



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 09:54 am GMT

PDB ID : 3K40
Title : Crystal structure of Drosophila 3,4-dihydroxyphenylalanine decarboxylase
Authors : Han, Q.; Ding, H.; Robinson, H.; Christensen, B.M.; Li, J.
Deposited on : 2009-10-05
Resolution : 1.75 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 : 1.7.2 (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

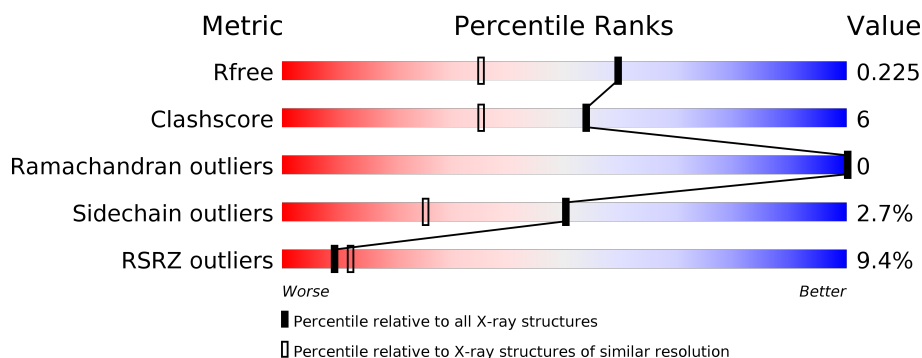
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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	1762 (1.76-1.76)
Clashscore	112137	1889 (1.76-1.76)
Ramachandran outliers	110173	1868 (1.76-1.76)
Sidechain outliers	110143	1868 (1.76-1.76)
RSRZ outliers	101464	1770 (1.76-1.76)

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	475	<div> <div>7%</div> <div> <div></div> <div>78%</div> <div>15%</div> <div>6%</div> </div> </div>
1	B	475	<div> <div>11%</div> <div> <div></div> <div>85%</div> <div>8%</div> <div>6%</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	GOL	A	476	-	-	-	X

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 7715 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 Aromatic-L-amino-acid decarboxylase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	448	Total	C	N	O	P	S	0	0	0
			3554	2272	610	643	1	28			
1	B	448	Total	C	N	O	P	S	0	0	0
			3554	2272	610	643	1	28			

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	344	Total	O	0	0
			344	344		

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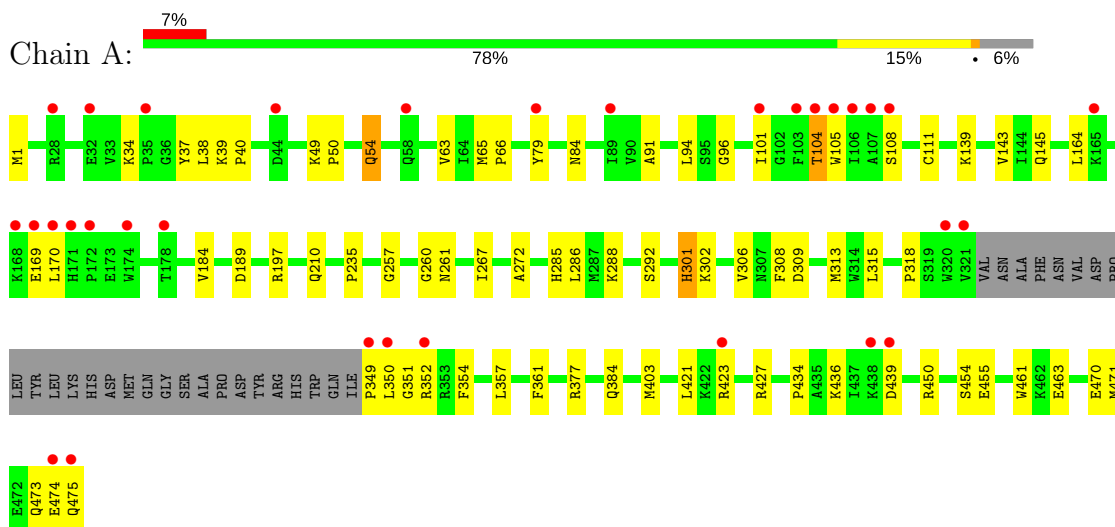
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	257	Total 257	O 257	0	0

