



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 03:20 PM GMT

PDB ID : 3P5R  
Title : Crystal Structure of Taxadiene Synthase from Pacific Yew (*Taxus brevifolia*)  
in complex with Mg<sup>2+</sup> and 2-fluorogeranylgeranyldiphosphate  
Authors : Koksai, M.; Christianson, D.W.  
Deposited on : 2010-10-10  
Resolution : 2.25 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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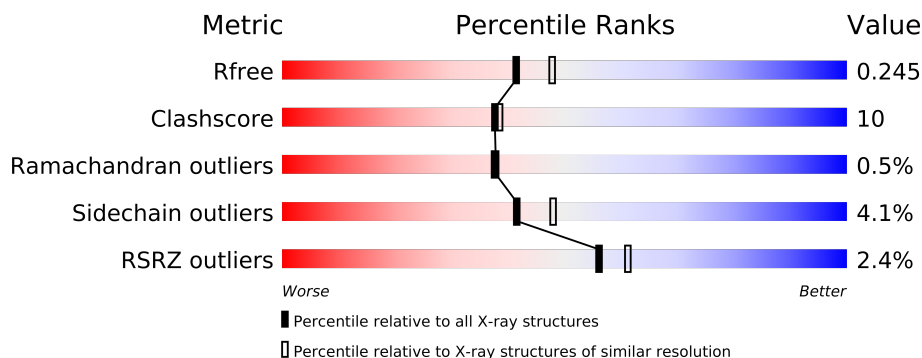
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.25 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))
$R_{free}$	66092	1112 (2.26-2.22)
Clashscore	79885	1317 (2.26-2.22)
Ramachandran outliers	78287	1282 (2.26-2.22)
Sidechain outliers	78261	1282 (2.26-2.22)
RSRZ outliers	66119	1112 (2.26-2.22)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	764	
1	B	764	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	MG	A	901	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13321 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Taxadiene synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	750	Total	C	N	O	S	0	4	0
			6078	3881	1022	1141	34			
1	B	738	Total	C	N	O	S	0	4	0
			5971	3812	999	1126	34			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	107	MET	-	INITIATING METHIONINE	UNP Q41594
A	863	GLY	-	EXPRESSION TAG	UNP Q41594
A	864	SER	-	EXPRESSION TAG	UNP Q41594
A	865	HIS	-	EXPRESSION TAG	UNP Q41594
A	866	HIS	-	EXPRESSION TAG	UNP Q41594
A	867	HIS	-	EXPRESSION TAG	UNP Q41594
A	868	HIS	-	EXPRESSION TAG	UNP Q41594
A	869	HIS	-	EXPRESSION TAG	UNP Q41594
A	870	HIS	-	EXPRESSION TAG	UNP Q41594
B	107	MET	-	INITIATING METHIONINE	UNP Q41594
B	863	GLY	-	EXPRESSION TAG	UNP Q41594
B	864	SER	-	EXPRESSION TAG	UNP Q41594
B	865	HIS	-	EXPRESSION TAG	UNP Q41594
B	866	HIS	-	EXPRESSION TAG	UNP Q41594
B	867	HIS	-	EXPRESSION TAG	UNP Q41594
B	868	HIS	-	EXPRESSION TAG	UNP Q41594
B	869	HIS	-	EXPRESSION TAG	UNP Q41594
B	870	HIS	-	EXPRESSION TAG	UNP Q41594

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

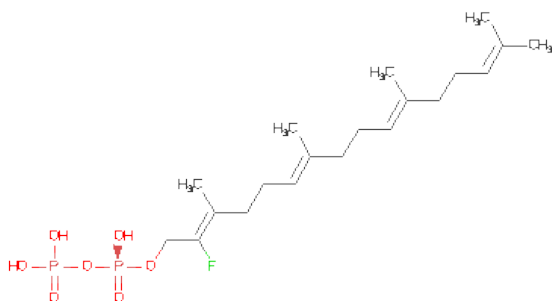
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	3	Total	Mg	0	0
			3	3		

