	N	O	S	0	0	0
			3025	1918	532	559	16			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	-1	GLY	-	cloning artifact	UNP Q66HM2
G	0	ALA	-	cloning artifact	UNP Q66HM2
B	-1	GLY	-	cloning artifact	UNP Q66HM2
B	0	ALA	-	cloning artifact	UNP Q66HM2
J	-1	GLY	-	cloning artifact	UNP Q66HM2
J	0	ALA	-	cloning artifact	UNP Q66HM2
A	-1	GLY	-	cloning artifact	UNP Q66HM2
A	0	ALA	-	cloning artifact	UNP Q66HM2

- Molecule 3 is a protein called Protein Nef.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	139	Total	C	N	O	S	0	0	0
			1134	732	197	202	3			
3	H	142	Total	C	N	O	S	0	0	0
			1143	737	199	204	3			
3	K	144	Total	C	N	O	S	0	0	0
			1157	745	202	207	3			
3	C	139	Total	C	N	O	S	0	0	0
			1134	732	197	202	3			

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	48	GLY	-	expression tag	UNP Q90VU7
E	49	UNK	-	expression tag	UNP Q90VU7
E	50	UNK	-	expression tag	UNP Q90VU7
E	51	UNK	-	expression tag	UNP Q90VU7
E	52	UNK	-	expression tag	UNP Q90VU7
E	53	UNK	-	expression tag	UNP Q90VU7
E	54	UNK	-	expression tag	UNP Q90VU7
E	55	UNK	-	expression tag	UNP Q90VU7
E	56	UNK	-	expression tag	UNP Q90VU7
E	57	UNK	-	expression tag	UNP Q90VU7
E	58	UNK	-	expression tag	UNP Q90VU7
E	59	UNK	-	expression tag	UNP Q90VU7
E	60	UNK	-	expression tag	UNP Q90VU7
E	61	UNK	-	expression tag	UNP Q90VU7
H	54	GLY	-	expression tag	UNP Q90VU7
H	55	UNK	-	expression tag	UNP Q90VU7
H	56	UNK	-	expression tag	UNP Q90VU7
H	57	UNK	-	expression tag	UNP Q90VU7
H	58	UNK	-	expression tag	UNP Q90VU7
H	59	UNK	-	expression tag	UNP Q90VU7
H	60	UNK	-	expression tag	UNP Q90VU7
H	61	UNK	-	expression tag	UNP Q90VU7
H	62	UNK	-	expression tag	UNP Q90VU7
H	63	UNK	-	expression tag	UNP Q90VU7
H	64	UNK	-	expression tag	UNP Q90VU7
H	65	UNK	-	expression tag	UNP Q90VU7
H	66	UNK	-	expression tag	UNP Q90VU7
H	67	UNK	-	expression tag	UNP Q90VU7
K	54	GLY	-	expression tag	UNP Q90VU7
K	55	UNK	-	expression tag	UNP Q90VU7

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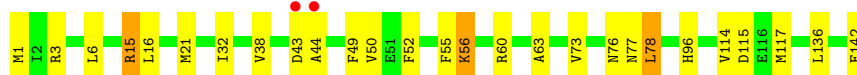
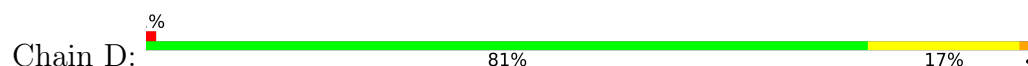
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Chain	Residue	Modelled	Actual	Comment	Reference
K	56	UNK	-	expression tag	UNP Q90VU7
K	57	UNK	-	expression tag	UNP Q90VU7
K	58	UNK	-	expression tag	UNP Q90VU7
K	59	UNK	-	expression tag	UNP Q90VU7
K	60	UNK	-	expression tag	UNP Q90VU7
K	61	UNK	-	expression tag	UNP Q90VU7
K	62	UNK	-	expression tag	UNP Q90VU7
K	63	UNK	-	expression tag	UNP Q90VU7
K	64	UNK	-	expression tag	UNP Q90VU7
K	65	UNK	-	expression tag	UNP Q90VU7
K	66	UNK	-	expression tag	UNP Q90VU7
K	67	UNK	-	expression tag	UNP Q90VU7
C	48	GLY	-	expression tag	UNP Q90VU7
C	49	UNK	-	expression tag	UNP Q90VU7
C	50	UNK	-	expression tag	UNP Q90VU7
C	51	UNK	-	expression tag	UNP Q90VU7
C	52	UNK	-	expression tag	UNP Q90VU7
C	53	UNK	-	expression tag	UNP Q90VU7
C	54	UNK	-	expression tag	UNP Q90VU7
C	55	UNK	-	expression tag	UNP Q90VU7
C	56	UNK	-	expression tag	UNP Q90VU7
C	57	UNK	-	expression tag	UNP Q90VU7
C	58	UNK	-	expression tag	UNP Q90VU7
C	59	UNK	-	expression tag	UNP Q90VU7
C	60	UNK	-	expression tag	UNP Q90VU7
C	61	UNK	-	expression tag	UNP Q90VU7

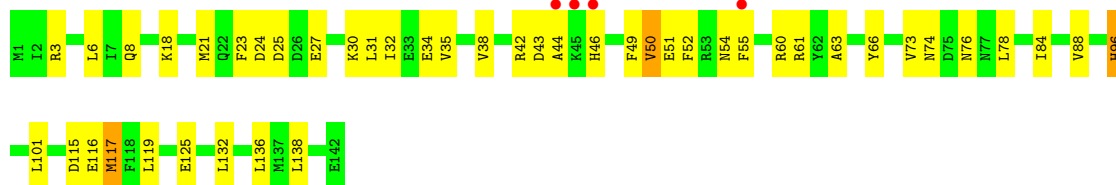
### 3 Residue-property plots

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.

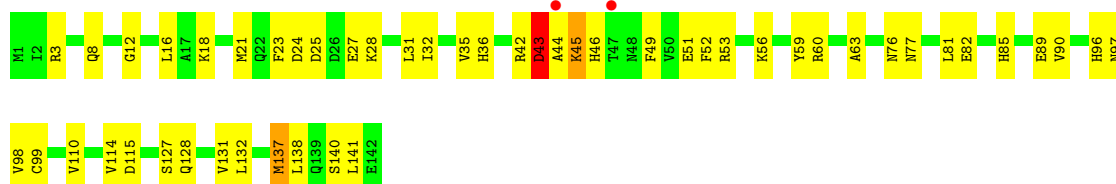
- Molecule 1: AP-2 complex subunit sigma



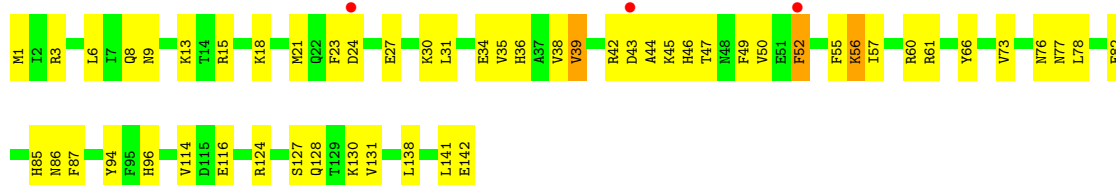
- Molecule 1: AP-2 complex subunit sigma



