



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 03:01 AM EDT

PDB ID : 1VQV / pdb_00001vqv
Title : Crystal Structure of Thiamine Monophosphate Kinase (thil) from Aquifex Aeolicus
Authors : Eswaramoorthy, S.; Swaminathan, S.; Burley, S.K.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2005-01-05
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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

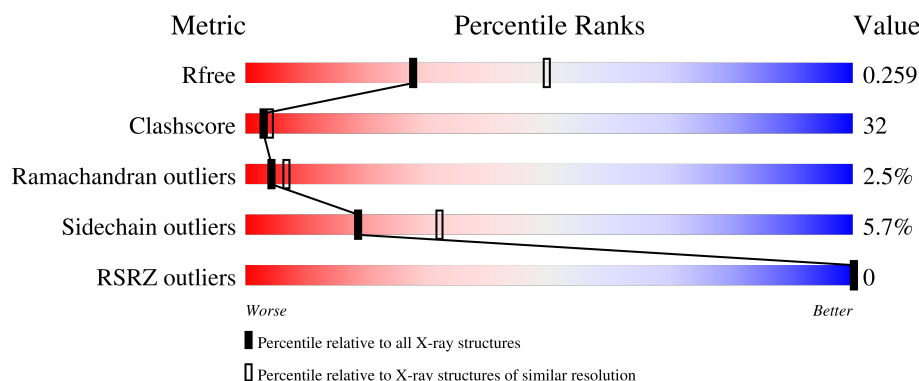
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

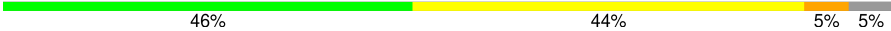
The reported resolution of this entry is 2.65 Å.

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}	180053	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	306	
1	B	306	

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	A	1002[B]	-	-	X	-
2	PO4	B	1004[B]	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4732 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.

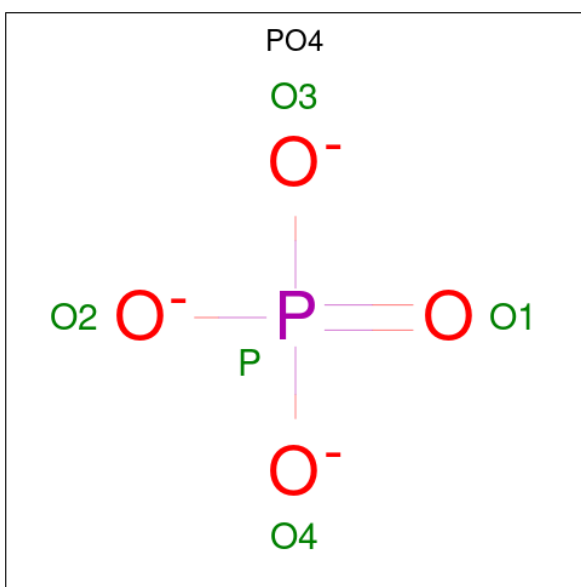
- Molecule 1 is a protein called thiamine monophosphate kinase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	292	Total	C	N	O	S	Se	0	1	0
			2285	1466	381	430	3	5			
1	B	295	Total	C	N	O	S	Se	0	1	0
			2303	1476	384	435	3	5			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	UNP O67883
A	169	MSE	MET	modified residue	UNP O67883
A	206	MSE	MET	modified residue	UNP O67883
A	242	MSE	MET	modified residue	UNP O67883
A	279	MSE	MET	modified residue	UNP O67883
B	1	MSE	MET	modified residue	UNP O67883
B	169	MSE	MET	modified residue	UNP O67883
B	206	MSE	MET	modified residue	UNP O67883
B	242	MSE	MET	modified residue	UNP O67883
B	279	MSE	MET	modified residue	UNP O67883

- Molecule 2 is PHOSPHATE ION (CCD ID: 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	1
			9	7	2		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	1
			9	7	2		

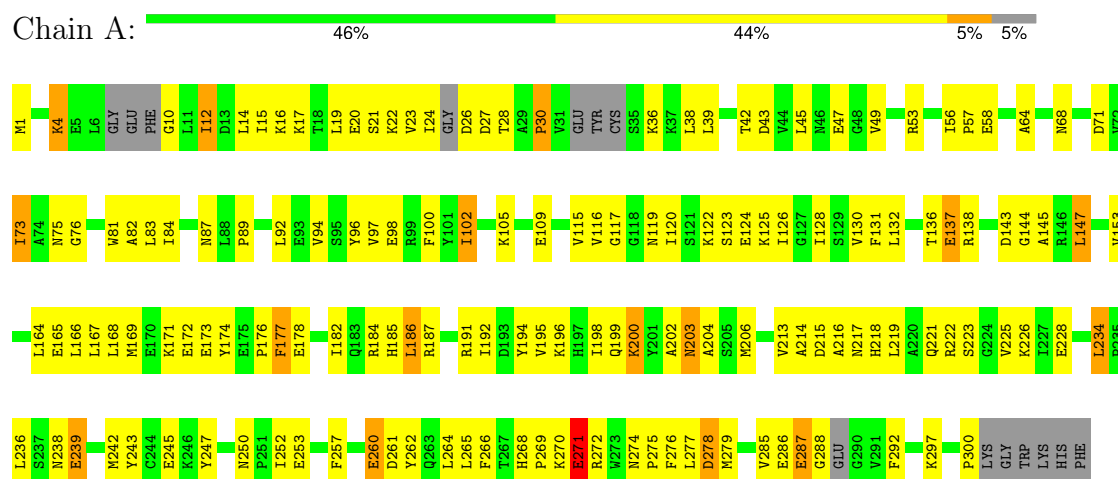
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	50	Total	O	0	0
			50	50		
3	B	66	Total	O	0	0
			66	66		

