



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:22 PM GMT

PDB ID : 4RLG
Title : The clear crystal structure of pyridoxal-dependent decarboxylase from sphaerobacter thermophilus dsm 20745
Authors : Wu, R.; Clancy, S.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2014-10-16
Resolution : 1.90 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

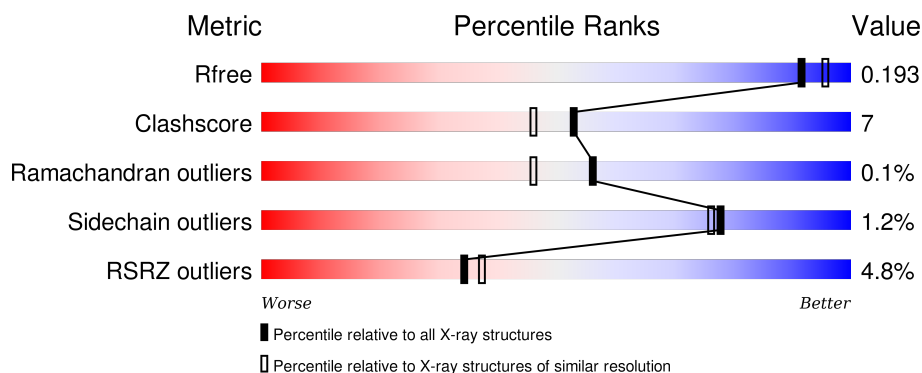
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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	486	<div> <div>6%</div> <div>86%10% . .</div> </div>
1	B	486	<div> <div>3%</div> <div>87%9% .</div> </div>
1	C	486	<div> <div>5%</div> <div>84%12% .</div> </div>
1	D	486	<div> <div>3%</div> <div>84%13% .</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
3	GOL	A	2002	-	-	-	X
3	GOL	B	503	-	-	-	X
3	GOL	C	502	-	-	-	X
5	ABU	D	502[A]	-	-	X	-
5	ABU	D	503[B]	-	-	X	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 15766 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.

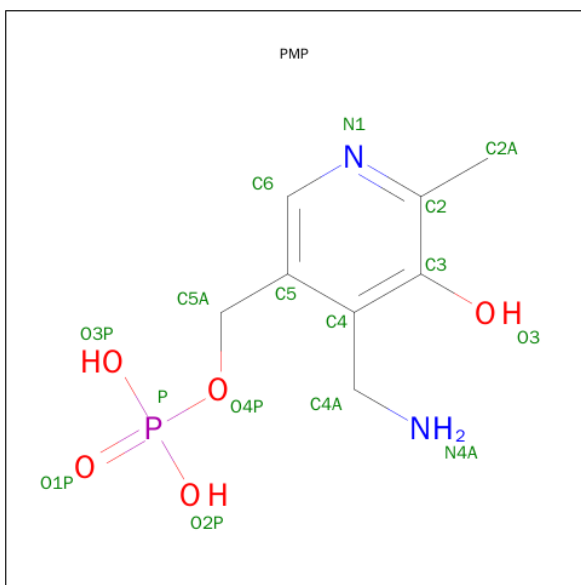
- Molecule 1 is a protein called Pyridoxal-dependent decarboxylase.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
1	A	472	Total	C	N	O	P	S	Se	0	3	0
			3640	2304	653	665	1	3	14			
1	B	470	Total	C	N	O	P	S	Se	0	5	0
			3625	2294	651	663	1	3	13			
1	D	470	Total	C	N	O	P	S	Se	0	3	0
			3628	2294	655	661	1	3	14			
1	C	469	Total	C	N	O	P	S	Se	0	2	0
			3612	2284	648	662	1	3	14			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	EXPRESSION TAG	UNP D1C7D8
A	-1	ASN	-	EXPRESSION TAG	UNP D1C7D8
A	0	ALA	-	EXPRESSION TAG	UNP D1C7D8
B	-2	SER	-	EXPRESSION TAG	UNP D1C7D8
B	-1	ASN	-	EXPRESSION TAG	UNP D1C7D8
B	0	ALA	-	EXPRESSION TAG	UNP D1C7D8
D	-2	SER	-	EXPRESSION TAG	UNP D1C7D8
D	-1	ASN	-	EXPRESSION TAG	UNP D1C7D8
D	0	ALA	-	EXPRESSION TAG	UNP D1C7D8
C	-2	SER	-	EXPRESSION TAG	UNP D1C7D8
C	-1	ASN	-	EXPRESSION TAG	UNP D1C7D8
C	0	ALA	-	EXPRESSION TAG	UNP D1C7D8

- Molecule 2 is 4'-DEOXY-4'-AMINOPYRIDOXAL-5'-PHOSPHATE (three-letter code: PMP) (formula: C₈H₁₃N₂O₅P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	1
			16	8	2	5	1		
2	B	1	Total	C	N	O	P	0	1
			16	8	2	5	1		
2	D	1	Total	C	N	O	P	0	1
			16	8	2	5	1		
2	C	1	Total	C	N	O	P	0	1
			16	8	2	5	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).

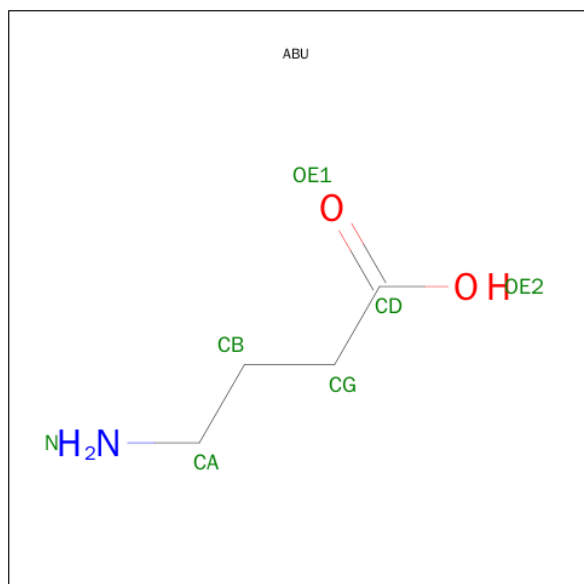


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	D	1	Total C O 6 3 3	0	0
3	C	1	Total C O 6 3 3	0	0