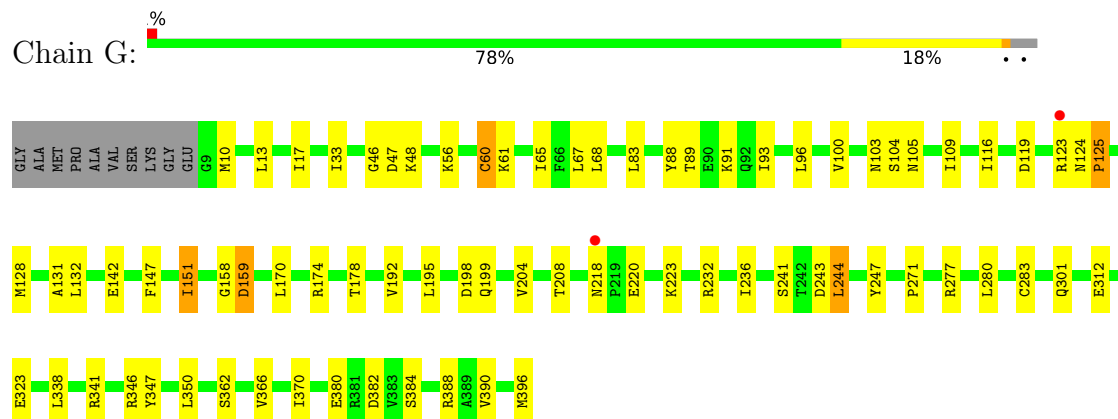
- Molecule 1: AP-2 complex subunit sigma



- Molecule 1: AP-2 complex subunit sigma



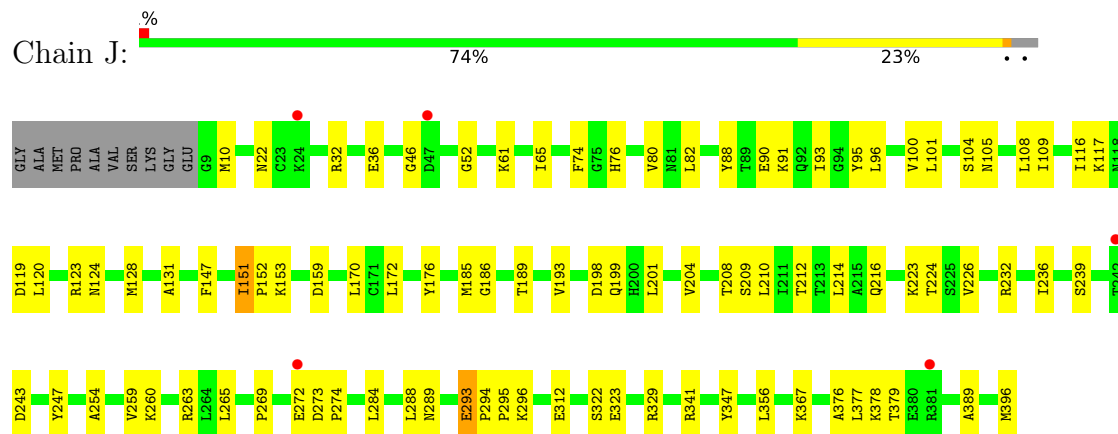
- Molecule 2: AP-2 complex subunit alpha-2



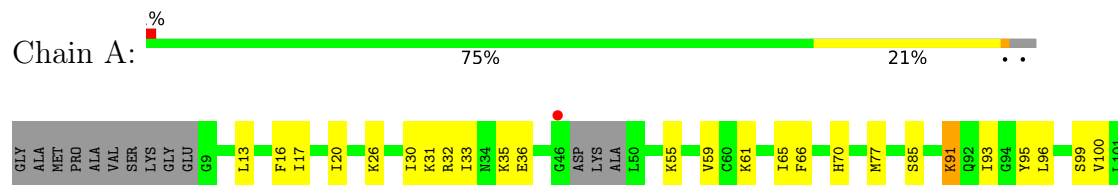
- Molecule 2: AP-2 complex subunit alpha-2

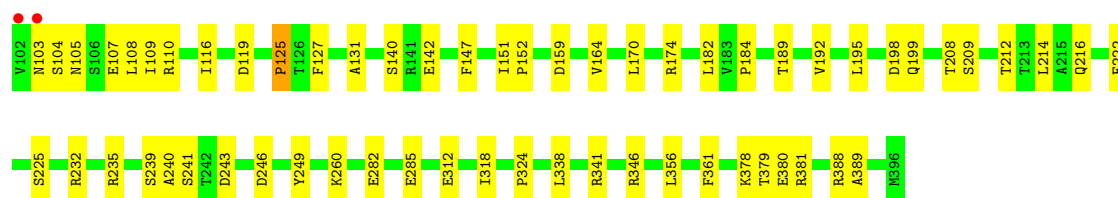


- Molecule 2: AP-2 complex subunit alpha-2



- Molecule 2: AP-2 complex subunit alpha-2

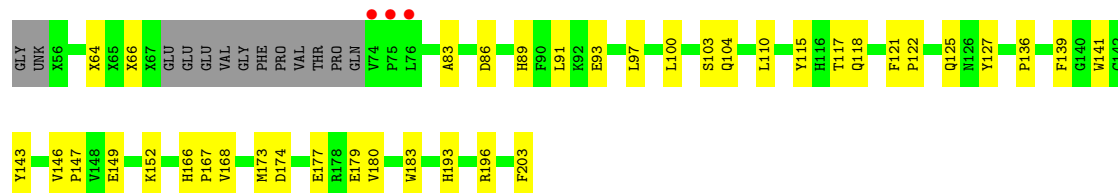




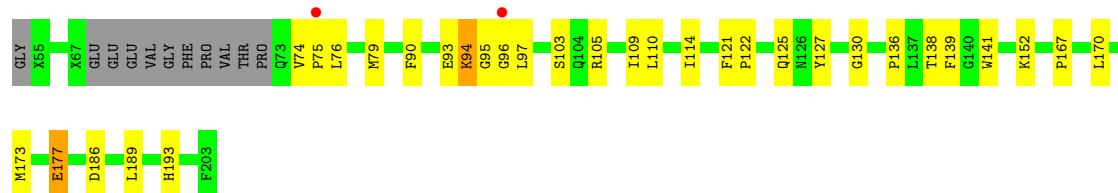
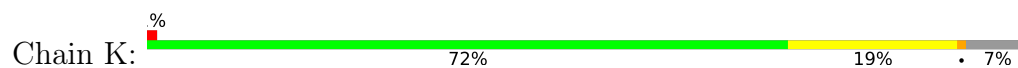
• Molecule 3: Protein Nef



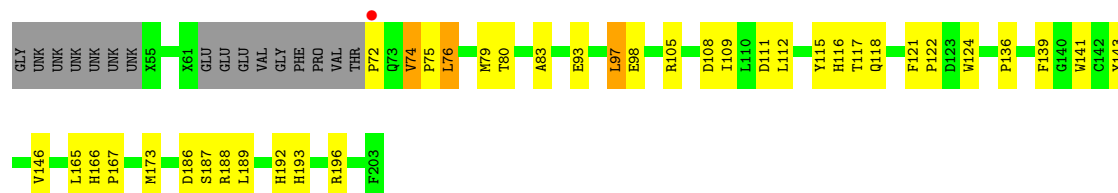
• Molecule 3: Protein Nef



• Molecule 3: Protein Nef



• Molecule 3: Protein Nef



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	109.56Å 168.03Å 200.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.88 – 2.88 48.88 – 2.88	Depositor EDS
% Data completeness (in resolution range)	95.5 (48.88-2.88) 94.6 (48.88-2.88)	Depositor EDS
$R_{merge}$	0.18	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.97 (at 2.91Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.3_1479)	Depositor
R, $R_{free}$	0.219 , 0.267 0.225 , 0.269	Depositor DCC
$R_{free}$ test set	4014 reflections (4.78%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	35.7	Xtriage
Anisotropy	0.169	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 44.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.42$ , $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	21530	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.36% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	D	1.07	0/1223	0.90	1/1650 (0.1%)
1	F	0.90	0/1223	0.90	5/1650 (0.3%)
1	I	1.12	0/1223	0.90	2/1650 (0.1%)
1	L	1.01	0/1223	0.97	2/1650 (0.1%)
2	A	0.87	0/3076	0.87	2/4161 (0.0%)
2	B	0.98	0/3099	0.90	10/4193 (0.2%)
2	G	1.00	0/3099	0.89	6/4193 (0.1%)
2	J	0.95	1/3099 (0.0%)	0.90	8/4193 (0.2%)
3	C	0.90	0/1139	0.89	2/1551 (0.1%)
3	E	0.91	0/1139	0.88	0/1551
3	H	0.83	0/1122	0.88	0/1528
3	K	0.81	0/1131	0.83	1/1540 (0.1%)
All	All	0.95	1/21796 (0.0%)	0.89	39/29510 (0.1%)

