



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:34 PM GMT

PDB ID : 3ZX6  
Title : Structure of Hamp(AF1503)-Tsr fusion - Hamp (A291V) mutant  
Authors : Zeth, K.; Ferris, H.U.; Lupas, A.L.  
Deposited on : 2011-08-08  
Resolution : 2.65 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

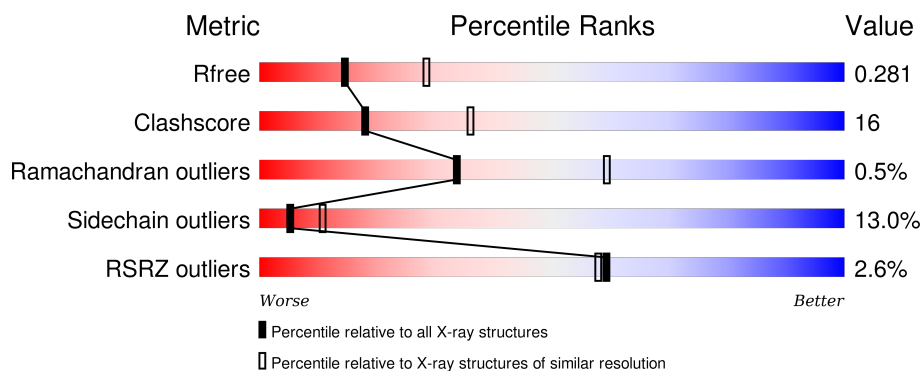
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

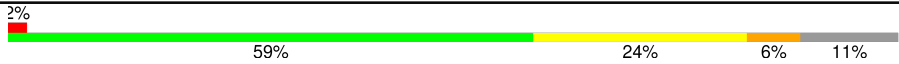

The reported resolution of this entry is 2.65 Å.

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	3152 (2.70-2.62)
Clashscore	102246	3524 (2.70-2.62)
Ramachandran outliers	100387	3469 (2.70-2.62)
Sidechain outliers	100360	3469 (2.70-2.62)
RSRZ outliers	91569	3161 (2.70-2.62)

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	341	 2% 59% 24% 6% 11%
1	B	341	 2% 62% 23% • 11%

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
2	CL	A	1306	-	-	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4514 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 HAMP, METHYL-ACCEPTING CHEMOTAXIS PROTEIN I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	303	Total	C	N	O	S	0	0	0
			2228	1343	410	469	6			
1	B	303	Total	C	N	O	S	0	0	0
			2208	1329	406	467	6			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	EXPRESSION TAG	UNP O28769
A	15	VAL	ALA	ENGINEERED MUTATION	UNP O28769
A	325	THR	ALA	CONFLICT	UNP P02942
A	335	GLY	GLU	CONFLICT	UNP P02942
B	1	MET	-	EXPRESSION TAG	UNP O28769
B	15	VAL	ALA	ENGINEERED MUTATION	UNP O28769
B	325	THR	ALA	CONFLICT	UNP P02942
B	335	GLY	GLU	CONFLICT	UNP P02942

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Cl	0	0
			1	1		
2	A	1	Total	Cl	0	0
			1	1		

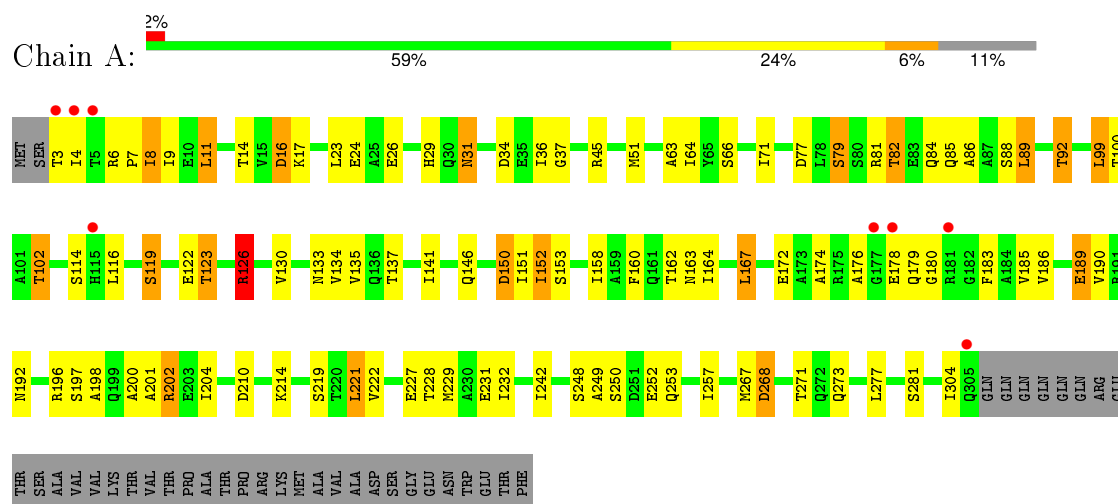
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	42	Total	O	0	0
			42	42		
3	B	34	Total	O	0	0
			34	34		