- Molecule 4 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total K 1 1	0	0
4	A	1	Total K 1 1	0	0
4	D	1	Total K 1 1	0	0
4	C	1	Total K 1 1	0	0

- Molecule 5 is GAMMA-AMINO-BUTANOIC ACID (three-letter code: ABU) (formula: C₄H₉NO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	D	1	Total	C	N	O	0	1
			7	4	1	2		
5	D	1	Total	C	N	O	0	1
			7	4	1	2		

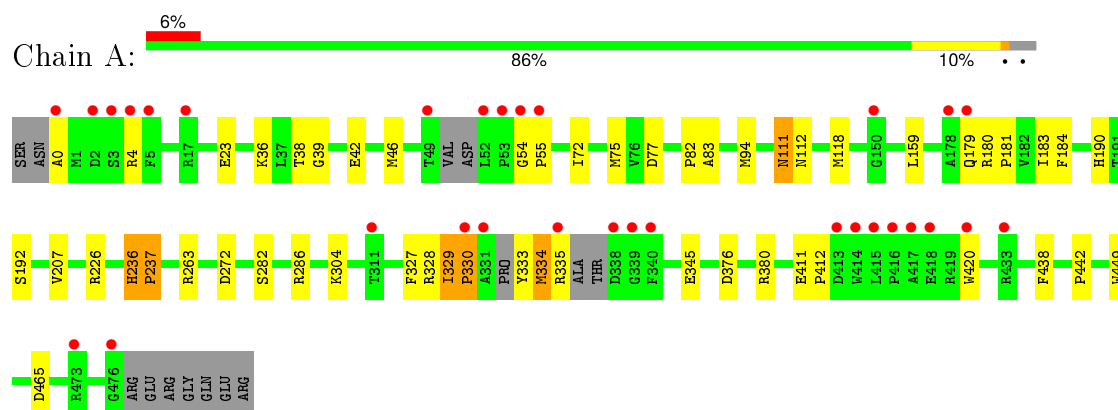
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	264	Total	O	0	0
			264	264		
6	B	305	Total	O	0	0
			305	305		
6	D	278	Total	O	0	0
			278	278		
6	C	302	Total	O	0	0
			302	302		

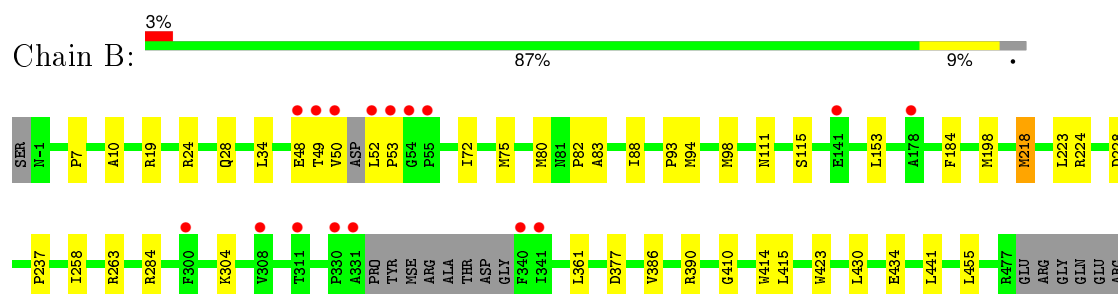
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

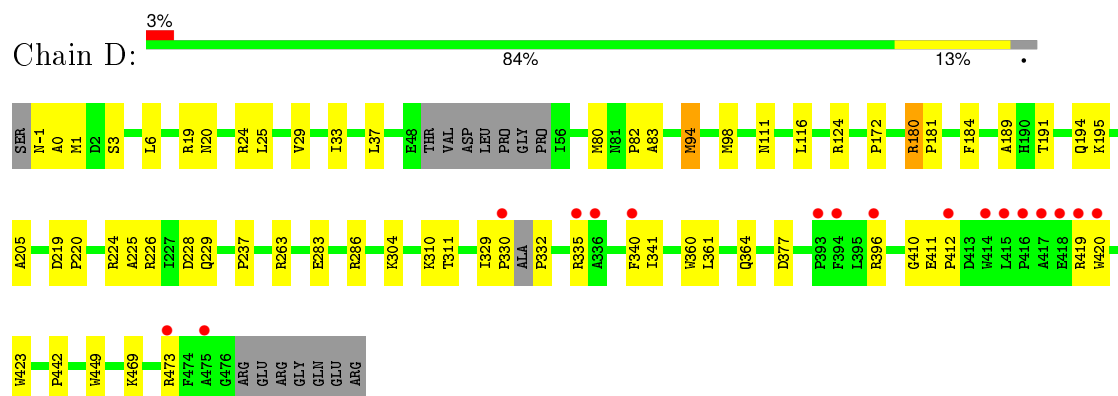
- Molecule 1: Pyridoxal-dependent decarboxylase



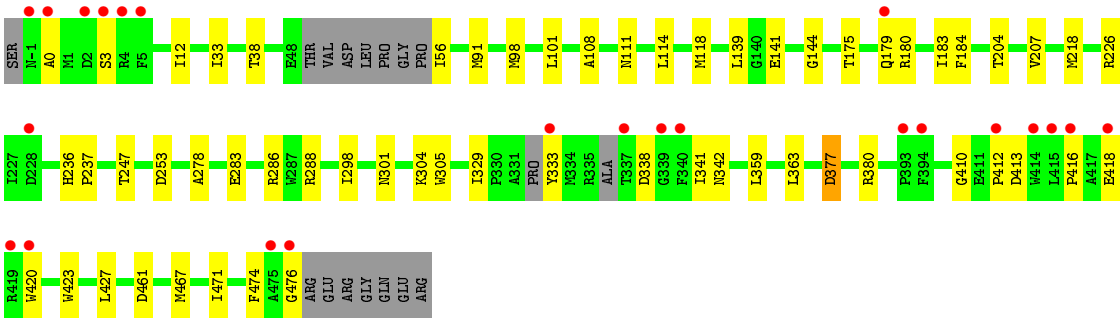
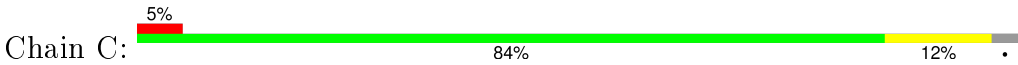
- Molecule 1: Pyridoxal-dependent decarboxylase