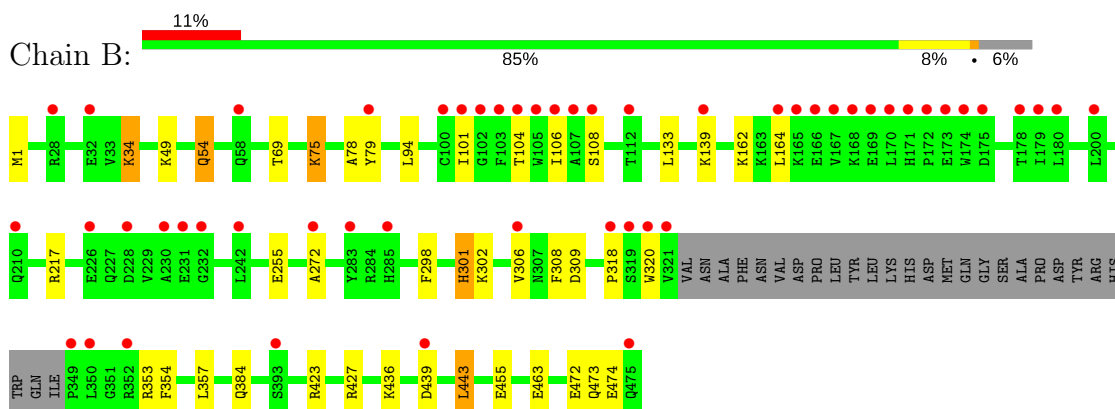
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Aromatic-L-amino-acid decarboxylase



• Molecule 1: Aromatic-L-amino-acid decarboxylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	105.81Å 108.60Å 86.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.88 – 1.75 29.87 – 1.75	Depositor EDS
% Data completeness (in resolution range)	98.9 (29.88-1.75) 98.9 (29.87-1.75)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.60 (at 1.75Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.197 , 0.226 0.196 , 0.225	Depositor DCC
R_{free} test set	4972 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	25.2	Xtriage
Anisotropy	0.051	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 49.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.009 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7715	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 5.04% 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: GOL, LLP

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	0.69	0/3618	0.70	1/4891 (0.0%)
1	B	0.61	0/3618	0.68	2/4891 (0.0%)
All	All	0.65	0/7236	0.69	3/9782 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	443	LEU	CA-CB-CG	5.44	127.81	115.30
1	B	443	LEU	CB-CG-CD2	5.25	119.92	111.00
1	A	461	TRP	CA-CB-CG	-5.23	103.77	113.70

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	3554	0	3499	62	0
1	B	3554	0	3499	31	0
2	A	6	0	8	3	0
3	A	344	0	0	9	0
3	B	257	0	0	4	0
All	All	7715	0	7006	86	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 6.

