



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:34 AM GMT

PDB ID : 3PD6  
Title : Crystal structure of mouse mitochondrial aspartate aminotransferase, a newly identified kynurenine aminotransferase-IV  
Authors : Han, Q.; Robinson, H.; Cai, T.; Tagle, D.A.; Li, J.  
Deposited on : 2010-10-22  
Resolution : 2.40 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 (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

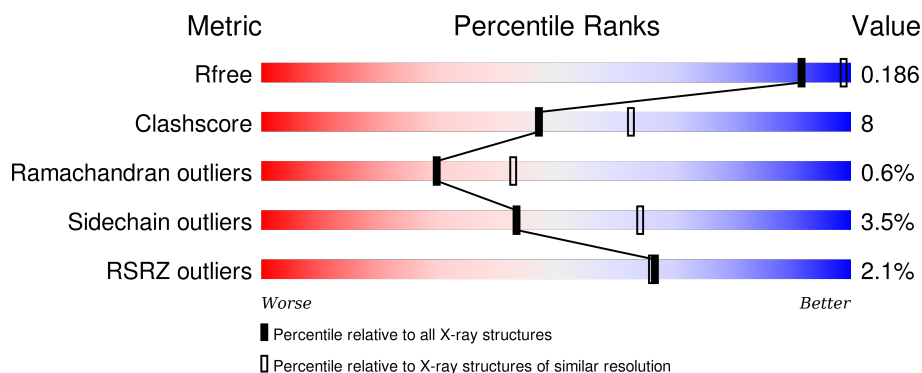
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	401	<div> <div>0%</div> <div> <div></div> <div>86%</div> <div>13%</div> </div> </div>
1	C	401	<div> <div></div> <div> <div></div> <div>86%</div> <div>13%</div> </div> </div>
2	B	401	<div> <div>2%</div> <div> <div></div> <div>81%</div> <div>18%</div> </div> </div>
2	D	401	<div> <div>5%</div> <div> <div></div> <div>80%</div> <div>19%</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
3	PMP	A	1	-	-	X	-
4	GOL	A	431	-	-	-	X
4	GOL	A	432	-	-	-	X
4	GOL	B	432	-	-	-	X
4	GOL	C	431	-	-	-	X
4	GOL	C	432	-	-	-	X
4	GOL	C	433	-	-	-	X
4	GOL	C	434	-	-	-	X
4	GOL	C	435	-	-	-	X
4	GOL	D	1	-	-	X	X
5	KYN	B	1	X	-	X	X



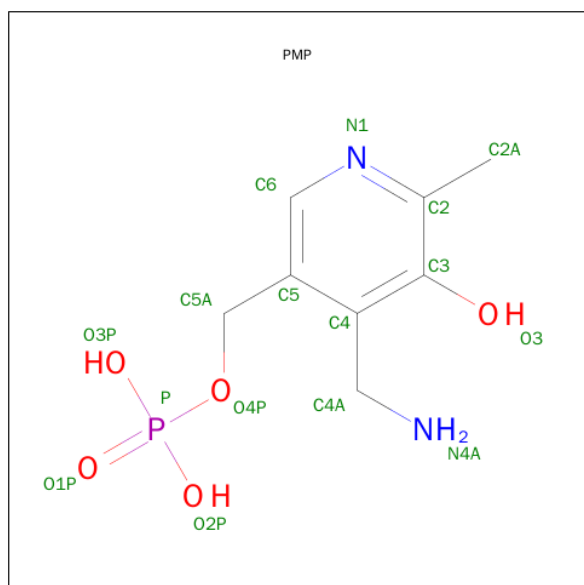
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 2 is a protein called Aspartate aminotransferase, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	401	Total 3140	C 1997	N 550	O 577	S 16	0	0	0
1	C	401	Total 3140	C 1997	N 550	O 577	S 16	0	0	0

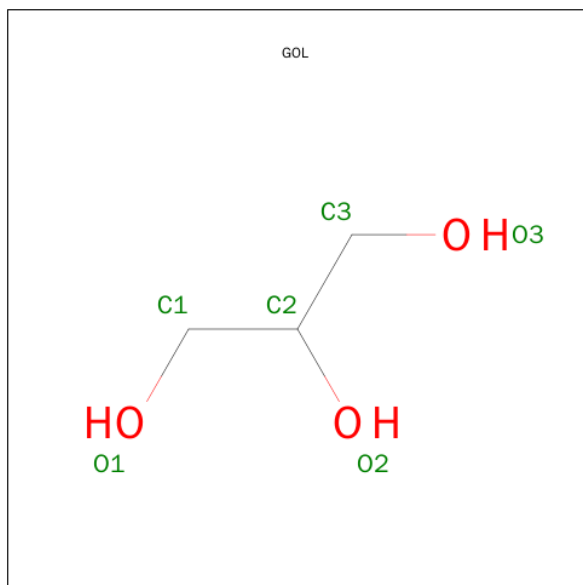
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	401	Total 3155	C 2005	N 551	O 582	P 1	S 16	0	0	0
2	D	401	Total 3155	C 2005	N 551	O 582	P 1	S 16	0	0	0

- Molecule 3 is 4'-DEOXY-4'-AMINOPYRIDOXAL-5'-PHOSPHATE (three-letter code: PMP) (formula:  $C_8H_{13}N_2O_5P$ ).