- Molecule 1: Pyridoxal-dependent decarboxylase



- Molecule 1: Pyridoxal-dependent decarboxylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	69.32Å 124.98Å 132.76Å 90.00° 99.49° 90.00°	Depositor
Resolution (Å)	39.70 – 1.90 39.70 – 1.90	Depositor EDS
% Data completeness (in resolution range)	88.1 (39.70-1.90) 88.1 (39.70-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.93 (at 1.89Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.162 , 0.191 0.166 , 0.193	Depositor DCC
R_{free} test set	7746 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	21.4	Xtriage
Anisotropy	0.322	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 48.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 154013 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	15766	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.36% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, K, LLP, PMP, ABU

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	0.78	3/3673 (0.1%)	0.76	3/4968 (0.1%)
1	B	0.77	0/3657	0.73	0/4951
1	C	0.80	0/3643	0.78	3/4925 (0.1%)
1	D	0.77	0/3661	0.75	1/4950 (0.0%)
All	All	0.78	3/14634 (0.0%)	0.76	7/19794 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	237	PRO	N-CD	5.24	1.55	1.47
1	A	55	PRO	N-CD	5.14	1.55	1.47
1	A	330	PRO	N-CD	5.02	1.54	1.47

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	195	LYS	CD-CE-NZ	5.59	124.55	111.70
1	A	329	ILE	C-N-CD	5.58	140.11	128.40
1	C	377	ASP	CB-CG-OD1	5.56	123.31	118.30
1	A	54	GLY	C-N-CD	5.52	140.00	128.40
1	C	288	ARG	NE-CZ-NH2	-5.43	117.58	120.30
1	A	236	HIS	C-N-CD	5.32	139.57	128.40
1	C	286	ARG	NE-CZ-NH2	-5.29	117.66	120.30

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	3640	0	3624	38	0
1	B	3625	0	3622	43	0
1	C	3612	0	3597	49	0
1	D	3628	0	3620	69	0
2	A	16	0	9	0	0
2	B	16	0	10	3	0
2	C	16	0	10	3	0
2	D	16	0	11	4	0
3	A	6	0	8	1	0
3	B	12	0	16	1	0
3	C	6	0	8	1	0
3	D	6	0	8	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	D	14	0	10	15	0
6	A	264	0	0	3	0
6	B	305	0	0	8	0
6	C	302	0	0	7	0
6	D	278	0	0	9	0
All	All	15766	0	14553	189	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 7.

