



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 11:56 am GMT

PDB ID : 4J0D  
Title : tannin acyl hydrolase from Lactobacillus plantarum (Cadmium)  
Authors : Ren, B.; Wu, M.; Wang, Q.; Peng, X.; Wen, H.; Chen, Q.; McKinsty, W.J.  
Deposited on : 2013-01-30  
Resolution : 1.60 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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

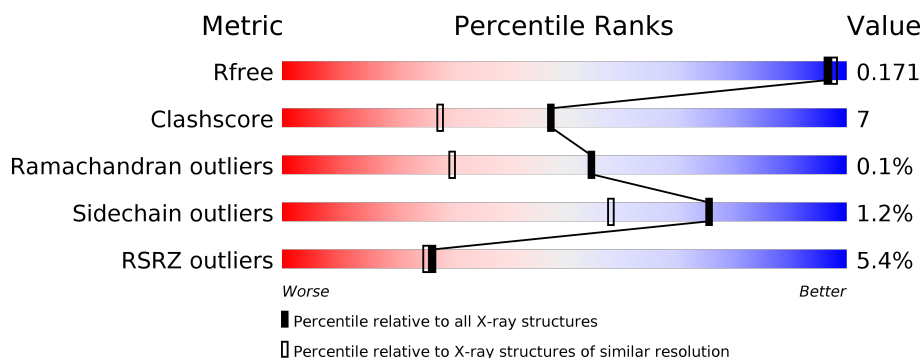
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.60 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))
$R_{free}$	100719	2696 (1.60-1.60)
Clashscore	112137	2967 (1.60-1.60)
Ramachandran outliers	110173	2887 (1.60-1.60)
Sidechain outliers	110143	2886 (1.60-1.60)
RSRZ outliers	101464	2714 (1.60-1.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	491	<div> <div>6%</div> <div> <div></div> <div>85%</div> <div>10%</div> <div>.</div> </div> </div>
1	B	491	<div> <div>4%</div> <div> <div></div> <div>86%</div> <div>9%</div> <div>.</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	CD	A	502	-	-	-	X
2	CD	B	502	-	-	-	X
3	PEG	A	503	-	-	X	X
3	PEG	A	505	-	-	X	-
3	PEG	B	505	-	-	-	X
3	PEG	B	506	-	-	-	X
3	PEG	B	507	-	-	-	X
3	PEG	B	508	-	-	X	X
3	PEG	B	510	-	-	-	X
3	PEG	B	511	-	-	X	X
5	PG4	B	503	-	-	X	-

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8177 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 Tannase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	470	Total	C	N	O	S	0	10	0
			3638	2307	630	690	11			
1	B	470	Total	C	N	O	S	0	9	0
			3635	2308	628	687	12			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	EXPRESSION TAG	UNP B3Y018
A	-20	HIS	-	EXPRESSION TAG	UNP B3Y018
A	-19	HIS	-	EXPRESSION TAG	UNP B3Y018
A	-18	HIS	-	EXPRESSION TAG	UNP B3Y018
A	-17	HIS	-	EXPRESSION TAG	UNP B3Y018
A	-16	HIS	-	EXPRESSION TAG	UNP B3Y018
A	-15	HIS	-	EXPRESSION TAG	UNP B3Y018
A	-14	SER	-	EXPRESSION TAG	UNP B3Y018
A	-13	SER	-	EXPRESSION TAG	UNP B3Y018
A	-12	GLY	-	EXPRESSION TAG	UNP B3Y018
A	-11	VAL	-	EXPRESSION TAG	UNP B3Y018
A	-10	ASP	-	EXPRESSION TAG	UNP B3Y018
A	-9	LEU	-	EXPRESSION TAG	UNP B3Y018
A	-8	GLY	-	EXPRESSION TAG	UNP B3Y018
A	-7	THR	-	EXPRESSION TAG	UNP B3Y018
A	-6	GLU	-	EXPRESSION TAG	UNP B3Y018
A	-5	ASN	-	EXPRESSION TAG	UNP B3Y018
A	-4	LEU	-	EXPRESSION TAG	UNP B3Y018
A	-3	TYR	-	EXPRESSION TAG	UNP B3Y018
A	-2	PHE	-	EXPRESSION TAG	UNP B3Y018
A	-1	GLN	-	EXPRESSION TAG	UNP B3Y018
A	0	SER	-	EXPRESSION TAG	UNP B3Y018
B	-21	MET	-	EXPRESSION TAG	UNP B3Y018
B	-20	HIS	-	EXPRESSION TAG	UNP B3Y018
B	-19	HIS	-	EXPRESSION TAG	UNP B3Y018

