



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:25 PM GMT

PDB ID : 1OSG  
Title : Complex between BAFF and a BR3 derived peptide presented in a beta-hairpin scaffold  
Authors : Gordon, N.C.; Pan, B.; Hymowitz, S.G.; Yin, J.P.; Kelley, R.F.; Cochran, A.G.; Yan, M.; Dixit, V.M.; Fairbrother, W.J.; Starovasnik, M.A.  
Deposited on : 2003-03-19  
Resolution : 3.00 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  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

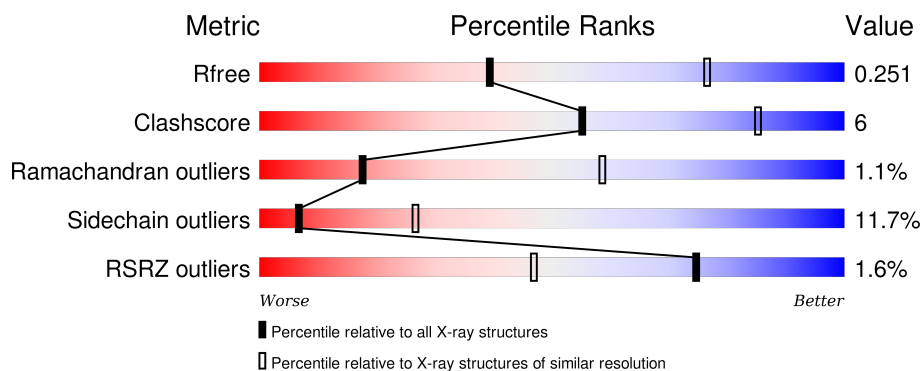
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.00 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.






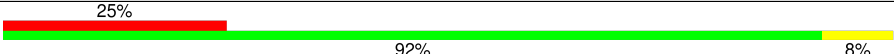
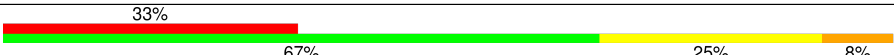
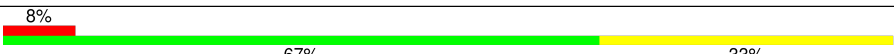

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	208	
1	B	208	
1	C	208	
1	D	208	
1	E	208	

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Mol	Chain	Length	Quality of chain
1	F	208	
2	G	12	
2	H	12	
2	I	12	
2	J	12	
2	K	12	
2	L	12	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	MG	A	701	-	-	-	X
3	MG	A	702	-	-	-	X
3	MG	D	802	-	-	-	X
3	MG	E	801	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7531 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 Tumor necrosis factor ligand superfamily member 13B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	144	Total	C	N	O	S	0	0	0
			1143	736	184	218	5			
1	B	144	Total	C	N	O	S	0	0	0
			1143	736	184	218	5			
1	C	144	Total	C	N	O	S	0	0	0
			1143	736	184	218	5			
1	D	144	Total	C	N	O	S	0	0	0
			1143	736	184	218	5			
1	E	144	Total	C	N	O	S	0	0	0
			1143	736	184	218	5			
1	F	144	Total	C	N	O	S	0	0	0
			1143	736	184	218	5			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	78	GLY	-	CLONING ARTIFACT	UNP Q9Y275
A	79	SER	-	CLONING ARTIFACT	UNP Q9Y275
A	80	HIS	-	CLONING ARTIFACT	UNP Q9Y275
A	81	MET	-	CLONING ARTIFACT	UNP Q9Y275
B	78	GLY	-	CLONING ARTIFACT	UNP Q9Y275
B	79	SER	-	CLONING ARTIFACT	UNP Q9Y275
B	80	HIS	-	CLONING ARTIFACT	UNP Q9Y275
B	81	MET	-	CLONING ARTIFACT	UNP Q9Y275
C	78	GLY	-	CLONING ARTIFACT	UNP Q9Y275
C	79	SER	-	CLONING ARTIFACT	UNP Q9Y275
C	80	HIS	-	CLONING ARTIFACT	UNP Q9Y275
C	81	MET	-	CLONING ARTIFACT	UNP Q9Y275
D	78	GLY	-	CLONING ARTIFACT	UNP Q9Y275
D	79	SER	-	CLONING ARTIFACT	UNP Q9Y275
D	80	HIS	-	CLONING ARTIFACT	UNP Q9Y275
D	81	MET	-	CLONING ARTIFACT	UNP Q9Y275
E	78	GLY	-	CLONING ARTIFACT	UNP Q9Y275