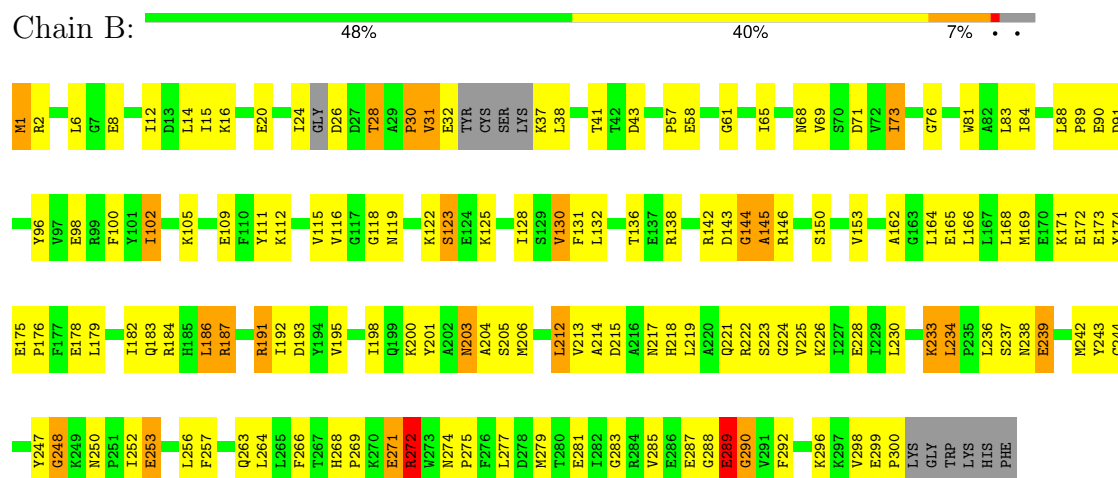
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: thiamine monophosphate kinase



- Molecule 1: thiamine monophosphate kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	65.58Å 65.58Å 192.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.65 50.00 – 2.65	Depositor EDS
% Data completeness (in resolution range)	95.7 (50.00-2.65) 95.7 (50.00-2.65)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.80 (at 2.65Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.238 , 0.260 0.237 , 0.259	Depositor DCC
R_{free} test set	870 reflections (3.71%)	wwPDB-VP
Wilson B-factor (Å ²)	59.9	Xtriage
Anisotropy	0.362	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 51.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.499 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4732	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.13% 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

5.1 Standard geometry

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	0.58	1/2325 (0.0%)	1.06	6/3129 (0.2%)
1	B	0.59	1/2345 (0.0%)	1.03	10/3159 (0.3%)
All	All	0.59	2/4670 (0.0%)	1.04	16/6288 (0.3%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	176	PRO	N-CD	5.40	1.55	1.47
1	A	176	PRO	N-CD	5.40	1.55	1.47

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	22	LYS	N-CA-C	-12.51	97.46	113.72
1	B	290	GLY	N-CA-C	7.97	122.74	111.42
1	A	203	ASN	N-CA-C	-7.59	104.15	113.41
1	A	36	LYS	N-CA-C	-7.47	98.21	109.94
1	A	137	GLU	N-CA-C	-6.95	103.80	112.90
1	B	203	ASN	N-CA-C	-6.79	105.12	113.41
1	B	272	ARG	N-CA-C	6.54	121.62	112.72
1	B	102	ILE	N-CA-C	-6.48	104.44	110.53
1	A	102	ILE	N-CA-C	-6.37	104.55	110.53
1	A	216	ALA	N-CA-C	-5.82	105.07	111.82
1	B	233	LYS	N-CA-C	-5.59	106.82	113.97
1	B	289	GLU	N-CA-C	5.44	122.38	110.80
1	B	31	VAL	N-CA-C	5.43	117.00	109.45
1	B	28	THR	N-CA-C	-5.30	100.27	108.90
1	B	175	GLU	N-CA-C	-5.29	102.58	110.20
1	B	142	ARG	N-CA-C	5.05	119.29	113.18

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	2285	0	2295	156	0
1	B	2303	0	2299	152	0
2	A	14	0	0	6	0
2	B	14	0	0	4	0
3	A	50	0	0	15	0
3	B	66	0	0	10	0
All	All	4732	0	4594	296	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 32.