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	HIS	-	EXPRESSION TAG	UNP B3Y018
B	-17	HIS	-	EXPRESSION TAG	UNP B3Y018
B	-16	HIS	-	EXPRESSION TAG	UNP B3Y018
B	-15	HIS	-	EXPRESSION TAG	UNP B3Y018
B	-14	SER	-	EXPRESSION TAG	UNP B3Y018
B	-13	SER	-	EXPRESSION TAG	UNP B3Y018
B	-12	GLY	-	EXPRESSION TAG	UNP B3Y018
B	-11	VAL	-	EXPRESSION TAG	UNP B3Y018
B	-10	ASP	-	EXPRESSION TAG	UNP B3Y018
B	-9	LEU	-	EXPRESSION TAG	UNP B3Y018
B	-8	GLY	-	EXPRESSION TAG	UNP B3Y018
B	-7	THR	-	EXPRESSION TAG	UNP B3Y018
B	-6	GLU	-	EXPRESSION TAG	UNP B3Y018
B	-5	ASN	-	EXPRESSION TAG	UNP B3Y018
B	-4	LEU	-	EXPRESSION TAG	UNP B3Y018
B	-3	TYR	-	EXPRESSION TAG	UNP B3Y018
B	-2	PHE	-	EXPRESSION TAG	UNP B3Y018
B	-1	GLN	-	EXPRESSION TAG	UNP B3Y018
B	0	SER	-	EXPRESSION TAG	UNP B3Y018

- Molecule 2 is CADMIUM ION (three-letter code: CD) (formula: Cd).

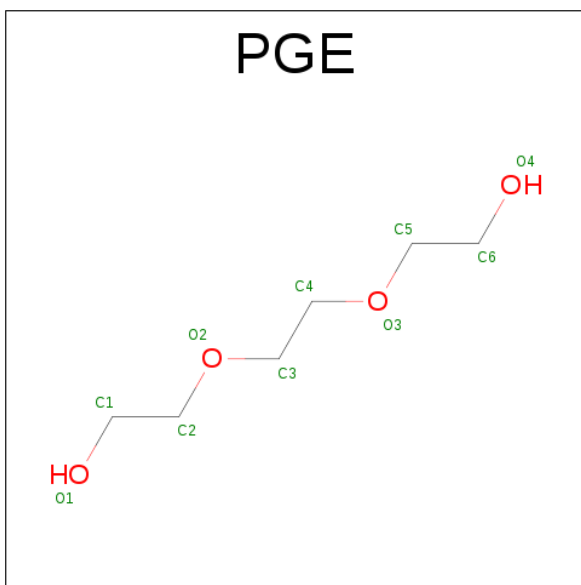
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	2	Total Cd 2 2	0	0
2	A	2	Total Cd 2 2	0	0

- Molecule 3 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C<sub>4</sub>H<sub>10</sub>O<sub>3</sub>).



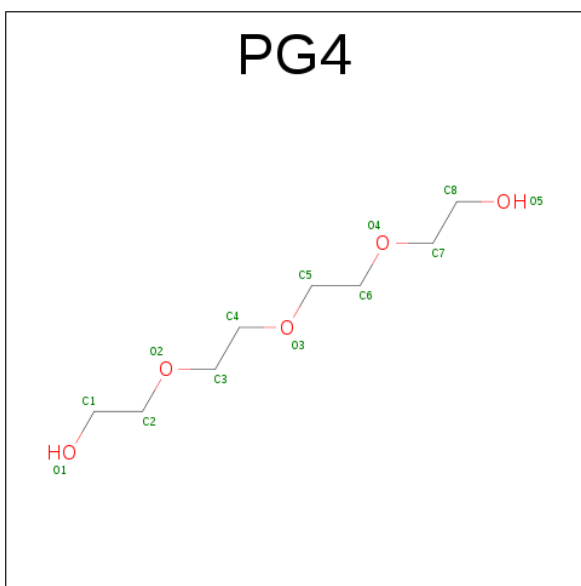
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	4	3		
3	A	1	Total	C	O	0	0
			7	4	3		
3	A	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 4 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C<sub>6</sub>H<sub>14</sub>O<sub>4</sub>).