All (189) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:191:THR:HG22	5:D:503[B]:ABU:CA	1.76	1.15
1:A:333:TYR:HB2	1:A:334:MSE:CE	1.80	1.10
1:D:191:THR:HA	5:D:503[B]:ABU:HB1	1.29	1.08
1:D:191:THR:HG22	5:D:502[A]:ABU:N	1.72	1.01
1:A:333:TYR:HB2	1:A:334:MSE:HE3	1.02	0.99
3:C:502:GOL:H11	6:C:810:HOH:O	1.64	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:191:THR:HA	5:D:503[B]:ABU:CB	1.94	0.95
1:A:333:TYR:CB	1:A:334:MSE:HE3	1.94	0.95
1:D:98:MSE:CE	1:D:361:LEU:HD13	1.95	0.95
1:D:80:MSE:HG2	1:D:82:PRO:HD3	1.52	0.91
1:D:191:THR:HA	5:D:502[A]:ABU:CA	1.99	0.91
1:A:0:ALA:HB2	1:B:377:ASP:OD2	1.71	0.88
1:D:283:GLU:HG2	1:D:286:ARG:NH2	1.94	0.82
1:A:82:PRO:HD2	6:A:2362:HOH:O	1.81	0.80
1:B:98:MSE:HE3	1:B:361:LEU:HB3	1.65	0.79
1:D:283:GLU:HG2	1:D:286:ARG:HH22	1.48	0.76
1:D:0:ALA:HB2	1:C:377:ASP:OD2	1.86	0.76
1:D:98:MSE:HE1	1:D:361:LEU:HD13	1.67	0.75
1:D:194:GLN:HG3	5:D:503[B]:ABU:OE2	1.87	0.74
1:C:416:PRO:HB2	1:C:418:GLU:OE2	1.87	0.74
1:D:98:MSE:HE2	1:D:361:LEU:HB3	1.69	0.72
1:A:39:GLY:O	1:A:42:GLU:HG2	1.88	0.72
1:D:20:ASN:O	1:D:24:ARG:HG3	1.89	0.71
1:D:25:LEU:HD21	1:C:98:MSE:CE	2.21	0.70
1:B:390:ARG:HG2	6:B:902:HOH:O	1.92	0.70
1:B:98:MSE:CE	1:B:361:LEU:HD13	2.22	0.70
1:C:380:ARG:NH2	1:C:461:ASP:OD1	2.24	0.70
1:C:183:ILE:HD11	1:C:207:VAL:HG22	1.75	0.68
1:D:-1:ASN:ND2	6:D:878:HOH:O	2.26	0.67
1:A:179:GLN:N	1:A:179:GLN:OE1	2.28	0.65
1:D:189:ALA:O	5:D:502[A]:ABU:HB1	1.97	0.63
1:D:25:LEU:HD21	1:C:98:MSE:HE3	1.80	0.63
1:B:98:MSE:CE	1:B:361:LEU:HB3	2.28	0.63
1:B:153:LEU:HG	6:B:896:HOH:O	1.98	0.63
1:C:474:PHE:C	1:C:476:GLY:N	2.52	0.62
1:B:390:ARG:CG	6:B:902:HOH:O	2.46	0.62
1:D:191:THR:CA	5:D:503[B]:ABU:CB	2.72	0.62
1:D:377:ASP:OD2	1:C:0:ALA:HB2	1.98	0.62
1:A:179:GLN:HG2	1:A:236:HIS:HB2	1.81	0.62
1:C:474:PHE:C	1:C:476:GLY:H	2.04	0.60
1:C:423:TRP:CE2	1:C:474:PHE:HE2	2.19	0.59
1:B:50:VAL:HG23	1:B:52:LEU:N	2.16	0.59
1:D:263:ARG:NH2	6:D:715:HOH:O	2.35	0.59
1:D:80:MSE:HE2	1:C:111:ASN:HA	1.84	0.59
1:D:184:PHE:HE2	1:D:237:PRO:HB3	1.67	0.59
1:B:98:MSE:HE3	1:B:361:LEU:HD13	1.83	0.59
1:C:179:GLN:N	1:C:179:GLN:OE1	2.35	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:ALA:HA	1:A:94:MSE:HE2	1.85	0.59
1:C:341:ILE:HG12	1:C:342:ASN:N	2.17	0.58
1:D:172:PRO:O	1:C:180:ARG:NH1	2.31	0.58
1:D:191:THR:HG22	5:D:503[B]:ABU:CB	2.32	0.58
1:C:410:GLY:HA3	1:C:423:TRP:CZ3	2.39	0.57
1:B:93:PRO:HB3	1:B:455:LEU:HD13	1.85	0.57
1:A:184:PHE:HE2	1:A:237:PRO:HB3	1.70	0.56
1:D:410:GLY:HA3	1:D:423:TRP:CZ3	2.41	0.56
1:A:46:MSE:HE2	1:A:438:PHE:CE1	2.40	0.56
1:B:80:MSE:HG2	1:B:82:PRO:HD3	1.89	0.55
1:D:304[A]:LLP:H5'1	6:D:660:HOH:O	2.05	0.55
1:B:98:MSE:HE1	1:B:361:LEU:HD13	1.87	0.55
1:D:124:ARG:HD2	6:D:842:HOH:O	2.05	0.55
1:D:191:THR:HG23	1:C:333:TYR:OH	2.05	0.55
1:D:329:ILE:HG23	1:C:204:THR:HG21	1.89	0.55
1:D:191:THR:HA	5:D:503[B]:ABU:HB2	1.87	0.55
1:D:469:LYS:HD3	1:D:469:LYS:N	2.21	0.55
1:C:56:ILE:HA	6:C:855:HOH:O	2.06	0.55
1:D:191:THR:CA	5:D:503[B]:ABU:HB2	2.37	0.54
1:D:224:ARG:NH1	1:D:228:ASP:OD1	2.27	0.54
1:D:98:MSE:CE	1:D:361:LEU:HB3	2.35	0.54
1:A:190:HIS:CD2	1:A:192:SER:H	2.26	0.54
1:C:184:PHE:HE2	1:C:237:PRO:HB3	1.73	0.53
1:D:191:THR:CG2	5:D:502[A]:ABU:N	2.62	0.53
1:B:410:GLY:HA3	1:B:423:TRP:CZ3	2.44	0.53
1:A:36:LYS:NZ	1:A:77:ASP:OD1	2.41	0.53
1:D:19[B]:ARG:HG3	1:C:38:THR:HG21	1.91	0.53
1:D:98:MSE:HE3	1:D:361:LEU:HD13	1.85	0.53
1:C:380:ARG:HH21	1:C:461:ASP:CG	2.11	0.52
1:D:191:THR:N	5:D:503[B]:ABU:HB2	2.24	0.52
1:D:83:ALA:HA	1:D:94:MSE:HE3	1.92	0.52
1:B:218:MSE:HE2	1:B:258[B]:ILE:HD12	1.92	0.52
1:C:412:PRO:HD2	1:C:420:TRP:CD1	2.45	0.52
1:A:183[B]:ILE:HD11	1:A:207:VAL:HG22	1.92	0.51
1:A:190:HIS:HD2	1:A:192:SER:OG	1.93	0.51
1:D:329:ILE:HG23	1:D:330:PRO:HD2	1.92	0.51
1:B:72:ILE:HA	1:B:75:MSE:HE2	1.93	0.51
1:C:329:ILE:HG12	6:C:870:HOH:O	2.11	0.50
1:D:473:ARG:HG3	1:D:473:ARG:HH11	1.74	0.50
1:A:335:ARG:HB2	1:B:441:LEU:HB3	1.94	0.50
2:D:501[B]:PMP:H5A1	6:D:660:HOH:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:224:ARG:HH12	1:B:228:ASP:CG	2.15	0.49