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

- Molecule 3 is (2Z,6E,10E)-2-FLUORO-3,7,11,15-TETRAMETHYLHEXADECA-2,6,10,14-TETRAEN-1-YLTRIHYDROGEN DIPHOSPHATE (three-letter code: FGG) (formula:  $C_{20}H_{35}FO_7P_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	O	P	0	0
			30	20	1	7	2		
3	B	1	Total	C	F	O	P	0	0
			30	20	1	7	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	569	Total	O	0	0
			569	569		
4	B	637	Total	O	0	0
			637	637		





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	54.05Å 201.98Å 81.43Å 90.00° 91.60° 90.00°	Depositor
Resolution (Å)	43.41 – 2.25 43.41 – 2.25	Depositor EDS
% Data completeness (in resolution range)	81.0 (43.41-2.25) 81.0 (43.41-2.25)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	0.13	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.41 (at 2.24Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, $R_{free}$	0.187 , 0.250 0.183 , 0.245	Depositor DCC
$R_{free}$ test set	1914 reflections (2.86%)	DCC
Wilson B-factor (Å <sup>2</sup> )	15.6	Xtriage
Anisotropy	0.142	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 32.4	EDS
Estimated twinning fraction	0.128 for h,-k,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 70530 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	13321	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.25% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

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

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	0.31	0/6239	0.45	0/8448
1	B	0.31	0/6127	0.46	0/8297
All	All	0.31	0/12366	0.45	0/16745

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6078	0	5956	121	0
1	B	5971	0	5854	112	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
3	A	30	0	32	7	0
3	B	30	0	32	7	0
4	A	569	0	0	7	0
4	B	637	0	0	8	0
All	All	13321	0	11874	241	0



Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 10.

