



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:08 AM GMT

PDB ID : 3L0Z
Title : Crystal structure of a putative Nicotinate-nucleotide-dimethylbenzimidazole phosphoribosyltransferase from *Methanocaldococcus jannaschii* DSM 2661
Authors : Nocek, B.; Hatzos, C.; Clancy, S.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2009-12-10
Resolution : 2.65 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

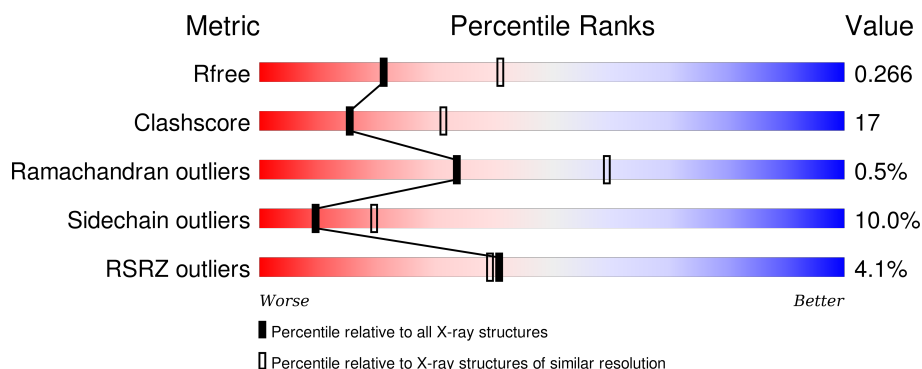
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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	3152 (2.70-2.62)
Clashscore	102246	3524 (2.70-2.62)
Ramachandran outliers	100387	3469 (2.70-2.62)
Sidechain outliers	100360	3469 (2.70-2.62)
RSRZ outliers	91569	3161 (2.70-2.62)

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	350	<div> <div>2%</div> <div>74%</div> <div>23%</div> <div>..</div> </div>
1	B	350	<div> <div>6%</div> <div>63%</div> <div>28%</div> <div>7%</div> <div>.</div> </div>
1	C	350	<div> <div>4%</div> <div>63%</div> <div>31%</div> <div>..</div> </div>

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PO4	B	351	-	-	X	-
2	PO4	C	352	-	-	-	X

2 Entry composition [i](#)

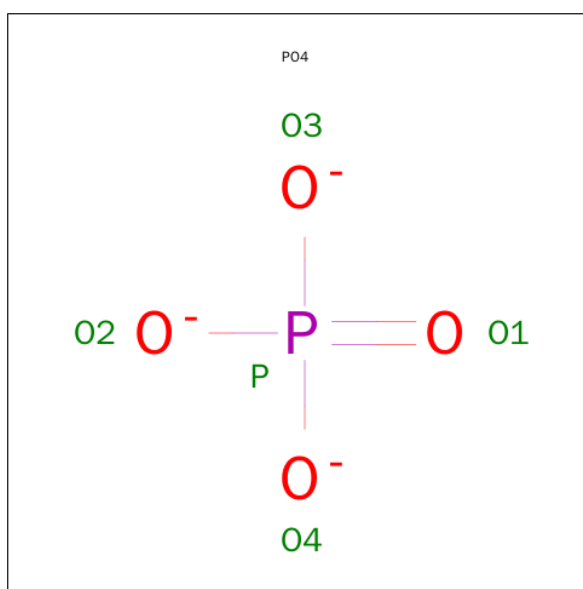
There are 3 unique types of molecules in this entry. The entry contains 7656 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 putative Nicotinate-nucleotide-dimethylbenzimidazole phosphoribosyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	346	Total	C	N	O	S	Se	0	0	0
			2606	1673	431	493	3	6			
1	B	342	Total	C	N	O	S	Se	0	0	0
			2542	1633	418	483	3	5			
1	C	340	Total	C	N	O	S	Se	0	0	0
			2464	1578	405	473	3	5			

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		

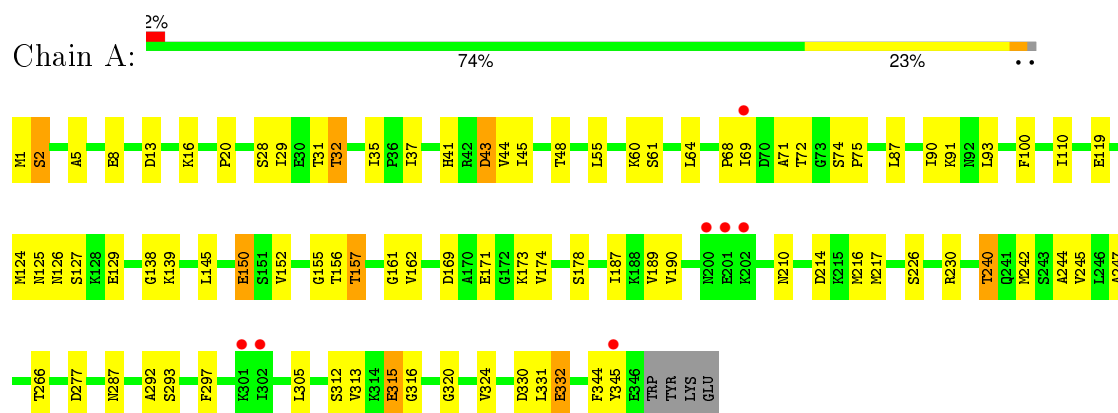
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	4	Total	O	0	0
			4	4		
3	B	3	Total	O	0	0
			3	3		
3	C	2	Total	O	0	0
			2	2		

