



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:57 AM GMT

PDB ID : 2VDU  
Title : STRUCTURE OF TRM8-TRM82, THE YEAST TRNA M7G METHYLATION COMPLEX  
Authors : Leulliot, N.; Chaillet, M.; Durand, D.; Ulryck, N.; Blondeau, K.; Van Tilbeurgh, H.  
Deposited on : 2007-10-11  
Resolution : 2.40 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

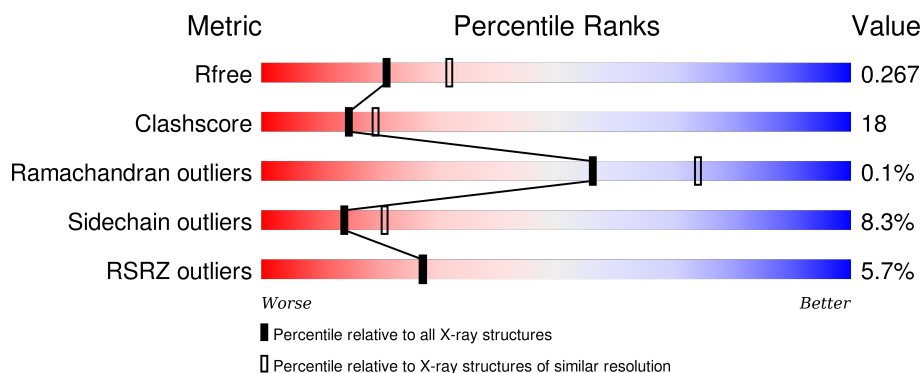
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	B	450	<div> <div>10%</div> <div> <div></div> <div>56%</div> <div>24%</div> <div>•</div> <div>17%</div> </div> </div>
1	D	450	<div> <div>3%</div> <div> <div></div> <div>58%</div> <div>22%</div> <div>•</div> <div>16%</div> </div> </div>
2	E	254	<div> <div>%</div> <div> <div></div> <div>50%</div> <div>25%</div> <div>•</div> <div>21%</div> </div> </div>
2	F	254	<div> <div>2%</div> <div> <div></div> <div>51%</div> <div>23%</div> <div>5%</div> <div>21%</div> </div> </div>

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9571 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.

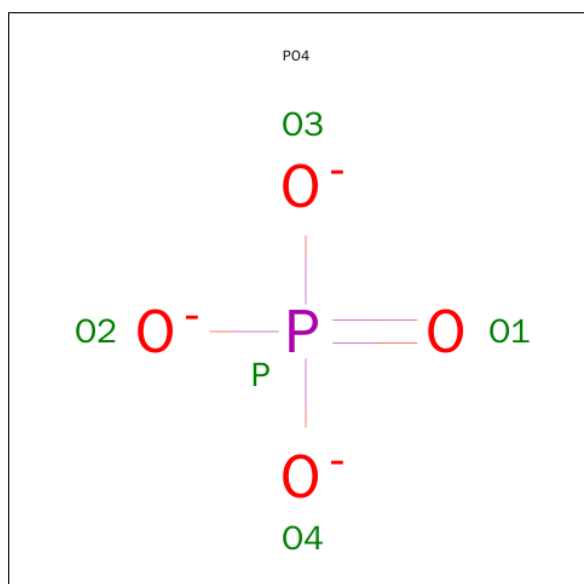
- Molecule 1 is a protein called TRNA (GUANINE-N(7)-)-METHYLTRANSFERASE-ASSOCIATED WD REPEAT PROTEIN TRM82.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	374	Total	C	N	O	S	0	0	0
			3013	1939	493	572	9			
1	D	376	Total	C	N	O	S	0	0	0
			3027	1949	495	574	9			

- Molecule 2 is a protein called TRNA (GUANINE-N(7)-)-METHYLTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	201	Total	C	N	O	S	0	0	0
			1631	1052	268	299	12			
2	F	201	Total	C	N	O	S	0	0	0
			1631	1052	268	299	12			

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).

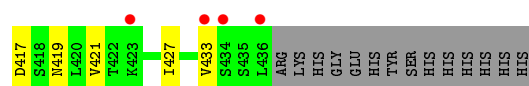


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	E	1	Total	O	P	0	0
			5	4	1		
3	F	1	Total	O	P	0	0
			5	4	1		

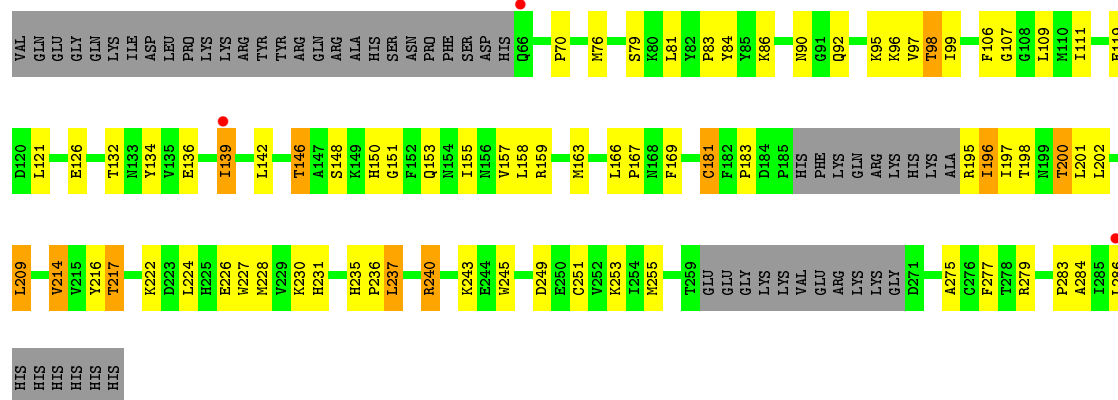
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	61	Total	O	0	0
			61	61		
4	D	102	Total	O	0	0
			102	102		
4	E	45	Total	O	0	0
			45	45		
4	F	51	Total	O	0	0
			51	51		