1:D:180:ARG:HH22	1:D:205:ALA:HB3	1.78	0.49
1:D:412:PRO:HD2	1:D:420:TRP:NE1	2.27	0.49
1:B:98:MSE:HE3	1:B:361:LEU:CB	2.41	0.48
1:A:329:ILE:HG23	1:A:330:PRO:HD2	1.96	0.48
1:D:335:ARG:HD2	6:D:871:HOH:O	2.14	0.48
1:B:24:ARG:O	1:B:28:GLN:HG3	2.14	0.47
1:C:278:ALA:CB	1:C:305:TRP:HB3	2.44	0.47
1:B:414:TRP:CD1	1:B:415:LEU:HD13	2.49	0.47
1:A:190:HIS:HD2	1:A:192:SER:CB	2.27	0.47
1:A:236:HIS:HE1	6:A:2337:HOH:O	1.97	0.47
1:C:141:GLU:OE1	1:C:141:GLU:HA	2.15	0.47
1:A:23:GLU:HG2	1:B:34:LEU:HD13	1.96	0.47
1:B:430:LEU:HD12	1:B:434:GLU:HG2	1.95	0.47
1:C:236:HIS:CE1	6:C:801:HOH:O	2.68	0.47
1:D:191:THR:CA	5:D:502[A]:ABU:CA	2.82	0.46
1:A:376:ASP:O	1:A:380:ARG:HG3	2.15	0.46
1:B:52:LEU:HD23	1:B:52:LEU:C	2.36	0.46
1:A:282:SER:O	1:A:286:ARG:HB2	2.15	0.46
1:C:218:MSE:HE3	1:C:253:ASP:OD2	2.15	0.46
1:D:116:LEU:HD21	1:D:124:ARG:NH2	2.30	0.46
1:B:7:PRO:HG2	1:B:10:ALA:HB2	1.97	0.46
1:B:284:ARG:NH2	6:B:721:HOH:O	2.36	0.46
1:D:473:ARG:HG3	1:D:473:ARG:NH1	2.31	0.46
1:B:263:ARG:NH2	6:B:762:HOH:O	2.49	0.46
1:B:83:ALA:HA	1:B:94:MSE:HE2	1.98	0.46
1:C:380:ARG:NH2	1:C:461:ASP:CG	2.68	0.45
1:C:91:MSE:HE1	1:C:247:THR:HA	1.97	0.45
1:A:411:GLU:HG2	1:A:420:TRP:CE3	2.52	0.45
1:D:33:ILE:HD12	1:D:33:ILE:HG23	1.77	0.45
1:A:442:PRO:HD2	1:A:449:TRP:O	2.16	0.45
1:D:225:ALA:O	1:D:229:GLN:HG3	2.17	0.45
1:A:46:MSE:HE2	1:A:438:PHE:CZ	2.53	0.44
1:B:115:SER:HB3	3:B:502:GOL:H2	1.98	0.44
1:D:37:LEU:HD13	1:C:12:ILE:HD12	1.99	0.44
1:B:184:PHE:HE2	1:B:237:PRO:HB3	1.81	0.44
1:D:1:MSE:SE	1:D:6:LEU:HD13	2.67	0.44
1:A:465:ASP:OD1	3:A:2002:GOL:C1	2.65	0.44
1:A:327:PHE:O	1:B:198:MSE:HE2	2.17	0.44
1:D:33:ILE:HD13	1:D:33:ILE:HA	1.85	0.44
1:D:29:VAL:HG21	1:C:101:LEU:HD13	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:179:GLN:CD	1:C:179:GLN:H	2.20	0.43
1:A:111:ASN:HA	1:B:80:MSE:HE2	2.00	0.43
1:D:360:TRP:O	1:D:364:GLN:HG3	2.19	0.43
1:A:72:ILE:HA	1:A:75:MSE:HE2	2.00	0.43
1:D:412:PRO:HA	6:D:716:HOH:O	2.18	0.42
1:C:33:ILE:HD12	1:C:33:ILE:HG23	1.78	0.42
1:B:284:ARG:NE	6:B:721:HOH:O	2.40	0.42
1:B:48:GLU:CD	1:B:49:THR:N	2.73	0.42
1:A:412:PRO:HD2	1:A:420:TRP:CD1	2.54	0.42
1:D:189:ALA:O	5:D:502[A]:ABU:OE1	2.37	0.42
1:C:412:PRO:HD2	1:C:420:TRP:NE1	2.35	0.42
1:C:139:LEU:HD11	1:C:298:ILE:HD12	2.02	0.42
1:A:263:ARG:NH2	6:A:2203:HOH:O	2.47	0.42
1:B:390:ARG:HG3	6:B:902:HOH:O	2.17	0.42
1:B:386:VAL:O	1:B:390:ARG:HG3	2.20	0.42
1:D:419:ARG:HA	6:D:767:HOH:O	2.20	0.42
1:C:236:HIS:HE1	6:C:801:HOH:O	2.01	0.41
1:C:33:ILE:HD13	1:C:33:ILE:HA	1.83	0.41
1:C:283:GLU:HG2	6:C:899:HOH:O	2.19	0.41
1:D:219:ASP:HA	1:D:220:PRO:HD3	1.80	0.41
1:A:159:LEU:HD22	1:A:183[A]:ILE:HD13	2.02	0.41
1:A:112:ASN:HB3	1:A:118:MSE:HB3	2.01	0.41
1:A:272:ASP:OD2	1:A:304[A]:LLP:N1	2.53	0.41
1:D:396:ARG:HD3	1:D:411:GLU:OE1	2.19	0.41
1:C:359:LEU:O	1:C:363:LEU:HG	2.20	0.41
1:C:144:GLY:HA3	6:C:876:HOH:O	2.20	0.41
1:D:310:LYS:HA	1:D:311:THR:HA	1.83	0.41
1:D:37:LEU:HD11	1:C:108:ALA:HB1	2.02	0.41
1:A:180:ARG:HA	1:A:181:PRO:HD3	1.89	0.41
1:D:442:PRO:HD2	1:D:449:TRP:O	2.21	0.41
1:A:328:ARG:NH2	1:A:345:GLU:OE1	2.54	0.41
1:A:38:THR:OG1	1:B:19:ARG:NH2	2.54	0.41
1:D:181:PRO:CD	1:C:175:THR:HG21	2.50	0.41
1:D:332:PRO:N	6:D:843:HOH:O	2.54	0.41
1:B:50:VAL:O	1:B:52:LEU:N	2.54	0.40
1:B:52:LEU:HG	1:B:53:PRO:HD2	2.04	0.40
1:B:218:MSE:CE	1:B:223:LEU:HD22	2.51	0.40
1:D:412:PRO:HD2	1:D:420:TRP:CD1	2.56	0.40
1:B:88:ILE:HG22	6:B:666:HOH:O	2.20	0.40
1:D:19[A]:ARG:NH1	1:C:38:THR:OG1	2.55	0.40
1:C:329:ILE:HG21	1:C:329:ILE:HD13	1.67	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:114:LEU:HD23	1:C:118:MSE:SE	2.71	0.40
1:C:427:LEU:HD21	1:C:471:ILE:HG12	2.03	0.40
1:A:335:ARG:HD2	1:B:441:LEU:CD2	2.52	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	466/486 (96%)	458 (98%)	8 (2%)	0	100	100
1	B	467/486 (96%)	456 (98%)	11 (2%)	0	100	100
1	C	461/486 (95%)	451 (98%)	9 (2%)	1 (0%)	52	42
1	D	465/486 (96%)	455 (98%)	10 (2%)	0	100	100
All	All	1859/1944 (96%)	1820 (98%)	38 (2%)	1 (0%)	56	46