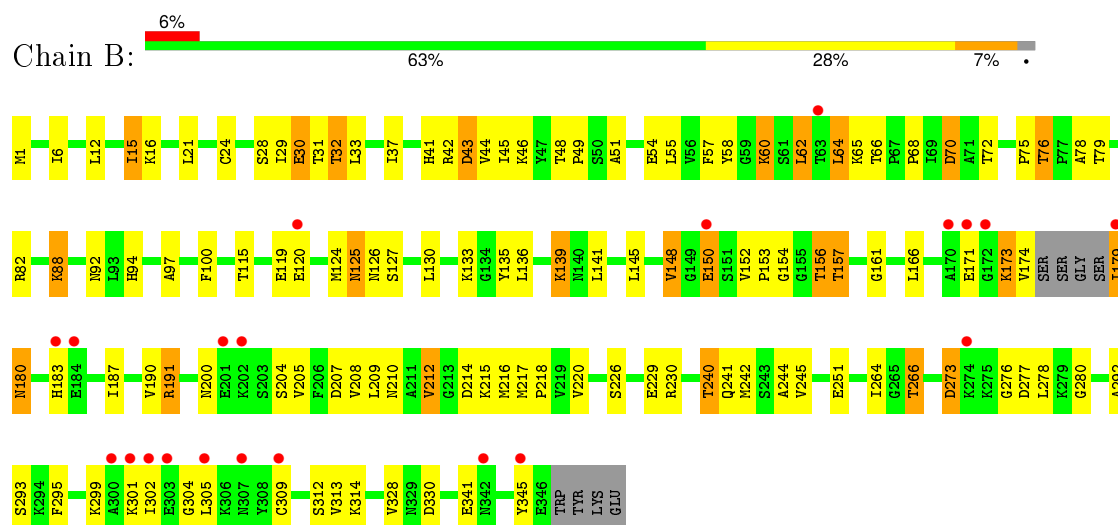
3 Residue-property plots

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: putative Nicotinate-nucleotide-dimethylbenzimidazole phosphoribosyltransferase

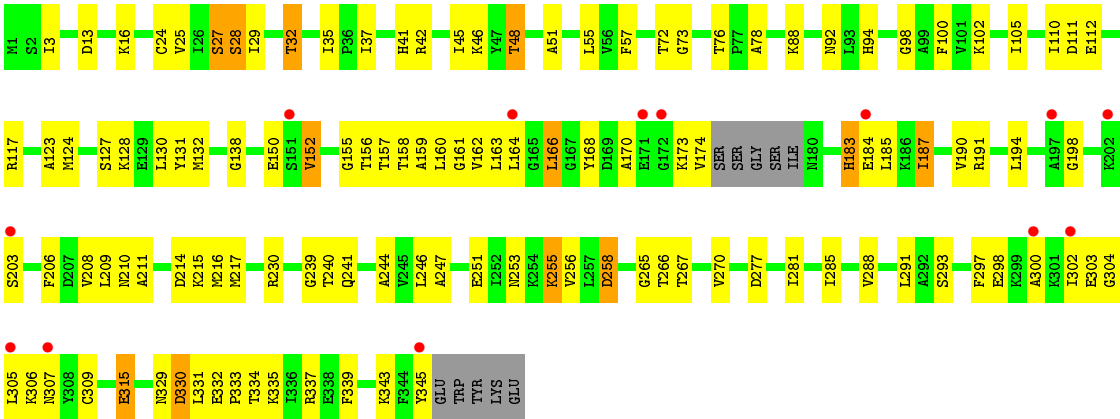


- Molecule 1: putative Nicotinate-nucleotide-dimethylbenzimidazole phosphoribosyltransferase



- Molecule 1: putative Nicotinate-nucleotide-dimethylbenzimidazole phosphoribosyltransferase





4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	119.31Å 119.31Å 170.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.00 – 2.65 39.05 – 2.65	Depositor EDS
% Data completeness (in resolution range)	99.4 (40.00-2.65) 99.4 (39.05-2.65)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.26 (at 2.65Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.202 , 0.262 0.212 , 0.266	Depositor DCC
R_{free} test set	2074 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	64.9	Xtriage
Anisotropy	0.098	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 69.0	EDS
Estimated twinning fraction	0.016 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 40986 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7656	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 24.15 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 4.0524e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² 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: PO4

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	0.94	0/2640	0.82	2/3556 (0.1%)
1	B	0.80	0/2575	0.77	0/3480
1	C	0.73	0/2497	0.77	0/3391
All	All	0.83	0/7712	0.79	2/10427 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	93	LEU	CB-CG-CD2	-6.07	100.68	111.00
1	A	125	ASN	CB-CA-C	-5.67	99.07	110.40

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	2606	0	2731	64	0
1	B	2542	0	2609	100	0
1	C	2464	0	2443	108	0
2	A	10	0	0	2	0
2	B	15	0	0	3	0
2	C	10	0	0	1	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	4	0	0	0	0
3	B	3	0	0	0	0
3	C	2	0	0	0	0
All	All	7656	0	7783	267	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 17.