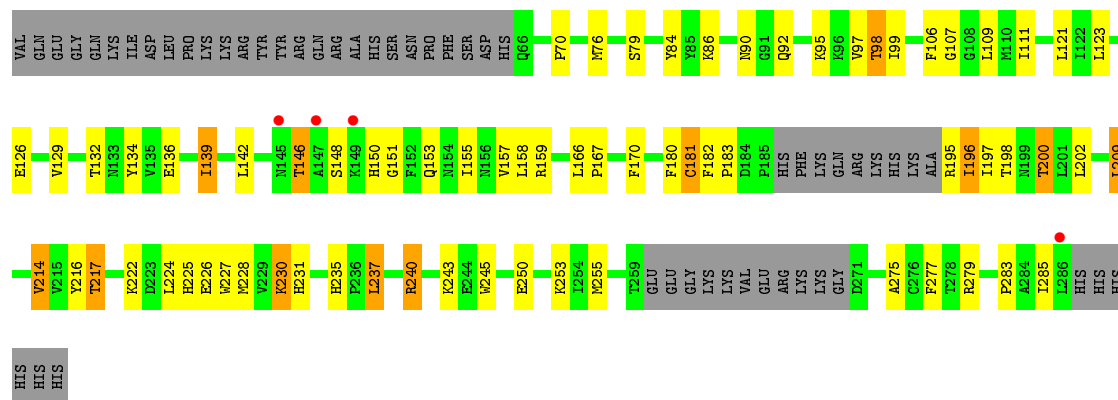




• Molecule 2: TRNA (GUANINE-N(7)-)-METHYLTRANSFERASE



• Molecule 2: TRNA (GUANINE-N(7)-)-METHYLTRANSFERASE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	53.38Å 107.68Å 127.75Å 90.00° 91.53° 90.00°	Depositor
Resolution (Å)	20.00 – 2.40 63.85 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.1 (20.00-2.40) 99.9 (63.85-2.40)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.82 (at 2.40Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.228 , 0.276 0.219 , 0.267	Depositor DCC
$R_{free}$ test set	2771 reflections (5.17%)	DCC
Wilson B-factor (Å <sup>2</sup> )	48.6	Xtriage
Anisotropy	0.099	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 49.9	EDS
Estimated twinning fraction	0.033 for h,-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 56477 reflections (0.002%)	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	9571	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 3.93% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PO4

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	B	0.38	0/3079	0.57	2/4161 (0.0%)
1	D	0.40	0/3095	0.58	1/4185 (0.0%)
2	E	0.40	0/1669	0.56	0/2256
2	F	0.41	0/1669	0.56	0/2256
All	All	0.40	0/9512	0.57	3/12858 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	192	LEU	CA-CB-CG	5.96	129.00	115.30
1	B	192	LEU	CA-CB-CG	5.81	128.66	115.30
1	B	294	ALA	N-CA-C	-5.02	97.44	111.00

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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	B	3013	0	2985	101	0
1	D	3027	0	2999	100	0
2	E	1631	0	1614	66	0
2	F	1631	0	1614	66	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	E	5	0	0	0	0
3	F	5	0	0	1	0
4	B	61	0	0	8	0
4	D	102	0	0	7	0
4	E	45	0	0	4	0
4	F	51	0	0	7	0
All	All	9571	0	9212	328	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 18.