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	F	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	J	254	ALA	CA-CB	-5.05	1.49	1.52

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	243	ASP	N-CA-C	9.99	122.25	111.36
2	B	158	GLY	N-CA-C	-8.64	98.67	112.66
1	L	25	ASP	N-CA-C	7.53	119.48	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	389	ALA	N-CA-C	-6.88	104.52	113.12
1	I	52	PHE	N-CA-C	6.69	119.85	109.07
2	G	244	LEU	N-CA-C	6.10	123.80	110.80
1	F	43	ASP	CA-C-N	6.06	132.60	121.70
1	F	43	ASP	C-N-CA	6.06	132.60	121.70
2	G	323	GLU	CA-C-N	5.98	125.19	118.97
2	G	323	GLU	C-N-CA	5.98	125.19	118.97
2	B	270	PRO	O-C-N	5.89	124.02	121.31
1	F	25	ASP	N-CA-C	-5.87	104.96	111.71
3	K	94	LYS	N-CA-C	-5.86	106.18	113.15
2	B	294	PRO	CA-C-N	5.62	125.80	119.90
2	B	294	PRO	C-N-CA	5.62	125.80	119.90
1	I	94	TYR	N-CA-C	5.59	117.46	111.36
2	J	151	ILE	CB-CA-C	-5.59	108.42	114.35
3	C	74	VAL	CA-C-N	5.58	125.53	119.78
3	C	74	VAL	C-N-CA	5.58	125.53	119.78
1	L	63	ALA	CB-CA-C	-5.58	110.13	116.54
1	F	63	ALA	CB-CA-C	-5.51	110.20	116.54
2	A	159	ASP	N-CA-C	-5.51	106.55	113.28
2	B	323	GLU	CA-C-N	5.48	124.67	118.97
2	B	323	GLU	C-N-CA	5.48	124.67	118.97
2	A	389	ALA	N-CA-C	-5.47	106.28	113.12
1	F	45	LYS	N-CA-C	5.46	117.23	111.28
2	B	270	PRO	CA-C-N	5.38	125.30	119.76
2	B	270	PRO	C-N-CA	5.38	125.30	119.76
2	J	74	PHE	N-CA-C	5.33	116.09	108.74
2	B	218	ASN	CA-C-N	5.32	124.96	119.05
2	B	218	ASN	C-N-CA	5.32	124.96	119.05
2	J	273	ASP	CA-C-N	5.20	124.82	119.05
2	J	273	ASP	C-N-CA	5.20	124.82	119.05
1	D	63	ALA	CB-CA-C	-5.15	110.62	116.54
2	J	269	PRO	CA-C-N	-5.13	115.91	119.66
2	J	269	PRO	C-N-CA	-5.13	115.91	119.66
2	G	218	ASN	CA-C-N	5.05	124.66	119.05
2	G	218	ASN	C-N-CA	5.05	124.66	119.05
2	G	151	ILE	CB-CA-C	-5.01	109.04	114.35