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Chain	Residue	Modelled	Actual	Comment	Reference
E	79	SER	-	CLONING ARTIFACT	UNP Q9Y275
E	80	HIS	-	CLONING ARTIFACT	UNP Q9Y275
E	81	MET	-	CLONING ARTIFACT	UNP Q9Y275
F	78	GLY	-	CLONING ARTIFACT	UNP Q9Y275
F	79	SER	-	CLONING ARTIFACT	UNP Q9Y275
F	80	HIS	-	CLONING ARTIFACT	UNP Q9Y275
F	81	MET	-	CLONING ARTIFACT	UNP Q9Y275

- Molecule 2 is a protein called BR3 derived PEPTIDE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	12	Total	C	N	O	S	0	0	0
			109	72	21	14	2			
2	J	12	Total	C	N	O	S	0	0	0
			109	72	21	14	2			
2	H	12	Total	C	N	O	S	0	0	0
			109	72	21	14	2			
2	I	12	Total	C	N	O	S	0	0	0
			104	69	19	14	2			
2	K	12	Total	C	N	O	S	0	0	0
			109	72	21	14	2			
2	L	12	Total	C	N	O	S	0	0	0
			109	72	21	14	2			

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	O	0	0
			3	3		
4	B	4	Total	O	0	0
			4	4		

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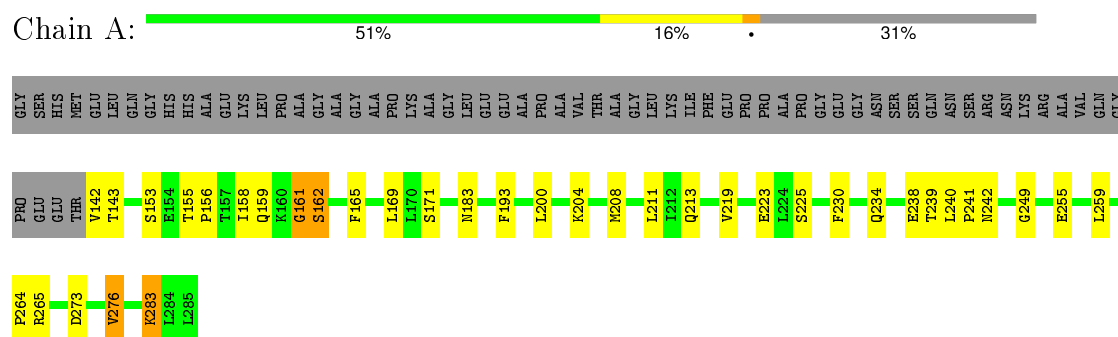
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	1	Total 1	O 1	0	0
4	E	3	Total 3	O 3	0	0
4	F	5	Total 5	O 5	0	0
4	G	1	Total 1	O 1	0	0
4	I	1	Total 1	O 1	0	0
4	K	1	Total 1	O 1	0	0
4	L	1	Total 1	O 1	0	0

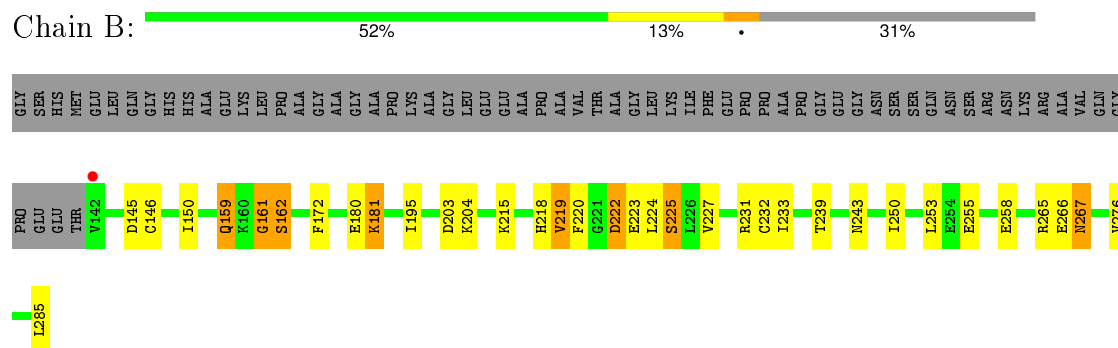
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