All (86) 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:350:LEU:HD21	3:A:584:HOH:O	1.65	0.97
1:A:313:MET:HE3	1:A:315:LEU:HD11	1.58	0.84
1:A:455:GLU:HG2	3:A:789:HOH:O	1.80	0.79
1:A:352:ARG:HB2	3:A:541:HOH:O	1.83	0.78
1:B:54:GLN:H	1:B:54:GLN:HE21	1.34	0.75
1:B:436:LYS:HE2	1:B:439:ASP:HA	1.71	0.73
1:A:105:TRP:HZ3	1:A:349:PRO:HD2	1.55	0.72
1:A:54:GLN:H	1:A:54:GLN:HE21	1.35	0.70
1:A:436:LYS:HE2	1:A:439:ASP:HA	1.73	0.69
3:A:585:HOH:O	1:B:49:LYS:HD3	1.93	0.69
1:A:423:ARG:HH21	1:A:470:GLU:HB3	1.58	0.69
1:A:143:VAL:CG1	1:A:313:MET:HE2	2.24	0.67
1:A:38:LEU:HD11	1:A:63:VAL:CG1	2.25	0.66
1:B:75:LYS:HE3	3:B:523:HOH:O	1.95	0.66
1:A:101:ILE:HD12	1:B:308:PHE:HB2	1.76	0.65
1:A:308:PHE:HE1	1:B:353:ARG:HE	1.42	0.64
1:A:104:THR:HB	1:B:79:TYR:HD2	1.63	0.63
1:A:38:LEU:CD1	1:A:63:VAL:HG13	2.30	0.62
1:A:79:TYR:HD2	1:B:104:THR:HG23	1.63	0.61
1:A:384:GLN:HE22	1:A:454:SER:HB2	1.66	0.61
1:A:1:MET:N	2:A:476:GOL:H31	2.16	0.60
1:B:384:GLN:NE2	3:B:702:HOH:O	2.34	0.60
1:B:217:ARG:NH1	1:B:255:GLU:OE1	2.35	0.59
1:A:104:THR:HG21	1:B:78:ALA:HB1	1.85	0.58
1:A:39:LYS:HD3	3:A:539:HOH:O	2.04	0.57
1:A:423:ARG:CD	1:A:471:MET:HG2	2.36	0.56
1:B:217:ARG:HH12	1:B:255:GLU:CD	2.10	0.55
1:A:49:LYS:HE3	1:A:50:PRO:HD2	1.88	0.54
1:A:94:LEU:HD23	1:A:357:LEU:HD21	1.90	0.54
1:A:1:MET:H3	2:A:476:GOL:H31	1.72	0.53
1:A:313:MET:CE	1:A:315:LEU:HD11	2.34	0.53
1:B:34:LYS:CE	1:B:34:LYS:H	2.21	0.53
1:A:377:ARG:HG3	3:A:603:HOH:O	2.09	0.53
1:A:84:ASN:HD22	1:A:450:ARG:HE	1.57	0.52
1:A:79:TYR:HB2	1:B:104:THR:OG1	2.09	0.52
1:A:1:MET:N	2:A:476:GOL:C3	2.72	0.52
1:A:139:LYS:O	1:A:318:PRO:HD2	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:261:ASN:HD21	1:A:292:SER:HB2	1.75	0.51
1:A:38:LEU:CD1	1:A:63:VAL:CG1	2.88	0.51
1:A:285:HIS:HD2	3:A:571:HOH:O	1.94	0.50
1:B:34:LYS:NZ	1:B:34:LYS:H	2.10	0.50
1:B:54:GLN:H	1:B:54:GLN:NE2	2.06	0.50
1:A:257:GLY:HA3	1:A:292:SER:OG	2.12	0.50
1:A:145:GLN:HE21	1:A:313:MET:HE1	1.78	0.49
1:A:260:GLY:HA3	1:A:267:ILE:HD11	1.94	0.49
1:A:421:LEU:HD22	1:A:434:PRO:HB3	1.94	0.49
1:A:272:ALA:HB3	1:A:302:LLP:C3	2.43	0.49
1:A:189:ASP:OD1	1:A:210:GLN:NE2	2.47	0.47
1:A:91:ALA:HB2	1:A:361:PHE:CD2	2.50	0.47
1:B:106:ILE:CG2	1:B:108:SER:HB3	2.44	0.47
1:A:427:ARG:HD3	1:A:463:GLU:OE1	2.14	0.47
1:A:423:ARG:CD	1:A:471:MET:CG	2.93	0.47
1:B:94:LEU:HD23	1:B:357:LEU:HD21	1.96	0.47
1:B:272:ALA:HB3	1:B:302:LLP:C3	2.46	0.46
1:A:261:ASN:ND2	1:A:292:SER:HB2	2.30	0.46
1:B:162:LYS:HG3	1:B:320:TRP:CZ2	2.51	0.46
1:B:301:HIS:HA	1:B:306:VAL:O	2.15	0.46
1:B:1:MET:N	3:B:530:HOH:O	2.42	0.46
1:B:308:PHE:HA	1:B:309:ASP:HA	1.65	0.46
1:A:302:LLP:O3	1:A:302:LLP:NZ	2.49	0.46
1:A:111:CYS:HB3	1:A:351:GLY:HA3	1.98	0.45
1:A:145:GLN:HE21	1:A:313:MET:CE	2.30	0.45
1:A:108:SER:HB2	1:B:69:THR:OG1	2.16	0.45
1:A:473:GLN:C	1:A:475:GLN:H	2.19	0.45
1:B:455:GLU:CD	1:B:455:GLU:H	2.20	0.45
1:A:39:LYS:HB3	1:A:40:PRO:HD3	1.99	0.44
1:A:423:ARG:HG3	3:A:719:HOH:O	2.16	0.44
1:A:423:ARG:HD3	1:A:471:MET:CG	2.47	0.44
1:A:54:GLN:H	1:A:54:GLN:NE2	2.09	0.44
1:A:261:ASN:HD21	1:A:292:SER:CB	2.31	0.43
1:A:308:PHE:HA	1:A:309:ASP:HA	1.78	0.43
1:A:105:TRP:HZ3	1:A:349:PRO:CD	2.29	0.43
1:B:423:ARG:HG3	3:B:508:HOH:O	2.17	0.43
1:A:286:LEU:CD1	1:A:403:MET:HE2	2.50	0.42
1:B:34:LYS:HE2	1:B:34:LYS:HB2	1.71	0.42
1:A:184:VAL:HG13	1:A:235:PRO:HA	2.01	0.42
1:B:427:ARG:HD3	1:B:463:GLU:OE1	2.20	0.42
1:A:423:ARG:NH2	1:A:470:GLU:HB3	2.31	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:286:LEU:CD1	1:A:403:MET:CE	2.99	0.41
1:A:96:GLY:HA3	3:A:736:HOH:O	2.20	0.41
1:A:34:LYS:O	1:A:37:TYR:HB3	2.20	0.41
1:B:472:GLU:C	1:B:474:GLU:H	2.24	0.41
1:A:301:HIS:HA	1:A:306:VAL:O	2.21	0.41
1:B:272:ALA:HA	1:B:298:PHE:HA	2.02	0.41
1:A:65:MET:N	1:A:66:PRO:CD	2.84	0.41
1:B:139:LYS:O	1:B:318:PRO:CD	2.69	0.41