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	F	43	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	D	1199	0	1195	20	0
1	F	1199	0	1195	38	0
1	I	1199	0	1195	39	0
1	L	1199	0	1195	34	0
2	A	3025	0	3105	54	0
2	B	3047	0	3128	55	0
2	G	3047	0	3128	48	0
2	J	3047	0	3128	62	0
3	C	1134	0	1058	26	0
3	E	1134	0	1058	28	0
3	H	1143	0	1047	31	0
3	K	1157	0	1056	20	0
All	All	21530	0	21488	405	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 (405) 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)
2:A:209:SER:OG	2:A:260:LYS:NZ	2.04	0.91
1:F:56:LYS:NZ	1:F:77:ASN:OD1	2.09	0.86
1:L:76:ASN:HD22	2:J:312:GLU:HG2	1.40	0.86
2:A:95:TYR:OH	2:A:119:ASP:OD2	1.95	0.83
3:E:56:UNK:HA	3:E:94:LYS:HE2	1.61	0.83
1:F:46:HIS:O	3:E:178:ARG:NH1	2.13	0.82
3:C:193:HIS:HD2	3:C:196:ARG:H	1.28	0.81
1:D:60:ARG:NH2	3:C:173:MET:SD	2.55	0.79
1:F:43:ASP:HB3	1:F:45:LYS:N	1.98	0.79
3:E:134:ARG:NH2	3:E:179:GLU:OE2	2.17	0.78
3:H:193:HIS:HD2	3:H:196:ARG:H	1.30	0.78
1:I:56:LYS:NZ	1:I:77:ASN:OD1	2.16	0.77
1:L:18:LYS:NZ	1:L:115:ASP:OD2	2.15	0.77
2:J:209:SER:OG	2:J:260:LYS:NZ	2.18	0.76
1:D:38:VAL:HG11	1:D:50:VAL:HG11	1.71	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:56:LYS:HG3	2:G:93:ILE:HG13	1.70	0.73
2:B:156:VAL:HG22	2:B:191:ARG:HB2	1.71	0.73
1:D:1:MET:SD	1:D:3:ARG:NH1	2.61	0.73
2:B:382:ASP:OD2	2:B:384:SER:OG	2.07	0.72
3:H:149:GLU:HB2	3:H:152:LYS:HG3	1.72	0.72
2:B:91:LYS:HE2	2:B:119:ASP:OD1	1.89	0.71
2:A:125:PRO:O	2:A:127:PHE:N	2.21	0.71
3:H:64:UNK:O	3:H:66:UNK:HA	1.91	0.71
2:A:318:ILE:HG12	2:A:356:LEU:HD13	1.75	0.69
1:L:38:VAL:HG11	1:L:50:VAL:HG11	1.74	0.69
3:C:80:THR:H	3:C:83:ALA:HB3	1.58	0.69
2:B:240:ALA:HB1	2:A:282:GLU:OE1	1.94	0.68
2:J:95:TYR:OH	2:J:119:ASP:OD1	2.04	0.67
2:A:235:ARG:O	2:A:239:SER:OG	2.12	0.67
1:D:16:LEU:HD21	1:D:114:VAL:HG21	1.77	0.67
1:D:43:ASP:OD1	1:D:44:ALA:N	2.26	0.67
1:I:60:ARG:NH2	3:H:173:MET:HE3	2.10	0.66
2:G:301:GLN:HE21	3:H:173:MET:HE2	1.59	0.66
2:J:116:ILE:HG23	2:J:131:ALA:HB1	1.76	0.66
2:A:85:SER:O	2:A:91:LYS:NZ	2.21	0.66
2:A:208:THR:O	2:A:212:THR:OG1	2.13	0.66
1:I:38:VAL:HG11	1:I:50:VAL:HG11	1.78	0.65
3:C:72:PRO:HD2	3:C:116:HIS:HA	1.78	0.65
2:G:91:LYS:HE2	2:G:119:ASP:OD1	1.96	0.65
2:A:174:ARG:O	2:A:174:ARG:NH1	2.30	0.65
2:G:116:ILE:HG23	2:G:131:ALA:HB1	1.77	0.64
1:I:1:MET:SD	1:I:3:ARG:NH2	2.71	0.64
1:L:76:ASN:ND2	2:J:312:GLU:HG2	2.11	0.64
2:A:198:ASP:OD1	2:A:199:GLN:N	2.31	0.64
3:C:105:ARG:O	3:C:109:ILE:HG13	1.98	0.64
2:J:236:ILE:O	2:J:239:SER:OG	2.10	0.63
2:B:377:LEU:O	2:B:386:ARG:NH2	2.31	0.63
1:D:56:LYS:HE2	1:D:73:VAL:HA	1.81	0.63
3:K:125:GLN:NE2	3:K:127:TYR:OH	2.26	0.63
2:B:95:TYR:OH	2:B:119:ASP:OD2	2.16	0.62
1:F:44:ALA:O	3:E:178:ARG:NH1	2.33	0.62
3:E:101:ILE:O	3:E:106:ARG:NH1	2.33	0.61
1:I:30:LYS:O	1:I:34:GLU:HG3	2.01	0.61
3:H:125:GLN:NE2	3:H:127:TYR:OH	2.25	0.61
1:D:76:ASN:ND2	2:B:312:GLU:OE1	2.31	0.60
2:A:346:ARG:NH1	2:A:380:GLU:OE1	2.35	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:192:VAL:O	2:A:195:LEU:HB2	2.01	0.60
2:A:240:ALA:HB3	2:A:243:ASP:OD2	2.02	0.59
1:F:3:ARG:HD2	1:F:21:MET:SD	2.43	0.59
1:F:21:MET:HE2	1:F:23:PHE:CZ	2.37	0.59
2:J:61:LYS:O	2:J:65:ILE:HG13	2.03	0.58
1:I:3:ARG:HD2	1:I:21:MET:SD	2.42	0.58
1:I:127:SER:O	1:I:131:VAL:HG23	2.03	0.58
2:G:346:ARG:NH1	2:G:380:GLU:OE1	2.37	0.58
3:K:103:SER:HB3	3:K:167:PRO:HB3	1.85	0.58
1:L:101:LEU:HD21	3:C:187:SER:OG	2.04	0.58
2:B:235:ARG:O	2:B:239:SER:OG	2.21	0.58
2:A:105:ASN:O	2:A:109:ILE:HG13	2.04	0.57
2:J:198:ASP:OD1	2:J:199:GLN:N	2.37	0.57
1:D:6:LEU:HD11	1:D:32:ILE:HD13	1.86	0.57
3:K:93:GLU:C	3:K:95:GLY:H	2.11	0.57
1:F:76:ASN:ND2	2:A:312:GLU:OE1	2.26	0.57
2:G:382:ASP:OD2	2:G:384:SER:OG	2.19	0.57
1:F:98:VAL:HG13	3:E:164:LEU:HD22	1.85	0.57
2:B:185:MET:HG3	2:B:218:ASN:HD22	1.69	0.57
3:C:193:HIS:CD2	3:C:196:ARG:H	2.18	0.57
1:I:46:HIS:ND1	2:G:382:ASP:OD1	2.38	0.57
2:G:220:GLU:O	2:G:223:LYS:HG3	2.05	0.57
2:A:147:PHE:O	2:A:151:ILE:HG12	2.05	0.57
3:H:166:HIS:HD2	3:H:168:VAL:HB	1.70	0.57
1:I:24:ASP:OD1	1:I:27:GLU:N	2.33	0.57
3:K:170:LEU:HD13	3:K:173:MET:HE2	1.87	0.57
1:I:86:ASN:OD1	1:I:128:GLN:NE2	2.28	0.56
2:G:390:VAL:HG21	2:J:378:LYS:HB2	1.87	0.56
2:B:240:ALA:HB3	2:B:243:ASP:OD1	2.04	0.56
1:L:42:ARG:NH2	1:L:46:HIS:HB3	2.20	0.56
3:K:130:GLY:HA3	3:K:177:GLU:HG3	1.86	0.56
2:G:204:VAL:O	2:G:208:THR:OG1	2.23	0.56
2:B:142:GLU:OE1	2:B:142:GLU:N	2.38	0.56
1:F:137:MET:O	1:F:141:LEU:HD13	2.05	0.56
1:F:18:LYS:NZ	1:F:115:ASP:OD1	2.39	0.56
1:F:43:ASP:HB3	1:F:44:ALA:C	2.31	0.56
1:I:42:ARG:NH1	1:I:46:HIS:HB3	2.21	0.56
1:I:85:HIS:CE1	3:H:173:MET:HE1	2.40	0.56
1:L:34:GLU:O	1:L:38:VAL:HG23	2.05	0.55
2:J:198:ASP:O	2:J:232:ARG:NH2	2.40	0.55
1:L:6:LEU:HD11	1:L:32:ILE:HD13	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:21:MET:HB3	1:L:23:PHE:CE2	2.41	0.55
1:L:138:LEU:HD22	2:J:88:TYR:CZ	2.41	0.55
1:F:45:LYS:HD3	2:A:381:ARG:HG2	1.88	0.55
1:F:99:CYS:HA	3:E:164:LEU:HD13	1.89	0.55
2:J:259:VAL:HG13	2:J:312:GLU:HG3	1.88	0.55
1:L:3:ARG:HD2	1:L:21:MET:SD	2.46	0.54
2:G:142:GLU:OE1	2:G:142:GLU:N	2.34	0.54
2:B:56:LYS:HG3	2:B:93:ILE:HG13	1.89	0.54
3:K:136:PRO:HB3	3:K:141:TRP:HD1	1.71	0.54
2:B:241:SER:HB3	2:A:282:GLU:OE2	2.07	0.54
2:J:289:ASN:OD1	2:J:329:ARG:NH2	2.39	0.54
2:A:32:ARG:HH12	2:A:36:GLU:HB2	1.71	0.54
1:L:61:ARG:HD2	1:L:66:TYR:CE1	2.42	0.54
2:G:271:PRO:HD2	2:G:277:ARG:HG3	1.89	0.54
1:F:16:LEU:HD21	1:F:114:VAL:HG21	1.89	0.54
2:A:142:GLU:OE1	2:A:142:GLU:N	2.36	0.54
1:D:78:LEU:HD13	2:B:348:LEU:HD11	1.90	0.54
1:F:42:ARG:CZ	1:F:46:HIS:HB3	2.38	0.54
3:K:121:PHE:CD1	3:K:122:PRO:HD2	2.42	0.54
2:A:61:LYS:O	2:A:65:ILE:HG13	2.08	0.53
3:C:186:ASP:HB3	3:C:189:LEU:HG	1.90	0.53
1:D:56:LYS:HD2	1:D:77:ASN:OD1	2.08	0.53
1:I:56:LYS:HG2	1:I:73:VAL:HA	1.90	0.53
2:J:100:VAL:HB	2:J:101:LEU:HD12	1.90	0.53
3:H:193:HIS:CD2	3:H:196:ARG:H	2.18	0.53
2:J:117:LYS:HG3	2:J:147:PHE:HE1	1.73	0.52
2:B:282:GLU:OE2	2:A:240:ALA:HB1	2.09	0.52
2:G:60:CYS:SG	2:G:93:ILE:HD12	2.50	0.52
2:G:147:PHE:O	2:G:151:ILE:HG12	2.10	0.52
1:F:24:ASP:N	1:F:27:GLU:OE1	2.34	0.52
1:F:97:ASN:ND2	3:E:163:SER:OG	2.43	0.52
2:G:198:ASP:O	2:G:232:ARG:NH2	2.43	0.52