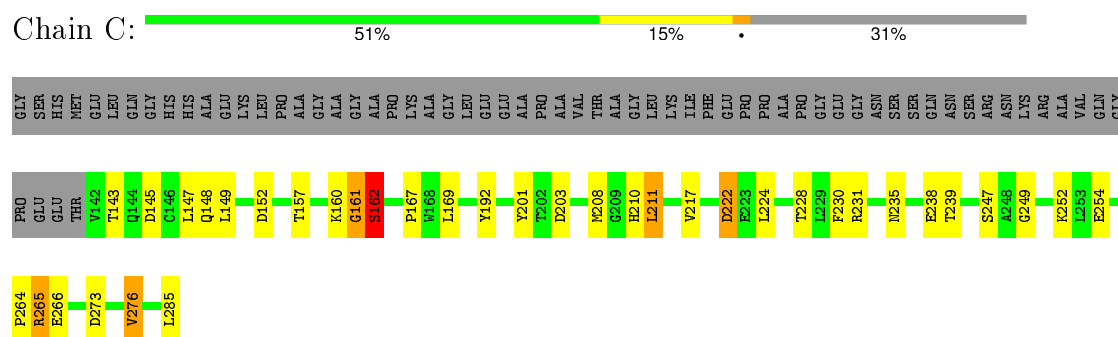
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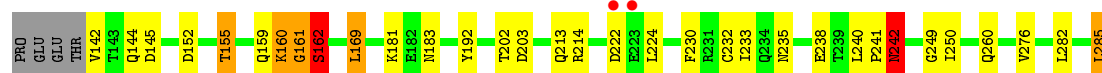
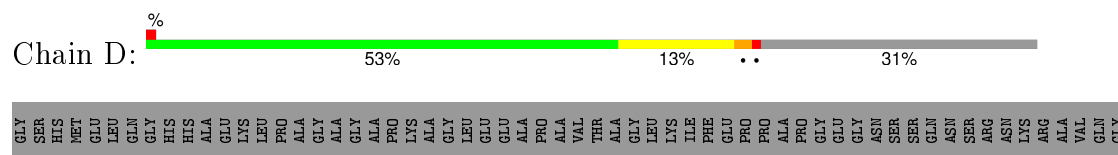
- Molecule 1: Tumor necrosis factor ligand superfamily member 13B



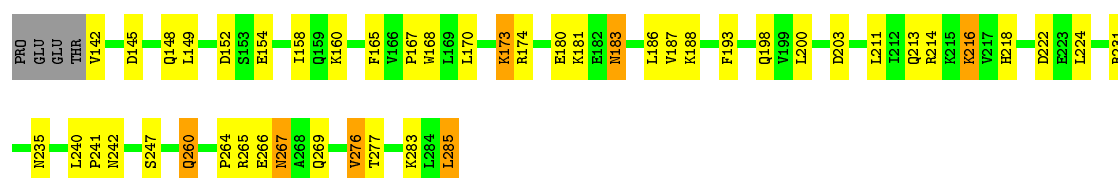
- Molecule 1: Tumor necrosis factor ligand superfamily member 13B



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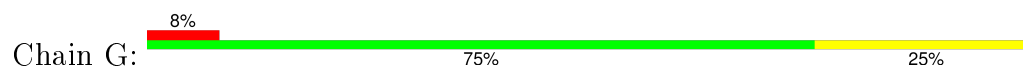
- Molecule 1: Tumor necrosis factor ligand superfamily member 13B



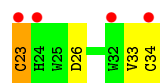
- Molecule 1: Tumor necrosis factor ligand superfamily member 13B



