



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:46 AM GMT

PDB ID : 2E8V
Title : S. cerevisiae geranylgeranyl pyrophosphate synthase in complex with product GGPP (P21)
Authors : Chen, C.K.-M.; Guo, R.T.; Ko, T.P.; Jeng, W.Y.; Chang, T.H.; Liang, P.H.; Oldfield, E.; Wang, A.H.-J.
Deposited on : 2007-01-23
Resolution : 1.80 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
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

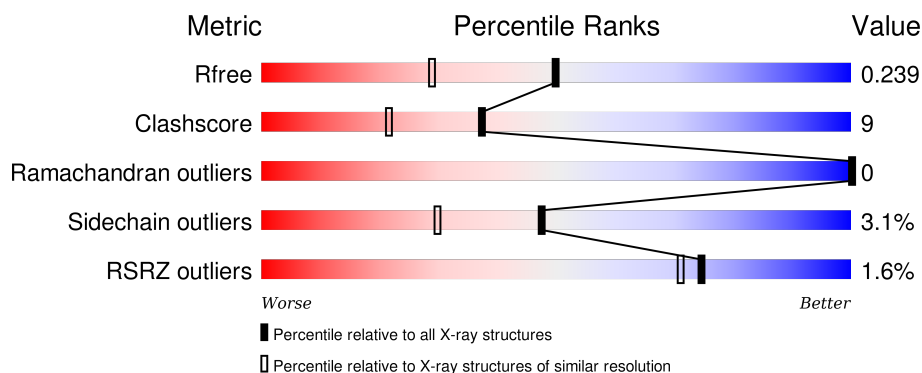
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

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 ≥ 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	340	<div> <div></div> <div>71% 14% • 14%</div> </div>
1	B	340	<div> <div></div> <div>68% 15% • 15%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5761 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.

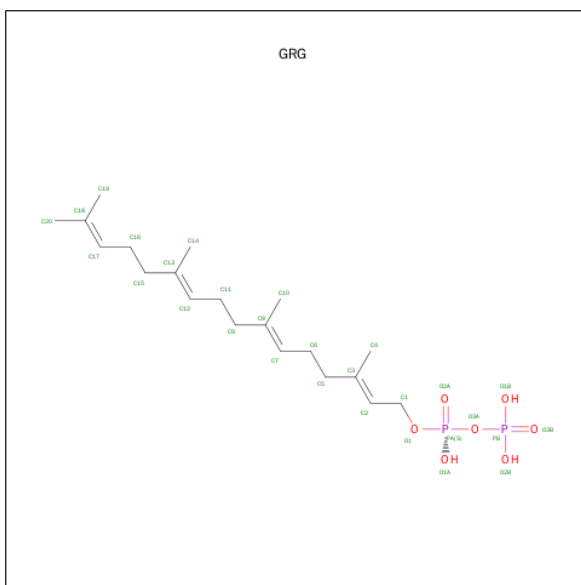
- Molecule 1 is a protein called Geranylgeranyl pyrophosphate synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	292	Total	C	N	O	S	0	0	0
			2378	1528	402	439	9			
1	B	288	Total	C	N	O	S	0	0	0
			2346	1508	397	432	9			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	CLONING ARTIFACT	UNP Q12051
A	2	THR	-	CLONING ARTIFACT	UNP Q12051
A	3	LYS	-	CLONING ARTIFACT	UNP Q12051
A	4	ASN	-	CLONING ARTIFACT	UNP Q12051
A	5	LYS	-	CLONING ARTIFACT	UNP Q12051
B	1	MET	-	CLONING ARTIFACT	UNP Q12051
B	2	THR	-	CLONING ARTIFACT	UNP Q12051
B	3	LYS	-	CLONING ARTIFACT	UNP Q12051
B	4	ASN	-	CLONING ARTIFACT	UNP Q12051
B	5	LYS	-	CLONING ARTIFACT	UNP Q12051

- Molecule 2 is GERANYLGERANYL DIPHOSPHATE (three-letter code: GRG) (formula: $C_{20}H_{36}O_7P_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 29	C 20	O 7	P 2	0	0
2	B	1	Total 29	C 20	O 7	P 2	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	532	Total O 532 532	0	0
3	B	447	Total O 447 447	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.35Å 48.82Å 93.23Å 90.00° 111.59° 90.00°	Depositor
Resolution (Å)	24.70 – 1.80 24.74 – 1.80	Depositor EDS
% Data completeness (in resolution range)	90.5 (24.70-1.80) 90.3 (24.74-1.80)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.49 (at 1.79Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.198 , 0.242 0.197 , 0.239	Depositor DCC
R_{free} test set	2945 reflections (5.06%)	DCC
Wilson B-factor (Å ²)	23.6	Xtriage
Anisotropy	0.388	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 69.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 58340 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5761	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.26% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: GRG