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

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



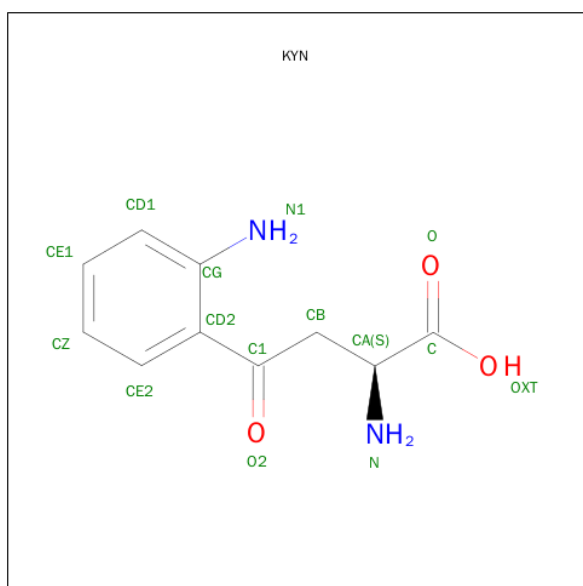
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total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

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

- Molecule 5 is (2S)-2-AMINO-4-(2-AMINOPHENYL)-4-OXOBUTANOIC ACID (three-letter code: KYN) (formula: C<sub>10</sub>H<sub>12</sub>N<sub>2</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	N	O	0	0
			15	10	2	3		

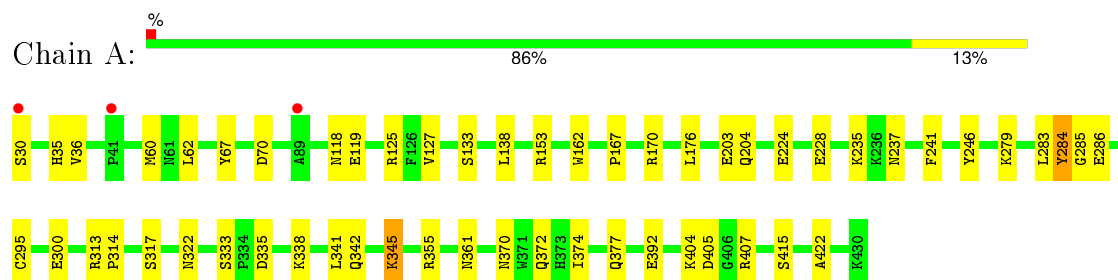
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	256	Total	O	0	0
			256	256		
6	B	172	Total	O	0	0
			172	172		
6	C	263	Total	O	0	0
			263	263		
6	D	140	Total	O	0	0
			140	140		

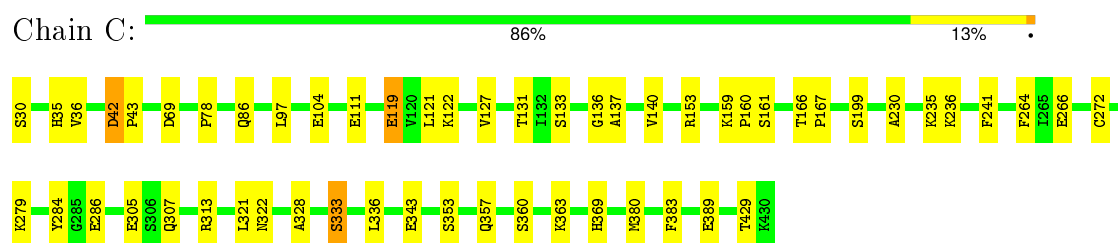
### 3 Residue-property plots

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

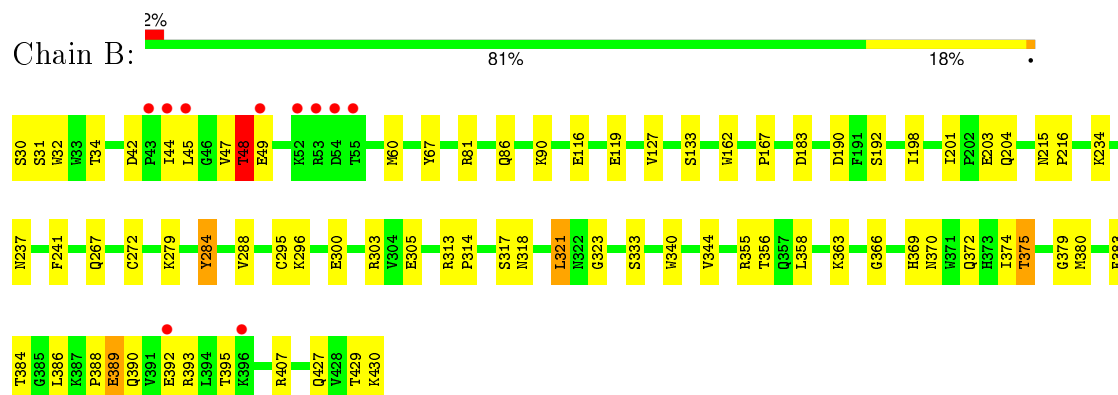
- Molecule 1: Aspartate aminotransferase, mitochondrial



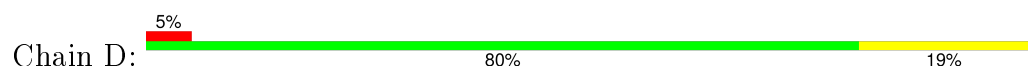
- Molecule 1: Aspartate aminotransferase, mitochondrial