All (267) 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:71:ALA:O	1:A:72:THR:HG22	1.32	1.22
1:A:1:MSE:CG	1:A:2:SER:H	1.62	1.12
1:A:1:MSE:HG3	1:A:2:SER:N	1.48	1.11
1:A:37:ILE:HD12	1:A:157:THR:CG2	1.85	1.05
1:A:37:ILE:HD12	1:A:157:THR:HG21	1.37	1.05
1:B:161:GLY:HA3	1:B:212:VAL:HG21	1.37	1.05
1:B:29:ILE:O	1:B:32:THR:HG22	1.61	1.01
1:B:173:LYS:O	1:B:174:VAL:HG13	1.62	0.97
1:A:71:ALA:O	1:A:72:THR:CG2	2.18	0.92
1:C:187:ILE:HD11	1:C:191:ARG:CZ	2.00	0.91
1:A:315:GLU:HG3	1:A:316:GLY:N	1.90	0.87
1:C:329:ASN:O	1:C:330:ASP:HB2	1.73	0.87
1:A:32:THR:HG23	1:A:45:ILE:HD13	1.57	0.85
1:C:124:MSE:CE	1:C:209:LEU:HB3	2.08	0.83
1:C:128:LYS:HA	1:C:206:PHE:CE1	2.13	0.83
1:C:166:LEU:HD22	1:C:208:VAL:HG21	1.62	0.81
1:C:183:HIS:O	1:C:187:ILE:HG22	1.81	0.80
1:C:124:MSE:HE2	1:C:209:LEU:HB3	1.63	0.80
1:C:190:VAL:O	1:C:194:LEU:HD22	1.82	0.79
1:C:302:ILE:HG22	1:C:304:GLY:H	1.48	0.79
1:C:166:LEU:CD2	1:C:208:VAL:HG11	2.12	0.79
1:C:255:LYS:HA	1:C:258:ASP:OD1	1.81	0.79
1:C:270:VAL:HG21	1:C:315:GLU:HG3	1.64	0.79
1:C:124:MSE:HE3	1:C:127:SER:CB	2.13	0.79
1:C:303:GLU:O	1:C:306:LYS:O	2.01	0.78
1:B:41:HIS:HB3	1:B:43:ASP:OD1	1.84	0.78
1:A:37:ILE:HD12	1:A:157:THR:HG23	1.66	0.78
1:A:37:ILE:CD1	1:A:157:THR:HG21	2.14	0.76
1:A:127:SER:H	1:A:210:ASN:HD21	1.38	0.72
1:B:30:GLU:O	1:B:31:THR:C	2.29	0.70