2:G:198:ASP:OD1	2:G:199:GLN:N	2.43	0.52
2:G:13:LEU:O	2:G:17:ILE:HG13	2.10	0.52
1:D:142:GLU:OE1	2:B:87:ARG:NH1	2.43	0.52
1:L:60:ARG:NH2	3:K:173:MET:SD	2.63	0.52
3:K:139:PHE:HB2	3:K:193:HIS:CD2	2.45	0.52
2:B:124:ASN:O	2:B:128:MET:HG3	2.10	0.51
3:E:72:PRO:HG3	3:E:116:HIS:CE1	2.46	0.51
2:A:16:PHE:O	2:A:20:ILE:HG12	2.11	0.51
3:E:139:PHE:HB2	3:E:193:HIS:CD2	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:96:LEU:O	2:G:100:VAL:HG23	2.10	0.51
2:J:22:ASN:OD1	3:C:188:ARG:HD3	2.10	0.51
3:K:186:ASP:HB3	3:K:189:LEU:HG	1.92	0.51
1:F:60:ARG:HH21	1:F:81:LEU:HG	1.75	0.51
1:F:43:ASP:CB	1:F:44:ALA:HB3	2.40	0.51
2:J:76:HIS:O	2:J:80:VAL:HG23	2.11	0.51
2:J:341:ARG:NH1	3:K:177:GLU:OE2	2.41	0.51
3:K:105:ARG:O	3:K:109:ILE:HG13	2.11	0.51
3:H:91:LEU:HD12	3:H:141:TRP:HH2	1.74	0.50
3:C:117:THR:OG1	3:C:118:GLN:HG3	2.12	0.50
2:A:378:LYS:HG3	2:A:379:THR:HG23	1.93	0.50
3:E:105:ARG:O	3:E:109:ILE:HG13	2.11	0.50
1:L:125:GLU:OE2	2:J:247:TYR:OH	2.22	0.50
1:I:61:ARG:HD3	1:I:66:TYR:CE1	2.46	0.50
2:G:124:ASN:O	2:G:128:MET:HG3	2.12	0.50
2:G:350:LEU:HB3	2:G:388:ARG:O	2.12	0.50
3:C:139:PHE:HB2	3:C:193:HIS:ND1	2.26	0.50
2:J:80:VAL:HG21	2:J:108:LEU:HD22	1.93	0.50
1:L:76:ASN:H	2:J:263:ARG:HH22	1.58	0.50
1:F:42:ARG:HD3	1:F:59:TYR:OH	2.11	0.50
1:D:3:ARG:HG3	1:D:55:PHE:CZ	2.47	0.49
2:B:318:ILE:HG12	2:B:356:LEU:HD13	1.95	0.49
3:C:115:TYR:CD1	3:C:122:PRO:HD3	2.47	0.49
2:A:31:LYS:O	2:A:35:LYS:HG3	2.13	0.49
1:D:115:ASP:HB3	2:B:133:HIS:CD2	2.47	0.49
1:I:138:LEU:HD23	1:I:141:LEU:HD12	1.94	0.49
3:E:123:ASP:O	3:E:126:ASN:ND2	2.45	0.49
3:E:172:GLY:O	3:E:178:ARG:HA	2.13	0.49
1:I:61:ARG:HD3	1:I:66:TYR:CZ	2.47	0.49
1:F:43:ASP:CG	1:F:44:ALA:HB3	2.38	0.48
1:F:43:ASP:HB3	1:F:45:LYS:H	1.75	0.48
1:F:82:GLU:OE1	2:A:249:TYR:OH	2.24	0.48
2:B:218:ASN:ND2	2:B:221:GLU:HB2	2.27	0.48
1:I:130:LYS:HE2	2:G:247:TYR:HE1	1.77	0.48
2:G:174:ARG:NH1	2:G:178:THR:OG1	2.46	0.48
2:J:376:ALA:HA	2:J:379:THR:HG22	1.95	0.48
3:E:79:MET:HE3	3:E:124:TRP:CD2	2.47	0.48
1:L:52:PHE:HB2	1:L:55:PHE:H	1.77	0.48
2:J:32:ARG:HH12	2:J:36:GLU:HG2	1.77	0.48
2:A:361:PHE:O	3:C:76:LEU:HB2	2.13	0.48
3:E:144:LYS:HE2	3:E:186:ASP:HB2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:138:LEU:HD22	2:G:88:TYR:CZ	2.48	0.48
2:A:55:LYS:O	2:A:59:VAL:HG23	2.14	0.48
2:A:66:PHE:HA	2:A:70:HIS:O	2.14	0.48
3:K:90:PHE:CE1	3:K:94:LYS:HE3	2.49	0.48
2:G:232:ARG:O	2:G:236:ILE:HG13	2.14	0.48
2:B:156:VAL:HG22	2:B:191:ARG:CB	2.43	0.48
2:A:96:LEU:O	2:A:100:VAL:HG23	2.14	0.47
3:H:103:SER:HB3	3:H:167:PRO:HB3	1.95	0.47
2:G:192:VAL:O	2:G:195:LEU:HB2	2.14	0.47
3:C:166:HIS:CD2	3:C:167:PRO:HD2	2.49	0.47
1:L:74:ASN:OD1	1:L:74:ASN:N	2.47	0.47
3:H:117:THR:HG22	3:H:118:GLN:HG2	1.95	0.47
2:J:367:LYS:HG2	2:J:396:MET:HB2	1.95	0.47
2:A:107:GLU:OE2	2:A:110:ARG:NH2	2.47	0.47
3:H:64:UNK:C	3:H:66:UNK:HA	2.44	0.47
2:J:147:PHE:O	2:J:151:ILE:HG12	2.14	0.47
2:J:185:MET:HB3	2:J:189:THR:HG21	1.95	0.47
2:A:95:TYR:O	2:A:99:SER:OG	2.13	0.47
1:L:119:LEU:HB3	2:J:170:LEU:HD13	1.97	0.47
1:I:116:GLU:O	1:I:124:ARG:HD3	2.14	0.47
2:A:222:PHE:O	2:A:225:SER:OG	2.27	0.47
1:L:27:GLU:HG3	1:L:30:LYS:HE3	1.96	0.47
2:G:103:ASN:OD1	2:G:104:SER:N	2.48	0.47
2:G:195:LEU:HA	2:G:195:LEU:HD23	1.80	0.47
1:F:21:MET:HB2	1:F:23:PHE:CE2	2.50	0.47
1:I:42:ARG:NH1	1:I:47:THR:O	2.45	0.47
2:B:100:VAL:HB	2:B:101:LEU:HD12	1.96	0.47
2:J:186:GLY:H	2:J:189:THR:HG23	1.80	0.47
2:G:301:GLN:HE21	3:H:173:MET:CE	2.28	0.47
1:I:21:MET:HE2	1:I:23:PHE:CZ	2.50	0.46
2:B:156:VAL:CG2	2:B:191:ARG:HB2	2.43	0.46
2:B:176:TYR:OH	2:B:217:LYS:HD2	2.15	0.46
2:J:341:ARG:HD2	3:K:177:GLU:OE2	2.15	0.46
2:G:347:TYR:CE2	2:G:388:ARG:HD2	2.50	0.46
2:A:140:SER:OG	2:A:142:GLU:OE1	2.33	0.46
1:L:96:HIS:HB2	3:C:93:GLU:HG2	1.98	0.46
1:I:8:GLN:NE2	1:I:36:HIS:HB2	2.30	0.46
2:G:105:ASN:O	2:G:109:ILE:HG13	2.16	0.46
3:K:74:VAL:HG13	3:K:75:PRO:HD2	1.98	0.46
3:C:136:PRO:HB3	3:C:141:TRP:HD1	1.80	0.46
2:B:392:LEU:HG	2:B:396:MET:HE3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:293:GLU:HG3	2:J:294:PRO:HD2	1.97	0.46
2:A:341:ARG:HD2	3:E:177:GLU:OE1	2.16	0.46
3:K:97:LEU:HD23	3:K:97:LEU:HA	1.73	0.46
3:C:97:LEU:HD23	3:C:143:TYR:CE2	2.51	0.46
3:E:117:THR:HG22	3:E:118:GLN:HG2	1.98	0.46
2:A:151:ILE:HB	2:A:152:PRO:HD3	1.97	0.46
2:A:170:LEU:HD23	2:A:170:LEU:HA	1.80	0.45
3:H:89:HIS:O	3:H:93:GLU:HG3	2.16	0.45
3:H:115:TYR:CD1	3:H:122:PRO:HD3	2.51	0.45
1:F:52:PHE:O	1:F:53:ARG:HB3	2.16	0.45
2:J:128:MET:HE2	2:J:128:MET:HB3	1.82	0.45
1:L:78:LEU:HD23	1:L:78:LEU:HA	1.73	0.45
2:G:123:ARG:HH22	2:G:159:ASP:HB2	1.81	0.45
3:E:115:TYR:CD1	3:E:122:PRO:HD3	2.51	0.45
3:E:121:PHE:CD1	3:E:122:PRO:HD2	2.51	0.45
1:I:34:GLU:O	1:I:38:VAL:HG23	2.16	0.45
2:B:16:PHE:CE1	2:B:36:GLU:HG3	2.51	0.45
2:J:105:ASN:O	2:J:109:ILE:HG13	2.16	0.45
1:L:30:LYS:O	1:L:34:GLU:HG3	2.16	0.45
2:J:193:VAL:HG11	2:J:224:THR:HB	1.98	0.45
3:E:56:UNK:CB	3:E:96:GLY:HA3	2.47	0.45
3:C:165:LEU:HD23	3:C:165:LEU:HA	1.73	0.45
1:I:56:LYS:HE3	1:I:73:VAL:O	2.16	0.45
2:G:170:LEU:HD23	2:G:170:LEU:HA	1.75	0.45
1:L:117:MET:HE3	1:L:117:MET:HB3	1.80	0.45
2:G:370:ILE:HA	2:G:396:MET:HE1	1.99	0.45
2:A:108:LEU:HD23	2:A:108:LEU:HA	1.78	0.45
3:H:100:LEU:HD23	3:H:183:TRP:CZ3	2.51	0.45
2:G:47:ASP:OD1	2:G:48:LYS:N	2.48	0.45
2:J:117:LYS:HG3	2:J:147:PHE:CE1	2.51	0.45
2:J:123:ARG:HH22	2:J:159:ASP:CG	2.25	0.45
2:A:125:PRO:C	2:A:127:PHE:N	2.74	0.45
3:C:166:HIS:CG	3:C:167:PRO:HD2	2.51	0.45
1:F:51:GLU:OE1	2:A:388:ARG:NH2	2.50	0.45
2:B:96:LEU:O	2:B:100:VAL:HG23	2.16	0.45
3:H:147:PRO:HA	3:H:180:VAL:O	2.17	0.45
3:C:111:ASP:HB3	3:C:122:PRO:HB3	1.99	0.45
2:B:324:PRO:HB3	2:B:361:PHE:CZ	2.52	0.45
2:J:212:THR:O	2:J:216:GLN:HG3	2.17	0.45
1:D:117:MET:HE2	1:D:117:MET:HB3	1.88	0.44
2:A:13:LEU:O	2:A:17:ILE:HG13	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:288:LEU:HD13	2:J:329:ARG:HB3	1.98	0.44
3:C:97:LEU:HD23	3:C:143:TYR:CZ	2.52	0.44
2:B:129:GLY:O	2:B:133:HIS:ND1	2.50	0.44
2:B:170:LEU:HD23	2:B:170:LEU:HA	1.70	0.44
2:B:185:MET:HE1	2:B:214:LEU:HD22	1.98	0.44
2:B:296:LYS:HE2	2:A:285:GLU:OE2	2.18	0.44
2:G:280:LEU:O	2:G:283:CYS:HB3	2.17	0.44
3:E:74:VAL:HG22	3:E:75:PRO:HD2	1.98	0.44