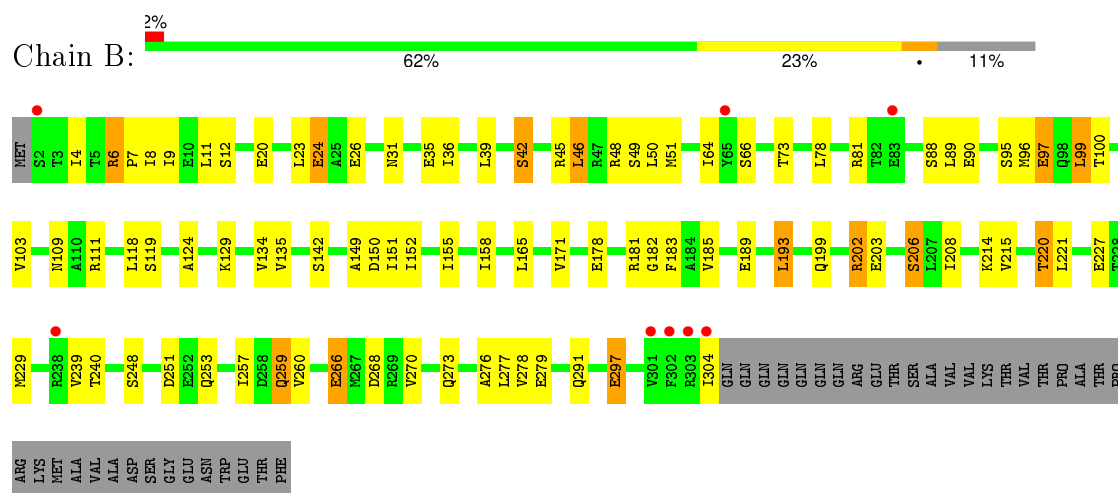
### 3 Residue-property plots

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.

#### • Molecule 1: HAMP, METHYL-ACCEPTING CHEMOTAXIS PROTEIN I



#### • Molecule 1: HAMP, METHYL-ACCEPTING CHEMOTAXIS PROTEIN I



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	112.22Å 80.02Å 139.17Å 90.00° 110.78° 90.00°	Depositor
Resolution (Å)	29.48 – 2.65 28.79 – 2.65	Depositor EDS
% Data completeness (in resolution range)	100.0 (29.48-2.65) 92.5 (28.79-2.65)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.58 (at 2.64Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.229 , 0.283 0.229 , 0.281	Depositor DCC
$R_{free}$ test set	1553 reflections (5.25%)	DCC
Wilson B-factor (Å <sup>2</sup> )	52.7	Xtriage
Anisotropy	0.277	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 64.2	EDS
Estimated twinning fraction	0.000 for h,-k,-h-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 31128 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4514	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.30% 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: CL

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	0.93	0/2235	0.93	5/3022 (0.2%)
1	B	1.03	5/2214 (0.2%)	0.99	2/2996 (0.1%)
All	All	0.98	5/4449 (0.1%)	0.96	7/6018 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	24	GLU	CG-CD	7.55	1.63	1.51
1	B	203	GLU	CG-CD	6.88	1.62	1.51
1	B	189	GLU	CG-CD	6.17	1.61	1.51
1	B	26	GLU	CG-CD	5.50	1.60	1.51
1	B	297	GLU	CG-CD	5.05	1.59	1.51