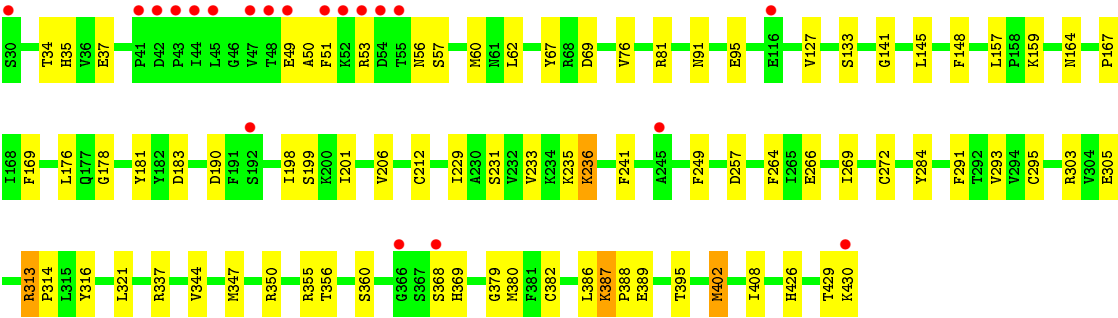


- Molecule 2: Aspartate aminotransferase, mitochondrial



- Molecule 2: Aspartate aminotransferase, mitochondrial







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	282.39 Å 78.11 Å 87.51 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.54 – 2.40 29.54 – 2.40	Depositor EDS
% Data completeness (in resolution range)	93.9 (29.54-2.40) 93.9 (29.54-2.40)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.59 (at 2.39 Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.177 , 0.189 0.173 , 0.186	Depositor DCC
$R_{free}$ test set	3603 reflections (5.28%)	DCC
Wilson B-factor (Å <sup>2</sup> )	27.6	Xtriage
Anisotropy	0.061	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 41.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 71871 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	13546	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.83% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.95	0/3213	0.81	0/4340
1	C	0.91	0/3213	0.81	1/4340 (0.0%)
2	B	0.87	2/3203 (0.1%)	0.81	0/4326
2	D	0.84	1/3203 (0.0%)	0.80	2/4326 (0.0%)
All	All	0.89	3/12832 (0.0%)	0.81	3/17332 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	284	TYR	CD2-CE2	-5.71	1.30	1.39
2	B	284	TYR	CD1-CE1	-5.42	1.31	1.39
2	D	95	GLU	CD-OE1	-5.30	1.19	1.25

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	337	ARG	NE-CZ-NH1	-5.88	117.36	120.30
1	C	69	ASP	CB-CG-OD1	5.87	123.58	118.30
2	D	321	LEU	CA-CB-CG	5.28	127.45	115.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3140	0	3125	41	0
1	C	3140	0	3125	33	0
2	B	3155	0	3130	55	0
2	D	3155	0	3130	60	0
3	A	16	0	10	8	0
3	C	16	0	10	1	0
4	A	18	0	24	4	0
4	B	18	0	24	5	0
4	C	36	0	48	4	0
4	D	6	0	8	6	0
5	B	15	0	11	15	0
6	A	256	0	0	10	0
6	B	172	0	0	4	0
6	C	263	0	0	4	0
6	D	140	0	0	2	0
All	All	13546	0	12645	191	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 8.