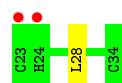
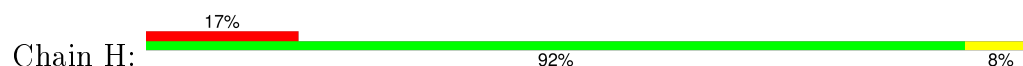
- Molecule 2: BR3 derived PEPTIDE



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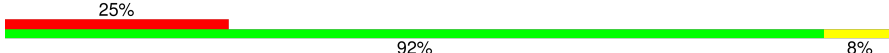


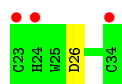
- Molecule 2: BR3 derived PEPTIDE





## ● Molecule 2: BR3 derived PEPTIDE

Chain I: 



## ● Molecule 2: BR3 derived PEPTIDE

Chain K: 



## ● Molecule 2: BR3 derived PEPTIDE

Chain L: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	121.63Å 121.63Å 157.21Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 3.00 29.85 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.6 (30.00-3.00) 99.7 (29.85-3.00)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.18 (at 3.00Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, $R_{free}$	0.213 , 0.286 0.193 , 0.251	Depositor DCC
$R_{free}$ test set	2536 reflections (10.67%)	DCC
Wilson B-factor (Å <sup>2</sup> )	53.3	Xtriage
Anisotropy	0.013	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 35.0	EDS
Estimated twinning fraction	0.056 for h,-h-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 26307 reflections (0.004%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	7531	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.67% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG

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	0.62	0/1165	0.80	1/1574 (0.1%)
1	B	0.59	0/1165	0.78	3/1574 (0.2%)
1	C	0.59	0/1165	0.77	4/1574 (0.3%)
1	D	0.58	0/1165	0.80	3/1574 (0.2%)
1	E	0.56	0/1165	0.82	3/1574 (0.2%)
1	F	0.59	0/1165	0.78	3/1574 (0.2%)
2	G	0.48	0/114	0.83	1/156 (0.6%)
2	H	0.50	0/114	0.71	0/156
2	I	0.54	0/108	0.80	1/148 (0.7%)
2	J	0.54	0/114	0.74	1/156 (0.6%)
2	K	0.48	0/114	0.85	1/156 (0.6%)
2	L	0.48	0/114	0.87	1/156 (0.6%)
All	All	0.58	0/7668	0.79	22/10372 (0.2%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	26	ASP	CB-CG-OD2	6.87	124.48	118.30
1	F	203	ASP	CB-CG-OD2	6.74	124.37	118.30
1	B	203	ASP	CB-CG-OD2	6.64	124.28	118.30
1	E	152	ASP	CB-CG-OD2	6.50	124.15	118.30
1	B	145	ASP	CB-CG-OD2	6.39	124.05	118.30
2	G	26	ASP	CB-CG-OD2	6.30	123.97	118.30
1	E	145	ASP	CB-CG-OD2	6.30	123.97	118.30
1	A	273	ASP	CB-CG-OD2	6.03	123.72	118.30
1	D	203	ASP	CB-CG-OD2	5.88	123.60	118.30
2	J	26	ASP	CB-CG-OD2	5.67	123.40	118.30
1	C	203	ASP	CB-CG-OD2	5.56	123.30	118.30
1	D	145	ASP	CB-CG-OD2	5.49	123.24	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	26	ASP	CB-CG-OD2	5.41	123.17	118.30
2	I	26	ASP	CB-CG-OD2	5.36	123.12	118.30
1	F	145	ASP	CB-CG-OD2	5.35	123.12	118.30
1	E	222	ASP	CB-CG-OD2	5.35	123.11	118.30
1	B	222	ASP	CB-CG-OD2	5.32	123.08	118.30
1	C	145	ASP	CB-CG-OD2	5.18	122.97	118.30
1	F	257	ASP	CB-CG-OD2	5.17	122.96	118.30
1	D	222	ASP	CB-CG-OD2	5.13	122.91	118.30
1	C	273	ASP	CB-CG-OD2	5.04	122.84	118.30
1	C	222	ASP	CB-CG-OD2	5.03	122.83	118.30

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	1143	0	1145	17	0
1	B	1143	0	1145	17	0
1	C	1143	0	1145	18	0
1	D	1143	0	1145	18	0
1	E	1143	0	1145	24	0
1	F	1143	0	1145	13	0
2	G	109	0	98	1	0
2	H	109	0	98	1	0
2	I	104	0	93	0	0
2	J	109	0	98	1	0
2	K	109	0	98	2	0
2	L	109	0	98	4	0
3	A	2	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
4	A	3	0	0	0	0
4	B	4	0	0	0	0
4	D	1	0	0	0	0
4	E	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	F	5	0	0	0	0
4	G	1	0	0	0	0
4	I	1	0	0	0	0
4	K	1	0	0	0	0
4	L	1	0	0	1	0
All	All	7531	0	7453	94	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 6.

