



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:21 AM UTC

PDB ID : 9C7I / pdb_00009c7i
Title : Crystal structure of caryolan-1-ol synthase from *S. griseus* with PEG molecule in the active site
Authors : Prem Kumar, R.; Matos, J.O.; Oprian, D.D.
Deposited on : 2024-06-10
Resolution : 2.33 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

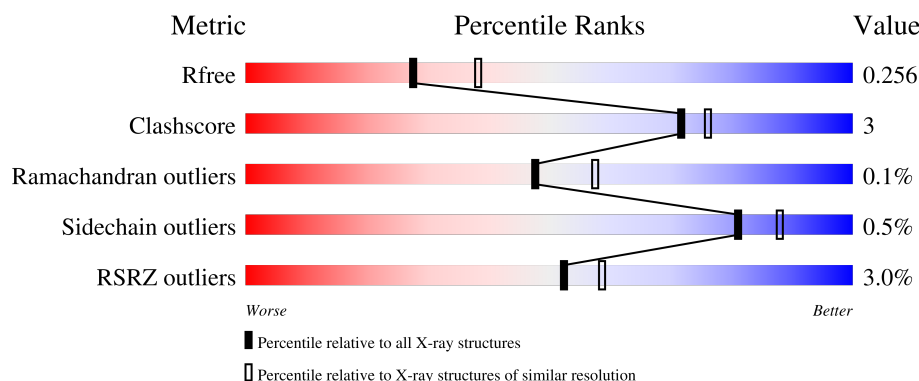
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.33 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}	180053	3031 (2.36-2.32)
Clashscore	190562	3127 (2.36-2.32)
Ramachandran outliers	187476	3095 (2.36-2.32)
Sidechain outliers	187428	3095 (2.36-2.32)
RSRZ outliers	180081	3033 (2.36-2.32)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	349	2% 78% 7% 15%
1	B	349	2% 77% 7% 16%
1	C	349	2% 78% 7% 15%
1	D	349	4% 80% 6% 15%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10212 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 (+)-caryolan-1-ol synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	297	Total	C	N	O	S	0	0	0
			2353	1470	442	428	13			
1	B	294	Total	C	N	O	S	0	0	0
			2337	1460	439	425	13			
1	C	295	Total	C	N	O	S	0	0	0
			2336	1460	437	426	13			
1	D	298	Total	C	N	O	S	0	0	0
			2355	1472	443	427	13			

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	initiating methionine	UNP B1W019
A	-12	GLY	-	expression tag	UNP B1W019
A	-11	HIS	-	expression tag	UNP B1W019
A	-10	HIS	-	expression tag	UNP B1W019
A	-9	HIS	-	expression tag	UNP B1W019
A	-8	HIS	-	expression tag	UNP B1W019
A	-7	HIS	-	expression tag	UNP B1W019
A	-6	HIS	-	expression tag	UNP B1W019
A	-5	GLU	-	expression tag	UNP B1W019
A	-4	ASN	-	expression tag	UNP B1W019
A	-3	LEU	-	expression tag	UNP B1W019
A	-2	TYR	-	expression tag	UNP B1W019
A	-1	PHE	-	expression tag	UNP B1W019
A	0	GLN	-	expression tag	UNP B1W019
A	1	GLY	-	expression tag	UNP B1W019
B	-13	MET	-	initiating methionine	UNP B1W019
B	-12	GLY	-	expression tag	UNP B1W019
B	-11	HIS	-	expression tag	UNP B1W019
B	-10	HIS	-	expression tag	UNP B1W019
B	-9	HIS	-	expression tag	UNP B1W019
B	-8	HIS	-	expression tag	UNP B1W019