All (296) 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:16:LYS:HA	1:A:24:ILE:HD12	1.36	1.06
1:B:206:MSE:HE1	1:B:219:LEU:HB2	1.41	1.01
1:B:136:THR:HG22	1:B:138:ARG:H	1.22	0.99
1:B:28:THR:O	1:B:30:PRO:HD3	1.70	0.91
1:B:30:PRO:HD2	1:B:38:LEU:O	1.70	0.90
1:A:225:VAL:HG22	1:A:226:LYS:H	1.35	0.89
1:A:105:LYS:O	1:A:109:GLU:HG3	1.73	0.88
1:A:225:VAL:HG21	1:A:285:VAL:CG1	2.07	0.83
1:A:136:THR:HG22	1:A:138:ARG:H	1.45	0.81
1:B:206:MSE:CE	1:B:219:LEU:HB2	2.11	0.81
1:A:166:LEU:HA	1:A:169:MSE:HE2	1.62	0.81
1:B:172:GLU:HG2	1:B:173:GLU:H	1.44	0.81
1:A:285:VAL:HG12	3:A:1036:HOH:O	1.80	0.80
1:B:125:LYS:HE2	3:B:1031:HOH:O	1.83	0.78
1:B:250:ASN:HD21	1:B:252:ILE:HB	1.48	0.78
1:A:27:ASP:OD2	1:A:39:LEU:HB3	1.84	0.78
1:A:1:MSE:N	3:A:1010:HOH:O	2.17	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:LEU:HD22	3:A:1036:HOH:O	1.83	0.77
1:A:225:VAL:HG21	1:A:285:VAL:HG12	1.65	0.77
1:B:225:VAL:HG21	1:B:285:VAL:CG1	2.16	0.75
1:A:238:ASN:O	1:A:242:MSE:HG3	1.87	0.75
1:A:43:ASP:OD1	1:A:130:VAL:HB	1.87	0.74
1:A:43:ASP:HB3	1:A:68:ASN:OD1	1.86	0.74
1:A:81:TRP:O	1:A:132:LEU:HD12	1.86	0.74
1:B:172:GLU:HG2	1:B:173:GLU:N	1.99	0.74
1:A:28:THR:O	1:A:30:PRO:HD3	1.88	0.74
1:B:144:GLY:HA3	1:B:203:ASN:O	1.88	0.74
1:A:250:ASN:HD21	1:A:252:ILE:HB	1.54	0.73
1:B:81:TRP:O	1:B:132:LEU:HD12	1.87	0.73
1:B:143:ASP:HA	1:B:223:SER:HB3	1.71	0.73
1:B:239:GLU:HA	1:B:242:MSE:HE3	1.70	0.72
1:B:238:ASN:O	1:B:242:MSE:HG3	1.89	0.72
1:A:30:PRO:HD2	1:A:38:LEU:O	1.88	0.72
1:A:166:LEU:HD13	1:A:178:GLU:HB3	1.72	0.71
1:B:68:ASN:ND2	1:B:130:VAL:HB	2.06	0.71
1:B:187:ARG:HH11	2:B:1004[B]:PO4:P	2.14	0.71
1:A:187:ARG:HH11	2:A:1002[B]:PO4:P	2.14	0.71
1:B:226:LYS:HB2	1:B:288:GLY:C	2.17	0.69
1:B:16:LYS:HA	1:B:24:ILE:HD12	1.73	0.69
1:A:165:GLU:O	1:A:169:MSE:HG3	1.93	0.69
1:A:206:MSE:HE3	1:A:265:LEU:HD23	1.74	0.69
1:B:178:GLU:O	1:B:182:ILE:HG13	1.93	0.69
1:A:217:ASN:HB3	1:A:221:GLN:NE2	2.08	0.68
1:A:225:VAL:HG22	1:A:226:LYS:N	2.07	0.68
1:B:26:ASP:N	3:B:1042:HOH:O	2.26	0.68
1:B:105:LYS:O	1:B:109:GLU:HG3	1.93	0.68
1:B:218:HIS:O	1:B:222:ARG:HG2	1.95	0.67
1:B:201:TYR:CD2	1:B:272:ARG:HG2	2.28	0.67
1:B:213:VAL:HG21	1:B:257:PHE:CE2	2.30	0.67
1:A:287:GLU:HG3	1:A:288:GLY:N	2.08	0.66
1:B:225:VAL:HG21	1:B:285:VAL:HG13	1.77	0.65
1:A:187:ARG:HD2	2:A:1002[B]:PO4:O3	1.96	0.65
1:A:136:THR:HG21	1:A:138:ARG:O	1.96	0.65
1:A:167:LEU:HD11	1:A:186:LEU:HD13	1.79	0.65
1:B:37:LYS:N	3:B:1040:HOH:O	2.29	0.65
1:B:206:MSE:HE1	1:B:219:LEU:CB	2.24	0.65
1:A:172:GLU:HG2	1:A:173:GLU:N	2.10	0.65
1:A:226:LYS:HD3	1:A:287:GLU:HB3	1.78	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:226:LYS:HE2	1:A:292:PHE:CE1	2.33	0.64
1:A:16:LYS:HG2	1:A:24:ILE:HB	1.79	0.64
1:A:84:ILE:HG12	1:A:130:VAL:HG22	1.79	0.63
1:A:143:ASP:HA	1:A:223:SER:OG	1.98	0.63
1:B:98:GLU:O	1:B:102:ILE:HG13	1.98	0.62
1:B:213:VAL:HG23	1:B:214:ALA:N	2.14	0.62
1:B:230:LEU:HD12	1:B:233:LYS:HD2	1.81	0.62
1:B:16:LYS:HG2	1:B:24:ILE:HB	1.81	0.62
1:A:268:HIS:HB2	1:A:269:PRO:HD2	1.81	0.62
1:B:31:VAL:HG12	1:B:32:GLU:N	2.14	0.62
1:A:206:MSE:CE	1:A:265:LEU:HD23	2.30	0.61
1:B:201:TYR:CE2	1:B:272:ARG:HG2	2.36	0.61
1:B:73:ILE:HG12	1:B:192:ILE:HA	1.83	0.60
1:A:239:GLU:CD	1:A:239:GLU:H	2.09	0.60