All (241) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:571:MET:HE1	1:B:650:CYS:HA	1.50	0.94
1:A:175:ASP:N	1:A:176:GLY:HA2	1.81	0.93
1:A:669:GLN:HE22	1:A:731:VAL:H	1.24	0.86
1:B:254:TYR:HA	1:B:259:ILE:HD11	1.58	0.84
1:A:609:GLN:HE22	1:A:719:CYS:HB3	1.43	0.84
1:B:669:GLN:HE22	1:B:731:VAL:H	1.23	0.83
3:A:911:FGG:H04	3:A:911:FGG:H17A	1.60	0.83
1:A:850:ASP:HB2	4:A:1193:HOH:O	1.78	0.82
3:B:911:FGG:H04	3:B:911:FGG:H17A	1.62	0.82
1:B:151:ILE:HG12	1:B:157:GLU:HG2	1.64	0.80
1:A:537:ASN:HD22	1:A:540:CYS:H	1.30	0.79
1:B:609:GLN:HE22	1:B:719:CYS:HB3	1.48	0.77
1:B:340:PHE:CZ	1:B:345:PRO:HG3	2.19	0.77
1:A:225:ASN:HD22	1:A:225:ASN:H	1.32	0.76
1:A:221:LEU:HD12	1:A:253:PRO:HD2	1.69	0.75
1:A:256:LEU:HB2	1:A:259:ILE:HG22	1.67	0.75
1:B:629:THR:HB	1:B:685:PHE:HB3	1.69	0.74
1:A:629:THR:HB	1:A:685:PHE:HB3	1.73	0.71
1:A:577:THR:HG22	1:A:578:ARG:H	1.54	0.71
1:A:174:GLN:HE21	1:A:174:GLN:HA	1.56	0.70
1:B:327:GLU:O	1:B:331:THR:HG22	1.93	0.69
1:A:150:THR:HG23	1:A:158:LYS:HG3	1.74	0.69
1:B:625:LEU:O	1:B:629:THR:HG23	1.92	0.69
1:A:664:ASP:HA	1:A:667:LYS:HE3	1.75	0.68
1:B:148:LEU:HD13	1:B:321:LEU:HD23	1.75	0.68
1:B:546:LEU:HD11	1:B:862:VAL:HG11	1.76	0.67
1:A:207:HIS:O	1:A:211:GLN:HG2	1.95	0.67
1:B:473:GLU:HB3	1:B:474:PRO:HD3	1.77	0.67
1:A:206:GLY:O	1:A:209:GLN:HG2	1.95	0.66
1:B:853:ARG:HD2	4:B:988:HOH:O	1.95	0.66
1:B:760:LYS:HG2	1:B:835:TYR:HB3	1.77	0.66
3:A:911:FGG:H16B	4:A:1182:HOH:O	1.95	0.66
1:A:119:LEU:HB3	1:A:543:LEU:HG	1.78	0.66
1:A:145:VAL:HG11	1:A:337:LEU:HD11	1.78	0.65
1:B:571:MET:CE	1:B:650:CYS:HA	2.26	0.65
1:A:797:VAL:HA	4:A:1074:HOH:O	1.97	0.65
1:A:580[B]:ARG:HH21	3:A:911:FGG:H19B	1.63	0.64
1:B:622:LEU:HD22	1:B:626:LYS:HE3	1.80	0.64
1:A:537:ASN:HD21	1:A:539:LYS:HB2	1.63	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:117:ASP:O	1:A:121:VAL:HG23	1.98	0.63
1:A:238:PHE:HB3	1:A:239:PRO:HD3	1.80	0.63
1:B:233:ASP:O	1:B:237:ILE:HG12	1.98	0.63
1:B:242:LEU:HD13	1:B:259:ILE:HD12	1.80	0.63
1:B:511:ALA:O	1:B:515:ILE:HG13	1.99	0.62
1:A:644:PRO:O	1:A:648:GLN:HG3	1.99	0.61
1:A:720:THR:O	1:A:724:ILE:HG12	1.99	0.61
1:B:625:LEU:HD12	1:B:692:ARG:HG3	1.83	0.60
1:B:235:GLN:CD	1:B:235:GLN:H	2.04	0.60
1:A:563:THR:OG1	1:A:578:ARG:HG3	2.01	0.60
1:B:792:LYS:O	1:B:796:ARG:HG2	2.02	0.60
1:A:280:ILE:HG21	1:A:300:ILE:HD13	1.84	0.59
1:A:177:SER:HB3	1:A:216:PHE:CE1	2.37	0.59
3:A:911:FGG:H04	3:A:911:FGG:C17	2.33	0.58
1:B:609:GLN:NE2	1:B:719:CYS:HB3	2.18	0.58
1:A:225:ASN:N	1:A:225:ASN:HD22	2.00	0.58
1:A:473:GLU:HB3	1:A:474:PRO:HD3	1.85	0.58
1:A:327:GLU:O	1:A:331:THR:HG22	2.03	0.58
1:A:826:ASN:HD22	1:A:828:ARG:HH12	1.50	0.57
1:B:611:LEU:HD21	1:B:650:CYS:SG	2.45	0.57
1:B:194:THR:O	1:B:198:ILE:HG13	2.04	0.57