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-7	HIS	-	expression tag	UNP B1W019
B	-6	HIS	-	expression tag	UNP B1W019
B	-5	GLU	-	expression tag	UNP B1W019
B	-4	ASN	-	expression tag	UNP B1W019
B	-3	LEU	-	expression tag	UNP B1W019
B	-2	TYR	-	expression tag	UNP B1W019
B	-1	PHE	-	expression tag	UNP B1W019
B	0	GLN	-	expression tag	UNP B1W019
B	1	GLY	-	expression tag	UNP B1W019
C	-13	MET	-	initiating methionine	UNP B1W019
C	-12	GLY	-	expression tag	UNP B1W019
C	-11	HIS	-	expression tag	UNP B1W019
C	-10	HIS	-	expression tag	UNP B1W019
C	-9	HIS	-	expression tag	UNP B1W019
C	-8	HIS	-	expression tag	UNP B1W019
C	-7	HIS	-	expression tag	UNP B1W019
C	-6	HIS	-	expression tag	UNP B1W019
C	-5	GLU	-	expression tag	UNP B1W019
C	-4	ASN	-	expression tag	UNP B1W019
C	-3	LEU	-	expression tag	UNP B1W019
C	-2	TYR	-	expression tag	UNP B1W019
C	-1	PHE	-	expression tag	UNP B1W019
C	0	GLN	-	expression tag	UNP B1W019
C	1	GLY	-	expression tag	UNP B1W019
D	-13	MET	-	initiating methionine	UNP B1W019
D	-12	GLY	-	expression tag	UNP B1W019
D	-11	HIS	-	expression tag	UNP B1W019
D	-10	HIS	-	expression tag	UNP B1W019
D	-9	HIS	-	expression tag	UNP B1W019
D	-8	HIS	-	expression tag	UNP B1W019
D	-7	HIS	-	expression tag	UNP B1W019
D	-6	HIS	-	expression tag	UNP B1W019
D	-5	GLU	-	expression tag	UNP B1W019
D	-4	ASN	-	expression tag	UNP B1W019
D	-3	LEU	-	expression tag	UNP B1W019
D	-2	TYR	-	expression tag	UNP B1W019
D	-1	PHE	-	expression tag	UNP B1W019
D	0	GLN	-	expression tag	UNP B1W019
D	1	GLY	-	expression tag	UNP B1W019

- Molecule 2 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: C₈H₁₈O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	8	5		
2	D	1	Total	C	O	0	0
			13	8	5		

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		
3	B	2	Total	Ca	0	0
			2	2		
3	C	2	Total	Ca	0	0
			2	2		
3	D	1	Total	Ca	0	0
			1	1		

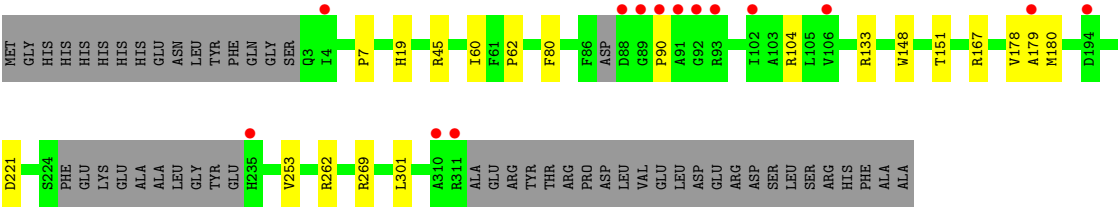
- Molecule 4 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: C₁₀H₂₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			16	10	6		
4	C	1	Total	C	O	0	0
			16	10	6		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	200	Total	O	0	0
			200	200		
5	B	191	Total	O	0	0
			191	191		
5	C	190	Total	O	0	0
			190	190		
5	D	186	Total	O	0	0
			186	186		



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	68.52Å 89.80Å 114.11Å 90.00° 90.82° 90.00°	Depositor
Resolution (Å)	70.57 – 2.33 70.57 – 2.33	Depositor EDS
% Data completeness (in resolution range)	100.0 (70.57-2.33) 100.0 (70.57-2.33)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.41 (at 2.32Å)	Xtriage
Refinement program	PHENIX 1.20_4444	Depositor
R, R_{free}	0.211 , 0.256 0.212 , 0.256	Depositor DCC
R_{free} test set	2862 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	26.0	Xtriage
Anisotropy	0.140	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 40.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.115 for h,-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	10212	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.06% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: PG4, CA, 1PE

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	0.07	0/2406	0.20	0/3263
1	B	0.07	0/2390	0.20	0/3241
1	C	0.07	0/2389	0.21	0/3241
1	D	0.08	0/2409	0.22	0/3268
All	All	0.07	0/9594	0.21	0/13013

There are no bond length outliers.