1:A:12:ILE:HG13	3:A:1035:HOH:O	2.01	0.60
1:A:76:GLY:O	1:A:136:THR:HG23	2.02	0.60
1:A:144:GLY:HA3	1:A:203:ASN:O	2.00	0.60
1:A:217:ASN:HB3	1:A:221:GLN:HE22	1.65	0.60
1:B:84:ILE:HG12	1:B:130:VAL:HG13	1.82	0.60
1:B:119:ASN:N	3:B:1006:HOH:O	2.35	0.60
1:A:239:GLU:HA	1:A:242:MSE:HE3	1.84	0.60
1:A:238:ASN:HB3	1:A:242:MSE:HE2	1.83	0.59
1:B:32:GLU:C	3:B:1041:HOH:O	2.45	0.59
1:A:136:THR:HG22	1:A:138:ARG:N	2.16	0.58
1:A:122:LYS:HD2	3:A:1020:HOH:O	2.04	0.58
1:B:171:LYS:HB2	1:B:174:TYR:CE2	2.39	0.58
1:A:23:VAL:O	1:A:24:ILE:HG13	2.03	0.58
1:B:143:ASP:O	1:B:223:SER:HB3	2.03	0.58
1:A:98:GLU:O	1:A:102:ILE:HG13	2.03	0.57
3:A:1043:HOH:O	1:B:105:LYS:HE3	2.03	0.57
1:A:19:LEU:HD12	1:A:24:ILE:HD13	1.86	0.57
1:B:166:LEU:HD13	1:B:178:GLU:HB3	1.86	0.57
1:B:250:ASN:ND2	1:B:252:ILE:HB	2.17	0.57
1:A:153:VAL:HA	1:A:264:LEU:O	2.05	0.57
1:A:213:VAL:HG21	1:A:257:PHE:CE2	2.40	0.57
1:B:225:VAL:HG22	1:B:226:LYS:H	1.70	0.57
1:B:226:LYS:HG3	1:B:290:GLY:O	2.05	0.57
1:A:286:GLU:N	3:A:1036:HOH:O	2.37	0.56
1:B:83:LEU:HB2	1:B:131:PHE:HB3	1.87	0.56
1:A:21:SER:HB2	1:A:24:ILE:HG13	1.88	0.56
1:B:43:ASP:OD1	1:B:68:ASN:ND2	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:171:LYS:HB2	1:B:174:TYR:CZ	2.40	0.56
1:B:238:ASN:HB3	1:B:242:MSE:HE2	1.88	0.56
1:A:27:ASP:HB2	1:A:75:ASN:ND2	2.20	0.56
1:B:239:GLU:CD	1:B:239:GLU:H	2.13	0.56
1:B:213:VAL:HG23	1:B:214:ALA:H	1.71	0.56
1:B:239:GLU:CA	1:B:242:MSE:HE3	2.36	0.56
1:B:73:ILE:HG13	1:B:195:VAL:HG23	1.88	0.55
1:A:172:GLU:HG2	1:A:173:GLU:H	1.71	0.55
1:A:28:THR:HG21	1:B:116:VAL:O	2.07	0.55
1:A:23:VAL:O	1:A:23:VAL:HG23	2.07	0.55
1:B:57:PRO:HG3	1:B:96:TYR:CE1	2.42	0.55
1:B:165:GLU:O	1:B:169:MSE:HG3	2.06	0.55
1:B:230:LEU:HB2	1:B:233:LYS:HD2	1.89	0.54
1:B:187:ARG:HD2	2:B:1004[B]:PO4:O3	2.08	0.54
1:B:272:ARG:NH1	3:B:1048:HOH:O	2.39	0.54
1:B:16:LYS:O	1:B:20:GLU:N	2.40	0.53
1:A:82:ALA:O	1:A:115:VAL:HA	2.07	0.53
1:A:184:ARG:HH12	1:A:239:GLU:CD	2.16	0.53
1:A:287:GLU:CG	1:A:288:GLY:N	2.72	0.52
1:A:23:VAL:O	1:A:24:ILE:CG1	2.57	0.52
1:A:16:LYS:HD2	3:A:1009:HOH:O	2.08	0.52
1:A:116:VAL:O	1:B:28:THR:HG21	2.09	0.52
1:B:65:ILE:O	1:B:69:VAL:HG23	2.09	0.52
1:B:68:ASN:HD21	1:B:130:VAL:HB	1.73	0.52
1:B:144:GLY:O	1:B:203:ASN:HB3	2.10	0.52
1:A:164:LEU:O	1:A:168:LEU:HG	2.09	0.52
1:B:112:LYS:HD2	3:B:1047:HOH:O	2.10	0.52
1:A:239:GLU:CA	1:A:242:MSE:HE3	2.40	0.52
1:A:214:ALA:HA	1:A:300:PRO:HG3	1.92	0.51
1:A:57:PRO:HG3	1:A:96:TYR:CD1	2.46	0.51
1:B:223:SER:O	1:B:225:VAL:N	2.43	0.51
1:A:47:GLU:OE2	1:A:53:ARG:HB2	2.10	0.51
1:A:206:MSE:HE1	1:A:215:ASP:O	2.11	0.51
1:A:225:VAL:CG2	3:A:1036:HOH:O	2.59	0.51
1:A:89:PRO:HA	1:A:123:SER:O	2.11	0.51
1:A:145:ALA:N	1:A:204:ALA:HB2	2.25	0.51
1:A:166:LEU:HD23	1:A:169:MSE:HE2	1.91	0.51
1:B:268:HIS:HB2	1:B:269:PRO:HD2	1.92	0.51
1:A:225:VAL:CG2	1:A:226:LYS:H	2.17	0.50
1:A:119:ASN:HB2	3:A:1027:HOH:O	2.11	0.50
1:A:42:THR:HG21	1:B:118:GLY:HA2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:250:ASN:ND2	1:A:252:ILE:HB	2.23	0.50
1:B:214:ALA:HA	1:B:300:PRO:HG3	1.94	0.50
1:B:217:ASN:HB3	1:B:221:GLN:HE22	1.77	0.50
1:A:4:LYS:HG3	1:B:90:GLU:O	2.11	0.50
1:A:234:LEU:O	1:A:236:LEU:HD13	2.11	0.50
1:B:299:GLU:CB	3:B:1034:HOH:O	2.59	0.50
1:B:143:ASP:CA	1:B:223:SER:HB3	2.41	0.50