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	443/475 (93%)	435 (98%)	8 (2%)	0	100	100
1	B	443/475 (93%)	430 (97%)	13 (3%)	0	100	100
All	All	886/950 (93%)	865 (98%)	21 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	371/395 (94%)	361 (97%)	10 (3%)	50	25

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	371/395 (94%)	361 (97%)	10 (3%)	50	25
All	All	742/790 (94%)	722 (97%)	20 (3%)	50	25

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	GLN
1	A	104	THR
1	A	164	LEU
1	A	169	GLU
1	A	170	LEU
1	A	197	ARG
1	A	288	LYS
1	A	301	HIS
1	A	354	PHE
1	A	474	GLU
1	B	34	LYS
1	B	54	GLN
1	B	75	LYS
1	B	101	ILE
1	B	133	LEU
1	B	164	LEU
1	B	301	HIS
1	B	354	PHE
1	B	443	LEU
1	B	473	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	GLN
1	A	77	HIS
1	A	84	ASN
1	A	145	GLN
1	A	210	GLN
1	A	261	ASN
1	A	384	GLN
1	A	473	GLN
1	B	54	GLN
1	B	77	HIS
1	B	84	ASN

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Mol	Chain	Res	Type
1	B	210	GLN
1	B	261	ASN
1	B	384	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 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	302	1	24,24,25	1.55	5 (20%)	28,32,34	1.73	6 (21%)
1	LLP	B	302	1	24,24,25	1.56	4 (16%)	28,32,34	1.44	5 (17%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	LLP	A	302	1	-	0/15/17/19	0/1/1/1
1	LLP	B	302	1	-	0/15/17/19	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	302	LLP	O3-C3	-4.81	1.25	1.37
1	A	302	LLP	O3-C3	-4.71	1.26	1.37
1	B	302	LLP	P-OP2	-2.32	1.45	1.54
1	A	302	LLP	P-OP2	-2.15	1.46	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	302	LLP	C6-N1	2.50	1.39	1.34
1	A	302	LLP	C4-C4'	2.50	1.51	1.46
1	A	302	LLP	C6-N1	2.58	1.40	1.34
1	A	302	LLP	C4'-NZ	2.89	1.35	1.27
1	B	302	LLP	C4-C4'	3.15	1.52	1.46