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:29:ILE:O	1:B:32:THR:CG2	2.38	0.70
1:B:212:VAL:O	1:B:212:VAL:CG2	2.40	0.69
1:C:131:TYR:HE2	1:C:132:MSE:HE2	1.55	0.69
1:C:163:LEU:HD13	1:C:281:ILE:HD13	1.73	0.69
1:A:41:HIS:CD2	1:A:43:ASP:H	2.11	0.68
1:A:242:MSE:HE1	1:A:245:VAL:HG21	1.74	0.68
1:B:187:ILE:HD11	1:B:191:ARG:NH1	2.09	0.68
1:B:242:MSE:HE1	1:B:245:VAL:HG21	1.76	0.67
1:B:173:LYS:HG2	1:B:277:ASP:HB3	1.75	0.67
1:C:166:LEU:HD23	1:C:208:VAL:HG11	1.77	0.67
1:A:91:LYS:NZ	2:A:352:PO4:O4	2.24	0.66
1:C:131:TYR:CE2	1:C:132:MSE:HE2	2.30	0.66
1:B:179:ILE:N	1:B:312:SER:HG	1.94	0.66
1:B:88:LYS:NZ	1:C:13:ASP:OD1	2.29	0.65
1:C:124:MSE:HE3	1:C:127:SER:HB2	1.77	0.65
1:B:148:VAL:HG21	1:B:220:VAL:HG13	1.78	0.65
1:B:242:MSE:CE	1:B:245:VAL:HG21	2.26	0.65
1:B:127:SER:H	1:B:210:ASN:HD21	1.44	0.65
1:B:124:MSE:HE3	1:B:209:LEU:HB3	1.79	0.64
1:A:345:TYR:CD2	1:A:345:TYR:N	2.62	0.64
1:B:273:ASP:OD1	1:B:276:GLY:N	2.30	0.64
1:A:5:ALA:HB1	1:A:8:GLU:HB2	1.79	0.64
1:A:242:MSE:HA	1:A:242:MSE:HE2	1.80	0.64
1:B:24:CYS:H	1:B:94:HIS:HD1	1.46	0.64
1:A:1:MSE:HG3	1:A:2:SER:H	0.67	0.63
1:C:124:MSE:HE3	1:C:127:SER:HB3	1.79	0.63
1:B:171:GLU:OE1	1:B:187:ILE:HD13	1.99	0.63
1:B:135:TYR:CZ	1:B:139:LYS:HE2	2.34	0.63
1:A:16:LYS:NZ	1:C:330:ASP:OD2	2.32	0.62
1:C:166:LEU:CD2	1:C:208:VAL:HG21	2.28	0.62
1:B:150:GLU:HG3	1:B:216:MSE:HE2	1.81	0.62
1:A:330:ASP:OD2	1:B:16:LYS:HE3	1.99	0.62
1:A:28:SER:HB3	1:A:100:PHE:CE2	2.35	0.62
1:B:97:ALA:HB1	1:B:215:LYS:HB3	1.81	0.62
1:C:214:ASP:OD1	1:C:216:MSE:HE3	2.00	0.62
1:C:270:VAL:HG21	1:C:315:GLU:CG	2.30	0.62
1:C:28:SER:OG	1:C:100:PHE:N	2.33	0.61
1:B:240:THR:HA	1:B:278:LEU:HD13	1.83	0.61
1:A:68:PRO:O	1:A:75:PRO:HA	2.01	0.61
1:C:297:PHE:CE1	1:C:305:LEU:HD22	2.36	0.61
1:B:173:LYS:O	1:B:174:VAL:CG1	2.43	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:242:MSE:CE	1:A:245:VAL:HG21	2.31	0.60
1:C:28:SER:HB2	1:C:100:PHE:CE2	2.36	0.60
1:B:31:THR:O	1:B:33:LEU:N	2.35	0.60
1:C:161:GLY:HA2	1:C:190:VAL:HG11	1.82	0.60
1:B:273:ASP:OD2	1:B:313:VAL:HG11	2.01	0.59
1:B:266:THR:O	1:B:292:ALA:HA	2.02	0.59
1:C:217:MSE:CE	1:C:244:ALA:HB1	2.32	0.59
1:B:156:THR:HG1	2:B:351:PO4:P	2.26	0.58
1:C:158:THR:HG21	1:C:214:ASP:HB3	1.86	0.58
1:B:75:PRO:HG3	1:B:305:LEU:CD2	2.32	0.58
1:B:173:LYS:CG	1:B:277:ASP:HB3	2.34	0.58
1:B:64:LEU:O	1:B:64:LEU:HD13	2.04	0.58
1:A:313:VAL:HG13	1:A:313:VAL:O	2.01	0.58
1:B:15:ILE:HB	1:B:145:LEU:CD1	2.33	0.57
1:C:152:VAL:CG1	1:C:152:VAL:O	2.52	0.57
1:C:173:LYS:NZ	1:C:277:ASP:OD2	2.28	0.57
1:C:191:ARG:HA	1:C:194:LEU:HD23	1.87	0.56
1:B:302:ILE:HD13	1:B:345:TYR:CG	2.41	0.56
1:C:194:LEU:O	1:C:198:GLY:N	2.36	0.56
1:C:339:PHE:O	1:C:343:LYS:HB2	2.05	0.56
1:C:57:PHE:HB3	1:C:105:ILE:HD13	1.86	0.56
1:A:156:THR:OG1	2:A:351:PO4:O1	2.24	0.56
1:C:246:LEU:HD11	1:C:288:VAL:HG21	1.87	0.56
1:C:190:VAL:HG12	1:C:194:LEU:HD21	1.88	0.55
1:C:163:LEU:CD1	1:C:281:ILE:HD13	2.34	0.55
1:C:28:SER:HB2	1:C:100:PHE:CD2	2.41	0.55
1:C:32:THR:HG23	1:C:45:ILE:HD13	1.88	0.55
1:C:27:SER:O	1:C:152:VAL:HG23	2.06	0.55
1:B:28:SER:HB3	1:B:100:PHE:CE2	2.41	0.55
1:B:125:ASN:C	1:B:125:ASN:HD22	2.10	0.55
1:C:124:MSE:HE2	1:C:209:LEU:CB	2.36	0.55
1:C:173:LYS:HD3	1:C:277:ASP:HB3	1.89	0.55
1:A:124:MSE:HE1	1:A:214:ASP:C	2.27	0.55
1:B:241:GLN:O	1:B:244:ALA:HB3	2.07	0.54
1:C:24:CYS:H	1:C:94:HIS:HD1	1.53	0.54
1:B:330:ASP:OD2	1:C:16:LYS:HE2	2.08	0.54
1:B:204:SER:O	1:B:207:ASP:N	2.41	0.54
1:C:123:ALA:HB2	1:C:211:ALA:O	2.07	0.54
1:C:157:THR:OG1	1:C:158:THR:N	2.40	0.54
1:B:70:ASP:OD1	1:B:72:THR:N	2.32	0.54