All (328) 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:23:ILE:HD11	1:D:28:LEU:HG	1.30	1.13
1:B:23:ILE:HD11	1:B:28:LEU:HG	1.34	1.02
2:F:79:SER:HB2	2:F:86:LYS:HE2	1.53	0.91
2:E:79:SER:HB2	2:E:86:LYS:HE2	1.52	0.91
2:F:198:THR:HG22	2:F:200:THR:H	1.37	0.90
1:B:289:ASN:H	1:B:292:HIS:HD2	1.20	0.88
2:E:198:THR:HG22	2:E:200:THR:H	1.35	0.88
1:D:289:ASN:H	1:D:292:HIS:HD2	1.21	0.87
1:D:133:ASP:HB2	1:D:141:LYS:HE3	1.58	0.86
1:B:158:ALA:HA	1:B:203:LEU:HD13	1.56	0.85
1:D:158:ALA:HA	1:D:203:LEU:HD13	1.56	0.85
1:D:271:THR:HG22	1:D:273:LYS:H	1.41	0.85
1:B:133:ASP:HB2	1:B:141:LYS:HE3	1.58	0.85
1:D:159:GLU:HG3	4:D:2018:HOH:O	1.77	0.85
1:B:271:THR:HG22	1:B:273:LYS:H	1.43	0.81
1:B:199:THR:HG21	1:B:244:VAL:O	1.79	0.80
2:E:183:PRO:HG2	2:E:224:LEU:HD21	1.63	0.79
1:D:199:THR:HG21	1:D:244:VAL:O	1.82	0.79
2:E:148:SER:HB2	2:E:150:HIS:CE1	2.19	0.78
2:F:148:SER:HB2	2:F:150:HIS:CE1	2.18	0.78
1:B:315:LYS:HD2	1:B:316:SER:O	1.84	0.78
2:F:183:PRO:HG2	2:F:224:LEU:HD21	1.66	0.76
1:D:210:HIS:HE1	4:D:2029:HOH:O	1.68	0.76
1:D:151:ARG:HH12	1:D:153:ASN:HD22	1.34	0.75
2:F:90:ASN:HB3	2:F:92:GLN:HG2	1.67	0.75
1:D:315:LYS:HD2	1:D:316:SER:O	1.87	0.75
2:E:90:ASN:HB3	2:E:92:GLN:HG2	1.68	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:7:LEU:O	1:B:370:THR:HG21	1.87	0.73
1:B:151:ARG:HH12	1:B:153:ASN:HD22	1.34	0.73
1:B:271:THR:HG22	1:B:273:LYS:N	2.05	0.72
1:D:7:LEU:O	1:D:370:THR:HG21	1.90	0.72
1:D:229:GLN:H	1:D:229:GLN:NE2	1.87	0.72
1:D:271:THR:HG22	1:D:273:LYS:N	2.05	0.71
1:D:240:HIS:HE1	1:D:257:SER:OG	1.72	0.71
1:B:289:ASN:H	1:B:292:HIS:CD2	2.07	0.70
1:B:8:GLN:N	1:B:8:GLN:HE21	1.91	0.69
4:D:2025:HOH:O	2:F:200:THR:HG23	1.91	0.69
1:B:224:ILE:HD12	1:B:235:LYS:HD3	1.73	0.69
1:B:229:GLN:H	1:B:229:GLN:NE2	1.90	0.68
2:F:90:ASN:HB3	2:F:92:GLN:CG	2.23	0.68
1:D:40:PHE:CD1	1:D:413:SER:HB2	2.29	0.68
2:E:146:THR:HG21	2:E:151:GLY:HA3	1.76	0.68
2:F:95:LYS:HG2	2:F:121:LEU:HG	1.75	0.68
1:B:40:PHE:CD1	1:B:413:SER:HB2	2.28	0.68
2:E:90:ASN:HB3	2:E:92:GLN:CG	2.24	0.67
1:B:240:HIS:HE1	1:B:257:SER:OG	1.76	0.67
2:E:95:LYS:HG2	2:E:121:LEU:HG	1.77	0.67
1:D:289:ASN:H	1:D:292:HIS:CD2	2.08	0.67
1:D:34:SER:CB	1:D:35:PRO:HA	2.25	0.66
2:E:109:LEU:HD22	2:E:181:CYS:SG	2.36	0.66
2:E:217:THR:HB	2:E:275:ALA:O	1.95	0.66
1:B:201:VAL:HG13	1:B:215:THR:HG22	1.78	0.66
1:D:224:ILE:HD12	1:D:235:LYS:HD3	1.78	0.66
2:F:146:THR:HG21	2:F:151:GLY:HA3	1.76	0.65
1:B:321:PHE:HE1	1:B:348:LYS:HG3	1.60	0.65
2:F:109:LEU:HD22	2:F:181:CYS:SG	2.37	0.65
2:F:240:ARG:HD3	4:F:2043:HOH:O	1.98	0.64
1:B:321:PHE:CE1	1:B:348:LYS:HG3	2.33	0.63
2:E:181:CYS:HB2	4:E:2043:HOH:O	1.98	0.63
1:D:34:SER:HB3	1:D:35:PRO:HA	1.81	0.62
1:D:271:THR:HG21	4:D:2057:HOH:O	1.98	0.62
2:E:136:GLU:O	2:E:139:ILE:HD13	2.00	0.62
1:D:201:VAL:HG13	1:D:215:THR:HG22	1.82	0.62
1:D:226:HIS:HD2	4:D:2030:HOH:O	1.83	0.62
1:D:359:SER:OG	1:D:370:THR:HB	2.01	0.61
2:F:227:TRP:CE2	2:F:231:HIS:HE1	2.18	0.61
2:F:70:PRO:HG3	2:F:76:MET:SD	2.41	0.61
2:F:214:VAL:HG22	2:F:216:TYR:CE1	2.34	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:230:LYS:HE2	4:F:2036:HOH:O	2.00	0.61
2:E:79:SER:CB	2:E:86:LYS:HE2	2.30	0.61
2:E:84:TYR:CB	2:E:153:GLN:HG3	2.31	0.61
2:F:98:THR:HB	4:F:2010:HOH:O	1.99	0.61
2:F:136:GLU:O	2:F:139:ILE:HD13	2.01	0.61
2:F:227:TRP:CE2	2:F:231:HIS:CE1	2.88	0.61
1:D:8:GLN:HE21	1:D:8:GLN:N	1.98	0.60
1:D:295:PRO:HB3	1:D:296:PRO:HD2	1.83	0.60
2:E:227:TRP:CE2	2:E:231:HIS:CE1	2.89	0.60
2:E:227:TRP:CE2	2:E:231:HIS:HE1	2.19	0.60
1:B:5:HIS:HB3	1:B:23:ILE:HG22	1.83	0.60
2:F:230:LYS:HG2	4:F:2036:HOH:O	2.01	0.60
2:E:196:ILE:HD12	2:E:197:ILE:HD13	1.83	0.59
2:E:70:PRO:HG3	2:E:76:MET:SD	2.41	0.59
1:D:23:ILE:HD11	1:D:28:LEU:CG	2.20	0.59