All (94) 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:D:250:ILE:H	1:E:148:GLN:HE22	1.27	0.82
1:A:242:ASN:H	1:C:235:ASN:HD21	1.35	0.73
1:D:235:ASN:HD21	1:E:242:ASN:H	1.40	0.70
1:C:161:GLY:O	1:C:162:SER:HB2	1.92	0.69
1:A:264:PRO:O	1:A:265:ARG:HD3	1.95	0.67
2:L:34:CYS:HA	4:L:106:HOH:O	1.96	0.66
1:E:200:LEU:HD22	1:E:276:VAL:HG21	1.80	0.63
1:E:193:PHE:CZ	1:E:283:LYS:HD2	2.34	0.62
1:E:216:LYS:HB3	1:E:218:HIS:O	2.00	0.62
1:C:208:MET:SD	1:C:265:ARG:HD2	2.41	0.61
1:A:219:VAL:HG21	1:A:225:SER:HB3	1.83	0.60
1:E:235:ASN:HD21	1:F:242:ASN:H	1.48	0.60
1:D:161:GLY:O	1:D:162:SER:HB2	2.03	0.59
1:C:201:TYR:OH	1:C:210:HIS:ND1	2.31	0.58
2:L:27:LEU:HD23	2:L:27:LEU:N	2.18	0.58
1:C:211:LEU:HD22	1:C:231:ARG:HG3	1.86	0.56
1:B:150:ILE:HD11	1:F:219:VAL:HG23	1.87	0.56
1:B:150:ILE:HD11	1:F:219:VAL:CG2	2.37	0.55
1:E:170:LEU:HD11	1:E:173:LYS:HB2	1.88	0.55
1:A:230:PHE:CE1	1:A:249:GLY:HA3	2.43	0.54
1:E:158:ILE:HB	1:E:165:PHE:HB2	1.89	0.54
2:L:27:LEU:H	2:L:27:LEU:HD23	1.74	0.52
1:D:235:ASN:ND2	1:E:242:ASN:H	2.06	0.52
1:B:220:PHE:O	1:B:223:GLU:HB2	2.10	0.51
1:D:230:PHE:CZ	1:D:249:GLY:HA3	2.45	0.51
1:B:219:VAL:HG21	1:B:225:SER:HA	1.93	0.51
1:E:168:TRP:H	1:E:183:ASN:HD21	1.58	0.50
1:A:161:GLY:O	1:A:162:SER:HB2	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:144:GLN:NE2	1:F:282:LEU:HD11	2.26	0.50
1:D:152:ASP:HB2	1:D:169:LEU:HB3	1.93	0.50
1:B:172:PHE:HA	1:F:218:HIS:CD2	2.47	0.50
1:C:264:PRO:O	1:C:265:ARG:HG2	2.12	0.50
1:C:192:TYR:CE1	1:C:252:LYS:HD2	2.47	0.50
1:B:181:LYS:NZ	1:B:258:GLU:OE2	2.45	0.49
1:A:240:LEU:N	1:A:241:PRO:HD3	2.27	0.49
1:A:234:GLN:HA	1:B:243:ASN:OD1	2.13	0.48
1:A:158:ILE:HB	1:A:165:PHE:HB2	1.95	0.48
1:E:213:GLN:OE1	1:E:260:GLN:NE2	2.46	0.48
1:E:167:PRO:HA	1:E:183:ASN:HD21	1.78	0.48
1:E:186:LEU:HD12	1:E:187:VAL:N	2.28	0.48
1:A:200:LEU:HD22	1:A:276:VAL:HG21	1.96	0.48
1:E:264:PRO:HG2	2:K:28:LEU:HB2	1.96	0.47
1:C:230:PHE:CZ	1:C:249:GLY:HA3	2.49	0.47
1:D:160:LYS:O	1:D:161:GLY:O	2.32	0.47
1:C:230:PHE:CE1	1:C:249:GLY:HA3	2.50	0.47
1:B:233:ILE:HD12	2:H:28:LEU:HG	1.96	0.47
1:D:192:TYR:CE2	1:E:174:ARG:HD2	2.50	0.47
1:F:167:PRO:HA	1:F:183:ASN:HD21	1.80	0.47