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

- Molecule 5 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula:  $C_8H_{18}O_5$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			13	8	5		
5	B	1	Total	C	O	0	0
			13	8	5		

- Molecule 6 is water.

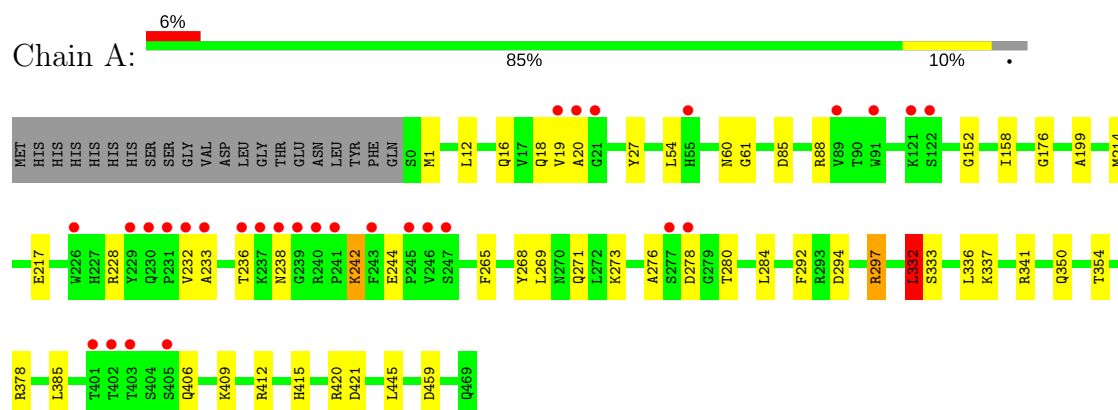
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	373	Total 373	O 373	0	0
6	B	414	Total 414	O 414	0	0



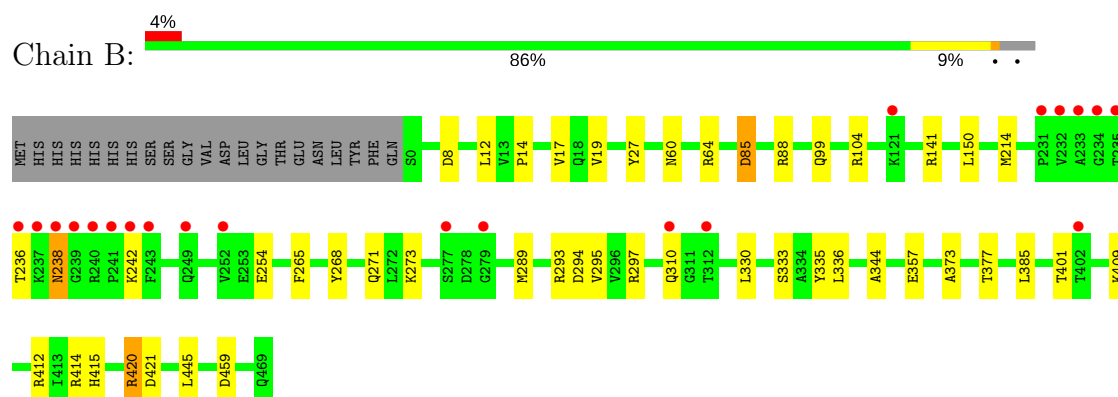
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