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	150	ASP	CB-CG-OD1	6.43	124.09	118.30
1	B	150	ASP	CB-CG-OD1	6.34	124.01	118.30
1	A	126	ARG	NE-CZ-NH1	5.75	123.17	120.30
1	A	45	ARG	NE-CZ-NH1	5.64	123.12	120.30
1	A	45	ARG	NE-CZ-NH2	-5.38	117.61	120.30
1	A	152	ILE	CG1-CB-CG2	-5.30	99.74	111.40
1	B	150	ASP	CB-CG-OD2	-5.02	113.78	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	2228	0	2232	99	0
1	B	2208	0	2196	87	0
2	A	1	0	0	1	0
2	B	1	0	0	0	0
3	A	42	0	0	4	0
3	B	34	0	0	8	0
All	All	4514	0	4428	139	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 16.

All (139) 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:63:ALA:CB	1:B:291:GLN:HE21	1.70	1.05
1:A:102:THR:HG22	1:B:253:GLN:HE21	1.22	1.04
1:A:102:THR:CG2	1:B:253:GLN:HE21	1.70	1.03
1:A:228:THR:HG21	1:B:229:MET:HE3	1.43	0.98
1:B:42:SER:HA	3:B:2006:HOH:O	1.62	0.98
1:A:8:ILE:HD13	1:B:36:ILE:HD13	1.49	0.93
1:B:97:GLU:HA	1:B:97:GLU:OE1	1.76	0.86
1:A:102:THR:HG21	1:B:253:GLN:CG	2.06	0.85
1:A:100:THR:HG23	1:A:257:ILE:HG21	1.59	0.84
1:A:102:THR:HG21	1:B:253:GLN:HG3	1.59	0.83
1:A:99:LEU:HB3	1:A:257:ILE:HD11	1.63	0.80
1:A:180:GLY:HA2	1:A:183:PHE:HD2	1.46	0.79
1:A:210:ASP:HB2	3:A:2036:HOH:O	1.82	0.79
1:A:228:THR:HG21	1:B:229:MET:CE	2.13	0.79
1:A:84:GLN:NE2	1:B:266:GLU:HG2	1.98	0.78
1:B:178:GLU:OE1	1:B:181:ARG:NH2	2.18	0.77
1:A:102:THR:CG2	1:B:253:GLN:NE2	2.47	0.76
1:A:119:SER:O	1:A:123:THR:HG23	1.86	0.74
1:A:162:THR:HG22	1:B:193:LEU:HD23	1.69	0.73
1:B:90:GLU:HA	1:B:90:GLU:OE1	1.87	0.73
1:B:97:GLU:O	1:B:100:THR:HG22	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:ALA:HB1	1:A:179:GLN:HB2	1.72	0.72
1:A:133:ASN:O	1:A:137:THR:HG23	1.89	0.71
1:A:81:ARG:HG2	1:B:270:VAL:HG22	1.72	0.71
1:A:84:GLN:HE21	1:B:266:GLU:CG	2.03	0.71
1:A:63:ALA:CB	1:B:291:GLN:NE2	2.51	0.69
1:B:9:ILE:O	1:B:12:SER:HB3	1.95	0.67
1:A:84:GLN:HE21	1:B:266:GLU:HG2	1.59	0.66
1:A:176:ALA:CB	1:A:179:GLN:HB2	2.26	0.66
1:A:146:GLN:OE1	1:A:146:GLN:HA	1.96	0.65
1:B:78:LEU:HD23	1:B:278:VAL:HG22	1.78	0.65
1:A:267:MET:SD	1:B:89:LEU:HD12	2.36	0.64
1:A:16:ASP:OD1	1:B:42:SER:OG	2.16	0.64
1:B:45:ARG:NH1	3:B:2006:HOH:O	2.31	0.63
1:A:8:ILE:HD11	1:B:35:GLU:C	2.19	0.63
1:A:227:GLU:HA	1:A:227:GLU:OE1	1.97	0.63