There are no bond angle outliers.

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	2353	0	2274	18	0
1	B	2337	0	2263	16	0
1	C	2336	0	2254	16	0
1	D	2355	0	2278	17	0
2	A	13	0	18	0	0
2	D	13	0	18	2	0
3	A	1	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	1	0	0	0	0
4	B	16	0	22	2	0
4	C	16	0	22	2	0
5	A	200	0	0	2	0
5	B	191	0	0	2	0
5	C	190	0	0	3	0
5	D	186	0	0	5	0
All	All	10212	0	9149	61	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 3.

All (61) 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:262:ARG:HH22	1:D:262:ARG:HH12	1.30	0.77
1:C:159:ARG:NH2	5:C:501:HOH:O	2.26	0.68
1:A:26:ARG:NH1	5:A:508:HOH:O	2.34	0.60
1:B:50:ARG:NH2	5:B:511:HOH:O	2.34	0.59
1:D:269:ARG:NH1	5:D:512:HOH:O	2.36	0.58
1:D:221:ASP:O	5:D:501:HOH:O	2.17	0.57
1:D:178:VAL:HG13	1:D:180:MET:H	1.70	0.56
1:B:262:ARG:NH1	1:C:255:GLU:OE2	2.38	0.56
1:C:148:TRP:O	1:C:151:THR:OG1	2.24	0.56
1:B:4:ILE:HD12	1:B:253:VAL:HG21	1.90	0.54
1:C:250:GLN:NE2	5:C:513:HOH:O	2.40	0.54
1:B:148:TRP:O	1:B:151:THR:OG1	2.25	0.53
1:D:45:ARG:NH2	5:D:518:HOH:O	2.41	0.53
1:B:220:ASN:HD21	4:B:401:1PE:H162	1.74	0.53
1:A:262:ARG:HH22	1:D:262:ARG:NH1	2.02	0.52
1:D:179:ALA:HB3	2:D:401:PG4:H62	1.90	0.51
1:A:178:VAL:HG13	1:A:180:MET:H	1.76	0.51
1:A:148:TRP:O	1:A:151:THR:OG1	2.28	0.51
1:C:155:GLU:OE1	5:C:501:HOH:O	2.20	0.50
1:B:178:VAL:HG13	1:B:180:MET:H	1.78	0.49
1:A:184:LEU:HD13	1:A:209:ARG:HG2	1.95	0.48
1:C:56:TRP:CE2	4:C:401:1PE:H242	2.48	0.47
1:A:51:THR:O	1:A:308:TYR:OH	2.22	0.47
1:D:80:PHE:HE1	2:D:401:PG4:H52	1.80	0.46
1:A:278:ILE:HA	1:A:283:ILE:HD12	1.97	0.46
1:A:95:PRO:HB3	1:A:157:VAL:HG22	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:172:LYS:HE3	1:B:172:LYS:HB3	1.62	0.46
1:B:148:TRP:CD1	1:B:179:ALA:HB2	2.50	0.46
1:C:284:ASP:HA	1:C:288:ARG:HD2	1.98	0.45
1:D:133:ARG:NH2	5:D:517:HOH:O	2.40	0.45
1:A:164:VAL:HB	1:A:243:ARG:HH12	1.82	0.44