1:B:153:VAL:HA	1:B:264:LEU:O	2.12	0.50
1:A:138:ARG:NH2	1:A:202:ALA:O	2.40	0.49
1:A:226:LYS:HB2	1:A:288:GLY:C	2.38	0.49
1:A:287:GLU:CG	1:A:288:GLY:H	2.25	0.49
1:B:263:GLN:HA	1:B:263:GLN:OE1	2.12	0.49
1:B:184:ARG:HD3	3:B:1028:HOH:O	2.12	0.49
1:B:30:PRO:CD	1:B:38:LEU:O	2.53	0.49
1:A:124:GLU:O	1:A:125:LYS:HD3	2.13	0.49
1:B:217:ASN:HB3	1:B:221:GLN:NE2	2.28	0.48
1:B:226:LYS:HE2	1:B:292:PHE:CE1	2.47	0.48
1:B:225:VAL:HG22	1:B:226:LYS:N	2.28	0.48
1:B:223:SER:C	1:B:225:VAL:H	2.20	0.48
1:A:30:PRO:CD	1:A:38:LEU:O	2.59	0.48
1:A:206:MSE:SE	1:A:215:ASP:HB3	2.63	0.48
1:B:274:ASN:CG	1:B:275:PRO:HD2	2.38	0.48
1:B:206:MSE:SE	1:B:215:ASP:HB3	2.64	0.48
1:B:266:PHE:HZ	1:B:279:MSE:HG3	1.79	0.48
1:A:286:GLU:CA	3:A:1036:HOH:O	2.60	0.48
1:B:228:GLU:O	1:B:283:GLY:HA3	2.14	0.48
1:A:73:ILE:HD11	1:A:192:ILE:HG13	1.96	0.47
1:A:195:VAL:O	1:A:199:GLN:HG2	2.14	0.47
1:A:292:PHE:CZ	1:A:297:LYS:HG3	2.48	0.47
1:A:292:PHE:CE2	1:A:297:LYS:HG3	2.49	0.47
1:A:17:LYS:O	1:A:20:GLU:N	2.47	0.47
1:A:277:LEU:C	1:A:279:MSE:H	2.22	0.47
1:B:119:ASN:C	1:B:119:ASN:OD1	2.57	0.47
1:B:244:CYS:O	1:B:248:GLY:N	2.48	0.47
1:A:187:ARG:HD2	2:A:1002[B]:PO4:P	2.54	0.47
1:A:26:ASP:CB	1:A:71:ASP:HB3	2.45	0.47
1:B:57:PRO:HG3	1:B:96:TYR:CD1	2.50	0.47
1:B:143:ASP:O	1:B:223:SER:CB	2.63	0.47
1:B:31:VAL:CG1	1:B:32:GLU:N	2.78	0.47
1:A:97:VAL:O	1:A:100:PHE:HB3	2.15	0.47
1:A:136:THR:HG22	1:A:137:GLU:N	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:187:ARG:NH1	2:A:1002[B]:PO4:O3	2.48	0.47
1:B:187:ARG:NH1	2:B:1004[B]:PO4:O3	2.47	0.47
1:A:266:PHE:HZ	1:A:279:MSE:HG3	1.79	0.46
1:B:73:ILE:CG1	1:B:192:ILE:HA	2.44	0.46
1:B:206:MSE:HE3	1:B:219:LEU:HD22	1.97	0.46
1:B:89:PRO:HA	1:B:123:SER:O	2.16	0.46
1:A:171:LYS:HB2	1:A:174:TYR:CE2	2.51	0.46
1:B:269:PRO:HB2	1:B:271:GLU:HG3	1.97	0.46
1:B:76:GLY:O	1:B:136:THR:HG23	2.15	0.46
1:A:21:SER:HA	3:A:1032:HOH:O	2.16	0.46
1:A:243:TYR:O	1:A:247:TYR:HD1	1.99	0.46
1:A:16:LYS:O	1:A:20:GLU:N	2.49	0.45
1:A:287:GLU:HG3	1:A:288:GLY:H	1.78	0.45
1:A:10:GLY:N	3:A:1035:HOH:O	2.50	0.45
1:A:119:ASN:OD1	1:A:119:ASN:C	2.59	0.45
1:B:84:ILE:HG12	1:B:130:VAL:CG1	2.45	0.45
1:B:150:SER:HB3	1:B:281:GLU:HG3	1.98	0.45
1:A:19:LEU:HD12	1:A:24:ILE:CD1	2.46	0.45
1:A:164:LEU:HD13	1:A:185:HIS:CE1	2.51	0.45
1:B:179:LEU:O	1:B:183:GLN:HG3	2.16	0.45
1:A:136:THR:CG2	1:A:138:ARG:O	2.63	0.45
1:B:166:LEU:HD21	1:B:243:TYR:CE1	2.52	0.45
1:A:285:VAL:C	3:A:1036:HOH:O	2.58	0.45
1:B:212:LEU:HD23	1:B:256:LEU:HD23	1.98	0.45
1:A:94:VAL:HG23	1:B:1:MSE:O	2.16	0.45
1:A:167:LEU:HD21	1:A:182:ILE:HG23	1.99	0.45
1:A:269:PRO:HB2	1:A:271:GLU:HG3	1.99	0.45
1:A:26:ASP:N	3:A:1019:HOH:O	2.51	0.44
1:A:204:ALA:O	1:A:266:PHE:HA	2.18	0.44
1:B:277:LEU:C	1:B:279:MSE:H	2.24	0.44
1:A:49:VAL:HG22	1:B:122:LYS:HE3	2.00	0.44
1:A:92:LEU:HD21	1:A:96:TYR:CD2	2.52	0.44
1:A:177:PHE:CD1	1:A:177:PHE:C	2.94	0.44
1:B:223:SER:C	1:B:225:VAL:N	2.75	0.44
1:A:15:ILE:O	1:A:16:LYS:C	2.59	0.44
1:B:145:ALA:N	1:B:204:ALA:HB2	2.32	0.44
1:A:120:ILE:HD11	1:B:15:ILE:CD1	2.47	0.44
1:B:164:LEU:O	1:B:168:LEU:HG	2.17	0.44
1:B:234:LEU:O	1:B:236:LEU:HD13	2.16	0.44
1:B:225:VAL:CG2	1:B:285:VAL:HG13	2.45	0.44
1:B:61:GLY:HA2	1:B:100:PHE:CE1	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:191:ARG:C	1:B:193:ASP:N	2.76	0.44