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	413	ASP

5.3.2 Protein sidechains [i](#)

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	365/362 (101%)	361 (99%)	4 (1%)	80	79
1	B	365/362 (101%)	363 (100%)	2 (0%)	92	92
1	C	363/362 (100%)	359 (99%)	4 (1%)	80	79
1	D	364/362 (101%)	357 (98%)	7 (2%)	65	59
All	All	1457/1448 (101%)	1440 (99%)	17 (1%)	78	76

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

Mol	Chain	Res	Type
1	A	4	ARG
1	A	111	ASN
1	A	226	ARG
1	A	334	MSE
1	B	111	ASN
1	B	218	MSE
1	D	3	SER
1	D	94	MSE
1	D	111	ASN
1	D	180	ARG
1	D	226	ARG
1	D	340	PHE
1	D	341	ILE
1	C	3	SER
1	C	226	ARG
1	C	338	ASP
1	C	467	MSE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	190	HIS
1	A	236	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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
1	LLP	A	304[A]	-	23,24,25	3.25	8 (34%)	28,32,34	2.65	9 (32%)
1	LLP	B	304[A]	-	23,24,25	4.62	8 (34%)	28,32,34	1.76	8 (28%)
1	LLP	C	304[A]	-	23,24,25	4.52	7 (30%)	28,32,34	2.18	10 (35%)
1	LLP	D	304[A]	-	23,24,25	4.76	9 (39%)	28,32,34	1.77	8 (28%)

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
1	LLP	A	304[A]	-	-	0/15/17/19	0/1/1/1
1	LLP	B	304[A]	-	-	0/15/17/19	0/1/1/1
1	LLP	C	304[A]	-	-	0/15/17/19	0/1/1/1
1	LLP	D	304[A]	-	-	0/15/17/19	0/1/1/1

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	304[A]	LLP	P-OP3	-3.82	1.41	1.54
1	D	304[A]	LLP	P-OP2	-3.59	1.41	1.54
1	D	304[A]	LLP	P-OP3	-3.05	1.43	1.54
1	D	304[A]	LLP	P-OP1	-3.01	1.41	1.51
1	A	304[A]	LLP	P-OP1	-2.69	1.42	1.51
1	B	304[A]	LLP	CE-NZ	-2.39	1.41	1.46
1	D	304[A]	LLP	CE-NZ	-2.35	1.41	1.46
1	A	304[A]	LLP	P-OP2	-2.12	1.47	1.54
1	C	304[A]	LLP	P-OP3	-2.10	1.47	1.54
1	B	304[A]	LLP	P-OP2	-2.08	1.47	1.54
1	C	304[A]	LLP	CE-NZ	-2.06	1.42	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	304[A]	LLP	P-OP3	-2.01	1.47	1.54
1	A	304[A]	LLP	C4-C4'	2.26	1.50	1.46
1	B	304[A]	LLP	C4-C4'	2.57	1.51	1.46
1	D	304[A]	LLP	C4-C4'	2.58	1.51	1.46
1	C	304[A]	LLP	C4-C4'	3.16	1.52	1.46
1	D	304[A]	LLP	C4-C5	4.04	1.47	1.42
1	A	304[A]	LLP	C4-C5	4.45	1.48	1.42
1	C	304[A]	LLP	C4-C3	5.10	1.47	1.40
1	D	304[A]	LLP	C4-C3	5.11	1.47	1.40
1	B	304[A]	LLP	C4-C5	5.13	1.48	1.42
1	C	304[A]	LLP	C4-C5	5.29	1.49	1.42
1	A	304[A]	LLP	C4-C3	5.64	1.47	1.40
1	B	304[A]	LLP	C4-C3	5.94	1.48	1.40
1	A	304[A]	LLP	C4'-NZ	7.62	1.50	1.27
1	A	304[A]	LLP	C3-C2	9.57	1.47	1.40
1	C	304[A]	LLP	C3-C2	10.40	1.48	1.40
1	D	304[A]	LLP	C3-C2	11.61	1.48	1.40
1	B	304[A]	LLP	C3-C2	11.72	1.48	1.40
1	B	304[A]	LLP	C4'-NZ	16.28	1.75	1.27
1	C	304[A]	LLP	C4'-NZ	16.55	1.76	1.27
1	D	304[A]	LLP	C4'-NZ	17.09	1.78	1.27

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	304[A]	LLP	C4-C4'-NZ	-4.45	100.27	125.06
1	B	304[A]	LLP	C4-C4'-NZ	-4.45	100.27	125.06
1	C	304[A]	LLP	C4-C4'-NZ	-4.42	100.44	125.06
1	A	304[A]	LLP	C4-C4'-NZ	-3.27	106.87	125.06
1	A	304[A]	LLP	C2'-C2-C3	-3.19	117.19	121.04
1	D	304[A]	LLP	C3-C4-C4'	-2.69	116.67	120.16
1	A	304[A]	LLP	C3-C4-C4'	-2.44	117.00	120.16
1	C	304[A]	LLP	O-C-CA	-2.36	119.34	125.49
1	C	304[A]	LLP	C3-C4-C4'	-2.27	117.22	120.16
1	B	304[A]	LLP	C3-C4-C4'	-2.27	117.22	120.16
1	C	304[A]	LLP	OP4-P-OP1	-2.25	101.42	107.14
1	C	304[A]	LLP	C2'-C2-C3	-2.08	118.53	121.04
1	C	304[A]	LLP	C3-C4-C5	-2.05	116.57	118.11
1	A	304[A]	LLP	OP2-P-OP1	2.06	117.20	110.58
1	A	304[A]	LLP	C5-C4-C4'	2.12	124.56	121.52
1	B	304[A]	LLP	C6-N1-C2	2.15	123.67	119.28
1	C	304[A]	LLP	O3-C3-C2	2.16	121.41	117.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	304[A]	LLP	CD-CE-NZ	2.24	114.64	110.98
1	D	304[A]	LLP	C6-N1-C2	2.28	123.93	119.28
1	D	304[A]	LLP	OP2-P-OP1	2.36	118.19	110.58
1	D	304[A]	LLP	OP4-C5'-C5	2.47	113.07	108.99
1	A	304[A]	LLP	C6-N1-C2	2.48	124.34	119.28
1	B	304[A]	LLP	C5-C4-C4'	2.60	125.25	121.52
1	B	304[A]	LLP	CE-NZ-C4'	2.62	126.55	118.97
1	D	304[A]	LLP	C5-C4-C4'	2.67	125.36	121.52
1	D	304[A]	LLP	CE-NZ-C4'	2.68	126.72	118.97
1	B	304[A]	LLP	O3-C3-C2	2.97	122.81	117.66
1	C	304[A]	LLP	C5-C4-C4'	3.22	126.14	121.52
1	C	304[A]	LLP	OP4-C5'-C5	3.82	115.31	108.99
1	B	304[A]	LLP	OP4-C5'-C5	4.11	115.79	108.99
1	D	304[A]	LLP	O3-C3-C2	4.15	124.87	117.66
1	A	304[A]	LLP	OP4-C5'-C5	4.32	116.14	108.99
1	A	304[A]	LLP	CD-CE-NZ	5.50	119.99	110.98
1	C	304[A]	LLP	CE-NZ-C4'	6.52	137.80	118.97
1	A	304[A]	LLP	CE-NZ-C4'	9.40	146.12	118.97