1:B:242:MSE:HE2	1:B:245:VAL:CG2	2.37	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:212:VAL:HG22	1:B:212:VAL:O	2.08	0.53
1:B:45:ILE:CG2	1:B:46:LYS:N	2.72	0.53
1:C:206:PHE:O	1:C:210:ASN:OD1	2.27	0.53
1:C:73:GLY:O	1:C:345:TYR:OH	2.23	0.53
1:A:32:THR:HG23	1:A:45:ILE:CD1	2.35	0.53
1:B:48:THR:HB	1:B:49:PRO:HD3	1.91	0.53
1:A:127:SER:OG	1:A:210:ASN:ND2	2.42	0.52
1:C:302:ILE:HG22	1:C:304:GLY:N	2.20	0.52
1:A:138:GLY:O	1:A:230:ARG:NH2	2.41	0.52
1:A:332:GLU:OE2	1:B:1:MSE:CB	2.58	0.52
1:A:37:ILE:HB	1:A:157:THR:HG21	1.90	0.52
1:A:173:LYS:NZ	1:A:277:ASP:OD2	2.43	0.52
1:A:29:ILE:O	1:A:32:THR:HB	2.09	0.51
1:C:166:LEU:HD12	1:C:251:GLU:HG3	1.91	0.51
1:B:54:GLU:O	1:B:58:TYR:HB3	2.10	0.51
1:B:161:GLY:CA	1:B:212:VAL:HG21	2.26	0.51
1:B:57:PHE:HA	1:B:92:ASN:HD21	1.75	0.51
1:A:266:THR:O	1:A:292:ALA:HA	2.09	0.51
1:C:217:MSE:CE	1:C:244:ALA:CB	2.88	0.51
1:C:247:ALA:HB2	1:C:285:ILE:HD13	1.93	0.51
1:B:12:LEU:HD22	1:B:328:VAL:HG21	1.92	0.51
1:A:55:LEU:HD12	1:A:60:LYS:HA	1.93	0.51
1:B:242:MSE:CE	1:B:245:VAL:CG2	2.89	0.51
1:C:247:ALA:HB2	1:C:285:ILE:CD1	2.41	0.51
1:B:37:ILE:HB	1:B:157:THR:HG21	1.92	0.51
1:B:212:VAL:HG23	1:B:212:VAL:O	2.11	0.51
1:B:179:ILE:HD13	1:B:180:ASN:H	1.77	0.50
1:B:55:LEU:HD13	1:B:60:LYS:HA	1.92	0.50
1:A:31:THR:HG23	1:A:35:ILE:HD12	1.94	0.50
1:C:247:ALA:CA	1:C:285:ILE:HD13	2.42	0.50
1:C:265:GLY:HA2	1:C:291:LEU:O	2.12	0.50
1:C:164:LEU:HD23	1:C:168:TYR:O	2.11	0.50
1:C:187:ILE:HD12	1:C:187:ILE:O	2.12	0.50
1:C:266:THR:HG23	1:C:267:THR:N	2.26	0.50
1:A:150:GLU:HG2	1:A:152:VAL:HG13	1.94	0.49
1:B:217:MSE:N	1:B:218:PRO:CD	2.75	0.49
1:C:27:SER:HB3	1:C:152:VAL:CG2	2.42	0.49
1:C:57:PHE:HA	1:C:92:ASN:HD21	1.76	0.49
1:C:88:LYS:HD3	1:C:331:LEU:HB2	1.93	0.49
1:A:345:TYR:HD2	1:A:345:TYR:N	2.09	0.49
1:C:300:ALA:HB2	1:C:337:ARG:HD2	1.93	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:51:ALA:HA	1:B:54:GLU:OE1	2.12	0.49
1:A:61:SER:HB2	1:A:64:LEU:O	2.13	0.49
1:C:161:GLY:HA2	1:C:190:VAL:CG1	2.42	0.49
1:C:217:MSE:HE1	1:C:244:ALA:HB1	1.95	0.49
1:C:239:GLY:O	1:C:240:THR:HB	2.12	0.49
1:A:126:ASN:HB2	1:A:129:GLU:HB2	1.95	0.48
1:B:31:THR:O	1:B:32:THR:C	2.49	0.48
1:A:244:ALA:O	1:A:247:ALA:HB3	2.13	0.48
1:C:166:LEU:HD21	1:C:208:VAL:HG11	1.95	0.48
1:C:191:ARG:HG3	1:C:191:ARG:HH11	1.78	0.48
1:B:302:ILE:HG22	1:B:304:GLY:H	1.77	0.48
1:C:98:GLY:HA2	1:C:102:LYS:NZ	2.28	0.48
1:A:242:MSE:HE2	1:A:245:VAL:CG2	2.44	0.48
1:C:127:SER:H	1:C:210:ASN:HD21	1.60	0.48
1:B:76:THR:HG22	1:B:79:THR:HG23	1.95	0.47
1:A:242:MSE:CE	1:A:245:VAL:CG2	2.92	0.47
1:C:55:LEU:HD23	1:C:78:ALA:CB	2.45	0.47
1:B:277:ASP:OD2	1:B:280:GLY:HA3	2.15	0.47
1:B:313:VAL:HG23	1:B:313:VAL:O	2.13	0.47
1:B:161:GLY:HA2	1:B:190:VAL:CG1	2.44	0.47
1:B:124:MSE:HE3	1:B:209:LEU:C	2.35	0.47
1:C:160:LEU:CD1	1:C:170:ALA:HB3	2.44	0.47
1:C:57:PHE:CB	1:C:105:ILE:HD13	2.45	0.47
1:A:20:PRO:HG2	1:A:90:ILE:HG23	1.96	0.47
1:B:295:PHE:N	1:B:295:PHE:CD2	2.83	0.47
1:B:207:ASP:O	1:B:208:VAL:C	2.54	0.46
1:A:171:GLU:HG3	1:A:187:ILE:CD1	2.45	0.46
1:B:301:LYS:HB3	1:B:341:GLU:OE2	2.15	0.46
1:B:215:LYS:O	1:B:218:PRO:HD2	2.15	0.46
1:C:35:ILE:HG22	1:C:185:LEU:HD21	1.97	0.46
1:A:320:GLY:O	1:A:324:VAL:HG23	2.15	0.46
1:A:72:THR:HG23	1:A:74:SER:OG	2.15	0.46
1:A:313:VAL:CG1	1:A:313:VAL:O	2.64	0.46
1:B:125:ASN:HD22	1:B:126:ASN:N	2.14	0.46
1:B:230:ARG:NH1	2:B:352:PO4:O3	2.49	0.46
1:B:130:LEU:CD1	1:B:215:LYS:HG3	2.46	0.46
1:C:29:ILE:O	1:C:32:THR:HG22	2.16	0.46
1:C:156:THR:OG1	2:C:351:PO4:O3	2.30	0.46
1:B:242:MSE:HE2	1:B:245:VAL:HG23	1.98	0.45