1:B:136:THR:HG22	1:B:138:ARG:N	2.07	0.44
1:B:166:LEU:HA	1:B:169:MSE:HE2	1.99	0.44
1:A:28:THR:CG2	1:B:116:VAL:O	2.66	0.43
1:A:64:ALA:HB1	1:A:130:VAL:HG21	1.99	0.43
1:A:38:LEU:HD11	1:B:38:LEU:HD11	2.00	0.43
1:A:166:LEU:CD2	1:A:169:MSE:HE2	2.48	0.43
1:A:196:LYS:HA	1:A:199:GLN:HG2	2.00	0.43
1:A:21:SER:CB	1:A:24:ILE:HG13	2.48	0.43
1:B:200:LYS:O	1:B:200:LYS:HG3	2.19	0.43
1:B:15:ILE:O	1:B:16:LYS:C	2.61	0.43
1:B:91:ASP:OD1	1:B:91:ASP:N	2.51	0.43
1:B:239:GLU:OE2	2:B:1003:PO4:O3	2.36	0.43
1:B:1:MSE:HG2	1:B:2:ARG:HE	1.83	0.43
1:A:191:ARG:HH22	1:A:278:ASP:CG	2.25	0.43
1:B:111:TYR:O	1:B:112:LYS:HB2	2.19	0.43
1:B:173:GLU:HG3	1:B:174:TYR:N	2.33	0.43
1:A:198:ILE:C	1:A:200:LYS:H	2.27	0.43
1:A:28:THR:HG21	1:B:116:VAL:C	2.44	0.42
1:A:214:ALA:O	1:A:217:ASN:HB2	2.19	0.42
1:A:239:GLU:N	1:A:242:MSE:HE3	2.34	0.42
1:A:239:GLU:OE2	2:A:1001:PO4:O1	2.37	0.42
1:B:162:ALA:O	1:B:166:LEU:HG	2.19	0.42
1:A:218:HIS:O	1:A:222:ARG:HB2	2.19	0.42
1:A:271:GLU:O	1:A:272:ARG:HG3	2.19	0.42
1:B:28:THR:C	1:B:30:PRO:HD3	2.40	0.42
1:A:45:LEU:HB2	1:A:128:ILE:HB	2.01	0.42
1:A:194:TYR:O	1:A:198:ILE:HG12	2.19	0.42
1:A:245:GLU:OE1	1:A:245:GLU:HA	2.20	0.42
1:B:288:GLY:O	1:B:289:GLU:CB	2.68	0.42
1:A:274:ASN:CG	1:A:275:PRO:HD2	2.44	0.42
1:B:73:ILE:HD11	1:B:192:ILE:HG13	2.02	0.42
1:B:83:LEU:HD12	1:B:131:PHE:CD2	2.55	0.42
1:B:205:SER:O	1:B:206:MSE:HB3	2.19	0.42
1:B:88:LEU:HA	1:B:89:PRO:HD3	1.91	0.42
1:B:296:LYS:O	1:B:298:VAL:HG23	2.20	0.42
1:B:184:ARG:HH12	1:B:239:GLU:CG	2.33	0.42
1:A:73:ILE:CG1	1:A:192:ILE:HA	2.49	0.42
1:A:217:ASN:C	1:A:219:LEU:N	2.77	0.42
1:B:26:ASP:CB	1:B:71:ASP:HB3	2.50	0.42
1:B:166:LEU:CD2	1:B:243:TYR:HE1	2.33	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:LEU:HD23	1:A:169:MSE:CE	2.50	0.42
1:B:213:VAL:CG2	1:B:214:ALA:N	2.82	0.42
1:A:169:MSE:HE3	1:A:171:LYS:HD2	2.01	0.41
1:A:83:LEU:HD12	1:B:131:PHE:CE1	2.54	0.41
1:A:199:GLN:HG3	1:A:199:GLN:O	2.20	0.41
1:A:275:PRO:HG2	1:A:276:PHE:CD1	2.55	0.41
1:A:260:GLU:OE1	1:A:262:TYR:CE2	2.73	0.41
1:A:94:VAL:HG21	1:B:6:LEU:HD22	2.02	0.41
1:B:191:ARG:C	1:B:193:ASP:H	2.28	0.41
1:B:146:ARG:NH1	3:B:1038:HOH:O	2.49	0.41
1:A:131:PHE:CE2	1:B:131:PHE:HB2	2.56	0.41
1:B:16:LYS:HE2	1:B:24:ILE:O	2.21	0.41
1:B:243:TYR:O	1:B:247:TYR:HD1	2.03	0.41
1:A:213:VAL:HG23	1:A:214:ALA:N	2.35	0.41
1:B:288:GLY:O	1:B:289:GLU:HB2	2.21	0.41
1:A:226:LYS:NZ	1:A:228:GLU:CG	2.84	0.40
1:B:73:ILE:HD12	1:B:73:ILE:HA	1.82	0.40
1:A:14:LEU:O	1:A:17:LYS:HB3	2.22	0.40
1:A:83:LEU:HA	1:A:117:GLY:O	2.21	0.40
1:B:198:ILE:C	1:B:200:LYS:H	2.29	0.40
1:B:250:ASN:ND2	1:B:253:GLU:H	2.19	0.40
1:A:14:LEU:HD12	1:A:17:LYS:CE	2.52	0.40
1:A:87:ASN:O	1:A:126:ILE:HA	2.22	0.40
1:B:41:THR:CG2	1:B:132:LEU:HB3	2.51	0.40
1:A:56:ILE:HA	1:A:57:PRO:HD3	1.91	0.40
1:A:144:GLY:O	1:A:203:ASN:HB3	2.21	0.40
1:B:43:ASP:OD1	1:B:130:VAL:HG23	2.20	0.40
1:B:100:PHE:CE1	1:B:128:ILE:HD13	2.57	0.40
1:B:186:LEU:HA	1:B:186:LEU:HD12	1.90	0.40
1:A:270:LYS:C	1:A:272:ARG:H	2.29	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	283/306 (92%)	256 (90%)	22 (8%)	5 (2%)	6	11
1	B	290/306 (95%)	263 (91%)	18 (6%)	9 (3%)	3	5
All	All	573/612 (94%)	519 (91%)	40 (7%)	14 (2%)	4	7