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	0.83	0/2419	0.82	1/3273 (0.0%)
1	B	0.80	0/2388	0.87	3/3235 (0.1%)
All	All	0.82	0/4807	0.85	4/6508 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	269	ARG	NE-CZ-NH2	-11.13	114.73	120.30
1	B	269	ARG	NE-CZ-NH1	10.21	125.41	120.30
1	B	154	ARG	NE-CZ-NH2	-6.61	116.99	120.30
1	A	36	LEU	CA-CB-CG	5.51	127.96	115.30

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	A	2378	0	2440	47	0
1	B	2346	0	2400	47	0
2	A	29	0	33	1	0
2	B	29	0	33	0	0
3	A	532	0	0	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	447	0	0	5	0
All	All	5761	0	4906	90	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

All (90) 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:B:284:GLU:HG3	1:B:290:LEU:HD12	1.37	1.04
1:B:52:ASN:HD21	1:B:57:LEU:H	1.09	1.01
1:A:116:MET:SD	3:B:1350:HOH:O	2.36	0.82
1:A:232:GLU:HG3	3:A:1364:HOH:O	1.80	0.81
1:A:6:MET:O	1:A:9:LYS:HG2	1.80	0.80
1:A:205:LEU:HD23	1:A:304:ILE:HD11	1.63	0.80
1:A:116:MET:HE3	1:B:137:ASN:CG	2.03	0.78
1:A:174:LYS:HE2	3:A:1348:HOH:O	1.84	0.76
1:B:85:ASN:ND2	1:B:96:HIS:HE1	1.83	0.76
1:A:235:THR:O	1:A:269:ARG:HD3	1.90	0.72
1:B:154:ARG:HD3	3:B:1343:HOH:O	1.91	0.71
1:B:85:ASN:HD22	1:B:96:HIS:HE1	1.37	0.71
1:A:236:GLU:OE2	1:A:238:LYS:HE3	1.93	0.69
1:A:174:LYS:CE	3:A:1348:HOH:O	2.39	0.68
1:A:154:ARG:HD3	1:A:236:GLU:O	1.93	0.68
1:B:269:ARG:HH11	1:B:269:ARG:HG2	1.58	0.68
1:B:147:GLN:HA	1:B:147:GLN:HE21	1.58	0.68
1:A:205:LEU:CD2	1:A:304:ILE:HD11	2.22	0.68
1:B:52:ASN:ND2	1:B:57:LEU:H	1.88	0.67
1:B:236:GLU:OE1	1:B:238:LYS:HE3	1.95	0.66
1:A:269:ARG:NH2	3:A:1093:HOH:O	2.27	0.66
1:A:113:PHE:HD2	1:A:116:MET:HE1	1.63	0.64
1:A:246:ALA:HB2	1:A:283:LEU:HD22	1.80	0.63
1:B:258:GLN:NE2	1:B:286:ASP:OD2	2.32	0.61
1:A:60:ASP:HB2	3:A:550:HOH:O	1.99	0.61
1:A:276:LYS:NZ	3:A:789:HOH:O	2.29	0.61
1:B:161:ILE:HD12	1:B:247:LEU:HD13	1.81	0.61
1:B:126:GLU:N	1:B:127:PRO:HD2	2.16	0.59
1:B:125:LYS:C	1:B:127:PRO:HD2	2.23	0.58
1:B:230:PHE:HB3	1:B:232:GLU:HG2	1.86	0.58
1:B:52:ASN:HD21	1:B:57:LEU:N	1.92	0.57
1:A:206:LEU:C	1:A:206:LEU:HD23	2.25	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:88:LEU:HD12	1:B:230:PHE:HE1	1.70	0.56
1:B:85:ASN:HD22	1:B:96:HIS:CE1	2.20	0.56
1:B:180:ARG:HH11	1:B:204:ASN:HD21	1.53	0.56
1:A:130:HIS:HE1	3:A:1323:HOH:O	1.87	0.56
1:A:130:HIS:CE1	3:A:1323:HOH:O	2.59	0.55
1:A:220:LYS:NZ	3:A:1292:HOH:O	2.26	0.55
1:A:116:MET:HE3	1:B:137:ASN:CB	2.37	0.54
1:A:144:HIS:HE1	3:A:1321:HOH:O	1.91	0.53