#### • Molecule 1: Tannase



#### • Molecule 1: Tannase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	46.94Å 62.92Å 84.07Å 70.58° 85.70° 78.96°	Depositor
Resolution (Å)	20.02 – 1.60 20.02 – 1.60	Depositor EDS
% Data completeness (in resolution range)	93.1 (20.02-1.60) 84.3 (20.02-1.60)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.64 (at 1.60Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.1_743)	Depositor
R, $R_{free}$	0.143 , 0.175 0.141 , 0.171	Depositor DCC
$R_{free}$ test set	1848 reflections (1.69%)	DCC
Wilson B-factor (Å <sup>2</sup> )	14.1	Xtriage
Anisotropy	0.551	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.41 , 69.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.012 for -h,-k,-k+l	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	8177	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.16% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PG4, PEG, PGE, CD

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	0.70	0/3743	0.79	5/5106 (0.1%)
1	B	0.78	1/3744 (0.0%)	0.84	7/5106 (0.1%)
All	All	0.74	1/7487 (0.0%)	0.82	12/10212 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	357	GLU	CG-CD	5.15	1.59	1.51

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	20	ALA	N-CA-C	-6.72	92.86	111.00
1	A	385	LEU	CA-CB-CG	6.29	129.76	115.30
1	B	104	ARG	NE-CZ-NH2	-6.17	117.21	120.30
1	B	85	ASP	CB-CG-OD1	5.98	123.68	118.30
1	B	150	LEU	CA-CB-CG	5.48	127.91	115.30
1	B	420	ARG	NE-CZ-NH1	5.40	123.00	120.30
1	B	141	ARG	NE-CZ-NH2	-5.22	117.69	120.30
1	A	297	ARG	NE-CZ-NH1	5.17	122.88	120.30
1	A	19	VAL	CB-CA-C	-5.12	101.68	111.40
1	A	332	LEU	CA-CB-CG	5.09	127.01	115.30
1	B	420	ARG	NE-CZ-NH2	-5.09	117.75	120.30
1	B	385	LEU	CA-CB-CG	5.08	126.99	115.30

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3638	0	3581	50	1
1	B	3635	0	3580	48	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	21	0	30	10	0
3	B	56	0	80	18	0
4	A	10	0	14	2	0
5	B	26	0	35	12	0
6	A	373	0	0	11	2
6	B	414	0	0	11	4
All	All	8177	0	7320	103	5

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 7.