1:B:609:GLN:HB3	3:B:911:FGG:H17B	1.87	0.57
1:B:238:PHE:HB3	1:B:239:PRO:HD3	1.86	0.56
1:A:340:PHE:CZ	1:A:345:PRO:HG3	2.40	0.56
1:A:453:ARG:O	1:A:456:ASP:HB2	2.06	0.56
1:B:198:ILE:O	1:B:202:VAL:HG23	2.06	0.56
1:B:537:ASN:HD22	1:B:540:CYS:H	1.52	0.56
1:B:222:ARG:HG2	1:B:256:LEU:HD11	1.87	0.55
1:B:587:SER:OG	3:B:911:FGG:H13	2.05	0.55
1:A:579:HIS:O	1:A:583:GLU:HG3	2.06	0.55
1:B:609:GLN:CB	3:B:911:FGG:H17B	2.36	0.55
3:B:911:FGG:H04	3:B:911:FGG:C17	2.35	0.55
1:A:178:TRP:CE3	1:A:193:THR:HG23	2.42	0.55
1:B:141:ASP:HB3	1:B:314:PRO:HB2	1.87	0.55
1:B:206:GLY:O	1:B:209:GLN:HG2	2.07	0.55
1:B:755:LEU:O	1:B:759:THR:HG23	2.07	0.55
1:B:847:GLU:HG2	1:B:848:ILE:H	1.72	0.55
1:B:788:GLU:HG2	1:B:792:LYS:HZ3	1.72	0.54
1:A:614:ASP:O	1:A:618:ILE:HB	2.07	0.54
1:B:644:PRO:O	1:B:648:GLN:HG3	2.07	0.54
1:A:340:PHE:CE1	1:A:345:PRO:HG3	2.43	0.53
1:A:220:ASN:HA	1:A:223:LEU:HD13	1.90	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:138:SER:HB2	1:A:346[A]:CYS:SG	2.48	0.53
1:A:571:MET:HE1	1:A:650:CYS:HA	1.90	0.53
1:A:113:GLN:OE1	1:A:550:ILE:HD13	2.09	0.53
1:A:389:ARG:NE	1:A:389:ARG:HA	2.23	0.53
1:A:450:ASN:ND2	1:A:453:ARG:HH11	2.07	0.53
1:B:847:GLU:HG2	1:B:848:ILE:N	2.24	0.53
1:B:414:MET:O	1:B:414:MET:HG3	2.09	0.53
1:A:178:TRP:CZ3	1:A:193:THR:HG23	2.43	0.52
1:A:622:LEU:HD22	1:A:626:LYS:HE3	1.92	0.52
1:A:225:ASN:ND2	1:A:225:ASN:N	2.57	0.52
1:B:562:LEU:HD13	1:B:603:THR:HG21	1.92	0.52
1:B:267:GLU:O	1:B:271:THR:HG23	2.09	0.52
1:B:788:GLU:CD	1:B:792:LYS:HZ1	2.12	0.52
1:B:519:ASP:HB3	1:B:522:TYR:HB2	1.92	0.51
1:A:609:GLN:NE2	1:A:719:CYS:HB3	2.20	0.51
1:A:347:MET:HE2	1:A:526:ARG:HD3	1.92	0.51
1:B:220:ASN:HA	1:B:223:LEU:HD12	1.92	0.51
1:A:175:ASP:N	1:A:176:GLY:CA	2.65	0.51
1:A:537:ASN:ND2	1:A:540:CYS:H	2.04	0.51
1:A:177:SER:HB3	1:A:216:PHE:CD1	2.46	0.51
1:A:316:SER:O	1:A:320:VAL:HG12	2.10	0.51
1:A:357:SER:CB	1:A:535:LEU:HD22	2.41	0.51
1:A:217:ILE:O	1:A:221:LEU:HG	2.12	0.50
1:B:402:LEU:HD11	1:B:451:LEU:HD13	1.93	0.50
1:A:254:TYR:HA	1:A:259:ILE:HD13	1.93	0.50
1:A:148:LEU:CD1	1:A:321:LEU:HD23	2.41	0.50
1:A:551:VAL:HG12	1:A:555:HIS:CE1	2.47	0.50
1:B:288:LEU:C	1:B:288:LEU:HD23	2.33	0.49
1:B:567:LYS:NZ	1:B:567:LYS:HB3	2.26	0.49
1:B:504:MET:CE	1:B:836:LYS:HD3	2.42	0.49
1:B:151:ILE:HA	1:B:157:GLU:HA	1.95	0.49
1:B:557:GLU:HG3	4:B:1058:HOH:O	2.13	0.49
1:B:714:VAL:HG21	1:B:753:TRP:HB3	1.93	0.49
1:A:224:LEU:HB3	1:A:258:PHE:CE2	2.48	0.49
1:B:314:PRO:HG2	1:B:345:PRO:O	2.12	0.49
1:B:460:PRO:O	1:B:461:ASP:HB2	2.13	0.49
1:B:546:LEU:CD1	1:B:862:VAL:HG11	2.43	0.49
1:A:622:LEU:CD2	1:A:626:LYS:HE3	2.43	0.49
1:A:307:ASP:O	1:A:332:PHE:HB2	2.12	0.49
1:A:474:PRO:O	1:A:478:GLU:HB2	2.13	0.49
1:A:826:ASN:ND2	1:A:828:ARG:HH12	2.12	0.48