1:D:23:ILE:CD1	1:D:28:LEU:HG	2.21	0.59
1:D:295:PRO:O	1:D:296:PRO:C	2.41	0.59
2:F:217:THR:HB	2:F:275:ALA:O	2.01	0.59
2:F:84:TYR:CB	2:F:153:GLN:HG3	2.32	0.59
2:F:84:TYR:HB2	2:F:153:GLN:CG	2.33	0.59
1:D:321:PHE:HE1	1:D:348:LYS:HG3	1.68	0.59
2:F:79:SER:CB	2:F:86:LYS:HE2	2.31	0.58
2:E:84:TYR:HB2	2:E:153:GLN:CG	2.33	0.58
2:F:196:ILE:HD12	2:F:197:ILE:HD13	1.84	0.58
1:D:321:PHE:CE1	1:D:348:LYS:HG3	2.38	0.58
2:E:214:VAL:HG22	2:E:216:TYR:CE1	2.37	0.58
2:F:227:TRP:CZ2	2:F:231:HIS:CE1	2.92	0.58
1:D:196:SER:HB2	1:D:219:ASP:OD1	2.04	0.58
1:D:159:GLU:HG2	1:D:205:LYS:NZ	2.18	0.58
1:D:106:ARG:HG2	1:D:120:CYS:HB2	1.86	0.58
1:D:5:HIS:HB3	1:D:23:ILE:HG22	1.86	0.58
1:B:158:ALA:HA	1:B:203:LEU:CD1	2.32	0.58
1:D:7:LEU:O	1:D:370:THR:CG2	2.50	0.58
1:D:14:ARG:HG2	1:D:114:GLU:OE2	2.04	0.58
1:B:7:LEU:O	1:B:370:THR:CG2	2.52	0.57
2:F:227:TRP:CD2	2:F:231:HIS:HE1	2.22	0.57
2:E:227:TRP:CZ2	2:E:231:HIS:CE1	2.92	0.57
2:E:283:PRO:HG3	2:F:245:TRP:CD1	2.39	0.57
1:D:4:ILE:HG12	1:D:421:VAL:HG22	1.86	0.57
1:B:23:ILE:CD1	1:B:28:LEU:HG	2.23	0.57
1:B:4:ILE:HG12	1:B:421:VAL:HG22	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:84:TYR:HB2	2:F:153:GLN:HG3	1.86	0.57
1:B:285:LYS:HB3	1:B:286:PRO:HD3	1.87	0.57
1:B:106:ARG:HG2	1:B:120:CYS:HB2	1.85	0.57
2:E:84:TYR:HB2	2:E:153:GLN:HG3	1.87	0.56
1:B:114:ARG:HG2	1:B:114:GLU:OE2	2.05	0.56
2:F:198:THR:HG22	2:F:200:THR:N	2.15	0.56
1:B:34:SER:CB	1:B:35:PRO:HA	2.34	0.56
1:B:159:GLU:HG2	1:B:205:LYS:NZ	2.21	0.56
2:E:198:THR:HG22	2:E:200:THR:N	2.13	0.56
2:E:227:TRP:CD2	2:E:231:HIS:HE1	2.23	0.56
1:D:285:LYS:HB3	1:D:286:PRO:HD3	1.86	0.56
1:D:216:SER:HB2	1:D:244:VAL:HG22	1.88	0.56
1:B:381:LYS:HG2	1:B:382:ASN:H	1.70	0.56
1:B:23:ILE:HD11	1:B:28:LEU:CG	2.24	0.56
2:E:196:ILE:CD1	2:E:197:ILE:HD13	2.36	0.56
2:E:132:THR:O	2:E:136:GLU:HG3	2.07	0.55
1:D:34:SER:HA	1:D:36:ASN:H	1.71	0.55
1:B:151:ARG:NH1	1:B:153:ASN:HD22	2.03	0.55
2:E:227:TRP:CZ2	2:E:231:HIS:HE1	2.25	0.55
1:B:196:SER:HB2	1:B:219:ASP:OD1	2.06	0.55
1:D:158:ALA:HA	1:D:203:LEU:CD1	2.32	0.54
1:B:8:GLN:HE22	1:B:23:ILE:HA	1.72	0.54
1:D:151:ARG:HH12	1:D:153:ASN:ND2	2.02	0.54
2:E:97:VAL:HA	2:E:121:LEU:HB2	1.90	0.54
2:F:227:TRP:CZ2	2:F:231:HIS:HE1	2.25	0.54
2:F:196:ILE:CD1	2:F:197:ILE:HD13	2.38	0.54
1:B:199:THR:HG22	1:B:216:SER:OG	2.08	0.54
1:B:34:SER:HA	1:B:36:ASN:H	1.72	0.54
2:F:107:GLY:O	2:F:111:ILE:HG12	2.08	0.54
1:B:151:ARG:HH12	1:B:153:ASN:ND2	2.04	0.54
1:D:296:PRO:HB2	4:D:2067:HOH:O	2.08	0.53
1:D:305:ILE:HD12	1:D:305:ILE:N	2.23	0.53
1:D:194:HIS:NE2	1:D:223:LYS:HE2	2.23	0.53
2:F:132:THR:O	2:F:136:GLU:HG3	2.09	0.53
1:D:151:ARG:NH1	1:D:153:ASN:HD22	2.05	0.53
1:B:41:ALA:HB1	1:B:139:VAL:HG23	1.89	0.53
1:B:359:SER:OG	1:B:370:THR:HB	2.08	0.53
2:E:111:ILE:HD12	2:E:142:LEU:CD1	2.39	0.52
1:D:256:LEU:CD2	1:D:322:VAL:HG21	2.39	0.52
2:F:97:VAL:HA	2:F:121:LEU:HB2	1.90	0.52
2:E:84:TYR:CB	2:E:153:GLN:CG	2.87	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:381:LYS:HG2	1:D:382:ASN:H	1.75	0.52
1:B:256:LEU:CD2	1:B:322:VAL:HG21	2.40	0.52
1:D:8:GLN:HE22	1:D:23:ILE:HA	1.74	0.52
1:D:199:THR:HG22	1:D:216:SER:OG	2.10	0.52
2:F:240:ARG:HG3	2:F:277:PHE:CZ	2.45	0.52
2:E:222:LYS:HE2	2:E:226:GLU:OE1	2.10	0.52
2:E:227:TRP:CH2	2:E:231:HIS:HE1	2.27	0.51
2:F:227:TRP:CE3	2:F:231:HIS:HE1	2.28	0.51
2:F:111:ILE:HD12	2:F:142:LEU:CD1	2.40	0.51
1:D:355:TYR:OH	1:D:376:SER:HB3	2.10	0.51
1:B:341:GLN:HG2	4:B:2045:HOH:O	2.08	0.51
2:E:227:TRP:CZ3	2:E:231:HIS:HE1	2.29	0.51
2:F:227:TRP:CH2	2:F:231:HIS:HE1	2.28	0.51
2:E:227:TRP:CE3	2:E:231:HIS:HE1	2.29	0.51
1:B:204:ILE:HG22	1:B:212:PHE:HB2	1.93	0.51
1:B:355:TYR:OH	1:B:376:SER:HB3	2.10	0.50
1:B:4:ILE:HG23	1:B:414:VAL:HG11	1.94	0.50
