



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:37 AM UTC

PDB ID : 1OTY / pdb_00001oty
Title : Native PNP +ALLO
Authors : Ealick, S.E.; Bennett, E.M.; Anand, R.; Secrist, J.A.; Parker, W.B.; Hassan, A.E.; Allan, P.W.; McPherson, D.T.; Sorscher, E.J.
Deposited on : 2003-03-23
Resolution : 2.50 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

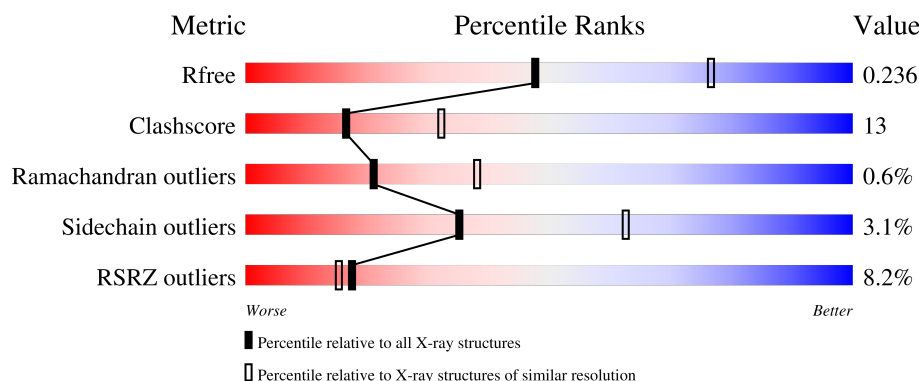
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	238	9% 72% 26% .
1	B	238	8% 77% 20% .
1	C	238	8% 71% 27% .

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	6MP	A	247	-	X	-	-
3	6MP	B	246	-	X	-	-
3	6MP	C	245	-	X	-	-

2 Entry composition [i](#)

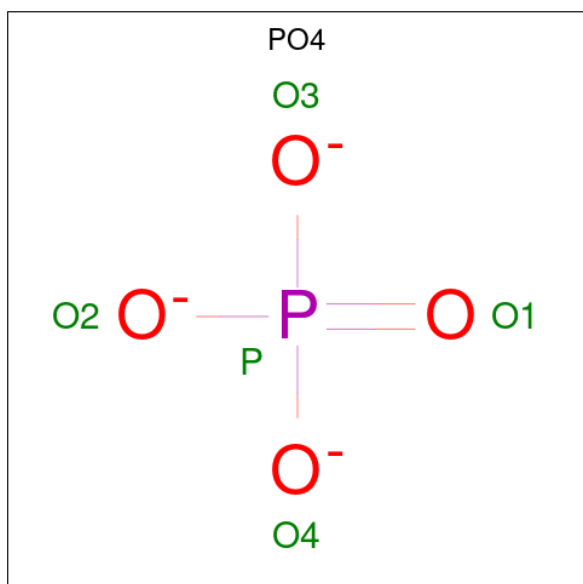
There are 4 unique types of molecules in this entry. The entry contains 5598 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Purine nucleoside phosphorylase.

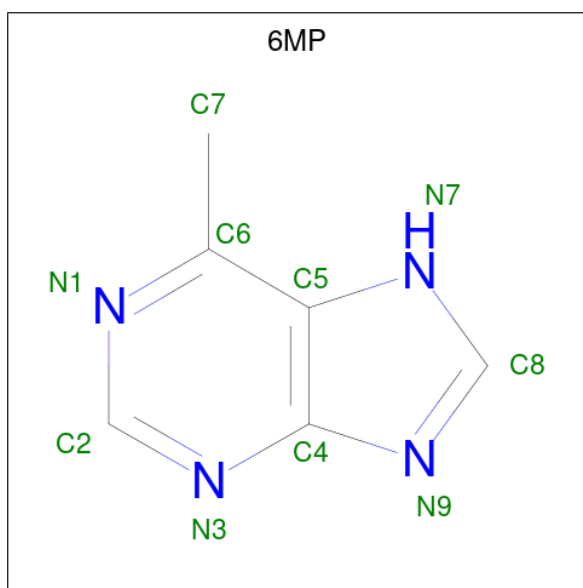
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	237	Total	C	N	O	S	0	0	0
			1793	1132	307	339	15			
1	B	237	Total	C	N	O	S	0	0	0
			1793	1132	307	339	15			
1	C	237	Total	C	N	O	S	0	0	0
			1793	1132	307	339	15			

- 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	B	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is 6-METHYLPURINE (CCD ID: 6MP) (formula: $C_6H_6N_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			10	6	4		
3	B	1	Total	C	N	0	0
			10	6	4		
3	C	1	Total	C	N	0	0
			10	6	4		

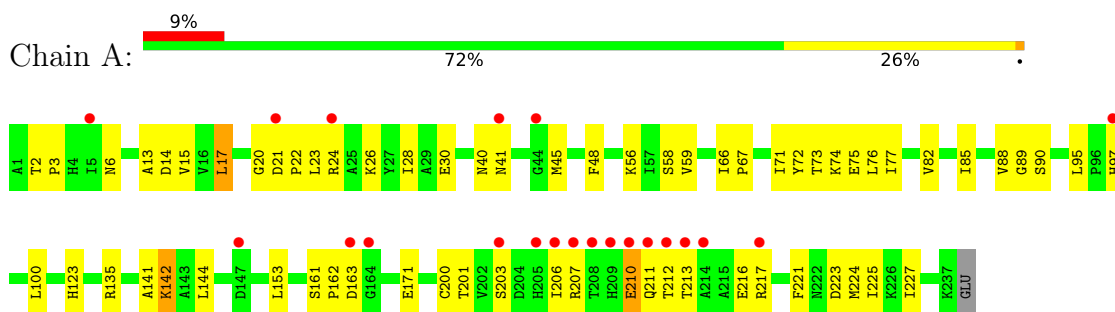
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	51	Total	O	0	0
			51	51		
4	B	74	Total	O	0	0
			74	74		
4	C	49	Total	O	0	0
			49	49		