1:B:241:PHE:HB3	1:B:242:PRO:HD3	1.92	0.52
1:B:150:ASP:OD2	1:B:174:LYS:NZ	2.42	0.52
1:A:116:MET:CE	1:B:137:ASN:CG	2.76	0.52
1:A:12:GLU:CD	3:A:1367:HOH:O	2.47	0.52
1:A:126:GLU:N	1:A:127:PRO:CD	2.73	0.51
1:A:281:GLN:HA	1:A:281:GLN:NE2	2.25	0.51
1:B:277:LEU:O	1:B:281:GLN:HG3	2.12	0.50
1:A:305:LYS:NZ	3:A:1020:HOH:O	2.45	0.49
1:A:140:LEU:O	1:A:144:HIS:HD2	1.95	0.49
1:A:30:LYS:HB2	1:A:31:PRO:HD3	1.94	0.49
1:A:61:GLN:OE1	1:A:189:LEU:HD22	2.13	0.49
1:B:88:LEU:HD12	1:B:230:PHE:CE1	2.48	0.48
1:B:180:ARG:HH11	1:B:204:ASN:ND2	2.11	0.48
1:A:241:PHE:HB3	1:A:242:PRO:HD3	1.95	0.48
1:A:304:ILE:N	1:A:304:ILE:HD12	2.29	0.48
1:A:113:PHE:HD2	1:A:116:MET:CE	2.26	0.47
1:B:126:GLU:N	1:B:127:PRO:CD	2.78	0.47
1:B:83:GLU:O	1:B:154:ARG:NH2	2.45	0.47
1:B:129:TYR:CZ	1:B:133:ILE:HG13	2.50	0.47
1:A:6:MET:HA	1:A:9:LYS:HD3	1.97	0.47
1:A:113:PHE:CD2	1:A:116:MET:HE1	2.47	0.46
1:A:228:LYS:HD2	3:A:581:HOH:O	2.16	0.46
1:A:232:GLU:CG	3:A:1364:HOH:O	2.52	0.45
1:A:116:MET:CE	1:B:137:ASN:OD1	2.65	0.45
1:A:125:LYS:C	1:A:127:PRO:HD2	2.35	0.45
1:B:171:VAL:HG22	1:B:211:GLN:HG2	1.99	0.45
1:B:20:TRP:CH2	1:B:24:ASN:HB3	2.52	0.45
1:A:91:GLY:HA2	1:A:227:GLU:HB3	1.99	0.45
1:B:180:ARG:HD3	1:B:204:ASN:HD21	1.83	0.44
1:A:147:GLN:OE1	1:A:174:LYS:HD2	2.18	0.44
1:A:176:GLY:HA3	3:A:1126:HOH:O	2.17	0.44
1:A:218:ASN:ND2	3:A:1340:HOH:O	2.48	0.44
1:B:39:PRO:HA	3:B:1007:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:253:LYS:HB2	1:B:255:GLN:HG3	2.01	0.43
1:B:125:LYS:NZ	3:B:1310:HOH:O	2.50	0.43
1:B:288:ASN:HD22	1:B:288:ASN:HA	1.59	0.43
1:A:117:GLN:N	1:B:137:ASN:HD21	2.17	0.43
1:A:246:ALA:CB	1:A:283:LEU:HD22	2.48	0.43
1:B:75:SER:OG	1:B:111:MET:HB2	2.19	0.43
1:B:127:PRO:HG2	1:B:128:LEU:H	1.85	0.42
1:A:19:VAL:HG12	1:A:20:TRP:N	2.34	0.42
1:B:37:LEU:HD23	3:B:933:HOH:O	2.20	0.41
1:A:125:LYS:C	1:A:127:PRO:CD	2.89	0.41
1:B:304:ILE:O	1:B:307:ASP:HB3	2.20	0.41
1:B:101:VAL:N	1:B:102:PRO:CD	2.84	0.41
1:B:147:GLN:HA	1:B:147:GLN:NE2	2.32	0.41
1:B:244:VAL:O	1:B:248:ASN:ND2	2.42	0.41
1:B:273:LYS:NZ	1:B:274:ASP:OD1	2.54	0.41
1:B:167:TYR:CZ	1:B:171:VAL:HG21	2.56	0.40
2:A:1501:GRG:HC62	2:A:1501:GRG:H101	1.87	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	286/340 (84%)	283 (99%)	3 (1%)	0	100	100
1	B	282/340 (83%)	281 (100%)	1 (0%)	0	100	100
All	All	568/680 (84%)	564 (99%)	4 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	272/317 (86%)	264 (97%)	8 (3%)	50	34
1	B	269/317 (85%)	260 (97%)	9 (3%)	45	27
All	All	541/634 (85%)	524 (97%)	17 (3%)	47	30