2:F:106:PHE:HD2	2:F:134:TYR:CD2	2.30	0.50
1:D:41:ALA:HB1	1:D:139:VAL:HG23	1.93	0.50
1:B:194:HIS:NE2	1:B:223:LYS:HE2	2.27	0.50
2:E:157:VAL:HG22	2:E:158:LEU:N	2.27	0.50
2:F:150:HIS:CE1	4:F:2017:HOH:O	2.65	0.50
2:F:227:TRP:CZ3	2:F:231:HIS:HE1	2.30	0.50
2:E:107:GLY:O	2:E:111:ILE:HG12	2.12	0.50
2:F:84:TYR:CB	2:F:153:GLN:CG	2.90	0.49
1:D:295:PRO:CB	1:D:296:PRO:HD2	2.42	0.49
1:B:195:VAL:HG12	1:B:195:VAL:O	2.12	0.49
2:E:209:LEU:HD13	2:E:279:ARG:HB2	1.93	0.49
1:D:223:LYS:HD3	1:D:236:TRP:CH2	2.47	0.49
2:F:222:LYS:HE2	2:F:226:GLU:OE1	2.12	0.49
1:B:133:ASP:HB3	1:B:136:SER:HB3	1.94	0.49
1:B:197:MET:HE3	4:B:2018:HOH:O	2.12	0.49
1:B:34:SER:HB3	1:B:35:PRO:HA	1.95	0.48
2:F:209:LEU:HD13	2:F:279:ARG:HB2	1.94	0.48
2:E:106:PHE:HD2	2:E:134:TYR:CD2	2.31	0.48
1:B:192:LEU:HD11	1:B:213:ILE:HD13	1.94	0.48
1:B:374:LYS:HG2	1:B:427:ILE:HD12	1.95	0.48
1:D:408:SER:O	1:D:412:GLN:HG2	2.13	0.48
2:E:163:MET:HE3	2:E:201:LEU:HD11	1.95	0.48
1:D:374:LYS:HG2	1:D:427:ILE:HD12	1.95	0.48
1:D:294:ALA:HB1	1:D:295:PRO:HD2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:240:HIS:CD2	1:B:244:VAL:HG22	2.49	0.48
1:B:290:ASP:HB2	1:B:293:LEU:HD12	1.94	0.48
1:B:226:HIS:HD2	4:B:2020:HOH:O	1.96	0.48
2:E:181:CYS:CB	4:E:2043:HOH:O	2.58	0.47
2:F:227:TRP:CH2	2:F:231:HIS:CE1	3.03	0.47
2:E:286:LEU:HB3	2:F:250:GLU:HG2	1.95	0.47
1:B:411:ILE:O	1:B:415:GLN:HG3	2.15	0.47
2:E:227:TRP:CD2	2:E:231:HIS:CE1	3.02	0.47
2:E:183:PRO:HG3	2:E:228:MET:CE	2.44	0.47
1:D:133:ASP:HB3	1:D:136:SER:HB3	1.96	0.47
2:F:150:HIS:HE1	4:F:2017:HOH:O	1.98	0.47
1:D:186:PHE:CZ	1:D:188:GLN:CG	2.98	0.47
1:B:8:GLN:H	1:B:8:GLN:HE21	1.62	0.47
1:D:34:SER:HA	1:D:36:ASN:N	2.30	0.47
2:E:227:TRP:CH2	2:E:231:HIS:CE1	3.02	0.47
1:B:157:ILE:HD12	4:B:2008:HOH:O	2.15	0.47
2:F:198:THR:O	2:F:202:LEU:HG	2.15	0.47
1:D:133:ASP:OD1	1:D:136:SER:HB2	2.14	0.47
2:F:126:GLU:O	2:F:159:ARG:HA	2.15	0.47
2:E:84:TYR:HB3	2:E:153:GLN:HG3	1.97	0.46
1:D:290:ASP:HB2	1:D:293:LEU:HD12	1.97	0.46
1:D:159:GLU:HG2	1:D:205:LYS:HZ3	1.80	0.46
1:B:398:VAL:HG23	4:B:2056:HOH:O	2.14	0.46
2:F:227:TRP:CD2	2:F:231:HIS:CE1	3.02	0.46
1:B:109:ARG:CG	1:B:157:ILE:HD11	2.46	0.46
1:D:411:ILE:O	1:D:415:GLN:HG3	2.15	0.46
1:B:408:SER:O	1:B:412:GLN:HG2	2.15	0.46
1:D:271:THR:CG2	1:D:273:LYS:CB	2.94	0.46
1:B:159:GLU:HG2	1:B:205:LYS:HZ2	1.80	0.46
1:B:31:LYS:HG3	1:B:41:ALA:HB2	1.98	0.46
2:E:126:GLU:O	2:E:159:ARG:HA	2.16	0.46
1:D:204:ILE:HG22	1:D:212:PHE:HB2	1.96	0.46
2:E:235:HIS:CE1	2:E:237:LEU:HB2	2.51	0.45
1:D:358:ILE:HG22	1:D:370:THR:HG22	1.98	0.45
1:D:229:GLN:H	1:D:229:GLN:HE21	1.64	0.45
2:E:240:ARG:HG3	2:E:277:PHE:CZ	2.51	0.45
1:B:171:GLY:C	1:B:198:LEU:HD22	2.37	0.45
1:B:8:GLN:NE2	1:B:23:ILE:HA	2.31	0.45
1:B:34:SER:HA	1:B:36:ASN:N	2.31	0.45
2:F:157:VAL:HG22	2:F:158:LEU:N	2.32	0.45
1:D:8:GLN:NE2	1:D:23:ILE:HA	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:271:THR:CG2	1:B:273:LYS:CB	2.94	0.45
1:B:186:PHE:CZ	1:B:188:GLN:CG	3.00	0.45
2:E:198:THR:O	2:E:202:LEU:HG	2.17	0.45
2:F:98:THR:HG22	2:F:99:ILE:HG13	1.98	0.45
1:D:186:PHE:CZ	1:D:188:GLN:HG3	2.51	0.45
1:B:150:LYS:HD2	1:B:170:PHE:CD1	2.51	0.45
1:D:109:ARG:CG	1:D:157:ILE:HD11	2.47	0.45
2:E:148:SER:CB	2:E:150:HIS:CE1	2.96	0.45
2:E:284:ALA:O	2:F:250:GLU:N	2.49	0.44
1:D:271:THR:HG21	1:D:273:LYS:HB3	2.00	0.44
1:B:358:ILE:HD12	1:B:358:ILE:HA	1.68	0.44
1:B:171:GLY:HA2	1:B:198:LEU:HD22	1.98	0.44
2:E:98:THR:HG22	2:E:99:ILE:HG13	1.99	0.44
1:D:192:LEU:HD11	1:D:213:ILE:HD13	1.98	0.44
1:D:358:ILE:HD12	1:D:358:ILE:HA	1.73	0.44
1:B:360:LEU:HD12	1:B:361:SER:N	2.32	0.44
2:E:166:LEU:N	2:E:167:PRO:CD	2.81	0.44
1:B:133:ASP:OD1	1:B:136:SER:HB2	2.17	0.44
2:F:235:HIS:CE1	2:F:237:LEU:HB2	2.53	0.44