1:A:252:GLU:HA	1:A:252:GLU:OE1	1.99	0.62
1:B:103:VAL:HG22	1:B:253:GLN:OE1	1.98	0.62
1:A:8:ILE:CD1	1:B:36:ILE:HD13	2.27	0.62
1:B:20:GLU:HG2	1:B:304:ILE:C	2.20	0.61
1:B:259:GLN:NE2	1:B:259:GLN:HA	2.16	0.61
1:A:63:ALA:HB3	1:B:291:GLN:HE21	1.59	0.60
3:A:2022:HOH:O	1:B:214:LYS:NZ	2.34	0.60
1:A:152:ILE:HD11	1:A:201:ALA:HA	1.81	0.60
1:B:11:LEU:HD13	1:B:39:LEU:HD23	1.84	0.59
1:A:8:ILE:HD13	1:B:36:ILE:CD1	2.29	0.59
1:B:45:ARG:CZ	3:B:2006:HOH:O	2.50	0.58
1:B:48:ARG:HD2	1:B:51:MET:CE	2.33	0.58
1:A:267:MET:SD	1:B:89:LEU:CD1	2.92	0.58
1:A:63:ALA:HB1	1:B:291:GLN:HE21	1.63	0.58
1:B:42:SER:CA	3:B:2006:HOH:O	2.34	0.57
1:A:197:SER:HB2	1:B:158:ILE:HG21	1.86	0.57
1:A:64:ILE:HD13	1:B:64:ILE:HD12	1.87	0.56
1:A:116:LEU:HD22	1:B:239:VAL:HG23	1.87	0.56
1:A:242:ILE:HG23	1:B:109:ASN:HB3	1.88	0.55
1:A:92:THR:HB	1:B:260:VAL:HG22	1.89	0.55
1:B:149:ALA:O	1:B:152:ILE:HG22	2.07	0.54
1:B:23:LEU:HD23	1:B:51:MET:HG2	1.88	0.54
1:A:89:LEU:CD2	1:A:268:ASP:HB3	2.38	0.54
1:A:158:ILE:HG23	1:B:193:LEU:HG	1.89	0.54
1:A:102:THR:CG2	1:B:253:GLN:CG	2.85	0.53
1:A:88:SER:O	1:A:92:THR:HG23	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:ARG:HG2	1:A:7:PRO:HD3	1.89	0.53
1:A:229:MET:HE1	1:B:229:MET:HG2	1.90	0.53
1:A:186:VAL:O	1:A:190:VAL:HG23	2.08	0.52
1:A:100:THR:CG2	1:A:257:ILE:HG21	2.37	0.52
1:B:48:ARG:HD2	1:B:51:MET:HE1	1.90	0.52
1:A:192:ASN:O	1:A:196:ARG:HG3	2.10	0.52
1:A:232:ILE:HD11	1:B:124:ALA:HB2	1.90	0.52
1:B:182:GLY:O	1:B:185:VAL:HG22	2.10	0.52
1:B:31:ASN:HB2	3:B:2003:HOH:O	2.10	0.52
1:A:228:THR:CG2	1:B:229:MET:HE3	2.31	0.51
1:A:135:VAL:HG23	1:A:222:VAL:HG12	1.92	0.51
1:B:11:LEU:CD1	1:B:39:LEU:HD23	2.40	0.51
1:B:142:SER:HB2	1:B:215:VAL:HG11	1.93	0.51
1:A:85:GLN:NE2	1:A:271:THR:OG1	2.35	0.50
1:A:198:ALA:O	1:A:202:ARG:HG2	2.10	0.50
1:A:84:GLN:HE21	1:B:266:GLU:HG3	1.74	0.50
1:A:102:THR:HG21	1:B:253:GLN:HG2	1.92	0.49
1:A:164:ILE:O	1:A:167:LEU:HB2	2.12	0.49
1:B:97:GLU:OE1	1:B:97:GLU:CA	2.56	0.49
1:A:130:VAL:O	1:A:134:VAL:HG23	2.12	0.49
1:B:206:SER:HB2	3:B:2026:HOH:O	2.13	0.49
1:A:79:SER:O	1:A:82:THR:N	2.45	0.49
1:A:135:VAL:CG2	1:A:222:VAL:HG12	2.43	0.49
1:A:277:LEU:O	1:A:277:LEU:HD23	2.12	0.49
1:A:126:ARG:HH11	1:A:126:ARG:HB2	1.77	0.49