1:A:126:ASN:OD1	1:A:126:ASN:N	2.49	0.45
1:B:156:THR:HG22	1:B:174:VAL:HB	1.98	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:72:THR:O	1:C:72:THR:CG2	2.64	0.45
1:A:287:ASN:C	1:A:287:ASN:OD1	2.54	0.45
1:C:130:LEU:HD11	1:C:215:LYS:HD3	1.99	0.45
1:C:162:VAL:O	1:C:166:LEU:HD23	2.17	0.45
1:B:15:ILE:HB	1:B:145:LEU:HD11	1.97	0.45
1:B:302:ILE:HG22	1:B:304:GLY:N	2.31	0.45
1:C:334:THR:HG23	1:C:337:ARG:NH2	2.32	0.45
1:B:21:LEU:CD2	1:B:141:LEU:HD13	2.47	0.45
1:A:169:ASP:O	1:A:173:LYS:HD2	2.17	0.44
1:B:153:PRO:HA	1:B:154:GLY:HA2	1.75	0.44
1:B:76:THR:CG2	1:B:78:ALA:HB3	2.48	0.44
1:A:87:LEU:HD13	1:A:331:LEU:HD11	1.98	0.44
1:C:187:ILE:HD11	1:C:191:ARG:NH1	2.30	0.44
1:B:150:GLU:HG2	1:B:152:VAL:HG13	1.98	0.44
1:C:174:VAL:HG22	1:C:240:THR:HG23	1.99	0.44
1:B:173:LYS:C	1:B:174:VAL:HG22	2.37	0.44
1:B:242:MSE:HE2	1:B:242:MSE:HA	1.98	0.44
1:B:64:LEU:HD22	1:B:64:LEU:C	2.37	0.44
1:C:160:LEU:O	1:C:160:LEU:HD12	2.18	0.44
1:C:306:LYS:O	1:C:307:ASN:HB2	2.17	0.44
1:B:55:LEU:O	1:B:82:ARG:HD3	2.18	0.44
1:A:315:GLU:HG3	1:A:316:GLY:H	1.74	0.44
1:B:240:THR:HA	1:B:278:LEU:CD1	2.47	0.44
1:A:71:ALA:O	1:A:72:THR:CB	2.63	0.43
1:C:35:ILE:CG2	1:C:185:LEU:HD21	2.49	0.43
1:C:72:THR:O	1:C:72:THR:HG22	2.18	0.43
1:A:119:GLU:N	1:A:119:GLU:OE1	2.41	0.43
1:C:190:VAL:HG12	1:C:194:LEU:CD2	2.47	0.43
1:C:45:ILE:CG2	1:C:46:LYS:N	2.82	0.43
1:B:32:THR:OG1	1:B:45:ILE:HD11	2.18	0.43
1:A:1:MSE:CG	1:A:2:SER:N	2.34	0.43
1:B:45:ILE:HG23	1:B:46:LYS:N	2.33	0.42
1:B:28:SER:OG	1:B:100:PHE:N	2.51	0.42
1:C:127:SER:OG	1:C:206:PHE:HD1	2.02	0.42
1:A:119:GLU:HB3	1:A:189:VAL:HG13	2.01	0.42
1:C:156:THR:O	1:C:159:ALA:HB3	2.20	0.42
1:C:48:THR:O	1:C:51:ALA:HB3	2.20	0.42
1:B:135:TYR:OH	1:B:229:GLU:OE1	2.21	0.42
1:C:239:GLY:O	1:C:241:GLN:N	2.50	0.42
1:B:124:MSE:HE3	1:B:209:LEU:O	2.19	0.42
1:A:152:VAL:HG22	1:A:155:GLY:HA3	2.02	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:28:SER:CB	1:C:100:PHE:CD2	3.02	0.42
1:B:28:SER:HB3	1:B:100:PHE:CD2	2.54	0.41
1:B:302:ILE:HD13	1:B:345:TYR:CD2	2.55	0.41
1:C:246:LEU:HD21	1:C:288:VAL:HG21	2.00	0.41
1:A:174:VAL:HG13	1:A:240:THR:CG2	2.50	0.41
1:A:72:THR:HG21	1:A:312:SER:OG	2.20	0.41
1:B:216:MSE:SE	1:B:242:MSE:HE1	2.70	0.41
1:B:124:MSE:CE	1:B:209:LEU:O	2.69	0.41
1:C:25:VAL:O	1:C:150:GLU:HA	2.20	0.41
1:A:161:GLY:HA2	1:A:190:VAL:CG1	2.51	0.41
1:B:179:ILE:HD13	1:B:180:ASN:N	2.34	0.41
1:C:37:ILE:HB	1:C:157:THR:HG21	2.03	0.41
1:C:111:ASP:OD2	1:C:215:LYS:NZ	2.35	0.41
1:B:166:LEU:HD22	1:B:251:GLU:HG3	2.03	0.41
1:C:332:GLU:O	1:C:335:LYS:N	2.40	0.41
1:C:253:ASN:C	1:C:255:LYS:N	2.73	0.41
1:C:138:GLY:O	1:C:230:ARG:NH2	2.54	0.41
1:B:156:THR:OG1	2:B:351:PO4:P	2.79	0.41
1:C:190:VAL:CG1	1:C:194:LEU:HD21	2.50	0.41
1:C:332:GLU:HA	1:C:333:PRO:HD3	1.95	0.41
1:B:173:LYS:O	1:B:174:VAL:HG22	2.21	0.41
1:C:216:MSE:HE3	1:C:216:MSE:HB3	1.92	0.41
1:C:297:PHE:CD1	1:C:305:LEU:HD22	2.55	0.41
1:C:152:VAL:CG1	1:C:155:GLY:HA3	2.51	0.41
1:B:115:THR:HG21	1:B:214:ASP:HA	2.03	0.41
1:A:297:PHE:CE1	1:A:305:LEU:HD13	2.56	0.41
1:C:27:SER:HB3	1:C:152:VAL:HG21	2.03	0.41
1:B:135:TYR:CE2	1:B:139:LYS:HE2	2.56	0.40
1:B:68:PRO:O	1:B:76:THR:HB	2.20	0.40
1:A:162:VAL:HG21	1:A:217:MSE:SE	2.71	0.40
1:A:150:GLU:HG3	1:A:216:MSE:HE2	2.04	0.40
1:B:62:LEU:HD22	1:B:62:LEU:HA	1.70	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	344/350 (98%)	323 (94%)	21 (6%)	0	100	100
1	B	338/350 (97%)	310 (92%)	25 (7%)	3 (1%)	21	44
1	C	336/350 (96%)	307 (91%)	27 (8%)	2 (1%)	30	54
All	All	1018/1050 (97%)	940 (92%)	73 (7%)	5 (0%)	34	59