1:D:60:ILE:HG23	1:D:301:LEU:HG	1.99	0.44
1:A:262:ARG:NH2	1:D:262:ARG:HH12	2.07	0.44
1:C:7:PRO:HD3	1:C:253:VAL:HG12	2.00	0.44
1:C:199:ALA:HB1	1:C:291:LEU:HD21	1.98	0.44
1:C:208:LEU:HG	1:C:298:TYR:HD2	1.82	0.44
1:C:147:ALA:O	1:C:151:THR:HG23	2.18	0.44
1:B:57:ILE:HG21	1:B:73:CYS:HB2	2.00	0.44
1:A:45:ARG:NH1	5:A:518:HOH:O	2.50	0.43
1:A:52:ARG:HE	1:A:55:LEU:HD22	1.84	0.43
1:D:133:ARG:NH1	5:D:514:HOH:O	2.37	0.43
1:C:115:ASN:N	1:C:119:GLU:OE1	2.44	0.43
1:C:308:TYR:HE2	4:C:401:1PE:H251	1.83	0.42
1:B:99:GLU:HB2	1:B:153:TYR:CE1	2.54	0.42
1:B:258:ILE:HG23	1:C:258:ILE:HG23	2.01	0.42
1:D:104:ARG:HD2	1:D:104:ARG:HA	1.89	0.42
1:A:147:ALA:O	1:A:151:THR:HG23	2.20	0.42
1:D:19:HIS:HB2	1:D:62:PRO:HA	2.02	0.42
1:A:148:TRP:CD1	1:A:179:ALA:HB2	2.54	0.42
1:B:56:TRP:CE2	4:B:401:1PE:H251	2.54	0.42
1:C:104:ARG:NH2	1:C:113:ALA:O	2.53	0.41
1:D:7:PRO:HD3	1:D:253:VAL:HG12	2.01	0.41
1:D:148:TRP:O	1:D:151:THR:OG1	2.36	0.41
1:B:52:ARG:HE	1:B:55:LEU:HD22	1.85	0.41
1:B:158:GLU:HB3	5:B:504:HOH:O	2.20	0.41
1:A:4:ILE:HG13	1:A:249:LEU:HD23	2.02	0.41
1:B:170:PHE:O	1:B:174:ARG:N	2.50	0.41
1:C:148:TRP:CD1	1:C:179:ALA:HB2	2.56	0.40
1:A:247:SER:HA	1:D:269:ARG:HH21	1.86	0.40
1:B:38:LEU:HD23	1:B:117:PRO:HB2	2.03	0.40
1:A:247:SER:HB2	1:A:251:GLU:HB2	2.04	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	291/349 (83%)	287 (99%)	4 (1%)	0	100	100
1	B	288/349 (82%)	282 (98%)	6 (2%)	0	100	100
1	C	289/349 (83%)	284 (98%)	5 (2%)	0	100	100
1	D	292/349 (84%)	288 (99%)	3 (1%)	1 (0%)	36	41
All	All	1160/1396 (83%)	1141 (98%)	18 (2%)	1 (0%)	48	57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	90	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	235/280 (84%)	234 (100%)	1 (0%)	84	89
1	B	235/280 (84%)	233 (99%)	2 (1%)	70	80
1	C	234/280 (84%)	233 (100%)	1 (0%)	84	89
1	D	235/280 (84%)	234 (100%)	1 (0%)	84	89
All	All	939/1120 (84%)	934 (100%)	5 (0%)	81	88