1:F:28:LYS:O	1:F:32:ILE:HG13	2.17	0.44
2:J:204:VAL:O	2:J:208:THR:OG1	2.29	0.44
2:A:324:PRO:HB3	2:A:361:PHE:CZ	2.53	0.44
2:J:209:SER:HG	2:J:260:LYS:NZ	2.15	0.44
2:J:377:LEU:HD12	2:J:377:LEU:HA	1.72	0.44
2:A:33:ILE:HG23	2:A:65:ILE:HD13	1.99	0.44
2:A:116:ILE:HG23	2:A:131:ALA:HB1	1.99	0.44
2:B:182:LEU:O	2:B:184:PRO:HD3	2.18	0.44
1:D:115:ASP:HB3	2:B:133:HIS:HD2	1.82	0.44
2:B:44:PHE:C	2:B:46:GLY:H	2.26	0.44
3:E:79:MET:HE3	3:E:124:TRP:CG	2.53	0.44
3:K:95:GLY:HA2	3:K:96:GLY:HA2	1.47	0.44
3:K:110:LEU:O	3:K:114:ILE:HG12	2.18	0.44
2:B:13:LEU:O	2:B:17:ILE:HG13	2.17	0.43
2:J:272:GLU:O	2:J:274:PRO:HD3	2.18	0.43
2:A:346:ARG:NH1	2:A:380:GLU:CD	2.76	0.43
3:C:108:ASP:O	3:C:112:LEU:HD23	2.18	0.43
1:L:52:PHE:C	1:L:54:ASN:H	2.25	0.43
1:F:85:HIS:NE2	1:F:89:GLU:OE2	2.52	0.43
1:L:24:ASP:OD1	1:L:27:GLU:HB3	2.18	0.43
3:H:100:LEU:HD23	3:H:183:TRP:CE3	2.54	0.43
3:H:196:ARG:HA	3:H:203:PHE:CE2	2.53	0.43
2:G:61:LYS:O	2:G:65:ILE:HG13	2.18	0.43
3:C:79:MET:HE3	3:C:124:TRP:CG	2.53	0.43
1:I:31:LEU:HD23	1:I:31:LEU:HA	1.78	0.43
2:J:52:GLY:HA3	2:J:90:GLU:OE2	2.18	0.43
2:A:26:LYS:O	2:A:30:ILE:HD12	2.18	0.43
3:H:166:HIS:CD2	3:H:168:VAL:H	2.37	0.43
3:K:152:LYS:HA	3:K:152:LYS:HD3	1.87	0.43
1:I:45:LYS:HB3	1:I:45:LYS:HE2	1.73	0.43
2:J:172:LEU:HD23	2:J:210:LEU:HD21	2.01	0.43
1:L:132:LEU:O	1:L:136:LEU:HD23	2.19	0.43
2:B:212:THR:O	2:B:216:GLN:HG3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:91:LYS:NZ	2:J:119:ASP:OD2	2.35	0.43
2:J:96:LEU:O	2:J:100:VAL:HG23	2.19	0.43
3:E:96:GLY:O	3:E:97:LEU:HB2	2.18	0.43
1:L:84:ILE:O	1:L:88:VAL:HG23	2.19	0.43
1:F:21:MET:HE3	1:F:21:MET:HB3	1.83	0.43
1:I:43:ASP:OD2	1:I:45:LYS:HE3	2.19	0.43
2:G:241:SER:HB3	2:G:244:LEU:HD13	2.00	0.43
2:G:341:ARG:HD2	3:H:177:GLU:OE2	2.19	0.43
2:B:11:ARG:O	2:B:15:VAL:HG23	2.19	0.42
2:B:271:PRO:O	2:B:277:ARG:NH2	2.51	0.42
1:F:110:VAL:O	1:F:114:VAL:HG23	2.19	0.42
2:A:77:MET:HE2	2:A:77:MET:HB2	1.92	0.42
1:F:12:GLY:HA2	1:F:36:HIS:CD2	2.54	0.42
1:F:127:SER:O	1:F:131:VAL:HG23	2.20	0.42
2:A:189:THR:O	2:A:192:VAL:HG22	2.19	0.42
1:I:87:PHE:CZ	1:I:114:VAL:HG22	2.53	0.42
2:G:83:LEU:O	2:G:91:LYS:HE3	2.19	0.42
3:E:115:TYR:CD1	3:E:121:PHE:HA	2.54	0.42
3:H:125:GLN:HG2	3:H:127:TYR:HE1	1.85	0.42
2:B:282:GLU:OE1	2:A:241:SER:HB3	2.20	0.42
3:H:97:LEU:HD11	3:H:110:LEU:HD21	2.01	0.42
1:I:9:ASN:OD1	1:I:13:LYS:N	2.52	0.42
1:I:60:ARG:NH1	1:I:85:HIS:HB2	2.35	0.42
2:B:317:ILE:HG22	2:B:327:LEU:HD23	2.02	0.42
1:F:98:VAL:CG1	3:E:164:LEU:HD22	2.48	0.42
1:F:138:LEU:HA	1:F:141:LEU:HD22	2.01	0.42
2:B:55:LYS:O	2:B:59:VAL:HG23	2.19	0.42
2:J:322:SER:OG	2:J:323:GLU:HG2	2.20	0.42
2:A:214:LEU:HD23	2:A:214:LEU:HA	1.81	0.42
3:E:196:ARG:HA	3:E:203:PHE:CE2	2.55	0.42
1:I:142:GLU:C	2:G:89:THR:HG23	2.45	0.42
2:J:284:LEU:HD23	2:J:284:LEU:HA	1.78	0.42
2:J:378:LYS:HE2	2:J:378:LYS:HB3	1.87	0.42
3:E:105:ARG:HH22	3:E:167:PRO:HG3	1.85	0.42
1:L:31:LEU:O	1:L:35:VAL:HG23	2.19	0.42
2:B:83:LEU:O	2:B:91:LYS:HE3	2.20	0.42
2:B:196:LEU:HD12	2:B:196:LEU:HA	1.80	0.42
1:I:43:ASP:OD1	1:I:44:ALA:N	2.50	0.42
3:H:139:PHE:HB2	3:H:193:HIS:ND1	2.34	0.42
1:L:8:GLN:NE2	1:L:32:ILE:HG23	2.34	0.41
2:B:116:ILE:HG23	2:B:131:ALA:HB1	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:147:PHE:O	2:B:151:ILE:HG12	2.19	0.41
3:K:79:MET:HE3	3:K:138:THR:HG21	2.02	0.41
1:D:43:ASP:CG	1:D:44:ALA:N	2.78	0.41
1:I:21:MET:HE3	1:I:21:MET:HB3	1.66	0.41
2:G:338:LEU:HD23	2:G:338:LEU:HA	1.87	0.41
2:B:377:LEU:HD12	2:B:377:LEU:HA	1.66	0.41
2:J:151:ILE:HB	2:J:152:PRO:HD3	2.02	0.41
2:G:362:SER:O	2:G:366:VAL:HG23	2.20	0.41
2:J:82:LEU:HD23	2:J:82:LEU:HA	1.89	0.41
2:A:103:ASN:OD1	2:A:104:SER:N	2.53	0.41
1:L:51:GLU:HG3	2:J:347:TYR:OH	2.21	0.41
2:J:123:ARG:NH2	2:J:159:ASP:OD2	2.51	0.41
1:L:30:LYS:HE3	1:L:30:LYS:HB3	1.83	0.41
2:G:125:PRO:O	2:G:128:MET:N	2.53	0.41
2:J:265:LEU:HA	2:J:265:LEU:HD23	1.84	0.41
1:D:15:ARG:HA	1:D:15:ARG:HD2	1.71	0.41
2:B:61:LYS:O	2:B:65:ILE:HG13	2.20	0.41
2:B:158:GLY:O	2:B:159:ASP:OD1	2.39	0.41
3:H:141:TRP:NE1	3:H:143:TYR:HB2	2.35	0.41
3:C:121:PHE:HA	3:C:122:PRO:HD3	1.91	0.41
1:F:90:VAL:CG2	1:F:128:GLN:HG2	2.51	0.41
3:H:104:GLN:N	3:H:174:ASP:OD2	2.23	0.41
3:C:74:VAL:HA	3:C:75:PRO:HD3	1.94	0.41
1:D:38:VAL:CG1	1:D:50:VAL:HG11	2.48	0.41
1:D:52:PHE:HB2	1:D:55:PHE:O	2.20	0.41
1:I:78:LEU:HD23	1:I:78:LEU:HA	1.78	0.41
2:G:67:LEU:HD23	2:G:67:LEU:HA	1.90	0.41
2:B:108:LEU:HD23	2:B:108:LEU:HA	1.92	0.41
2:B:366:VAL:HG12	2:B:396:MET:HE2	2.03	0.41
2:J:91:LYS:HZ1	2:J:119:ASP:CG	2.21	0.41
2:A:212:THR:O	2:A:216:GLN:HG2	2.21	0.41
1:F:8:GLN:HE21	1:F:8:GLN:HB3	1.68	0.41
1:I:50:VAL:HB	1:I:57:ILE:HB	2.03	0.41
2:G:68:LEU:HD23	2:G:68:LEU:HA	1.90	0.41
2:G:132:LEU:HD23	2:G:132:LEU:HA	1.95	0.41
2:B:31:LYS:O	2:B:35:LYS:HG3	2.20	0.41
2:B:47:ASP:OD1	2:B:48:LYS:N	2.54	0.41
2:B:193:VAL:O	2:B:196:LEU:HB2	2.21	0.41
2:J:294:PRO:HA	2:J:295:PRO:HD3	1.82	0.41
3:H:83:ALA:O	3:H:86:ASP:HB2	2.21	0.41
1:L:76:ASN:N	2:J:263:ARG:HH22	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:31:LEU:O	1:F:35:VAL:HG23	2.20	0.41
1:I:3:ARG:HG3	1:I:55:PHE:CE2	2.56	0.41
1:I:76:ASN:ND2	2:G:312:GLU:OE1	2.48	0.41
2:B:32:ARG:HH12	2:B:36:GLU:HG2	1.85	0.41
2:J:201:LEU:HA	2:J:201:LEU:HD23	1.75	0.41
3:H:136:PRO:HB3	3:H:141:TRP:HD1	1.84	0.41
1:I:35:VAL:O	1:I:39:VAL:HB	2.21	0.40
2:A:182:LEU:O	2:A:184:PRO:HD3	2.21	0.40
3:H:121:PHE:HA	3:H:122:PRO:HD3	1.88	0.40
1:D:3:ARG:HD2	1:D:21:MET:SD	2.61	0.40
1:L:43:ASP:OD1	1:L:44:ALA:N	2.48	0.40
1:I:3:ARG:HG3	1:I:55:PHE:CZ	2.56	0.40
2:G:301:GLN:NE2	3:H:173:MET:HE2	2.30	0.40
2:J:124:ASN:O	2:J:128:MET:HG3	2.21	0.40
1:L:76:ASN:H	2:J:263:ARG:NH2	2.19	0.40
2:J:223:LYS:O	2:J:226:VAL:HG23	2.22	0.40
3:E:77:ARG:N	3:E:119:GLY:O	2.51	0.40
1:F:132:LEU:HD23	1:F:132:LEU:HA	1.97	0.40
2:J:108:LEU:HA	2:J:108:LEU:HD23	1.78	0.40
2:J:176:TYR:CE2	2:J:214:LEU:HD23	2.57	0.40
3:C:74:VAL:HG13	3:C:75:PRO:HD2	2.03	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	D	140/142 (99%)	137 (98%)	2 (1%)	1 (1%)	18	44
1	F	140/142 (99%)	135 (96%)	4 (3%)	1 (1%)	18	44
1	I	140/142 (99%)	137 (98%)	2 (1%)	1 (1%)	18	44
1	L	140/142 (99%)	137 (98%)	2 (1%)	1 (1%)	18	44