2:E:237:LEU:HA	2:E:237:LEU:HD12	1.82	0.43
1:B:186:PHE:CZ	1:B:188:GLN:HG3	2.52	0.43
1:B:253:TYR:CZ	1:B:270:LYS:HD2	2.53	0.43
1:D:330:LYS:HG3	4:D:2081:HOH:O	2.18	0.43
1:B:139:VAL:O	1:B:141:LYS:HE2	2.19	0.43
2:E:236:PRO:HA	4:E:2037:HOH:O	2.18	0.43
1:D:253:TYR:CZ	1:D:270:LYS:HD2	2.53	0.43
1:D:417:ASP:OD2	1:D:419:ASN:HB3	2.18	0.43
2:F:183:PRO:HG3	2:F:228:MET:CE	2.48	0.43
2:F:214:VAL:HG22	2:F:216:TYR:CZ	2.53	0.43
1:B:381:LYS:HB3	1:B:381:LYS:HE3	1.76	0.43
1:B:40:PHE:CE1	1:B:413:SER:HB2	2.54	0.43
2:F:84:TYR:HB3	2:F:153:GLN:HG3	2.00	0.43
1:D:196:SER:OG	1:D:217:ASP:HB2	2.19	0.43
1:D:381:LYS:HB3	1:D:381:LYS:HE3	1.75	0.43
2:E:245:TRP:CD1	2:F:283:PRO:HG3	2.54	0.43
2:F:166:LEU:N	2:F:167:PRO:CD	2.82	0.43
1:B:271:THR:CG2	1:B:273:LYS:HB2	2.48	0.43
1:B:28:LEU:HD21	1:B:43:LYS:HD3	2.00	0.42
1:D:271:THR:CG2	1:D:273:LYS:HB2	2.49	0.42
1:B:271:THR:HG21	1:B:273:LYS:HB3	2.01	0.42
1:B:223:LYS:HD3	1:B:236:TRP:CH2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:40:PHE:CE1	1:D:413:SER:HB2	2.53	0.42
1:D:374:LYS:HB2	1:D:374:LYS:HE3	1.86	0.42
1:B:196:SER:OG	1:B:217:ASP:HB2	2.20	0.42
2:E:235:HIS:HA	2:E:236:PRO:HD3	1.91	0.42
1:B:360:LEU:HD12	1:B:361:SER:H	1.84	0.42
1:B:160:ASP:CG	1:B:162:THR:HG22	2.39	0.42
1:B:315:LYS:NZ	1:B:315:LYS:HB3	2.34	0.42
1:D:34:SER:HB3	1:D:35:PRO:CA	2.48	0.42
1:B:417:ASP:OD2	1:B:419:ASN:HB3	2.19	0.42
2:E:249:ASP:OD1	2:E:251:CYS:HB2	2.19	0.42
1:D:160:ASP:CG	1:D:162:THR:HG22	2.40	0.42
1:D:169:LYS:O	1:D:197:MET:HG3	2.19	0.42
1:D:171:GLY:C	1:D:198:LEU:HD22	2.40	0.42
1:D:139:VAL:O	1:D:141:LYS:HE2	2.20	0.42
1:D:294:ALA:HA	1:D:295:PRO:HD3	1.56	0.42
1:B:285:LYS:CB	4:B:2037:HOH:O	2.67	0.42
1:D:31:LYS:HG3	1:D:41:ALA:HB2	2.01	0.42
1:B:374:LYS:HE3	1:B:374:LYS:HB2	1.85	0.42
1:D:235:LYS:HG2	1:D:236:TRP:N	2.35	0.42
1:B:221:HIS:HA	4:B:2022:HOH:O	2.18	0.42
1:D:22:ILE:HD11	1:D:108:LEU:HG	2.02	0.42
1:D:240:HIS:CE1	1:D:257:SER:OG	2.63	0.41
1:B:204:ILE:HD13	1:B:204:ILE:HA	1.92	0.41
2:F:182:PHE:HA	4:F:2030:HOH:O	2.20	0.41
2:E:70:PRO:HD3	2:E:169:PHE:CE1	2.55	0.41
1:B:348:LYS:O	1:B:349:GLN:HB2	2.20	0.41
1:B:195:VAL:O	1:B:195:VAL:CG1	2.69	0.41
2:F:180:PHE:CD2	2:F:196:ILE:HD13	2.55	0.41
1:D:106:ARG:HG3	1:D:107:ASN:ND2	2.35	0.41
1:B:121:ALA:HB2	1:B:128:LEU:HD11	2.03	0.41
1:D:229:GLN:HG2	2:F:285:ILE:HD11	2.01	0.41
2:E:158:LEU:HD12	2:E:158:LEU:O	2.20	0.41
2:E:81:LEU:C	2:E:83:PRO:HD3	2.40	0.41
2:F:148:SER:CB	2:F:150:HIS:CE1	2.96	0.41
1:D:32:TYR:CZ	1:D:402:LYS:HE3	2.55	0.41
2:F:225:HIS:NE2	3:F:1287:PO4:O1	2.51	0.41
1:B:231:PHE:CD1	1:B:232:ILE:HG13	2.55	0.41
1:B:169:LYS:O	1:B:197:MET:HG3	2.21	0.41
1:B:124:ASP:HB3	1:B:126:SER:OG	2.21	0.41
2:E:119:GLU:HB3	4:E:2010:HOH:O	2.20	0.41
2:F:123:LEU:HD23	2:F:170:PHE:HZ	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:148:PHE:HZ	1:B:176:ILE:HG13	1.85	0.41
1:D:4:ILE:HG23	1:D:414:VAL:HG11	2.03	0.40
1:B:282:SER:C	4:B:2037:HOH:O	2.58	0.40
2:F:166:LEU:HD23	2:F:166:LEU:HA	1.88	0.40
1:D:155:ILE:HA	1:D:165:ILE:O	2.21	0.40
1:B:235:LYS:HG2	1:B:236:TRP:N	2.36	0.40
1:B:283:LEU:O	1:B:286:PRO:HD2	2.22	0.40
2:E:119:GLU:O	2:E:119:GLU:HG2	2.22	0.40
2:E:96:LYS:HD3	2:E:96:LYS:HA	1.84	0.40
1:B:106:ARG:HG3	1:B:107:ASN:ND2	2.36	0.40
1:D:271:THR:HG21	1:D:273:LYS:CB	2.52	0.40
2:E:196:ILE:HG13	2:E:196:ILE:H	1.69	0.40
1:D:109:ARG:HG3	1:D:157:ILE:HD11	2.04	0.40
1:D:198:LEU:HA	1:D:198:LEU:HD12	1.88	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	368/450 (82%)	344 (94%)	24 (6%)	0	100	100
1	D	370/450 (82%)	348 (94%)	21 (6%)	1 (0%)	46	63
2	E	195/254 (77%)	190 (97%)	5 (3%)	0	100	100
2	F	195/254 (77%)	189 (97%)	6 (3%)	0	100	100
All	All	1128/1408 (80%)	1071 (95%)	56 (5%)	1 (0%)	56	74