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	126	ARG
1	B	25	THR
1	B	299	ARG
1	C	220	ASN
1	D	167	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	GLN
1	B	220	ASN
1	B	309	HIS
1	C	163	GLN
1	C	181	GLN
1	D	163	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 6 are monoatomic - leaving 4 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
4	1PE	C	401	-	15,15,15	0.11	0	14,14,14	0.10	0
2	PG4	A	401	-	12,12,12	0.10	0	11,11,11	0.64	0
4	1PE	B	401	-	15,15,15	0.11	0	14,14,14	0.16	0
2	PG4	D	401	-	12,12,12	0.11	0	11,11,11	0.65	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
4	1PE	C	401	-	-	5/13/13/13	-
2	PG4	A	401	-	-	4/10/10/10	-
4	1PE	B	401	-	-	7/13/13/13	-
2	PG4	D	401	-	-	7/10/10/10	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	401	PG4	O2-C3-C4-O3
4	C	401	1PE	OH6-C15-C25-OH5
2	D	401	PG4	O4-C7-C8-O5
2	A	401	PG4	O3-C5-C6-O4
2	D	401	PG4	O3-C5-C6-O4
2	D	401	PG4	O1-C1-C2-O2
4	B	401	1PE	OH6-C15-C25-OH5
4	B	401	1PE	OH7-C16-C26-OH6
4	C	401	1PE	OH5-C14-C24-OH4
2	A	401	PG4	O2-C3-C4-O3
4	B	401	1PE	OH4-C13-C23-OH3
4	C	401	1PE	OH7-C16-C26-OH6
4	C	401	1PE	C25-C15-OH6-C26
4	B	401	1PE	C13-C23-OH3-C22
4	B	401	1PE	C16-C26-OH6-C15
2	D	401	PG4	C8-C7-O4-C6
4	B	401	1PE	OH5-C14-C24-OH4
4	B	401	1PE	C12-C22-OH3-C23

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Mol	Chain	Res	Type	Atoms
2	A	401	PG4	C1-C2-O2-C3
2	D	401	PG4	C3-C4-O3-C5
2	A	401	PG4	O4-C7-C8-O5
2	D	401	PG4	C5-C6-O4-C7
4	C	401	1PE	C23-C13-OH4-C24

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	401	1PE	2	0
4	B	401	1PE	2	0
2	D	401	PG4	2	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, 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	297/349 (85%)	0.37	7 (2%) 59 65	16, 27, 46, 64	0
1	B	294/349 (84%)	0.35	7 (2%) 59 65	20, 28, 44, 63	0
1	C	295/349 (84%)	0.27	7 (2%) 59 65	19, 27, 43, 63	0
1	D	298/349 (85%)	0.43	14 (4%) 36 42	18, 28, 49, 71	0
All	All	1184/1396 (84%)	0.36	35 (2%) 52 59	16, 28, 46, 71	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	91	ALA	5.2
1	B	92	GLY	4.9
1	C	87	ASP	3.4
1	C	92	GLY	3.2
1	D	90	PRO	3.1
1	A	225	PHE	2.9
1	D	91	ALA	2.9
1	D	235	HIS	2.8
1	D	311	ARG	2.7
1	C	235	HIS	2.6
1	B	96	LEU	2.6
1	C	310	ALA	2.5
1	B	235	HIS	2.4
1	D	106	VAL	2.4
1	D	4	ILE	2.3
1	D	89	GLY	2.3
1	C	88	ASP	2.3
1	A	92	GLY	2.3
1	A	93	ARG	2.2
1	C	97	MET	2.2
1	D	93	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	310	ALA	2.2
1	B	88	ASP	2.2
1	B	115	ASN	2.2
1	D	179	ALA	2.2
1	D	88	ASP	2.2
1	B	39	ASP	2.2
1	D	92	GLY	2.2
1	C	93	ARG	2.1
1	D	102	ILE	2.1
1	B	4	ILE	2.1
1	A	115	ASN	2.0
1	D	194	ASP	2.0
1	A	311	ARG	2.0
1	A	249	LEU	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 oligosaccharides 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. 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	B-factors(Å ²)	Q<0.9
2	PG4	D	401	13/13	0.78	0.17	32,44,48,50	0
2	PG4	A	401	13/13	0.84	0.13	27,35,46,46	0
4	1PE	B	401	16/16	0.84	0.13	26,36,45,46	0
4	1PE	C	401	16/16	0.90	0.11	24,32,46,50	0
3	CA	B	403	1/1	0.97	0.04	33,33,33,33	0
3	CA	C	403	1/1	0.97	0.04	35,35,35,35	0
3	CA	A	402	1/1	0.98	0.05	32,32,32,32	0
3	CA	B	402	1/1	0.98	0.03	33,33,33,33	0
3	CA	C	402	1/1	0.99	0.03	23,23,23,23	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	CA	D	402	1/1	0.99	0.03	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.