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	32	THR
1	B	299	LYS
1	C	330	ASP
1	C	255	LYS
1	B	205	VAL

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	286/286 (100%)	267 (93%)	19 (7%)	21	42
1	B	272/286 (95%)	233 (86%)	39 (14%)	4	8
1	C	254/286 (89%)	231 (91%)	23 (9%)	12	24
All	All	812/858 (95%)	731 (90%)	81 (10%)	9	19

All (81) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	SER
1	A	13	ASP
1	A	32	THR
1	A	43	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	44	VAL
1	A	48	THR
1	A	69	ILE
1	A	110	ILE
1	A	139	LYS
1	A	145	LEU
1	A	150	GLU
1	A	157	THR
1	A	178	SER
1	A	226	SER
1	A	240	THR
1	A	293	SER
1	A	315	GLU
1	A	332	GLU
1	A	344	PHE
1	B	6	ILE
1	B	15	ILE
1	B	30	GLU
1	B	42	ARG
1	B	43	ASP
1	B	44	VAL
1	B	60	LYS
1	B	62	LEU
1	B	64	LEU
1	B	65	LYS
1	B	66	THR
1	B	70	ASP
1	B	76	THR
1	B	88	LYS
1	B	119	GLU
1	B	120	GLU
1	B	125	ASN
1	B	133	LYS
1	B	136	LEU
1	B	139	LYS
1	B	148	VAL
1	B	150	GLU
1	B	156	THR
1	B	157	THR
1	B	173	LYS
1	B	179	ILE
1	B	180	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	183	HIS
1	B	191	ARG
1	B	200	ASN
1	B	212	VAL
1	B	226	SER
1	B	240	THR
1	B	264	ILE
1	B	266	THR
1	B	273	ASP
1	B	293	SER
1	B	309	CYS
1	B	314	LYS
1	C	3	ILE
1	C	27	SER
1	C	28	SER
1	C	32	THR
1	C	41	HIS
1	C	42	ARG
1	C	48	THR
1	C	76	THR
1	C	110	ILE
1	C	112	GLU
1	C	117	ARG
1	C	152	VAL
1	C	166	LEU
1	C	183	HIS
1	C	184	GLU
1	C	187	ILE
1	C	203	SER
1	C	256	VAL
1	C	258	ASP
1	C	293	SER
1	C	298	GLU
1	C	309	CYS
1	C	315	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	HIS
1	A	210	ASN
1	A	260	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	92	ASN
1	B	125	ASN
1	B	180	ASN
1	B	183	HIS
1	B	210	ASN
1	B	241	GLN
1	B	342	ASN
1	C	7	ASN
1	C	92	ASN
1	C	260	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