All (103) 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:60:ASN:HB3	6:B:987:HOH:O	1.57	1.05
1:A:273:LYS:NZ	6:A:940:HOH:O	1.90	1.03
1:B:459:ASP:HB3	3:B:511:PEG:H41	1.34	1.01
1:A:409:LYS:NZ	6:A:905:HOH:O	1.92	1.00
1:A:294:ASP:OD1	1:A:297:ARG:NH2	1.98	0.96
1:B:294:ASP:OD1	1:B:297[A]:ARG:NH2	1.99	0.95
1:A:278:ASP:OD1	1:A:280[A]:THR:HG22	1.66	0.94
1:B:289[A]:MET:CE	6:B:981:HOH:O	2.21	0.89
5:B:504:PG4:H12	6:B:801:HOH:O	1.74	0.86
1:B:293:ARG:HE	5:B:503:PG4:H62	1.41	0.85
1:A:269[A]:LEU:HD21	1:A:292:PHE:CD2	2.13	0.84
5:B:504:PG4:O1	6:B:824:HOH:O	1.97	0.83
1:A:459:ASP:HB3	3:A:503:PEG:H21	1.61	0.82
1:A:61:GLY:H	3:A:505:PEG:H32	1.43	0.82
1:B:293:ARG:HH21	5:B:503:PG4:H41	1.42	0.82
1:B:336:LEU:HD12	5:B:503:PG4:H32	1.62	0.82
1:A:412:ARG:HH21	3:A:503:PEG:H21	1.47	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:64:ARG:HD2	3:B:509:PEG:H21	1.65	0.78
1:B:409:LYS:HB3	3:B:508:PEG:H41	1.67	0.77
1:A:61:GLY:H	3:A:505:PEG:C3	1.99	0.74
1:A:350[A]:GLN:OE1	1:A:354:THR:OG1	2.06	0.74
6:A:749:HOH:O	1:B:271:GLN:HG3	1.90	0.71
1:A:269[A]:LEU:HD21	1:A:292:PHE:CG	2.25	0.71
1:B:293:ARG:HH21	5:B:503:PG4:C4	2.04	0.70
1:B:238:ASN:OD1	1:B:238:ASN:N	2.25	0.70
1:B:236:THR:HG1	1:B:238:ASN:CG	1.95	0.69
1:B:293:ARG:NE	5:B:503:PG4:H62	2.07	0.69
1:A:61:GLY:N	3:A:505:PEG:H32	2.11	0.66
1:A:27:TYR:CG	1:A:54[B]:LEU:HD12	2.30	0.66
1:A:269[A]:LEU:HD23	1:A:284:LEU:CD1	2.26	0.66
1:A:332:LEU:HD13	4:A:504:PGE:H3	1.77	0.65
1:A:271:GLN:HG3	6:A:949:HOH:O	1.96	0.65
1:A:459:ASP:HB3	3:A:503:PEG:C2	2.28	0.64
1:B:409:LYS:NZ	3:B:508:PEG:H21	2.12	0.64
1:A:61:GLY:HA2	3:A:505:PEG:H12	1.80	0.64
1:A:214[B]:MET:HE3	1:A:265:PHE:HA	1.81	0.63
1:A:16:GLN:NE2	6:A:875:HOH:O	2.05	0.63
1:A:228:ARG:NH1	6:A:873:HOH:O	2.32	0.62
1:A:333:SER:O	1:A:337:LYS:HG3	1.99	0.62
1:B:273:LYS:NZ	6:B:973:HOH:O	2.07	0.60
1:A:27:TYR:CD2	1:A:54[B]:LEU:HD12	2.38	0.58
1:B:459:ASP:CB	3:B:511:PEG:H41	2.22	0.58
1:B:293:ARG:HE	5:B:503:PG4:C6	2.16	0.57
1:B:412:ARG:HH21	3:B:511:PEG:H42	1.68	0.57
1:B:214[B]:MET:HE3	1:B:265:PHE:HA	1.87	0.57
1:A:336:LEU:CD1	4:A:504:PGE:H32	2.35	0.57
1:A:158:ILE:HG12	1:A:199:ALA:HB3	1.86	0.56
1:B:214[B]:MET:HE1	1:B:268:TYR:CB	2.35	0.55
1:B:409:LYS:HZ1	3:B:508:PEG:H42	1.72	0.55
1:B:333:SER:OG	5:B:503:PG4:H31	2.06	0.55
1:A:214[B]:MET:HE1	1:A:268:TYR:CB	2.37	0.54