1:A:427:PHE:O	1:A:434:PHE:HA	2.13	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:679:LYS:HB2	1:A:680:PRO:HD3	1.95	0.48
1:B:679:LYS:HB2	1:B:680:PRO:HD3	1.95	0.48
1:A:161:PHE:HE1	1:A:334:ASN:HD22	1.62	0.48
1:A:814:ASN:OD1	1:B:425:ASN:HB3	2.13	0.48
1:A:755:LEU:HB3	1:A:794:ILE:HG23	1.95	0.48
1:B:213:GLY:O	1:B:217:ILE:HG13	2.13	0.48
1:A:357:SER:HB2	1:A:535:LEU:HD22	1.95	0.47
1:B:763:GLN:HG3	4:B:992:HOH:O	2.13	0.47
1:B:242:LEU:O	1:B:245:ALA:HB3	2.15	0.47
1:B:148:LEU:CD1	1:B:321:LEU:HD23	2.44	0.47
1:A:148:LEU:HD13	1:A:321:LEU:HD23	1.95	0.47
1:A:722:GLN:HB2	1:A:723:PRO:HD3	1.95	0.47
3:A:911:FGG:H05A	3:A:911:FGG:H16	1.56	0.47
1:A:625:LEU:O	1:A:629:THR:HG23	2.14	0.47
1:A:818:MET:HE1	4:A:1047:HOH:O	2.13	0.47
1:A:239:PRO:HB2	1:A:266:ARG:HD2	1.97	0.47
1:B:788:GLU:HB2	4:B:1105:HOH:O	2.14	0.47
1:A:167:TRP:CD1	1:A:343:CYS:HB3	2.50	0.47
1:B:664:ASP:HA	1:B:667:LYS:HE3	1.97	0.47
1:B:113:GLN:NE2	1:B:113:GLN:HA	2.30	0.46
1:A:788:GLU:HG2	4:A:885:HOH:O	2.15	0.46
1:A:147:ARG:HH11	1:A:195:ASN:HD21	1.63	0.46
3:B:911:FGG:H09	3:B:911:FGG:H17	1.78	0.46
1:B:115:ARG:HH22	1:B:862:VAL:HG12	1.81	0.46
1:B:119:LEU:HB3	1:B:543:LEU:HG	1.96	0.46
1:A:288:LEU:HD23	1:A:288:LEU:C	2.35	0.46
1:A:279:ASN:CG	1:A:280:ILE:H	2.19	0.46
1:A:150:THR:HG22	1:A:160:ARG:HA	1.97	0.46
1:B:754:ARG:HD3	1:B:754:ARG:O	2.15	0.46
1:B:720:THR:O	1:B:724:ILE:HG12	2.15	0.46
1:A:235:GLN:HB3	1:A:262:LEU:HD22	1.98	0.46
1:B:132:ASP:HB3	1:B:348:TYR:CE1	2.50	0.46
1:A:111:THR:N	4:A:1158:HOH:O	2.49	0.46
1:A:740:HIS:CG	1:A:741:TYR:H	2.32	0.46
1:A:357:SER:HB2	1:A:535:LEU:CD2	2.46	0.45
1:A:571:MET:CE	1:A:650:CYS:HA	2.46	0.45
1:B:504:MET:HE2	1:B:836:LYS:HD3	1.98	0.45
1:A:745:MET:O	1:A:749:VAL:HG23	2.16	0.45
1:A:204:LYS:HA	1:A:204:LYS:HD2	1.64	0.45
1:B:147:ARG:HH11	1:B:195:ASN:HD21	1.65	0.45
1:B:197:VAL:CG1	1:B:250:ILE:HD13	2.46	0.45
1:A:174:GLN:C	1:A:176:GLY:HA2	2.37	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:197:VAL:HG12	1:B:250:ILE:HD13	1.99	0.45
1:B:340:PHE:CE1	1:B:345:PRO:HG3	2.50	0.45
1:A:677:ILE:O	1:A:680:PRO:HD2	2.16	0.45
1:B:504:MET:HE1	1:B:836:LYS:HB2	1.99	0.45
1:B:348:TYR:CG	1:B:349:SER:HA	2.51	0.45
1:A:190:LEU:HA	1:A:190:LEU:HD12	1.83	0.45
1:B:150:THR:HG23	1:B:158:LYS:HG3	1.98	0.45
1:A:232:PRO:HB3	1:A:442:HIS:CE1	2.52	0.45
1:A:236:ILE:HD11	1:A:269:ARG:NH1	2.32	0.45
1:A:410:ARG:HB2	1:A:454:ALA:HA	1.99	0.44
1:A:206:GLY:O	1:A:210:VAL:HG23	2.17	0.44
1:B:328:LYS:HE2	1:B:328:LYS:HB3	1.82	0.44
1:A:546:LEU:HD12	1:A:862:VAL:HG21	1.98	0.44
1:A:564:ARG:HD2	1:A:568:GLU:OE2	2.16	0.44
1:A:119:LEU:HD12	1:A:119:LEU:HA	1.81	0.44
1:B:800:ARG:HG2	4:B:1235:HOH:O	2.16	0.44
1:A:622:LEU:HD12	1:A:696:GLU:OE2	2.17	0.44
1:B:662:ASN:ND2	1:B:677:ILE:HD13	2.33	0.44
1:A:633:LYS:NZ	1:A:633:LYS:HB2	2.32	0.44