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	295	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	B	344/411 (84%)	318 (92%)	26 (8%)	16	25
1	D	346/411 (84%)	320 (92%)	26 (8%)	17	26
2	E	180/228 (79%)	163 (91%)	17 (9%)	11	16
2	F	180/228 (79%)	162 (90%)	18 (10%)	9	14
All	All	1050/1278 (82%)	963 (92%)	87 (8%)	14	21

All (87) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	3	VAL
1	B	8	GLN
1	B	9	ASN
1	B	11	LEU
1	B	23	ILE
1	B	100	PRO
1	B	127	LEU
1	B	141	LYS
1	B	187	THR
1	B	192	LEU
1	B	196	SER
1	B	198	LEU
1	B	199	THR
1	B	201	VAL
1	B	204	ILE
1	B	213	ILE
1	B	220	GLU
1	B	229	GLN
1	B	235	LYS
1	B	248	CYS
1	B	262	ASP
1	B	315	LYS
1	B	358	ILE
1	B	370	THR

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Mol	Chain	Res	Type
1	B	414	VAL
1	B	433	VAL
1	D	3	VAL
1	D	8	GLN
1	D	9	ASN
1	D	11	LEU
1	D	23	ILE
1	D	100	PRO
1	D	127	LEU
1	D	141	LYS
1	D	187	THR
1	D	192	LEU
1	D	196	SER
1	D	198	LEU
1	D	199	THR
1	D	201	VAL
1	D	204	ILE
1	D	213	ILE
1	D	220	GLU
1	D	229	GLN
1	D	235	LYS
1	D	248	CYS
1	D	262	ASP
1	D	315	LYS
1	D	358	ILE
1	D	370	THR
1	D	414	VAL
1	D	433	VAL
2	E	98	THR
2	E	139	ILE
2	E	146	THR
2	E	155	ILE
2	E	181	CYS
2	E	195	ARG
2	E	196	ILE
2	E	200	THR
2	E	209	LEU
2	E	214	VAL
2	E	217	THR
2	E	230	LYS
2	E	237	LEU
2	E	240	ARG