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	LEU
1	A	60	ASP
1	A	61	GLN
1	A	126	GLU
1	A	158	PRO
1	A	191	PRO
1	A	273	LYS
1	A	288	ASN
1	B	37	LEU
1	B	41	LYS
1	B	85	ASN
1	B	147	GLN
1	B	149	LEU
1	B	158	PRO
1	B	269	ARG
1	B	283	LEU
1	B	288	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	ASN
1	A	85	ASN
1	A	109	ASN
1	A	130	HIS
1	A	137	ASN

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Mol	Chain	Res	Type
1	A	144	HIS
1	A	255	GLN
1	A	260	ASN
1	A	281	GLN
1	A	288	ASN
1	B	46	ASN
1	B	52	ASN
1	B	85	ASN
1	B	96	HIS
1	B	109	ASN
1	B	121	GLN
1	B	137	ASN
1	B	147	GLN
1	B	204	ASN
1	B	288	ASN
1	B	298	ASN
1	B	299	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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
2	GRG	A	1501	-	26,28,28	2.06	8 (30%)	33,37,37	1.46	5 (15%)
2	GRG	B	1502	-	26,28,28	2.06	8 (30%)	33,37,37	1.23	5 (15%)

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
2	GRG	A	1501	-	-	0/31/31/31	0/0/0/0
2	GRG	B	1502	-	-	0/31/31/31	0/0/0/0

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1501	GRG	C11-C12	-3.99	1.39	1.50
2	A	1501	GRG	C6-C7	-3.97	1.39	1.50
2	B	1502	GRG	C11-C12	-3.94	1.39	1.50
2	B	1502	GRG	C6-C7	-3.90	1.39	1.50
2	A	1501	GRG	C16-C17	-3.84	1.39	1.50
2	B	1502	GRG	C16-C17	-3.83	1.39	1.50
2	A	1501	GRG	C1-C2	-2.91	1.39	1.49
2	B	1502	GRG	C1-C2	-2.82	1.39	1.49
2	B	1502	GRG	C17-C18	2.34	1.39	1.32
2	A	1501	GRG	C17-C18	2.43	1.39	1.32
2	A	1501	GRG	C12-C13	3.13	1.39	1.33
2	A	1501	GRG	C7-C8	3.21	1.39	1.33
2	A	1501	GRG	C2-C3	3.28	1.39	1.33
2	B	1502	GRG	C12-C13	3.31	1.39	1.33
2	B	1502	GRG	C7-C8	3.31	1.39	1.33
2	B	1502	GRG	C2-C3	3.31	1.39	1.33

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1501	GRG	C9-C11-C12	-2.72	104.57	111.69
2	B	1502	GRG	PA-O3A-PB	-2.29	124.99	132.67
2	A	1501	GRG	PA-O3A-PB	-2.18	125.36	132.67
2	B	1502	GRG	C19-C18-C20	2.20	120.05	114.64
2	B	1502	GRG	C14-C13-C15	2.37	119.03	115.41
2	B	1502	GRG	C10-C8-C9	2.59	119.36	115.41
2	A	1501	GRG	C10-C8-C9	2.88	119.81	115.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1502	GRG	C4-C3-C5	3.15	120.22	115.41
2	A	1501	GRG	C4-C3-C5	3.35	120.53	115.41
2	A	1501	GRG	C14-C13-C15	3.67	121.01	115.41

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
2	A	1501	GRG	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 [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, 95th 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 ²)	Q<0.9
1	A	292/340 (85%)	-0.35	4 (1%) 78 74	16, 28, 52, 66	0
1	B	288/340 (84%)	-0.31	5 (1%) 73 69	16, 28, 49, 69	0
All	All	580/680 (85%)	-0.33	9 (1%) 74 71	16, 28, 50, 69	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	230	PHE	5.0
1	B	126	GLU	2.9
1	B	307	ASP	2.7
1	B	306	ASN	2.6
1	A	39	PRO	2.4
1	B	124	THR	2.3
1	A	126	GLU	2.2
1	A	5	LYS	2.0
1	A	56	ASN	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, 95th 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(Å ²)	Q<0.9
2	GRG	A	1501	29/29	0.95	0.11	0.39	26,37,40,41	0
2	GRG	B	1502	29/29	0.96	0.09	0.15	19,26,32,34	0

6.5 Other polymers

There are no such residues in this entry.