1:A:233:ASP:O	1:A:237:ILE:HG12	2.17	0.44
1:B:719:CYS:O	1:B:723:PRO:HG2	2.18	0.44
1:B:389:ARG:NE	1:B:389:ARG:HA	2.32	0.44
1:B:547:ASP:O	1:B:551:VAL:HG23	2.18	0.43
1:B:259:ILE:O	1:B:259:ILE:HG13	2.16	0.43
1:B:484:ILE:HD13	1:B:494:ILE:HD12	2.00	0.43
1:A:559:LEU:HB3	1:A:578:ARG:HH21	1.84	0.43
1:B:384:ARG:NH1	4:B:1032:HOH:O	2.52	0.43
1:B:226:GLU:HG3	4:B:1076:HOH:O	2.19	0.43
1:B:255:ASP:O	1:B:260:LYS:HD3	2.19	0.43
1:A:348:TYR:CG	1:A:349:SER:HA	2.54	0.42
3:A:911:FGG:C04	3:A:911:FGG:C17	2.97	0.42
1:A:184:PHE:CE2	1:A:189:ARG:HG3	2.54	0.42
1:A:270:LEU:HD13	1:A:293:GLU:OE2	2.18	0.42
1:A:633:LYS:HZ3	1:A:633:LYS:HB2	1.84	0.42
1:B:244:LYS:HA	1:B:244:LYS:HD2	1.85	0.42
1:B:253:PRO:HB2	1:B:256:LEU:HG	2.01	0.42
1:A:775:ILE:N	1:A:775:ILE:HD12	2.35	0.42
1:B:537:ASN:HD21	1:B:539:LYS:HB2	1.85	0.41
1:B:270:LEU:O	1:B:294:VAL:HG11	2.20	0.41
1:A:164:ALA:O	1:A:168:VAL:HG23	2.20	0.41
1:A:225:ASN:HA	1:A:226:GLU:HA	1.68	0.41
1:B:479:ALA:HA	1:B:483:LYS:HG2	2.02	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:150:THR:O	1:B:151:ILE:C	2.59	0.41
1:B:504:MET:HE3	1:B:504:MET:HB2	1.92	0.41
1:A:491:PHE:CD2	1:A:491:PHE:C	2.94	0.41
1:A:159:PRO:HA	1:A:203:TRP:CZ2	2.55	0.41
1:A:225:ASN:ND2	1:A:225:ASN:H	2.06	0.41
1:B:484:ILE:CD1	1:B:494:ILE:HD12	2.51	0.41
1:B:554:LEU:O	1:B:558:GLU:HG3	2.20	0.41
1:A:172:GLN:OE1	1:A:209:GLN:HB3	2.19	0.41
1:B:722:GLN:HB2	1:B:723:PRO:HD3	2.02	0.41
3:B:911:FGG:H05A	3:B:911:FGG:H16	1.65	0.41
1:B:480:LEU:HD21	1:B:494:ILE:HB	2.02	0.41
1:B:427:PHE:O	1:B:434:PHE:HA	2.20	0.41
1:A:140:TYR:HB2	1:A:192:ASN:OD1	2.21	0.41
3:A:911:FGG:H12A	3:A:911:FGG:H20B	2.03	0.41
1:A:344:VAL:HA	1:A:345:PRO:HD3	1.88	0.41
1:B:123:ILE:HD12	1:B:543:LEU:HD12	2.02	0.41
1:B:556:GLN:NE2	4:B:886:HOH:O	2.54	0.41
1:B:565:TRP:CZ3	1:B:566:TRP:HE3	2.39	0.41
1:A:523:VAL:HG21	1:A:529:LEU:HB3	2.03	0.41
1:B:811:LYS:HA	1:B:812:PRO:HD3	1.94	0.41
1:B:201:SER:HB2	1:B:248:LEU:HD13	2.02	0.41
1:A:617:ASP:OD1	1:A:618:ILE:HG13	2.20	0.40
1:B:459:PHE:HB3	1:B:460:PRO:CD	2.52	0.40
1:B:809:TYR:HA	1:B:821:LYS:HE2	2.03	0.40
1:B:759:THR:OG1	1:B:760:LYS:HD3	2.20	0.40
1:A:178:TRP:CE3	1:A:193:THR:HA	2.56	0.40
1:A:310:PHE:CE2	1:A:320:VAL:HG11	2.56	0.40
1:A:730:LEU:HB3	1:B:389:ARG:HD2	2.03	0.40
1:A:399:VAL:HA	1:A:400:PRO:HD3	1.89	0.40
1:A:578:ARG:NE	4:A:1299:HOH:O	2.55	0.40
1:B:635:TRP:CE3	1:B:678:ARG:HG2	2.56	0.40
1:B:230:LEU:HD22	1:B:234:PHE:CD1	2.57	0.40
1:A:127:PHE:CG	1:A:374:GLU:HB3	2.57	0.40
1:B:254:TYR:HA	1:B:259:ILE:CD1	2.40	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	750/764 (98%)	712 (95%)	35 (5%)	3 (0%)	43	46
1	B	736/764 (96%)	701 (95%)	31 (4%)	4 (0%)	38	37
All	All	1486/1528 (97%)	1413 (95%)	66 (4%)	7 (0%)	38	37