1:A:160:PHE:O	1:A:163:ASN:HB3	2.13	0.48
1:A:64:ILE:HD13	1:B:64:ILE:CD1	2.43	0.48
1:A:250:SER:OG	2:A:1306:CL:CL	2.65	0.48
1:A:77:ASP:HB3	1:A:81:ARG:HH21	1.77	0.48
1:B:199:GLN:OE1	1:B:202:ARG:NH2	2.47	0.48
1:A:86:ALA:HB2	1:A:271:THR:HG21	1.95	0.48
1:A:249:ALA:O	1:A:253:GLN:HG3	2.13	0.48
1:A:273:GLN:OE1	1:B:81:ARG:HD3	2.13	0.48
1:B:273:GLN:O	1:B:276:ALA:HB3	2.14	0.48
1:A:152:ILE:HG13	1:A:152:ILE:O	2.13	0.48
1:A:137:THR:O	1:A:141:ILE:HG13	2.14	0.47
1:A:11:LEU:HD13	1:A:29:HIS:HD2	1.78	0.47
1:A:31:ASN:HA	3:A:2013:HOH:O	2.15	0.47
1:B:95:SER:O	1:B:99:LEU:HB2	2.15	0.46
1:A:277:LEU:C	1:A:277:LEU:HD23	2.36	0.46
1:A:158:ILE:O	1:A:162:THR:HG23	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:172:GLU:OE2	1:B:182:GLY:HA3	2.16	0.46
1:A:172:GLU:HG2	1:B:183:PHE:N	2.31	0.46
1:B:134:VAL:O	1:B:135:VAL:C	2.53	0.45
1:B:46:LEU:HD22	1:B:50:LEU:CD1	2.47	0.44
1:B:6:ARG:N	1:B:7:PRO:CD	2.81	0.44
1:A:17:LYS:HB3	1:A:17:LYS:HE2	1.54	0.44
1:A:99:LEU:HB3	1:A:257:ILE:CD1	2.42	0.44
1:B:24:GLU:OE1	1:B:24:GLU:N	2.50	0.44
1:A:26:GLU:HA	3:A:2009:HOH:O	2.17	0.44
1:B:96:MET:HE3	1:B:257:ILE:HG13	1.99	0.44
1:A:167:LEU:HD12	1:A:167:LEU:HA	1.79	0.44
1:A:8:ILE:HD11	1:B:35:GLU:CB	2.47	0.44
1:B:291:GLN:OE1	1:B:291:GLN:HA	2.18	0.44
1:B:48:ARG:HD2	1:B:51:MET:HE2	2.00	0.43
1:A:180:GLY:HA2	1:A:183:PHE:CD2	2.37	0.43
1:A:227:GLU:OE1	1:A:227:GLU:CA	2.67	0.43
1:B:220:THR:HG22	1:B:221:LEU:HD23	2.01	0.43
1:B:8:ILE:HD13	1:B:8:ILE:HA	1.87	0.43
1:B:73:THR:HG22	3:B:2012:HOH:O	2.19	0.43
1:A:8:ILE:HG22	1:A:36:ILE:HD12	2.01	0.42
1:A:8:ILE:HD11	1:B:35:GLU:O	2.19	0.42
1:A:34:ASP:O	1:A:37:GLY:N	2.47	0.42
1:B:208:ILE:HD13	1:B:208:ILE:HA	1.93	0.42
1:A:189:GLU:HB3	1:B:165:LEU:HD21	2.01	0.42
1:B:42:SER:N	3:B:2006:HOH:O	2.48	0.42
1:A:11:LEU:O	1:A:14:THR:HB	2.20	0.42
1:A:102:THR:HG22	1:B:253:GLN:NE2	2.07	0.41
1:A:273:GLN:OE1	1:B:81:ARG:CD	2.68	0.41
1:A:151:ILE:O	1:A:152:ILE:C	2.56	0.41
1:A:77:ASP:CB	1:A:81:ARG:HH21	2.33	0.41
1:A:221:LEU:HB3	1:B:134:VAL:CG2	2.51	0.41
1:A:51:MET:HG2	1:A:304:ILE:HG23	2.03	0.41
1:A:221:LEU:HD12	1:A:221:LEU:HA	1.94	0.41
1:A:150:ASP:O	1:A:153:SER:OG	2.33	0.41
1:A:200:ALA:CB	1:B:155:ILE:HD11	2.51	0.41
1:A:204:ILE:HD11	1:B:151:ILE:HB	2.03	0.41