1:B:409:LYS:HZ2	3:B:508:PEG:H21	1.73	0.54
1:A:152:GLY:O	6:A:639:HOH:O	2.18	0.54
1:A:236:THR:HG22	1:A:238:ASN:H	1.73	0.54
1:B:64:ARG:HH11	3:B:509:PEG:H21	1.72	0.54
1:B:17:VAL:HG22	1:B:19:VAL:HG13	1.90	0.53
1:A:265:PHE:CE2	1:A:269[A]:LEU:HD22	2.44	0.53
1:B:373:ALA:O	1:B:377[B]:THR:HG23	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:293:ARG:NH2	5:B:503:PG4:H41	2.20	0.53
1:A:378:ARG:NH2	6:A:908:HOH:O	2.38	0.52
1:A:61:GLY:N	3:A:505:PEG:C3	2.69	0.51
1:A:214[A]:MET:HG3	6:A:638:HOH:O	2.11	0.51
1:A:378:ARG:NE	6:A:908:HOH:O	2.41	0.50
1:B:310:GLN:HG3	6:B:932:HOH:O	2.11	0.49
1:B:64:ARG:CD	3:B:509:PEG:H21	2.39	0.49
1:B:415:HIS:O	1:B:445:LEU:HA	2.13	0.48
1:A:85:ASP:HB3	1:A:88:ARG:HB2	1.96	0.48
1:A:278:ASP:OD1	1:A:280[B]:THR:HG23	2.14	0.47
1:A:269[A]:LEU:CD2	1:A:292:PHE:CD2	2.93	0.47
1:B:295:VAL:HG23	3:B:512:PEG:H21	1.95	0.47
1:A:18:GLN:NE2	6:A:875:HOH:O	2.34	0.47
1:B:8:ASP:OD2	6:B:873:HOH:O	2.21	0.46
1:B:414:ARG:HH21	3:B:511:PEG:H42	1.81	0.46
1:B:295:VAL:CG2	3:B:512:PEG:H21	2.45	0.46
1:A:12:LEU:HB3	1:A:27:TYR:HB3	1.97	0.46
1:B:214[A]:MET:HG3	6:B:632:HOH:O	2.15	0.46
1:A:415:HIS:O	1:A:445:LEU:HA	2.16	0.46
1:A:276:ALA:HB3	1:A:280[A]:THR:HG22	1.97	0.45
1:A:412:ARG:HH21	3:A:503:PEG:C2	2.23	0.45
1:A:232:VAL:HG22	1:A:242:LYS:HB2	1.96	0.45
1:A:176:GLY:O	1:A:406:GLN:HB3	2.16	0.45
1:B:330:LEU:HD11	1:B:335:TYR:HB2	1.99	0.45
1:B:85:ASP:HB3	1:B:88:ARG:HB2	1.99	0.44
5:B:503:PG4:H41	5:B:503:PG4:H61	1.60	0.44
1:B:19:VAL:CG1	3:B:510:PEG:H42	2.48	0.44
1:A:236:THR:OG1	1:A:242:LYS:HD3	2.18	0.43
1:A:269[A]:LEU:HD23	1:A:284:LEU:HD11	2.01	0.43
1:A:276:ALA:HB3	1:A:280[A]:THR:CG2	2.49	0.43
1:B:401:THR:OG1	6:B:938:HOH:O	2.21	0.43
1:B:409:LYS:HZ1	3:B:508:PEG:H21	1.84	0.43
3:B:508:PEG:H42	3:B:508:PEG:H21	1.76	0.42
1:B:289[A]:MET:HE2	6:B:981:HOH:O	2.04	0.42
1:B:14:PRO:HB2	3:B:507:PEG:H32	2.02	0.42
1:A:217:GLU:OE2	1:A:341[B]:ARG:NH2	2.37	0.41
5:B:503:PG4:H61	5:B:503:PG4:H81	1.70	0.41
1:A:60:ASN:N	3:A:505:PEG:H32	2.35	0.41
1:A:269[A]:LEU:HD23	1:A:284:LEU:HD12	2.02	0.41
1:A:214[B]:MET:HE2	1:A:214[B]:MET:HB2	1.88	0.41
1:B:17:VAL:HG11	1:B:99:GLN:HG2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:254:GLU:OE2	1:B:344:ALA:HB1	2.21	0.41
1:B:14:PRO:HB2	3:B:507:PEG:H11	2.03	0.40
1:B:289[A]:MET:HE1	6:B:981:HOH:O	2.09	0.40
1:B:12:LEU:HB3	1:B:27:TYR:HB3	2.03	0.40