All (191) 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)
2:B:375:THR:HG23	6:B:591:HOH:O	1.59	1.02
3:A:1:PMP:HNA2	5:B:1:KYN:HB	1.20	1.01
3:A:1:PMP:N4A	5:B:1:KYN:HB	1.77	0.99
3:C:1:PMP:H4A1	6:C:786:HOH:O	1.68	0.94
2:B:318:ASN:HD21	5:B:1:KYN:HD1	1.33	0.93
2:B:318:ASN:ND2	5:B:1:KYN:HD1	1.85	0.92
3:A:1:PMP:H4A1	6:A:785:HOH:O	1.68	0.92
2:D:198:ILE:HA	2:D:201:ILE:HD12	1.51	0.91
2:D:402:MET:HE1	2:D:408:ILE:HG23	1.56	0.87
2:D:164:ASN:O	2:D:167:PRO:HD2	1.76	0.86
2:D:350:ARG:HD3	4:D:1:GOL:H31	1.59	0.85
2:D:387:LYS:HB2	2:D:387:LYS:HZ3	1.41	0.85
1:C:119:GLU:HG2	6:C:497:HOH:O	1.79	0.82
1:C:357:GLN:HG2	6:C:553:HOH:O	1.82	0.79
1:C:279:LYS:HE2	1:C:380:MET:HE3	1.63	0.79
2:D:402:MET:CE	2:D:408:ILE:CG2	2.61	0.79
2:B:318:ASN:HD21	5:B:1:KYN:CD1	1.97	0.77
1:A:162:TRP:HE1	5:B:1:KYN:HBA	1.50	0.76
2:B:67:TYR:OH	4:B:432:GOL:H12	1.85	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:402:MET:HE1	2:D:408:ILE:CG2	2.15	0.75
2:D:426:HIS:CE1	2:D:430:LYS:HB3	2.23	0.73
3:A:1:PMP:HNA2	5:B:1:KYN:CB	2.01	0.73
1:A:170:ARG:HD3	6:A:601:HOH:O	1.89	0.71
2:B:198:ILE:HA	2:B:201:ILE:HD12	1.73	0.70
3:A:1:PMP:N4A	5:B:1:KYN:HE2	2.07	0.70
1:A:246:TYR:CE1	1:A:279:LYS:HG3	2.28	0.69
2:D:313:ARG:CB	2:D:314:PRO:HD3	2.22	0.68
2:D:387:LYS:NZ	2:D:387:LYS:HB2	2.08	0.68
2:B:318:ASN:ND2	5:B:1:KYN:CD1	2.54	0.68
1:C:313:ARG:HD3	6:C:611:HOH:O	1.95	0.67
1:C:127:VAL:HG21	1:C:305:GLU:HB2	1.78	0.66
1:C:30:SER:O	4:C:435:GOL:H12	1.95	0.66
2:D:402:MET:HE2	2:D:408:ILE:HG22	1.79	0.65
1:C:279:LYS:CE	1:C:380:MET:HE3	2.27	0.64
1:A:370:ASN:HD21	1:A:372:GLN:HE21	1.43	0.64
1:A:237:ASN:OD1	4:A:433:GOL:H12	1.97	0.64
2:B:375:THR:CG2	6:B:591:HOH:O	2.31	0.63
2:D:60:MET:HE2	2:D:62:LEU:HD21	1.81	0.63
2:D:69:ASP:HB2	4:D:1:GOL:H32	1.81	0.63
1:A:67:TYR:OH	4:A:431:GOL:H12	2.00	0.62
1:C:199:SER:O	1:C:236:LYS:HE2	1.99	0.62
2:D:231:SER:O	2:D:235:LYS:HG2	2.00	0.61
2:B:183:ASP:HB2	2:B:190:ASP:HB2	1.82	0.61
1:A:203:GLU:HG3	1:A:204:GLN:HG3	1.81	0.61
1:C:35:HIS:CD2	1:C:35:HIS:H	2.17	0.61
1:C:307:GLN:NE2	2:D:37:GLU:O	2.31	0.60
2:B:369:HIS:HE1	2:B:429:THR:O	1.85	0.59
2:B:234:LYS:HD2	2:B:267:GLN:O	2.02	0.59
2:D:34:THR:HG22	2:D:35:HIS:HD2	1.66	0.59
2:B:370:ASN:HD21	2:B:372:GLN:HE21	1.52	0.58
2:D:389:GLU:OE2	2:D:389:GLU:N	2.29	0.57
1:C:97:LEU:HG	1:C:321:LEU:HD13	1.86	0.57
2:B:389:GLU:H	2:B:389:GLU:CD	2.08	0.57
2:D:60:MET:CE	2:D:62:LEU:HD21	2.35	0.56
1:C:104:GLU:HG2	1:C:328:ALA:HB1	1.86	0.56
1:A:355:ARG:HD3	1:A:374:ILE:O	2.06	0.55
1:A:370:ASN:OD1	1:A:372:GLN:HG2	2.07	0.55
2:D:382:CYS:HB3	2:D:408:ILE:HG13	1.88	0.55
2:B:67:TYR:OH	4:B:432:GOL:C1	2.52	0.55
2:D:183:ASP:HB2	2:D:190:ASP:HB2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:342:GLN:HG2	6:A:540:HOH:O	2.06	0.55
1:A:246:TYR:HE1	1:A:279:LYS:HG3	1.70	0.55
1:A:279:LYS:HD3	1:A:284:TYR:CE1	2.42	0.54
1:C:166:THR:HB	1:C:167:PRO:CD	2.37	0.54
2:B:241:PHE:CD1	2:B:272:CYS:SG	3.00	0.54
2:D:387:LYS:HB3	2:D:389:GLU:OE2	2.08	0.54
1:C:136:GLY:O	1:C:140:VAL:HG23	2.08	0.53
1:C:153:ARG:HA	1:C:153:ARG:HE	1.74	0.53
1:C:333:SER:OG	1:C:336:LEU:HB2	2.08	0.53
3:A:1:PMP:C4A	6:A:785:HOH:O	2.39	0.53
2:D:69:ASP:CB	4:D:1:GOL:H32	2.38	0.53
2:B:313:ARG:HB3	2:B:314:PRO:HD3	1.90	0.53
2:B:389:GLU:OE2	2:B:389:GLU:N	2.39	0.52
2:B:295:CYS:HB3	2:B:300:GLU:OE1	2.09	0.52
2:D:402:MET:HE2	2:D:408:ILE:CG2	2.32	0.52
1:C:161:SER:H	4:C:432:GOL:H2	1.74	0.52
4:C:434:GOL:H12	2:D:76:VAL:HG23	1.90	0.52
1:A:153:ARG:HD2	6:A:668:HOH:O	2.10	0.52
1:A:118:ASN:HA	6:A:507:HOH:O	2.10	0.52
2:D:355:ARG:NH2	2:D:379:GLY:O	2.43	0.52
1:C:78:PRO:HB2	1:C:343:GLU:OE1	2.09	0.52
2:D:313:ARG:HB3	2:D:314:PRO:HD3	1.89	0.52
2:B:45:LEU:HA	2:B:48:THR:HG22	1.92	0.52
2:B:31:SER:HB3	2:B:34:THR:HG23	1.93	0.50
3:A:1:PMP:P	5:B:1:KYN:HZ	2.51	0.50
2:B:31:SER:HB3	2:B:34:THR:CG2	2.42	0.50
1:C:35:HIS:HE1	2:D:148:PHE:O	1.94	0.50
4:A:433:GOL:H31	2:B:32:TRP:CE3	2.47	0.49
2:D:50:ALA:HA	2:D:53:ARG:NH2	2.28	0.49
2:D:49:GLU:HG2	2:D:53:ARG:HH12	1.78	0.49
2:B:215:ASN:OD1	2:B:216:PRO:HA	2.11	0.49
2:D:388:PRO:HD3	6:D:462:HOH:O	2.13	0.49
2:B:318:ASN:HD21	5:B:1:KYN:CE1	2.25	0.49
1:C:389:GLU:H	1:C:389:GLU:CD	2.16	0.49
2:D:127:VAL:HG21	2:D:305:GLU:HB2	1.95	0.49
5:B:1:KYN:HE2	5:B:1:KYN:HB	1.68	0.49
1:A:279:LYS:HD3	1:A:284:TYR:HE1	1.77	0.49
2:D:91:ASN:HB2	6:D:599:HOH:O	2.12	0.49
2:B:47:VAL:C	2:B:49:GLU:H	2.15	0.49