1:D:144:GLN:HG3	1:D:285:LEU:HD23	1.98	0.46
1:B:161:GLY:O	1:B:162:SER:CB	2.64	0.46
1:B:159:GLN:HE22	1:B:267:ASN:HA	1.81	0.45
1:B:285:LEU:HD12	1:C:285:LEU:CD1	2.46	0.45
1:A:242:ASN:H	1:C:235:ASN:ND2	2.09	0.45
1:C:152:ASP:HB2	1:C:169:LEU:HG	1.98	0.45
1:D:152:ASP:OD2	1:D:155:THR:CG2	2.64	0.45
1:A:213:GLN:O	1:A:259:LEU:HA	2.16	0.45
2:J:23:CYS:HA	2:J:33:VAL:O	2.17	0.45
1:C:147:LEU:HD21	1:C:149:LEU:HD21	1.99	0.44
1:A:155:THR:HG22	1:A:156:PRO:HD2	1.98	0.44
1:D:282:LEU:HD11	1:F:284:LEU:HD22	2.00	0.43
1:E:203:ASP:HA	1:E:269:GLN:HB2	1.99	0.43
1:D:242:ASN:HD22	1:D:242:ASN:HA	1.65	0.43
1:B:231:ARG:O	1:C:276:VAL:HG23	2.18	0.43
1:A:223:GLU:HG3	1:F:227:VAL:HG11	2.01	0.43
1:E:240:LEU:N	1:E:241:PRO:CD	2.81	0.43
1:F:157:THR:HA	1:F:167:PRO:HD3	2.01	0.43
1:D:240:LEU:N	1:D:241:PRO:HD3	2.33	0.43
1:E:149:LEU:HB3	1:E:168:TRP:HB3	2.00	0.42
1:D:213:GLN:HB2	1:D:260:GLN:HG3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:193:PHE:CZ	1:A:283:LYS:HD3	2.55	0.42
2:K:29:VAL:HG23	2:K:31:HIS:HB3	2.01	0.42
1:B:219:VAL:HG13	1:B:223:GLU:HB3	2.01	0.42
1:D:213:GLN:OE1	1:D:260:GLN:NE2	2.53	0.41
1:E:198:GLN:O	1:E:277:THR:HA	2.20	0.41
2:G:25:TRP:CZ3	2:G:31:HIS:HA	2.55	0.41
1:F:264:PRO:HG2	2:L:28:LEU:HB2	2.01	0.41
1:E:285:LEU:HA	1:E:285:LEU:HD22	1.98	0.41
1:D:202:THR:HG22	1:D:242:ASN:ND2	2.35	0.41
1:A:193:PHE:CE2	1:A:283:LYS:HD3	2.56	0.41
1:B:195:ILE:HD11	1:B:253:LEU:HD11	2.01	0.41
1:D:233:ILE:HG21	1:E:242:ASN:HB3	2.02	0.41
1:D:192:TYR:CZ	1:E:174:ARG:HD2	2.56	0.41
1:C:157:THR:HA	1:C:167:PRO:HD3	2.02	0.41
1:E:211:LEU:HD21	1:E:231:ARG:HD3	2.03	0.41
1:E:266:GLU:O	1:E:267:ASN:C	2.59	0.41
1:B:250:ILE:H	1:C:148:GLN:NE2	2.19	0.40
1:F:146:CYS:HB3	1:F:282:LEU:HA	2.03	0.40
1:A:240:LEU:N	1:A:241:PRO:CD	2.84	0.40
1:A:208:MET:SD	1:A:265:ARG:HG3	2.62	0.40
1:C:211:LEU:HB3	1:C:228:THR:HG23	2.03	0.40
1:F:148:GLN:HB2	1:F:172:PHE:CE2	2.56	0.40
1:F:211:LEU:HB3	1:F:228:THR:HG23	2.04	0.40
1:B:285:LEU:HD12	1:C:285:LEU:HD11	2.02	0.40
1:B:267:ASN:C	1:B:267:ASN:HD22	2.25	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	A	142/208 (68%)	134 (94%)	7 (5%)	1 (1%)	26 70