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B:956:HOH:O	6:B:974:HOH:O[1_455]	1.89	0.31
6:B:816:HOH:O	6:B:817:HOH:O[1_565]	1.97	0.23
6:B:872:HOH:O	6:B:907:HOH:O[1_545]	2.17	0.03
6:A:959:HOH:O	6:B:767:HOH:O[1_556]	2.19	0.01
1:A:244:GLU:OE1	6:A:918:HOH:O[1_455]	2.19	0.01

## 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	478/491 (97%)	467 (98%)	10 (2%)	1 (0%)	51	27
1	B	477/491 (97%)	466 (98%)	11 (2%)	0	100	100
All	All	955/982 (97%)	933 (98%)	21 (2%)	1 (0%)	55	30

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	233	ALA

### 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	379/389 (97%)	374 (99%)	5 (1%)	73	55
1	B	378/389 (97%)	374 (99%)	4 (1%)	78	61
All	All	757/778 (97%)	748 (99%)	9 (1%)	75	58

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	242	LYS
1	A	332	LEU
1	A	420	ARG
1	A	421	ASP
1	B	238	ASN
1	B	242	LYS
1	B	420	ARG
1	B	421	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.



## 5.6 Ligand geometry

Of 18 ligands modelled in this entry, 4 are monoatomic - leaving 14 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	PEG	A	503	-	6,6,6	0.38	0	5,5,5	2.68	4 (80%)
4	PGE	A	504	-	9,9,9	0.67	0	8,8,8	1.32	1 (12%)
3	PEG	A	505	-	6,6,6	0.71	0	5,5,5	1.68	2 (40%)
3	PEG	A	506	-	6,6,6	0.55	0	5,5,5	1.43	1 (20%)
5	PG4	B	503	-	12,12,12	0.70	0	11,11,11	1.57	2 (18%)
5	PG4	B	504	-	12,12,12	0.73	0	11,11,11	1.36	0
3	PEG	B	505	-	6,6,6	0.54	0	5,5,5	0.93	0
3	PEG	B	506	-	6,6,6	0.37	0	5,5,5	1.36	1 (20%)
3	PEG	B	507	-	6,6,6	0.44	0	5,5,5	1.70	1 (20%)
3	PEG	B	508	-	6,6,6	0.64	0	5,5,5	1.29	1 (20%)
3	PEG	B	509	-	6,6,6	0.61	0	5,5,5	1.40	0
3	PEG	B	510	-	6,6,6	0.52	0	5,5,5	1.64	1 (20%)
3	PEG	B	511	-	6,6,6	0.51	0	5,5,5	1.48	1 (20%)
3	PEG	B	512	-	6,6,6	0.63	0	5,5,5	1.36	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	PEG	A	503	-	-	0/4/4/4	0/0/0/0
4	PGE	A	504	-	-	0/7/7/7	0/0/0/0
3	PEG	A	505	-	-	0/4/4/4	0/0/0/0
3	PEG	A	506	-	-	0/4/4/4	0/0/0/0
5	PG4	B	503	-	-	0/10/10/10	0/0/0/0
5	PG4	B	504	-	-	0/10/10/10	0/0/0/0
3	PEG	B	505	-	-	0/4/4/4	0/0/0/0
3	PEG	B	506	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PEG	B	507	-	-	0/4/4/4	0/0/0/0
3	PEG	B	508	-	-	0/4/4/4	0/0/0/0
3	PEG	B	509	-	-	0/4/4/4	0/0/0/0
3	PEG	B	510	-	-	0/4/4/4	0/0/0/0
3	PEG	B	511	-	-	0/4/4/4	0/0/0/0
3	PEG	B	512	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	508	PEG	O2-C2-C1	2.01	119.42	110.15
3	A	505	PEG	O2-C3-C4	2.05	119.60	110.15
3	A	503	PEG	O1-C1-C2	2.26	124.84	111.89
3	A	506	PEG	O2-C2-C1	2.32	120.87	110.15
3	B	510	PEG	O2-C2-C1	2.32	120.88	110.15
3	A	505	PEG	O2-C2-C1	2.44	121.41	110.15
3	B	511	PEG	C3-O2-C2	2.59	124.51	113.30
5	B	503	PG4	O3-C4-C3	2.60	122.36	110.41
4	A	504	PGE	O2-C2-C1	2.62	122.22	110.15
3	A	503	PEG	C3-O2-C2	2.62	124.65	113.30
3	B	507	PEG	O2-C3-C4	2.70	122.59	110.15
5	B	503	PG4	O2-C3-C4	2.77	123.14	110.41
3	A	503	PEG	O2-C3-C4	2.81	123.11	110.15
3	B	506	PEG	C3-O2-C2	2.83	125.57	113.30
3	A	503	PEG	O2-C2-C1	3.63	126.91	110.15