2:B:237:ASN:ND2	4:B:433:GOL:O2	2.42	0.48
1:A:167:PRO:HB2	2:B:314:PRO:HG3	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:145:LEU:HD13	2:D:206:VAL:HG21	1.94	0.48
2:B:81:ARG:O	6:B:504:HOH:O	2.20	0.48
2:D:293:VAL:HG12	2:D:295:CYS:SG	2.53	0.48
2:D:199:SER:O	2:D:236:LYS:HD2	2.14	0.48
1:C:286:GLU:CD	1:C:322:ASN:ND2	2.67	0.48
2:B:390:GLN:HE22	2:B:429:THR:HG22	1.78	0.47
2:B:162:TRP:CD1	6:B:747:HOH:O	2.67	0.47
1:A:224:GLU:OE1	1:A:228:GLU:OE2	2.31	0.47
1:A:285:GLY:HA3	2:B:321:LEU:HD22	1.96	0.47
2:D:402:MET:CE	2:D:408:ILE:HG22	2.37	0.47
2:D:169:PHE:HB2	2:D:176:LEU:HD11	1.97	0.47
2:B:384:THR:OG1	2:B:386:LEU:HD12	2.15	0.47
2:D:241:PHE:CD1	2:D:272:CYS:SG	3.07	0.47
1:A:162:TRP:HE1	5:B:1:KYN:CB	2.25	0.46
2:D:76:VAL:HG21	2:D:81:ARG:NH2	2.30	0.46
1:A:119:GLU:HB2	6:A:479:HOH:O	2.14	0.46
2:B:234:LYS:CD	2:B:267:GLN:O	2.64	0.46
2:D:157:LEU:O	2:D:178:GLY:HA2	2.16	0.46
2:D:387:LYS:HZ3	2:D:387:LYS:CB	2.17	0.45
1:A:176:LEU:HD22	1:A:176:LEU:N	2.31	0.45
1:A:60:MET:CE	1:A:62:LEU:HD21	2.46	0.45
1:A:138:LEU:HD23	1:A:241:PHE:HE2	1.80	0.45
1:A:341:LEU:HD23	1:A:341:LEU:HA	1.84	0.45
1:C:266:GLU:O	1:C:266:GLU:HG2	2.15	0.45
1:C:230:ALA:HB2	1:C:264:PHE:CE1	2.52	0.45
2:B:215:ASN:HD21	4:B:431:GOL:H32	1.81	0.45
1:C:166:THR:HG1	4:C:432:GOL:C3	2.30	0.45
1:A:286:GLU:CD	1:A:322:ASN:HD22	2.20	0.45
2:D:313:ARG:CB	2:D:314:PRO:CD	2.95	0.45
1:C:36:VAL:O	2:D:303:ARG:HD2	2.17	0.45
1:A:313:ARG:HG3	1:A:313:ARG:HH11	1.82	0.44
1:A:392:GLU:HG3	1:A:404:LYS:HE2	1.99	0.44
2:B:389:GLU:O	2:B:393:ARG:HG3	2.18	0.44
2:D:49:GLU:HG2	2:D:53:ARG:NH1	2.31	0.44
2:B:388:PRO:O	2:B:392:GLU:HG3	2.18	0.44
2:D:344:VAL:HA	2:D:347:MET:HE3	2.00	0.44
2:B:313:ARG:HA	2:B:317:SER:HA	2.00	0.44
2:D:56:ASN:OD1	2:D:56:ASN:C	2.56	0.44
1:C:86:GLN:HE21	1:C:86:GLN:HB3	1.53	0.43
2:B:127:VAL:HG21	2:B:305:GLU:HB2	1.99	0.43
2:B:358:LEU:HD23	2:B:374:ILE:HG21	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:69:ASP:HA	4:D:1:GOL:H32	2.00	0.43
2:D:249:PHE:CZ	2:D:380:MET:CE	3.02	0.43
2:D:369:HIS:HE1	2:D:429:THR:O	2.02	0.43
1:C:369:HIS:HE1	1:C:429:THR:O	2.02	0.43
1:C:241:PHE:CD1	1:C:272:CYS:SG	3.12	0.43
1:C:131:THR:HG21	1:C:137:ALA:HA	2.01	0.43
2:B:427:GLN:O	2:B:430:LYS:HG2	2.19	0.43
1:A:407:ARG:NH1	5:B:1:KYN:O	2.49	0.43
2:B:198:ILE:HA	2:B:201:ILE:CD1	2.47	0.42
1:C:35:HIS:CE1	2:D:148:PHE:O	2.71	0.42
2:D:159:LYS:O	2:D:181:TYR:HB3	2.18	0.42
2:B:369:HIS:CE1	2:B:429:THR:O	2.68	0.42
2:B:288:VAL:HG22	2:B:323:GLY:HA3	2.01	0.42
2:D:141:GLY:HA2	2:D:291:PHE:CZ	2.54	0.42
1:C:42:ASP:HA	1:C:43:PRO:HD3	1.81	0.42
2:D:212:CYS:SG	2:D:257:ASP:HB3	2.60	0.42
2:B:372:GLN:O	2:B:375:THR:HG22	2.20	0.42
1:A:283:LEU:O	1:A:284:TYR:C	2.58	0.42
2:B:288:VAL:CG2	2:B:323:GLY:HA3	2.49	0.42
2:B:363:LYS:O	2:B:366:GLY:N	2.52	0.42
1:A:36:VAL:O	2:B:303:ARG:HD2	2.19	0.42
3:A:1:PMP:N4A	5:B:1:KYN:CB	2.66	0.41
1:A:67:TYR:OH	4:A:431:GOL:C1	2.66	0.41
1:A:35:HIS:CD2	1:A:35:HIS:H	2.38	0.41
2:B:203:GLU:O	2:B:204:GLN:HB2	2.21	0.41
2:D:229:ILE:HG22	2:D:264:PHE:HZ	1.85	0.41
2:D:67:TYR:OH	4:D:1:GOL:H12	2.20	0.41
2:D:344:VAL:HA	2:D:347:MET:CE	2.50	0.41
2:B:279:LLP:OP4	2:B:279:LLP:H4'2	2.20	0.41
2:B:355:ARG:NH2	2:B:379:GLY:O	2.30	0.41
1:C:111:GLU:HA	1:C:121:LEU:HD11	2.02	0.41
2:D:233:VAL:CG1	2:D:269:ILE:HD13	2.50	0.41
1:C:159:LYS:HA	1:C:160:PRO:HA	1.74	0.41
1:A:153:ARG:NH1	6:A:668:HOH:O	2.39	0.41
2:B:86:GLN:O	2:B:90:LYS:HG3	2.21	0.41
2:D:386:LEU:HA	2:D:386:LEU:HD23	1.94	0.41
1:A:295:CYS:HB3	1:A:300:GLU:OE1	2.21	0.41
2:B:340:TRP:O	2:B:344:VAL:HG23	2.21	0.41
1:A:314:PRO:CG	2:B:167:PRO:HB2	2.51	0.41
1:A:345:LYS:HE3	6:A:761:HOH:O	2.20	0.41
2:B:162:TRP:HE1	4:B:431:GOL:H31	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:56:ASN:OD1	2:D:57:SER:N	2.54	0.40
1:A:338:LYS:NZ	6:A:580:HOH:O	2.54	0.40
2:B:47:VAL:C	2:B:49:GLU:N	2.75	0.40
1:A:361:ASN:HB3	1:A:422:ALA:CB	2.51	0.40
2:D:69:ASP:HB2	4:D:1:GOL:C3	2.49	0.40
1:A:355:ARG:NH1	1:A:377:GLN:HB2	2.36	0.40
1:A:125:ARG:O	1:A:125:ARG:HG3	2.21	0.40
1:A:405:ASP:OD1	1:A:405:ASP:N	2.54	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	399/401 (100%)	388 (97%)	9 (2%)	2 (0%)	34	48
1	C	399/401 (100%)	385 (96%)	13 (3%)	1 (0%)	46	63
2	B	398/401 (99%)	379 (95%)	15 (4%)	4 (1%)	19	28
2	D	398/401 (99%)	376 (94%)	20 (5%)	2 (0%)	34	48
All	All	1594/1604 (99%)	1528 (96%)	57 (4%)	9 (1%)	30	43