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	173	LEU
1	B	151	ILE
1	A	204	LYS
1	B	225	ASN
1	B	224	LEU
1	A	577	THR
1	B	644	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	A	665/674 (99%)	637 (96%)	28 (4%)	40	46
1	B	656/674 (97%)	629 (96%)	27 (4%)	41	48
All	All	1321/1348 (98%)	1266 (96%)	55 (4%)	41	46

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	119	LEU
1	A	150	THR

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Mol	Chain	Res	Type
1	A	166	ASN
1	A	174	GLN
1	A	190	LEU
1	A	191	LEU
1	A	209	GLN
1	A	225	ASN
1	A	235	GLN
1	A	251	ASN
1	A	259	ILE
1	A	320	VAL
1	A	331	THR
1	A	373	GLN
1	A	509	LEU
1	A	543	LEU
1	A	546	LEU
1	A	557	GLU
1	A	577	THR
1	A	580[A]	ARG
1	A	580[B]	ARG
1	A	581	VAL
1	A	611	LEU
1	A	622	LEU
1	A	625	LEU
1	A	629	THR
1	A	671	ARG
1	A	683	LEU
1	B	119	LEU
1	B	144	TRP
1	B	166	ASN
1	B	190	LEU
1	B	191	LEU
1	B	195	ASN
1	B	209	GLN
1	B	251	ASN
1	B	259	ILE
1	B	320	VAL
1	B	331	THR
1	B	393	TRP
1	B	492	LYS
1	B	509	LEU
1	B	513	SER
1	B	517	SER

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Mol	Chain	Res	Type
1	B	543	LEU
1	B	546	LEU
1	B	573	ASP
1	B	594	GLU
1	B	622	LEU
1	B	625	LEU
1	B	629	THR
1	B	671	ARG
1	B	683	LEU
1	B	766	LYS
1	B	788	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	166	ASN
1	A	174	GLN
1	A	183	HIS
1	A	195	ASN
1	A	207	HIS
1	A	212	GLN
1	A	220	ASN
1	A	225	ASN
1	A	235	GLN
1	A	286	ASN
1	A	334	ASN
1	A	373	GLN
1	A	403	ASN
1	A	450	ASN
1	A	537	ASN
1	A	556	GLN
1	A	579	HIS
1	A	609	GLN
1	A	662	ASN
1	A	669	GLN
1	A	826	ASN
1	B	166	ASN
1	B	170	ASN
1	B	174	GLN
1	B	183	HIS
1	B	195	ASN
1	B	212	GLN

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Mol	Chain	Res	Type
1	B	220	ASN
1	B	251	ASN
1	B	286	ASN
1	B	298	ASN
1	B	334	ASN
1	B	403	ASN
1	B	450	ASN
1	B	537	ASN
1	B	556	GLN
1	B	609	GLN
1	B	662	ASN
1	B	669	GLN
1	B	826	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 8 ligands modelled in this entry, 6 are monoatomic - leaving 2 for Mogul analysis.