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	287	GLU
1	B	287	GLU
1	B	289	GLU
1	A	30	PRO
1	B	224	GLY
1	A	278	ASP
1	B	8	GLU
1	B	144	GLY
1	A	271	GLU
1	B	30	PRO
1	A	200	LYS
1	B	212	LEU
1	B	145	ALA
1	B	248	GLY

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	246/257 (96%)	234 (95%)	12 (5%)	22	38
1	B	245/257 (95%)	229 (94%)	16 (6%)	15	27
All	All	491/514 (96%)	463 (94%)	28 (6%)	18	32

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LYS
1	A	12	ILE
1	A	73	ILE
1	A	147	LEU
1	A	177	PHE
1	A	186	LEU
1	A	234	LEU
1	A	239	GLU
1	A	253	GLU
1	A	260	GLU
1	A	261	ASP
1	A	271	GLU
1	B	1	MSE
1	B	12	ILE
1	B	14	LEU
1	B	73	ILE
1	B	115	VAL
1	B	123	SER
1	B	130	VAL
1	B	186	LEU
1	B	187	ARG
1	B	191	ARG
1	B	234	LEU
1	B	237	SER
1	B	239	GLU
1	B	253	GLU
1	B	271	GLU
1	B	272	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	75	ASN
1	A	183	GLN
1	A	185	HIS
1	A	199	GLN
1	A	221	GLN
1	A	250	ASN
1	B	68	ASN
1	B	185	HIS
1	B	217	ASN
1	B	221	GLN
1	B	250	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