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	304[A]	LLP	1	0
1	B	304[A]	LLP	2	0
1	C	304[A]	LLP	2	0
1	D	304[A]	LLP	3	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 4 are monoatomic - leaving 11 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
2	PMP	A	2001[B]	1	16,16,16	3.54	6 (37%)	20,23,23	1.72	5 (25%)
3	GOL	A	2002	-	5,5,5	0.35	0	5,5,5	0.45	0
2	PMP	B	501[B]	1	16,16,16	3.82	6 (37%)	20,23,23	1.65	4 (20%)
3	GOL	B	502	-	5,5,5	0.38	0	5,5,5	0.57	0
3	GOL	B	503	-	5,5,5	0.44	0	5,5,5	0.60	0
2	PMP	C	501[B]	1	16,16,16	3.35	5 (31%)	20,23,23	1.73	4 (20%)
3	GOL	C	502	-	5,5,5	0.58	0	5,5,5	0.71	0
2	PMP	D	501[B]	1	16,16,16	3.75	6 (37%)	20,23,23	1.54	5 (25%)
5	ABU	D	502[A]	-	3,6,6	0.21	0	3,6,6	1.33	0
5	ABU	D	503[B]	-	3,6,6	0.15	0	3,6,6	1.01	0
3	GOL	D	504	-	5,5,5	0.24	0	5,5,5	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
2	PMP	A	2001[B]	1	-	0/8/8/8	0/1/1/1
3	GOL	A	2002	-	-	0/4/4/4	0/0/0/0
2	PMP	B	501[B]	1	-	0/8/8/8	0/1/1/1
3	GOL	B	502	-	-	0/4/4/4	0/0/0/0
3	GOL	B	503	-	-	0/4/4/4	0/0/0/0
2	PMP	C	501[B]	1	-	0/8/8/8	0/1/1/1
3	GOL	C	502	-	-	0/4/4/4	0/0/0/0
2	PMP	D	501[B]	1	-	0/8/8/8	0/1/1/1
5	ABU	D	502[A]	-	-	0/2/4/4	0/0/0/0
5	ABU	D	503[B]	-	-	0/2/4/4	0/0/0/0
3	GOL	D	504	-	-	0/4/4/4	0/0/0/0

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	501[B]	PMP	P-O3P	-4.11	1.40	1.54
2	D	501[B]	PMP	P-O2P	-3.74	1.41	1.54
2	D	501[B]	PMP	P-O3P	-3.59	1.41	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501[B]	PMP	P-O2P	-3.49	1.42	1.54
2	A	2001[B]	PMP	P-O2P	-3.46	1.42	1.54
2	B	501[B]	PMP	P-O3P	-3.18	1.43	1.54
2	C	501[B]	PMP	P-O2P	-3.10	1.43	1.54
2	A	2001[B]	PMP	P-O1P	-3.08	1.41	1.51
2	B	501[B]	PMP	P-O1P	-2.54	1.42	1.51
2	D	501[B]	PMP	P-O1P	-2.24	1.43	1.51
2	A	2001[B]	PMP	P-O3P	-2.11	1.47	1.54
2	D	501[B]	PMP	C5-C4	4.43	1.46	1.40
2	C	501[B]	PMP	C3-C4	4.65	1.47	1.40
2	C	501[B]	PMP	C5-C4	5.09	1.47	1.40
2	D	501[B]	PMP	C3-C4	5.16	1.48	1.40
2	A	2001[B]	PMP	C5-C4	5.22	1.47	1.40
2	B	501[B]	PMP	C5-C4	5.60	1.48	1.40
2	A	2001[B]	PMP	C3-C4	5.63	1.48	1.40
2	B	501[B]	PMP	C3-C4	5.80	1.49	1.40
2	C	501[B]	PMP	C3-C2	9.70	1.47	1.40
2	A	2001[B]	PMP	C3-C2	10.39	1.48	1.40
2	B	501[B]	PMP	C3-C2	11.54	1.48	1.40
2	D	501[B]	PMP	C3-C2	11.85	1.49	1.40

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501[B]	PMP	C2A-C2-C3	-3.49	116.82	121.04
2	A	2001[B]	PMP	C2A-C2-C3	-2.86	117.59	121.04
2	D	501[B]	PMP	C4-C3-C2	-2.59	115.35	120.03
2	C	501[B]	PMP	O3P-P-O4P	-2.39	99.70	106.56
2	C	501[B]	PMP	C4-C3-C2	-2.24	115.99	120.03
2	A	2001[B]	PMP	C4-C3-C2	-2.06	116.32	120.03
2	A	2001[B]	PMP	C6-N1-C2	2.38	124.13	119.28
2	B	501[B]	PMP	O3P-P-O2P	2.38	116.46	107.38
2	D	501[B]	PMP	C6-N1-C2	2.40	124.17	119.28
2	B	501[B]	PMP	C6-N1-C2	2.42	124.21	119.28
2	D	501[B]	PMP	O4P-C5A-C5	2.54	113.19	108.99
2	A	2001[B]	PMP	O3P-P-O2P	2.58	117.21	107.38
2	D	501[B]	PMP	O3P-P-O2P	2.83	118.14	107.38
2	B	501[B]	PMP	O3-C3-C2	2.97	122.82	117.66
2	C	501[B]	PMP	O4P-C5A-C5	3.33	114.50	108.99
2	D	501[B]	PMP	O3-C3-C2	3.64	124.00	117.66
2	A	2001[B]	PMP	O4P-C5A-C5	4.23	115.98	108.99
2	B	501[B]	PMP	O4P-C5A-C5	4.56	116.52	108.99