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	142/208 (68%)	132 (93%)	8 (6%)	2 (1%)	14	51
1	C	142/208 (68%)	132 (93%)	8 (6%)	2 (1%)	14	51
1	D	142/208 (68%)	133 (94%)	6 (4%)	3 (2%)	9	40
1	E	142/208 (68%)	132 (93%)	9 (6%)	1 (1%)	26	70
1	F	142/208 (68%)	134 (94%)	7 (5%)	1 (1%)	26	70
2	G	10/12 (83%)	10 (100%)	0	0	100	100
2	H	10/12 (83%)	9 (90%)	1 (10%)	0	100	100
2	I	10/12 (83%)	9 (90%)	1 (10%)	0	100	100
2	J	10/12 (83%)	10 (100%)	0	0	100	100
2	K	10/12 (83%)	10 (100%)	0	0	100	100
2	L	10/12 (83%)	10 (100%)	0	0	100	100
All	All	912/1320 (69%)	855 (94%)	47 (5%)	10 (1%)	17	58

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	162	SER
1	D	161	GLY
1	F	161	GLY
1	A	161	GLY
1	C	161	GLY
1	D	242	ASN
1	B	162	SER
1	E	267	ASN
1	B	161	GLY
1	D	162	SER

### 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	126/171 (74%)	111 (88%)	15 (12%)	6	26

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	126/171 (74%)	107 (85%)	19 (15%)	3	17
1	C	126/171 (74%)	112 (89%)	14 (11%)	8	29
1	D	126/171 (74%)	111 (88%)	15 (12%)	6	26
1	E	126/171 (74%)	110 (87%)	16 (13%)	5	23
1	F	126/171 (74%)	113 (90%)	13 (10%)	9	33
2	G	12/12 (100%)	12 (100%)	0	100	100
2	H	12/12 (100%)	12 (100%)	0	100	100
2	I	11/12 (92%)	11 (100%)	0	100	100
2	J	12/12 (100%)	10 (83%)	2 (17%)	3	13
2	K	12/12 (100%)	12 (100%)	0	100	100
2	L	12/12 (100%)	9 (75%)	3 (25%)	1	3
All	All	827/1098 (75%)	730 (88%)	97 (12%)	7	27

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	142	VAL
1	A	143	THR
1	A	153	SER
1	A	159	GLN
1	A	162	SER
1	A	169	LEU
1	A	171	SER
1	A	183	ASN
1	A	204	LYS
1	A	211	LEU
1	A	238	GLU
1	A	239	THR
1	A	255	GLU
1	A	276	VAL
1	A	283	LYS
1	B	146	CYS
1	B	159	GLN
1	B	180	GLU
1	B	181	LYS
1	B	204	LYS
1	B	215	LYS
1	B	218	HIS

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Mol	Chain	Res	Type
1	B	219	VAL
1	B	222	ASP
1	B	224	LEU
1	B	225	SER
1	B	227	VAL
1	B	232	CYS
1	B	239	THR
1	B	255	GLU
1	B	265	ARG
1	B	266	GLU
1	B	267	ASN
1	B	276	VAL
1	C	143	THR
1	C	160	LYS
1	C	162	SER
1	C	211	LEU
1	C	217	VAL
1	C	222	ASP
1	C	224	LEU
1	C	238	GLU
1	C	239	THR
1	C	247	SER
1	C	254	GLU
1	C	265	ARG
1	C	266	GLU
1	C	276	VAL
1	D	142	VAL
1	D	155	THR
1	D	159	GLN
1	D	160	LYS
1	D	162	SER
1	D	169	LEU
1	D	181	LYS
1	D	183	ASN
1	D	214	ARG
1	D	224	LEU
1	D	232	CYS
1	D	238	GLU
1	D	242	ASN
1	D	276	VAL
1	D	285	LEU
1	E	142	VAL