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Mol	Chain	Res	Type
2	E	243	LYS
2	E	253	LYS
2	E	255	MET
2	F	98	THR
2	F	129	VAL
2	F	139	ILE
2	F	146	THR
2	F	155	ILE
2	F	181	CYS
2	F	195	ARG
2	F	196	ILE
2	F	200	THR
2	F	209	LEU
2	F	214	VAL
2	F	217	THR
2	F	230	LYS
2	F	237	LEU
2	F	240	ARG
2	F	243	LYS
2	F	253	LYS
2	F	255	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	8	GLN
1	B	33	GLN
1	B	37	HIS
1	B	153	ASN
1	B	211	GLN
1	B	226	HIS
1	B	229	GLN
1	B	240	HIS
1	B	274	ASN
1	B	292	HIS
1	B	356	ASN
1	D	8	GLN
1	D	33	GLN
1	D	37	HIS
1	D	153	ASN
1	D	210	HIS
1	D	211	GLN

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Mol	Chain	Res	Type
1	D	226	HIS
1	D	229	GLN
1	D	240	HIS
1	D	292	HIS
1	D	356	ASN
2	E	90	ASN
2	E	92	GLN
2	E	133	ASN
2	E	199	ASN
2	E	231	HIS
2	E	257	ASN
2	F	90	ASN
2	F	92	GLN
2	F	133	ASN
2	F	150	HIS
2	F	168	ASN
2	F	199	ASN
2	F	231	HIS
2	F	257	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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	PO4	E	1287	-	4,4,4	0.39	0	6,6,6	0.27	0
3	PO4	F	1287	-	4,4,4	0.39	0	6,6,6	0.30	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PO4	E	1287	-	-	0/0/0/0	0/0/0/0
3	PO4	F	1287	-	-	0/0/0/0	0/0/0/0

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	1287	PO4	1	0

## 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	B	374/450 (83%)	0.73	45 (12%) 6 5	11, 34, 73, 100	0
1	D	376/450 (83%)	0.42	14 (3%) 45 46	10, 29, 72, 100	0
2	E	201/254 (79%)	0.13	3 (1%) 76 75	8, 26, 70, 90	0
2	F	201/254 (79%)	0.26	4 (1%) 68 68	8, 25, 70, 90	0
All	All	1152/1408 (81%)	0.44	66 (5%) 27 27	8, 30, 72, 100	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	186	PHE	10.3
1	B	159	GLU	6.8
1	D	187	THR	6.6
1	B	433	VAL	5.8
1	B	2	SER	5.7
1	B	185	LYS	5.6
1	B	305	ILE	5.5
2	E	286	LEU	5.3
2	F	145	ASN	5.2
1	B	294	ALA	5.0
1	B	3	VAL	4.9
1	B	158	ALA	4.7
1	B	291	GLN	4.1
1	B	187	THR	3.9
1	B	48	PHE	3.9
1	B	292	HIS	3.8
1	B	421	VAL	3.8
1	B	396	PHE	3.7
1	D	186	PHE	3.5
1	D	100	PRO	3.5
1	B	157	ILE	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	141	LYS	3.2
1	B	103	SER	3.1
1	D	434	SER	3.1
1	B	262	ASP	3.1
1	B	420	LEU	3.0
1	D	305	ILE	3.0
1	B	389	TYR	2.9
1	B	148	PHE	2.9
1	B	431	TYR	2.8
2	E	139	ILE	2.8
1	B	161	ASP	2.8
1	B	414	VAL	2.8
1	D	37	HIS	2.8
1	B	284	ILE	2.8
1	D	433	VAL	2.7
1	B	47	ASP	2.7
1	B	430	LEU	2.7
1	B	293	LEU	2.6
2	F	286	LEU	2.6
1	B	375	GLU	2.6
1	B	425	GLU	2.6
1	D	415	GLN	2.6
1	D	436	LEU	2.6
1	B	123	SER	2.5
1	B	5	HIS	2.5
1	B	290	ASP	2.5
1	B	8	GLN	2.4
2	F	149	LYS	2.4
1	D	423	LYS	2.3
1	B	308	PHE	2.3
2	E	66	GLN	2.3
1	B	134	LYS	2.3
1	B	355	TYR	2.3
1	B	184	GLU	2.2
1	D	378	GLY	2.2
1	B	394	ASN	2.2
1	D	191	ILE	2.1
2	F	147	ALA	2.1
1	D	204	ILE	2.1
1	D	138	ASN	2.1
1	B	4	ILE	2.1
1	B	422	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	138	ASN	2.1
1	B	178	ILE	2.0
1	B	415	GLN	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	PO4	F	1287	5/5	0.94	0.15	-	25,33,81,88	0
3	PO4	E	1287	5/5	0.91	0.19	-	30,45,76,80	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.