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

11 monomers are involved in 42 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	503	PEG	4	0
4	A	504	PGE	2	0
3	A	505	PEG	6	0
5	B	503	PG4	10	0
5	B	504	PG4	2	0
3	B	507	PEG	2	0
3	B	508	PEG	6	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	509	PEG	3	0
3	B	510	PEG	1	0
3	B	511	PEG	4	0
3	B	512	PEG	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, 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	470/491 (95%)	-0.25	30 (6%) 20 19	10, 17, 45, 136	0
1	B	470/491 (95%)	-0.34	21 (4%) 34 32	7, 14, 38, 124	0
All	All	940/982 (95%)	-0.29	51 (5%) 26 25	7, 16, 41, 136	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	234	GLY	9.7
1	B	237	LYS	8.4
1	B	235	THR	8.4
1	B	232	VAL	8.0
1	B	233	ALA	6.8
1	A	243	PHE	6.2
1	A	237	LYS	6.1
1	A	236	THR	5.9
1	B	236	THR	5.9
1	B	243	PHE	5.6
1	B	238	ASN	5.4
1	A	231	PRO	5.3
1	B	241	PRO	5.1
1	A	229	TYR	5.0
1	A	233	ALA	4.5
1	A	238	ASN	4.5
1	A	246	VAL	4.5
1	A	232	VAL	4.4
1	A	247	SER	3.9
1	A	241	PRO	3.8
1	A	402	THR	3.8
1	B	231	PRO	3.6
1	A	277	SER	3.5
1	A	403	THR	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	245	PRO	3.3
1	B	239	GLY	3.2
1	A	230	GLN	2.9
1	A	122	SER	2.8
1	B	310	GLN	2.8
1	A	91	TRP	2.8
1	B	242	LYS	2.7
1	B	277	SER	2.7
1	A	401	THR	2.5
1	B	249	GLN	2.4
1	A	226	TRP	2.4
1	A	405	SER	2.4
1	A	19	VAL	2.3
1	A	239	GLY	2.3
1	A	21	GLY	2.3
1	B	240	ARG	2.3
1	A	89	VAL	2.2
1	B	402	THR	2.2
1	B	279	GLY	2.2
1	A	121	LYS	2.2
1	B	121	LYS	2.2
1	A	20	ALA	2.2
1	B	252	VAL	2.1
1	A	55	HIS	2.1
1	A	240	ARG	2.0
1	A	278	ASP	2.0
1	B	312	THR	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(Å <sup>2</sup> )	Q<0.9
2	CD	B	502	1/1	0.51	1.18	60.83	76,76,76,76	1
3	PEG	B	508	7/7	0.93	0.17	10.42	22,29,35,44	0
3	PEG	A	503	7/7	0.88	0.12	6.07	20,28,38,39	0
2	CD	A	502	1/1	0.93	0.17	5.75	27,27,27,27	1
3	PEG	B	507	7/7	0.71	0.18	4.50	33,42,46,46	0
3	PEG	B	505	7/7	0.94	0.13	3.96	19,22,26,27	0
3	PEG	B	511	7/7	0.92	0.12	3.11	18,32,37,41	0
3	PEG	B	510	7/7	0.82	0.12	2.23	38,40,46,52	0
3	PEG	B	506	7/7	0.91	0.10	2.08	19,23,31,33	0
5	PG4	B	503	13/13	0.90	0.11	1.79	20,32,46,49	0
3	PEG	A	505	7/7	0.94	0.11	1.44	28,35,39,44	0
3	PEG	B	512	7/7	0.90	0.15	0.90	41,43,47,48	0
5	PG4	B	504	13/13	0.92	0.09	0.74	27,36,47,48	0
3	PEG	B	509	7/7	0.87	0.09	0.59	37,45,48,51	0
3	PEG	A	506	7/7	0.89	0.13	0.03	35,39,45,49	0
4	PGE	A	504	10/10	0.93	0.09	0.02	25,33,45,49	0
2	CD	B	501	1/1	0.98	0.06	-	31,31,31,31	1
2	CD	A	501	1/1	0.96	0.06	-	35,35,35,35	1

## 6.5 Other polymers [i](#)

There are no such residues in this entry.