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	284	TYR
1	C	284	TYR
2	D	284	TYR
1	A	317	SER
2	B	44	ILE
2	B	48	THR
2	B	119	GLU

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Mol	Chain	Res	Type
2	B	284	TYR
2	D	316	TYR

### 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	334/334 (100%)	325 (97%)	9 (3%)	52	73
1	C	334/334 (100%)	324 (97%)	10 (3%)	48	70
2	B	333/333 (100%)	316 (95%)	17 (5%)	29	46
2	D	333/333 (100%)	322 (97%)	11 (3%)	45	66
All	All	1334/1334 (100%)	1287 (96%)	47 (4%)	43	64

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	SER
1	A	70	ASP
1	A	127	VAL
1	A	133	SER
1	A	235	LYS
1	A	333	SER
1	A	335	ASP
1	A	345	LYS
1	A	415	SER
2	B	30	SER
2	B	42	ASP
2	B	48	THR
2	B	60	MET
2	B	116	GLU
2	B	133	SER
2	B	192	SER
2	B	296	LYS
2	B	321	LEU
2	B	333	SER

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Mol	Chain	Res	Type
2	B	356	THR
2	B	375	THR
2	B	380	MET
2	B	383	PHE
2	B	389	GLU
2	B	395	THR
2	B	407	ARG
1	C	42	ASP
1	C	119	GLU
1	C	122	LYS
1	C	133	SER
1	C	235	LYS
1	C	333	SER
1	C	353	SER
1	C	360	SER
1	C	363	LYS
1	C	383	PHE
2	D	51	PHE
2	D	133	SER
2	D	236	LYS
2	D	266	GLU
2	D	313	ARG
2	D	356	THR
2	D	360	SER
2	D	368	SER
2	D	387	LYS
2	D	395	THR
2	D	402	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	HIS
1	A	117	ASN
1	A	204	GLN
1	A	275	GLN
1	A	322	ASN
1	A	369	HIS
1	A	372	GLN
2	B	86	GLN
2	B	91	ASN
2	B	237	ASN

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Mol	Chain	Res	Type
2	B	318	ASN
2	B	361	ASN
2	B	369	HIS
2	B	372	GLN
1	C	35	HIS
1	C	86	GLN
1	C	322	ASN
1	C	361	ASN
1	C	369	HIS
2	D	35	HIS
2	D	86	GLN
2	D	91	ASN
2	D	237	ASN
2	D	369	HIS
2	D	372	GLN
2	D	426	HIS
2	D	427	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	$\# Z  > 2$
2	LLP	B	279	2	23,24,25	3.50	11 (47%)	28,32,34	2.84	8 (28%)
2	LLP	D	279	2	23,24,25	2.66	9 (39%)	28,32,34	2.99	9 (32%)