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	A	381/398 (96%)	378 (99%)	2 (0%)	1 (0%)	36	62
2	B	386/398 (97%)	383 (99%)	3 (1%)	0	100	100
2	G	386/398 (97%)	376 (97%)	6 (2%)	4 (1%)	12	35
2	J	386/398 (97%)	380 (98%)	5 (1%)	1 (0%)	36	62
3	C	130/155 (84%)	128 (98%)	1 (1%)	1 (1%)	16	41
3	E	130/155 (84%)	128 (98%)	2 (2%)	0	100	100
3	H	128/155 (83%)	128 (100%)	0	0	100	100
3	K	129/155 (83%)	128 (99%)	1 (1%)	0	100	100
All	All	2616/2780 (94%)	2575 (98%)	30 (1%)	11 (0%)	30	56

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	G	159	ASP
1	L	96	HIS
2	G	46	GLY
1	D	96	HIS
1	F	96	HIS
1	I	96	HIS
2	G	158	GLY
3	C	97	LEU
2	G	125	PRO
2	J	46	GLY
2	A	125	PRO

### 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	D	131/131 (100%)	126 (96%)	5 (4%)	29	61
1	F	131/131 (100%)	128 (98%)	3 (2%)	44	74
1	I	131/131 (100%)	123 (94%)	8 (6%)	17	43