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	FGG	A	911	2	29,29,29	1.24	2 (6%)	39,39,39	1.47	7 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	FGG	B	911	2	29,29,29	1.30	3 (10%)	39,39,39	1.64	7 (17%)

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	FGG	A	911	2	-	0/29/34/34	0/0/0/0
3	FGG	B	911	2	-	0/29/34/34	0/0/0/0

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	911	FGG	O01-C01	-3.88	1.40	1.43
3	A	911	FGG	F01-C02	-3.87	1.32	1.37
3	B	911	FGG	F01-C02	-3.66	1.32	1.37
3	A	911	FGG	O01-C01	-3.63	1.40	1.43
3	B	911	FGG	P04-O08	-2.08	1.47	1.54

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	911	FGG	C18-C11-C12	4.05	121.55	115.39
3	B	911	FGG	F01-C02-C01	3.97	118.92	111.42
3	A	911	FGG	F01-C02-C01	3.76	118.52	111.42
3	B	911	FGG	C09-C10-C11	-3.66	119.91	127.80
3	A	911	FGG	C18-C11-C12	3.19	120.23	115.39
3	A	911	FGG	C09-C10-C11	-3.00	121.33	127.80
3	B	911	FGG	C05-C06-C07	-2.58	122.24	127.80
3	B	911	FGG	P02-O03-P04	-2.49	124.38	131.68
3	B	911	FGG	O01-C01-C02	-2.44	106.63	110.77
3	B	911	FGG	C18-C11-C10	-2.42	118.72	123.52
3	A	911	FGG	C05-C06-C07	-2.36	122.72	127.80
3	A	911	FGG	P02-O03-P04	-2.27	125.02	131.68
3	A	911	FGG	C17-C07-C08	2.20	118.73	115.39
3	A	911	FGG	C19-C15-C20	2.18	120.17	114.62

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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

### 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	A	750/764 (98%)	-0.20	28 (3%)	39 44	9, 19, 64, 128	0
1	B	738/764 (96%)	-0.33	8 (1%)	77 81	9, 18, 48, 113	0
All	All	1488/1528 (97%)	-0.26	36 (2%)	56 61	9, 18, 57, 128	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	153	SER	6.4
1	B	154	ASP	4.5
1	B	152	SER	4.4
1	A	152	SER	4.3
1	A	258	PHE	4.0
1	A	227	GLU	3.9
1	A	264	THR	3.8
1	A	259	ILE	3.7
1	A	577	THR	3.6
1	B	227	GLU	3.5
1	B	155	GLY	3.5
1	A	225	ASN	3.5
1	A	273	VAL	3.4
1	A	257	PRO	3.2
1	A	226	GLU	3.1
1	A	262	LEU	3.0
1	A	579	HIS	3.0
1	A	260	LYS	2.9
1	B	264	THR	2.9
1	A	224	LEU	2.8
1	A	268	ALA	2.8
1	A	256	LEU	2.8
1	A	154	ASP	2.7
1	A	153	SER	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	265	THR	2.5
1	A	255	ASP	2.4
1	B	431	ASN	2.4
1	A	576	PHE	2.4
1	A	261	TYR	2.3
1	A	228	ASP	2.3
1	A	578	ARG	2.3
1	A	263	SER	2.3
1	A	272	ASP	2.1
1	A	573	ASP	2.1
1	A	155	GLY	2.0
1	B	575	ASN	2.0

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

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	MG	A	901	1/1	0.14	3.75	13,13,13,13	0
3	FGG	B	911	30/30	0.14	1.26	14,29,38,40	0
3	FGG	A	911	30/30	0.13	0.70	13,28,39,45	0
2	MG	B	901	1/1	0.09	0.41	12,12,12,12	0
2	MG	A	903	1/1	0.07	-1.07	13,13,13,13	0
2	MG	A	902	1/1	0.08	-1.83	16,16,16,16	0
2	MG	B	903	1/1	0.07	-3.01	18,18,18,18	0
2	MG	B	902	1/1	0.07	-6.17	12,12,12,12	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.