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 28 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	2002	GOL	1	0
2	B	501[B]	PMP	3	0
3	B	502	GOL	1	0
2	C	501[B]	PMP	3	0
3	C	502	GOL	1	0
2	D	501[B]	PMP	4	0
5	D	502[A]	ABU	6	0
5	D	503[B]	ABU	9	0

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, 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	457/486 (94%)	0.07	31 (6%)	20 23	14, 25, 53, 70	7 (1%)
1	B	456/486 (93%)	-0.12	16 (3%)	48 51	13, 24, 48, 65	1 (0%)
1	C	454/486 (93%)	-0.13	23 (5%)	32 35	12, 21, 50, 73	5 (1%)
1	D	455/486 (93%)	-0.09	17 (3%)	45 49	12, 23, 54, 68	8 (1%)
All	All	1822/1944 (93%)	-0.07	87 (4%)	34 37	12, 23, 51, 73	21 (1%)

All (87) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	52	LEU	7.0
1	A	52	LEU	6.9
1	A	4	ARG	6.4
1	A	340	PHE	6.2
1	C	476	GLY	6.0
1	A	54	GLY	5.9
1	B	50	VAL	5.5
1	B	49	THR	5.5
1	C	4	ARG	5.5
1	B	340	PHE	5.4
1	B	331	ALA	5.4
1	A	5	PHE	5.4
1	B	54	GLY	4.7
1	D	414	TRP	4.5
1	D	340	PHE	4.5
1	D	419	ARG	4.5
1	B	55	PRO	4.4
1	D	394	PHE	4.2
1	D	415	LEU	4.2
1	C	0	ALA	4.2
1	D	416	PRO	4.0

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Mol	Chain	Res	Type	RSRZ
1	D	473	ARG	4.0
1	C	414	TRP	3.9
1	A	49	THR	3.9
1	D	475	ALA	3.8
1	C	415	LEU	3.7
1	C	475	ALA	3.6
1	D	418	GLU	3.6
1	A	0	ALA	3.6
1	C	-1	ASN	3.5
1	C	2	ASP	3.4
1	C	5	PHE	3.4
1	B	48	GLU	3.3
1	C	179	GLN	3.3
1	A	2	ASP	3.3
1	A	331	ALA	3.3
1	B	141	GLU	3.3
1	A	414	TRP	3.2
1	C	3	SER	3.2
1	A	55	PRO	3.2
1	B	53	PRO	3.2
1	C	416	PRO	3.2
1	A	417	ALA	3.1
1	A	473	ARG	3.1
1	A	330	PRO	3.0
1	A	339	GLY	3.0
1	D	393	PRO	3.0
1	D	412	PRO	3.0
1	A	476	GLY	3.0
1	C	339	GLY	2.9
1	A	338	ASP	2.9
1	D	417	ALA	2.8
1	A	3	SER	2.8
1	A	53	PRO	2.8
1	A	413	ASP	2.8
1	D	420	TRP	2.8
1	A	179	GLN	2.8
1	C	333	TYR	2.7
1	B	341	ILE	2.7
1	D	335	ARG	2.6
1	A	418	GLU	2.6
1	D	396	ARG	2.6
1	C	337	THR	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	412	PRO	2.5
1	A	415	LEU	2.5
1	A	416	PRO	2.5
1	C	419	ARG	2.4
1	C	340	PHE	2.4
1	D	336	ALA	2.3
1	D	330	PRO	2.3
1	A	17	ARG	2.3
1	A	335	ARG	2.3
1	A	311	THR	2.3
1	B	311	THR	2.2
1	A	420	TRP	2.2
1	C	420	TRP	2.1
1	C	418	GLU	2.1
1	B	308	VAL	2.1
1	C	228	ASP	2.1
1	C	393	PRO	2.1
1	B	178	ALA	2.1
1	B	300	PHE	2.1
1	A	150	GLY	2.1
1	B	330	PRO	2.1
1	C	394	PHE	2.1
1	A	178	ALA	2.1
1	A	433	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	LLP	A	304[A]	24/25	0.94	0.17	-	14,25,31,33	15
1	LLP	C	304[A]	24/25	0.96	0.13	-	11,21,25,27	15
1	LLP	B	304[A]	24/25	0.97	0.17	-	13,24,29,31	15
1	LLP	D	304[A]	24/25	0.96	0.16	-	12,23,30,30	15

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, 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
3	GOL	B	503	6/6	0.85	0.28	6.42	59,62,65,65	0
3	GOL	A	2002	6/6	0.82	0.19	3.58	50,59,62,67	0
3	GOL	C	502	6/6	0.86	0.23	2.93	56,58,60,63	0
5	ABU	D	503[B]	7/7	0.69	0.22	2.74	27,30,33,35	7
3	GOL	D	504	6/6	0.86	0.24	1.73	62,63,65,67	0
5	ABU	D	502[A]	7/7	0.87	0.16	1.53	24,25,33,33	6
2	PMP	D	501[B]	16/16	0.96	0.15	1.08	16,27,30,30	16
2	PMP	B	501[B]	16/16	0.96	0.16	0.93	19,25,30,31	16
2	PMP	A	2001[B]	16/16	0.96	0.15	0.74	21,27,31,33	16
2	PMP	C	501[B]	16/16	0.96	0.12	0.14	10,22,26,27	16
4	K	D	505	1/1	0.97	0.14	-1.19	49,49,49,49	0
4	K	A	2003	1/1	0.99	0.04	-1.60	44,44,44,44	0
4	K	B	504	1/1	0.98	0.03	-1.76	38,38,38,38	0
4	K	C	503	1/1	0.98	0.06	-1.88	40,40,40,40	0
3	GOL	B	502	6/6	0.90	0.15	-	41,47,55,57	0

6.5 Other polymers ⓘ

There are no such residues in this entry.