7 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	$\# Z > 2$
2	PO4	A	351	-	4,4,4	0.77	0	6,6,6	0.34	0
2	PO4	A	352	-	4,4,4	0.77	0	6,6,6	0.34	0
2	PO4	B	351	-	4,4,4	0.49	0	6,6,6	0.30	0
2	PO4	B	352	-	4,4,4	0.44	0	6,6,6	0.33	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PO4	B	353	-	4,4,4	0.38	0	6,6,6	0.29	0
2	PO4	C	351	-	4,4,4	0.50	0	6,6,6	0.34	0
2	PO4	C	352	-	4,4,4	0.43	0	6,6,6	0.27	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	PO4	A	351	-	-	0/0/0/0	0/0/0/0
2	PO4	A	352	-	-	0/0/0/0	0/0/0/0
2	PO4	B	351	-	-	0/0/0/0	0/0/0/0
2	PO4	B	352	-	-	0/0/0/0	0/0/0/0
2	PO4	B	353	-	-	0/0/0/0	0/0/0/0
2	PO4	C	351	-	-	0/0/0/0	0/0/0/0
2	PO4	C	352	-	-	0/0/0/0	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.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	351	PO4	1	0
2	A	352	PO4	1	0
2	B	351	PO4	2	0
2	B	352	PO4	1	0
2	C	351	PO4	1	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	340/350 (97%)	-0.19	7 (2%) 67 66	15, 29, 40, 59	0
1	B	336/350 (96%)	0.08	21 (6%) 23 21	19, 33, 54, 63	0
1	C	334/350 (95%)	0.05	13 (3%) 43 41	19, 45, 63, 69	0
All	All	1010/1050 (96%)	-0.02	41 (4%) 41 39	15, 34, 58, 69	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	303	GLU	4.7
1	A	201	GLU	4.4
1	B	301	LYS	4.3
1	B	171	GLU	4.2
1	B	120	GLU	4.0
1	C	171	GLU	3.8
1	A	202	LYS	3.8
1	A	302	ILE	3.6
1	C	197	ALA	3.5
1	B	300	ALA	3.5
1	B	201	GLU	3.5
1	C	345	TYR	3.4
1	B	309	CYS	3.3
1	B	302	ILE	2.8
1	C	300	ALA	2.8
1	C	307	ASN	2.8
1	C	302	ILE	2.7
1	B	179	ILE	2.7
1	B	305	LEU	2.6
1	B	345	TYR	2.6
1	C	184	GLU	2.6
1	A	345	TYR	2.5
1	A	69	ILE	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	172	GLY	2.5
1	B	342	ASN	2.4
1	B	63	THR	2.4
1	B	150	GLU	2.3
1	B	184	GLU	2.3
1	C	164	LEU	2.3
1	B	172	GLY	2.3
1	B	274	LYS	2.3
1	C	202	LYS	2.2
1	B	202	LYS	2.1
1	B	307	ASN	2.1
1	C	203	SER	2.1
1	A	200	ASN	2.1
1	A	301	LYS	2.1
1	B	183	HIS	2.0
1	B	170	ALA	2.0
1	C	151	SER	2.0
1	C	305	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no 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(Å ²)	Q<0.9
2	PO4	C	352	5/5	0.29	0.83	13.33	235,235,235,235	0
2	PO4	A	351	5/5	0.99	0.20	0.93	44,46,48,55	0
2	PO4	C	351	5/5	0.93	0.19	-0.07	81,81,83,85	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	PO4	B	351	5/5	0.95	0.20	-0.09	66,68,70,72	0
2	PO4	B	352	5/5	0.95	0.09	-2.30	69,73,76,76	0
2	PO4	A	352	5/5	0.98	0.06	-2.90	63,67,67,69	0
2	PO4	B	353	5/5	0.97	0.10	-	73,77,78,78	0

6.5 Other polymers [i](#)

There are no such residues in this entry.