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	LLP	B	279	2	-	0/15/17/19	0/1/1/1
2	LLP	D	279	2	-	0/15/17/19	0/1/1/1

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	279	LLP	O3-C3	-7.58	1.19	1.37
2	D	279	LLP	O3-C3	-7.33	1.19	1.37
2	B	279	LLP	P-OP3	-5.19	1.36	1.54
2	B	279	LLP	P-OP2	-5.08	1.36	1.54
2	D	279	LLP	P-OP3	-4.45	1.38	1.54
2	B	279	LLP	C3-C2	-4.44	1.37	1.40
2	D	279	LLP	P-OP2	-4.31	1.39	1.54
2	D	279	LLP	P-OP1	-4.14	1.37	1.51
2	B	279	LLP	P-OP1	-4.10	1.37	1.51
2	B	279	LLP	P-OP4	-3.55	1.48	1.60
2	D	279	LLP	C3-C2	-3.26	1.38	1.40
2	D	279	LLP	P-OP4	-2.86	1.50	1.60
2	B	279	LLP	C4-C3	-2.66	1.37	1.40
2	D	279	LLP	C4-C3	-2.42	1.37	1.40
2	D	279	LLP	CE-NZ	2.65	1.52	1.46
2	D	279	LLP	C4-C4'	3.04	1.52	1.46
2	B	279	LLP	C4'-NZ	3.23	1.37	1.27
2	B	279	LLP	CD-CE	3.93	1.63	1.51
2	B	279	LLP	C4-C4'	4.69	1.54	1.46
2	B	279	LLP	CE-NZ	7.45	1.62	1.46

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	279	LLP	C3-C4-C5	-6.17	113.49	118.11
2	D	279	LLP	OP4-P-OP1	-5.57	92.97	107.14
2	D	279	LLP	C3-C4-C5	-4.02	115.10	118.11
2	D	279	LLP	C5'-C5-C6	-3.29	113.06	119.28
2	B	279	LLP	OP2-P-OP4	-3.11	97.60	106.56
2	B	279	LLP	O-C-CA	-3.07	117.48	125.49
2	B	279	LLP	C5'-C5-C6	-2.77	114.05	119.28
2	D	279	LLP	O-C-CA	-2.61	118.70	125.49
2	D	279	LLP	OP3-P-OP4	-2.40	99.67	106.56
2	B	279	LLP	C6-C5-C4	2.20	123.04	118.17
2	D	279	LLP	OP3-P-OP2	2.25	115.93	107.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	279	LLP	OP3-P-OP1	2.34	118.10	110.58
2	D	279	LLP	C5-C4-C4'	2.40	124.97	121.52
2	B	279	LLP	OP3-P-OP2	3.41	120.35	107.38
2	B	279	LLP	OP4-C5'-C5	4.90	117.10	108.99
2	B	279	LLP	CD-CE-NZ	9.96	127.29	110.98
2	D	279	LLP	OP4-C5'-C5	11.88	128.63	108.99

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	279	LLP	1	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

16 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	PMP	A	1	-	16,16,16	1.07	1 (6%)	20,23,23	2.16	7 (35%)
4	GOL	A	431	-	5,5,5	0.27	0	5,5,5	0.81	0
4	GOL	A	432	-	5,5,5	0.28	0	5,5,5	0.34	0
4	GOL	A	433	-	5,5,5	0.53	0	5,5,5	0.76	0
5	KYN	B	1	-	12,15,15	1.44	2 (16%)	11,20,20	1.18	1 (9%)
4	GOL	B	431	-	5,5,5	0.49	0	5,5,5	0.88	0
4	GOL	B	432	-	5,5,5	0.65	0	5,5,5	0.71	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	GOL	B	433	-	5,5,5	0.33	0	5,5,5	0.41	0
3	PMP	C	1	-	16,16,16	2.92	7 (43%)	20,23,23	1.44	3 (15%)
4	GOL	C	431	-	5,5,5	0.34	0	5,5,5	0.38	0
4	GOL	C	432	-	5,5,5	0.45	0	5,5,5	1.08	1 (20%)
4	GOL	C	433	-	5,5,5	1.04	1 (20%)	5,5,5	1.19	0
4	GOL	C	434	-	5,5,5	0.30	0	5,5,5	0.94	0
4	GOL	C	435	-	5,5,5	0.49	0	5,5,5	0.70	0
4	GOL	C	436	-	5,5,5	0.31	0	5,5,5	0.25	0
4	GOL	D	1	-	5,5,5	0.52	0	5,5,5	0.49	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	PMP	A	1	-	-	0/8/8/8	0/1/1/1
4	GOL	A	431	-	-	0/4/4/4	0/0/0/0
4	GOL	A	432	-	-	0/4/4/4	0/0/0/0
4	GOL	A	433	-	-	0/4/4/4	0/0/0/0
5	KYN	B	1	-	1/1/3/3	0/8/12/12	0/1/1/1
4	GOL	B	431	-	-	0/4/4/4	0/0/0/0
4	GOL	B	432	-	-	0/4/4/4	0/0/0/0
4	GOL	B	433	-	-	0/4/4/4	0/0/0/0
3	PMP	C	1	-	-	0/8/8/8	0/1/1/1
4	GOL	C	431	-	-	0/4/4/4	0/0/0/0
4	GOL	C	432	-	-	0/4/4/4	0/0/0/0
4	GOL	C	433	-	-	0/4/4/4	0/0/0/0
4	GOL	C	434	-	-	0/4/4/4	0/0/0/0
4	GOL	C	435	-	-	0/4/4/4	0/0/0/0
4	GOL	C	436	-	-	0/4/4/4	0/0/0/0
4	GOL	D	1	-	-	0/4/4/4	0/0/0/0