6 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	1002[A]	-	4,4,4	1.61	0	6,6,6	0.45	0
2	PO4	B	1004[B]	-	4,4,4	2.01	2 (50%)	6,6,6	0.47	0
2	PO4	B	1004[A]	-	4,4,4	1.66	0	6,6,6	0.44	0
2	PO4	A	1002[B]	-	4,4,4	1.98	2 (50%)	6,6,6	0.46	0
2	PO4	A	1001	-	4,4,4	0.98	0	6,6,6	0.53	0
2	PO4	B	1003	-	4,4,4	2.00	3 (75%)	6,6,6	0.45	0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1002[B]	PO4	P-O2	-2.89	1.46	1.54
2	B	1004[B]	PO4	P-O2	-2.72	1.46	1.54
2	B	1003	PO4	P-O3	-2.43	1.47	1.54
2	B	1004[B]	PO4	P-O3	-2.31	1.47	1.54
2	B	1003	PO4	P-O2	-2.19	1.48	1.54
2	A	1002[B]	PO4	P-O3	-2.08	1.48	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1003	PO4	P-O4	-2.04	1.48	1.54

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1004[B]	PO4	3	0
2	A	1002[B]	PO4	5	0
2	A	1001	PO4	1	0
2	B	1003	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	287/306 (93%)	-1.27	0 100 100	33, 57, 70, 74	1 (0%)
1	B	290/306 (94%)	-1.26	0 100 100	32, 58, 70, 76	1 (0%)
All	All	577/612 (94%)	-1.27	0 100 100	32, 57, 71, 76	2 (0%)

There are no RSRZ outliers to report.

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 oligosaccharides 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. 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	B-factors(Å ²)	Q<0.9
2	PO4	B	1004[A]	5/5	0.97	0.07	72,72,73,74	4
2	PO4	B	1004[B]	5/5	0.97	0.07	70,70,70,74	4
2	PO4	A	1002[A]	5/5	0.98	0.06	70,70,71,71	4
2	PO4	A	1002[B]	5/5	0.98	0.06	66,66,66,70	4
2	PO4	A	1001	5/5	0.99	0.04	88,88,89,92	0
2	PO4	B	1003	5/5	0.99	0.04	103,103,103,108	0

6.5 Other polymers [i](#)

There are no such residues in this entry.