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	302	LLP	CE-NZ-C4'	-3.81	107.98	119.03
1	A	302	LLP	C4-C3-C2	-3.70	117.88	120.15
1	A	302	LLP	C5-C6-N1	-3.63	117.74	123.87
1	A	302	LLP	C5'-C5-C6	-2.90	114.35	119.33
1	B	302	LLP	C5-C6-N1	-2.61	119.46	123.87
1	B	302	LLP	C4-C4'-NZ	-2.55	112.26	124.66
1	A	302	LLP	C4-C4'-NZ	-2.45	112.78	124.66
1	B	302	LLP	C4-C3-C2	-2.27	118.76	120.15
1	B	302	LLP	OP2-P-OP4	-2.25	100.73	106.73
1	A	302	LLP	C6-C5-C4	2.10	122.33	118.20
1	B	302	LLP	OP3-P-OP2	3.21	120.55	107.61

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	302	LLP	2	0
1	B	302	LLP	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	GOL	A	476	-	5,5,5	0.23	0	5,5,5	0.56	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	GOL	A	476	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	476	GOL	3	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	447/475 (94%)	0.47	32 (7%) 16 22	14, 24, 46, 69	0
1	B	447/475 (94%)	0.73	52 (11%) 5 7	16, 29, 55, 75	0
All	All	894/950 (94%)	0.60	84 (9%) 9 12	14, 27, 50, 75	0

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	105	TRP	15.7
1	A	105	TRP	12.1
1	B	174	TRP	11.9
1	B	475	GLN	8.5
1	B	106	ILE	7.8
1	A	107	ALA	7.0
1	A	475	GLN	6.9
1	A	349	PRO	6.8
1	B	350	LEU	6.1
1	B	107	ALA	5.8
1	B	168	LYS	5.7
1	A	172	PRO	5.6
1	B	178	THR	5.6
1	A	174	TRP	5.3
1	B	173	GLU	5.3
1	A	106	ILE	5.2
1	B	352	ARG	5.1
1	B	104	THR	5.0
1	A	321	VAL	4.8
1	B	319	SER	4.7
1	A	104	THR	4.7
1	B	230	ALA	4.6
1	B	320	TRP	4.6
1	A	350	LEU	4.6

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Mol	Chain	Res	Type	RSRZ
1	B	175	ASP	4.6
1	B	101	ILE	4.5
1	A	438	LYS	4.3
1	B	167	VAL	4.3
1	B	103	PHE	4.3
1	A	352	ARG	4.2
1	A	439	ASP	4.1
1	B	164	LEU	4.0
1	B	171	HIS	3.9
1	B	232	GLY	3.8
1	A	32	GLU	3.8
1	B	349	PRO	3.8
1	B	79	TYR	3.7
1	B	32	GLU	3.5
1	B	139	LYS	3.3
1	A	474	GLU	3.1
1	A	168	LYS	3.1
1	B	318	PRO	3.1
1	B	102	GLY	3.1
1	A	79	TYR	3.0
1	B	179	ILE	3.0
1	B	28	ARG	2.9
1	B	170	LEU	2.9
1	B	180	LEU	2.9
1	B	321	VAL	2.9
1	A	169	GLU	2.8
1	B	172	PRO	2.8
1	A	320	TRP	2.8
1	B	231	GLU	2.8
1	A	101	ILE	2.7
1	A	35	PRO	2.7
1	A	170	LEU	2.7
1	A	103	PHE	2.6
1	B	200	LEU	2.6
1	A	28	ARG	2.6
1	B	108	SER	2.5
1	B	393	SER	2.5
1	B	439	ASP	2.5
1	B	285	HIS	2.4
1	A	178	THR	2.4
1	B	100	CYS	2.4
1	B	242	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	165	LYS	2.3
1	B	166	GLU	2.3
1	A	171	HIS	2.3
1	B	169	GLU	2.3
1	B	112	THR	2.3
1	A	44	ASP	2.3
1	B	58	GLN	2.2
1	B	228	ASP	2.2
1	B	226	GLU	2.2
1	A	165	LYS	2.2
1	A	58	GLN	2.1
1	B	306	VAL	2.1
1	A	108	SER	2.1
1	B	283	TYR	2.1
1	A	423	ARG	2.1
1	B	210	GLN	2.1
1	B	272	ALA	2.1
1	A	89	ILE	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	B	302	24/25	0.96	0.11	-	21,24,26,26	0
1	LLP	A	302	24/25	0.97	0.10	-	14,18,21,22	0

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(\AA^2)	Q<0.9
2	GOL	A	476	6/6	0.71	0.19	2.00	54,55,56,56	0

6.5 Other polymers [i](#)

There are no such residues in this entry.