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	301/341 (88%)	287 (95%)	12 (4%)	2 (1%)	26	51
1	B	301/341 (88%)	281 (93%)	19 (6%)	1 (0%)	46	72
All	All	602/682 (88%)	568 (94%)	31 (5%)	3 (0%)	34	59

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	79	SER
1	A	174	ALA
1	B	4	ILE

### 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	233/266 (88%)	200 (86%)	33 (14%)	4	8
1	B	228/266 (86%)	201 (88%)	27 (12%)	6	13
All	All	461/532 (87%)	401 (87%)	60 (13%)	5	11

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	THR
1	A	4	ILE
1	A	8	ILE

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Mol	Chain	Res	Type
1	A	9	ILE
1	A	11	LEU
1	A	16	ASP
1	A	23	LEU
1	A	24	GLU
1	A	31	ASN
1	A	66	SER
1	A	71	ILE
1	A	82	THR
1	A	89	LEU
1	A	92	THR
1	A	99	LEU
1	A	102	THR
1	A	114	SER
1	A	119	SER
1	A	122	GLU
1	A	123	THR
1	A	126	ARG
1	A	167	LEU
1	A	178	GLU
1	A	185	VAL
1	A	189	GLU
1	A	202	ARG
1	A	214	LYS
1	A	219	SER
1	A	221	LEU
1	A	231	GLU
1	A	248	SER
1	A	268	ASP
1	A	281	SER
1	B	6	ARG
1	B	42	SER
1	B	46	LEU
1	B	49	SER
1	B	66	SER
1	B	88	SER
1	B	97	GLU
1	B	99	LEU
1	B	111	ARG
1	B	118	LEU
1	B	119	SER
1	B	129	LYS

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Mol	Chain	Res	Type
1	B	171	VAL
1	B	193	LEU
1	B	202	ARG
1	B	206	SER
1	B	220	THR
1	B	227	GLU
1	B	240	THR
1	B	248	SER
1	B	251	ASP
1	B	259	GLN
1	B	266	GLU
1	B	268	ASP
1	B	277	LEU
1	B	279	GLU
1	B	297	GLU

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

Mol	Chain	Res	Type
1	A	13	ASN
1	A	84	GLN
1	A	192	ASN
1	B	30	GLN
1	B	253	GLN
1	B	259	GLN
1	B	273	GLN
1	B	291	GLN

### 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 2 ligands modelled in this entry, 2 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	303/341 (88%)	-0.10	8 (2%) 59 58	28, 61, 93, 111	0
1	B	303/341 (88%)	-0.13	8 (2%) 59 58	26, 60, 89, 124	0
All	All	606/682 (88%)	-0.11	16 (2%) 59 58	26, 61, 91, 124	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	304	ILE	5.3
1	A	3	THR	4.9
1	A	305	GLN	4.0
1	B	301	VAL	3.7
1	A	5	THR	3.4
1	B	302	PHE	2.8
1	A	115	HIS	2.7
1	A	4	ILE	2.6
1	A	177	GLY	2.5
1	A	178	GLU	2.4
1	B	238	ARG	2.3
1	B	83	GLU	2.2
1	A	181	ARG	2.1
1	B	303	ARG	2.1
1	B	65	TYR	2.1
1	B	2	SER	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
2	CL	A	1306	1/1	0.94	0.44	13.18	111,111,111,111	0
2	CL	B	1305	1/1	0.91	0.06	-	70,70,70,70	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.