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1	PMP	C3-C2	-7.25	1.35	1.40
3	C	1	PMP	P-O3P	-4.51	1.38	1.54
3	C	1	PMP	P-O2P	-4.20	1.39	1.54
3	C	1	PMP	P-O1P	-3.84	1.38	1.51
5	B	1	KYN	CD2-CG	-3.58	1.36	1.41
3	C	1	PMP	C5-C4	-3.04	1.36	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1	PMP	C3-C4	-2.72	1.35	1.40
5	B	1	KYN	CB-C1	-2.64	1.47	1.51
4	C	433	GOL	O2-C2	-2.14	1.37	1.43
3	C	1	PMP	O3-C3	-2.10	1.32	1.37
3	A	1	PMP	C6-N1	2.17	1.39	1.34

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1	PMP	C4A-C4-C3	-4.52	113.59	120.40
5	B	1	KYN	O2-C1-CB	-2.30	118.18	120.75
3	A	1	PMP	C3-C2-N1	-2.21	117.56	120.61
3	A	1	PMP	C5-C6-N1	-2.07	120.26	123.86
3	C	1	PMP	O4P-P-O1P	2.14	112.60	107.14
4	C	432	GOL	O3-C3-C2	2.27	121.22	110.18
3	A	1	PMP	C3-C4-C5	2.57	121.52	118.82
3	A	1	PMP	O3-C3-C2	2.66	122.29	117.66
3	C	1	PMP	O2P-P-O4P	2.81	114.66	106.56
3	A	1	PMP	O3P-P-O4P	4.19	118.63	106.56
3	A	1	PMP	O4P-C5A-C5	4.61	116.61	108.99
3	C	1	PMP	O4P-C5A-C5	4.64	116.66	108.99

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	B	1	KYN	CA

There are no torsion outliers.

There are no ring outliers.

12 monomers are involved in 37 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1	PMP	8	0
4	A	431	GOL	2	0
4	A	433	GOL	2	0
5	B	1	KYN	15	0
4	B	431	GOL	2	0
4	B	432	GOL	2	0
4	B	433	GOL	1	0
3	C	1	PMP	1	0
4	C	432	GOL	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	434	GOL	1	0
4	C	435	GOL	1	0
4	D	1	GOL	6	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	401/401 (100%)	-0.57	3 (0%) 89 88	11, 19, 33, 44	0
1	C	401/401 (100%)	-0.57	0 100 100	12, 22, 37, 46	0
2	B	400/401 (99%)	-0.29	10 (2%) 61 60	14, 31, 64, 91	0
2	D	400/401 (99%)	-0.03	20 (5%) 32 33	16, 35, 70, 89	0
All	All	1602/1604 (99%)	-0.37	33 (2%) 67 66	11, 26, 55, 91	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	44	ILE	4.8
2	D	41	PRO	3.7
2	B	44	ILE	3.6
2	D	55	THR	3.5
2	B	53	ARG	3.5
2	D	52	LYS	3.4
2	D	45	LEU	3.4
2	D	42	ASP	3.4
2	B	45	LEU	3.3
2	D	30	SER	3.3
2	B	49	GLU	3.2
2	D	53	ARG	3.1
2	D	49	GLU	3.1
2	D	43	PRO	2.9
2	D	48	THR	2.8
2	B	55	THR	2.8
2	D	47	VAL	2.7
1	A	30	SER	2.7
2	D	430	LYS	2.5
2	D	366	GLY	2.4
1	A	41	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
2	D	54	ASP	2.3
2	D	192	SER	2.3
2	B	43	PRO	2.3
2	B	396	LYS	2.2
2	B	54	ASP	2.2
2	D	245	ALA	2.2
2	D	368	SER	2.2
1	A	89	ALA	2.1
2	B	52	LYS	2.0
2	D	51	PHE	2.0
2	B	392	GLU	2.0
2	D	116	GLU	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, 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	LLP	B	279	24/25	0.94	0.15	-	23,28,35,39	0
2	LLP	D	279	24/25	0.95	0.17	-	27,39,45,46	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, 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
4	GOL	C	432	6/6	0.94	0.26	7.73	20,26,34,36	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
4	GOL	C	434	6/6	0.71	0.35	7.62	54,56,57,58	0
5	KYN	B	1	15/15	0.83	0.30	5.83	62,64,65,67	0
4	GOL	C	433	6/6	0.92	0.28	4.28	48,49,49,51	0
4	GOL	A	432	6/6	0.82	0.17	3.34	59,59,60,60	0
4	GOL	A	431	6/6	0.87	0.15	2.98	48,49,49,51	0
4	GOL	C	435	6/6	0.82	0.28	2.93	61,62,63,63	0
4	GOL	C	431	6/6	0.90	0.15	2.48	25,30,32,37	0
4	GOL	B	432	6/6	0.96	0.15	2.26	38,39,40,43	0
4	GOL	D	1	6/6	0.84	0.15	2.20	40,43,44,45	0
4	GOL	C	436	6/6	0.79	0.26	1.19	59,61,61,61	0
4	GOL	B	431	6/6	0.85	0.22	0.61	48,49,49,51	0
4	GOL	B	433	6/6	0.91	0.20	0.56	46,47,47,48	0
4	GOL	A	433	6/6	0.92	0.12	0.03	40,42,43,45	0
3	PMP	A	1	16/16	0.98	0.13	-0.56	14,20,30,36	0
3	PMP	C	1	16/16	0.97	0.13	-0.70	17,25,36,43	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.