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Mol	Chain	Res	Type
1	E	154	GLU
1	E	160	LYS
1	E	173	LYS
1	E	180	GLU
1	E	181	LYS
1	E	183	ASN
1	E	188	LYS
1	E	214	ARG
1	E	216	LYS
1	E	224	LEU
1	E	247	SER
1	E	260	GLN
1	E	265	ARG
1	E	276	VAL
1	E	285	LEU
1	F	142	VAL
1	F	143	THR
1	F	146	CYS
1	F	169	LEU
1	F	180	GLU
1	F	225	SER
1	F	227	VAL
1	F	232	CYS
1	F	239	THR
1	F	254	GLU
1	F	265	ARG
1	F	276	VAL
1	F	282	LEU
2	J	23	CYS
2	J	34	CYS
2	L	24	HIS
2	L	27	LEU
2	L	30	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	144	GLN
1	A	159	GLN
1	A	210	HIS
1	A	218	HIS
1	A	242	ASN

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Mol	Chain	Res	Type
1	A	267	ASN
1	B	144	GLN
1	B	267	ASN
1	C	148	GLN
1	C	235	ASN
1	C	260	GLN
1	C	267	ASN
1	D	159	GLN
1	D	234	GLN
1	D	235	ASN
1	D	242	ASN
1	D	267	ASN
1	E	148	GLN
1	E	183	ASN
1	E	213	GLN
1	E	235	ASN
1	E	260	GLN
1	F	210	HIS
1	F	218	HIS
2	G	31	HIS

### 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 carbohydrates in this entry.

### 5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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, 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	A	144/208 (69%)	-0.49	0 <span>100</span> <span>100</span>	19, 29, 44, 51	0
1	B	144/208 (69%)	-0.43	1 (0%) <span>89</span> <span>70</span>	20, 31, 49, 61	0
1	C	144/208 (69%)	-0.47	0 <span>100</span> <span>100</span>	20, 31, 44, 51	0
1	D	144/208 (69%)	-0.39	2 (1%) <span>78</span> <span>51</span>	20, 31, 59, 79	0
1	E	144/208 (69%)	-0.47	0 <span>100</span> <span>100</span>	20, 31, 44, 51	0
1	F	144/208 (69%)	-0.49	0 <span>100</span> <span>100</span>	17, 29, 44, 51	0
2	G	12/12 (100%)	0.32	1 (8%) <span>14</span> <span>5</span>	38, 54, 63, 64	0
2	H	12/12 (100%)	0.70	2 (16%) <span>2</span> <span>1</span>	38, 54, 62, 64	0
2	I	12/12 (100%)	0.66	3 (25%) <span>1</span> <span>1</span>	39, 54, 63, 64	0
2	J	12/12 (100%)	0.98	4 (33%) <span>0</span> <span>0</span>	38, 54, 63, 64	0
2	K	12/12 (100%)	0.68	1 (8%) <span>14</span> <span>5</span>	38, 54, 63, 65	0
2	L	12/12 (100%)	0.60	1 (8%) <span>14</span> <span>5</span>	38, 54, 63, 64	0
All	All	936/1320 (70%)	-0.37	15 (1%) <span>74</span> <span>47</span>	17, 31, 56, 79	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	222	ASP	3.6
2	J	23	CYS	3.1
2	K	34	CYS	3.1
2	H	23	CYS	2.6
2	I	24	HIS	2.6
2	J	32	TRP	2.5
1	B	142	VAL	2.5
2	J	34	CYS	2.5
2	I	34	CYS	2.5
2	I	23	CYS	2.4
2	G	23	CYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	223	GLU	2.2
2	H	24	HIS	2.1
2	L	23	CYS	2.1
2	J	24	HIS	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	MG	A	702	1/1	0.76	0.79	34.96	58,58,58,58	0
3	MG	A	701	1/1	0.84	0.51	20.38	56,56,56,56	0
3	MG	E	801	1/1	0.84	0.62	16.56	75,75,75,75	0
3	MG	D	802	1/1	0.89	0.46	11.05	48,48,48,48	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.