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	L	131/131 (100%)	126 (96%)	5 (4%)	29	61
2	A	338/346 (98%)	332 (98%)	6 (2%)	51	78
2	B	340/346 (98%)	333 (98%)	7 (2%)	47	75
2	G	340/346 (98%)	336 (99%)	4 (1%)	63	84
2	J	340/346 (98%)	332 (98%)	8 (2%)	43	73
3	C	118/126 (94%)	114 (97%)	4 (3%)	32	64
3	E	118/126 (94%)	115 (98%)	3 (2%)	42	72
3	H	116/126 (92%)	114 (98%)	2 (2%)	53	80
3	K	117/126 (93%)	115 (98%)	2 (2%)	53	80
All	All	2351/2412 (98%)	2294 (98%)	57 (2%)	43	73

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	15	ARG
1	D	49	PHE
1	D	56	LYS
1	D	78	LEU
1	D	136	LEU
1	L	49	PHE
1	L	50	VAL
1	L	73	VAL
1	L	116	GLU
1	L	117	MET
1	F	49	PHE
1	F	137	MET
1	F	140	SER
1	I	6	LEU
1	I	15	ARG
1	I	18	LYS
1	I	39	VAL
1	I	49	PHE
1	I	52	PHE
1	I	56	LYS
1	I	82	GLU
2	G	10	MET
2	G	33	ILE
2	G	60	CYS
2	G	243	ASP

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Mol	Chain	Res	Type
2	B	10	MET
2	B	60	CYS
2	B	92	GLN
2	B	156	VAL
2	B	159	ASP
2	B	196	LEU
2	B	332	ASN
2	J	10	MET
2	J	93	ILE
2	J	104	SER
2	J	120	LEU
2	J	153	LYS
2	J	293	GLU
2	J	296	LYS
2	J	356	LEU
2	A	91	LYS
2	A	93	ILE
2	A	164	VAL
2	A	232	ARG
2	A	246	ASP
2	A	338	LEU
3	E	74	VAL
3	E	146	VAL
3	E	164	LEU
3	H	146	VAL
3	H	179	GLU
3	K	76	LEU
3	K	177	GLU
3	C	76	LEU
3	C	98	GLU
3	C	146	VAL
3	C	192	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	77	ASN
1	F	8	GLN
2	G	70	HIS
2	G	114	ASN
2	G	332	ASN
2	G	363	HIS

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Mol	Chain	Res	Type
2	B	70	HIS
2	B	76	HIS
2	B	92	GLN
2	B	218	ASN
2	B	320	HIS
2	B	333	GLN
2	J	70	HIS
2	J	103	ASN
2	J	166	GLN
2	J	218	ASN
2	J	333	GLN
2	A	266	GLN
3	E	73	GLN
3	E	116	HIS
3	E	126	ASN
3	H	126	ASN
3	H	166	HIS
3	K	125	GLN
3	C	118	GLN
3	C	161	ASN

### 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 ⓘ

There are no oligosaccharides in this entry.

### 5.6 Ligand geometry ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	D	142/142 (100%)	-0.28	2 (1%) 73 66	18, 27, 52, 92	0
1	F	142/142 (100%)	-0.05	2 (1%) 73 66	27, 37, 67, 98	0
1	I	142/142 (100%)	-0.19	3 (2%) 63 55	15, 24, 53, 81	0
1	L	142/142 (100%)	-0.10	4 (2%) 55 46	20, 30, 69, 104	0
2	A	385/398 (96%)	0.11	3 (0%) 82 77	26, 45, 70, 96	0
2	B	388/398 (97%)	-0.08	0 100 100	23, 36, 56, 87	0
2	G	388/398 (97%)	-0.16	2 (0%) 87 83	20, 34, 55, 78	0
2	J	388/398 (97%)	-0.06	5 (1%) 75 68	24, 38, 63, 88	0
3	C	132/155 (85%)	-0.01	1 (0%) 82 77	23, 36, 61, 77	0
3	E	132/155 (85%)	-0.05	1 (0%) 82 77	19, 37, 56, 69	0
3	H	130/155 (83%)	0.26	3 (2%) 61 52	24, 45, 76, 92	0
3	K	131/155 (84%)	0.13	2 (1%) 72 64	22, 39, 72, 89	0
All	All	2642/2780 (95%)	-0.04	28 (1%) 78 72	15, 36, 66, 104	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	44	ALA	4.0
3	H	76	LEU	4.0
2	A	46	GLY	3.7
3	K	96	GLY	3.5
1	L	44	ALA	3.3
3	K	75	PRO	3.2
1	F	44	ALA	3.0
3	H	74	VAL	2.8
2	A	102	VAL	2.7
1	I	43	ASP	2.7
2	J	242	THR	2.6

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Mol	Chain	Res	Type	RSRZ
3	E	72	PRO	2.5
1	D	43	ASP	2.5
1	L	46	HIS	2.5
1	L	45	LYS	2.4
3	H	75	PRO	2.3
1	I	52	PHE	2.2
2	J	272	GLU	2.2
3	C	72	PRO	2.2
2	J	24	LYS	2.2
1	I	24	ASP	2.2
1	L	55	PHE	2.1
2	G	123	ARG	2.1
2	J	47	ASP	2.1
2	A	103	ASN	2.1
2	G	218	ASN	2.0
2	J	381	ARG	2.0
1	F	47	THR	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.