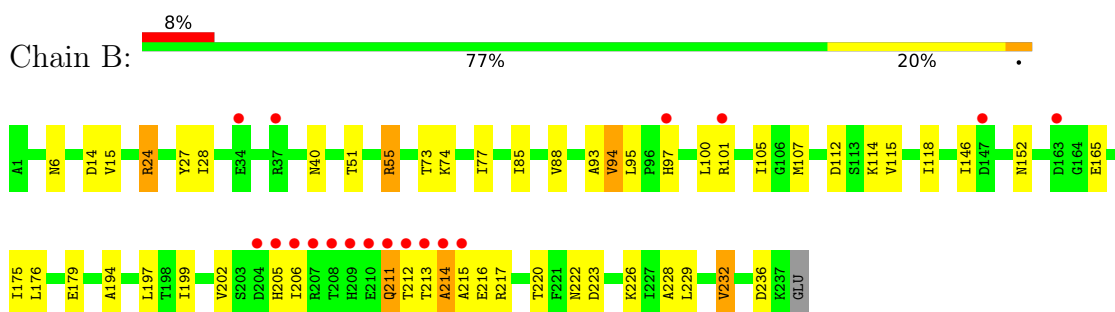
3 Residue-property plots [i](#)

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.

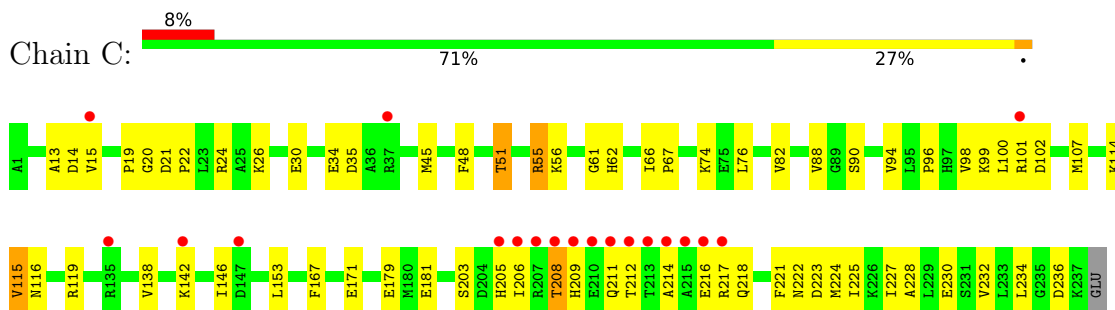
• Molecule 1: Purine nucleoside phosphorylase



• Molecule 1: Purine nucleoside phosphorylase



• Molecule 1: Purine nucleoside phosphorylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	119.97Å 119.97Å 240.46Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	26.20 – 2.50 26.20 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (26.20-2.50) 99.8 (26.20-2.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.56 (at 2.50Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.203 , 0.236 0.203 , 0.236	Depositor DCC
R_{free} test set	3603 reflections (9.98%)	wwPDB-VP
Wilson B-factor (Å ²)	31.3	Xtriage
Anisotropy	0.309	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 44.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5598	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.72% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, 6MP

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	0.42	0/1822	0.91	1/2457 (0.0%)
1	B	0.41	0/1822	0.94	3/2457 (0.1%)
1	C	0.40	0/1822	0.94	3/2457 (0.1%)
All	All	0.41	0/5466	0.93	7/7371 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	115	VAL	N-CA-C	6.80	117.56	110.62
1	A	153	LEU	N-CA-C	-6.17	101.07	110.14
1	C	115	VAL	N-CA-C	5.83	116.48	110.36
1	C	208	THR	N-CA-C	5.31	117.82	111.71
1	C	153	LEU	N-CA-C	-5.24	102.44	110.14
1	B	107	MET	N-CA-C	-5.20	105.78	111.82
1	B	215	ALA	N-CA-C	-5.13	105.87	111.82

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	1793	0	1791	50	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1793	0	1791	38	0
1	C	1793	0	1791	52	1
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
3	A	10	0	6	0	0
3	B	10	0	6	0	0
3	C	10	0	6	0	0
4	A	51	0	0	1	0
4	B	74	0	0	0	0
4	C	49	0	0	0	0
All	All	5598	0	5391	139	1

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 13.

All (139) 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:21:ASP:HB3	1:A:24:ARG:HD3	1.59	0.84
1:A:142:LYS:HE3	1:A:142:LYS:HA	1.59	0.84
1:A:210:GLU:HG2	1:A:211:GLN:H	1.44	0.80
1:C:94:VAL:HG12	1:C:171:GLU:HB2	1.65	0.79
1:B:101:ARG:HA	1:B:220:THR:HG21	1.66	0.77
1:A:223:ASP:O	1:A:227:ILE:HG12	1.85	0.76
1:A:73:THR:O	1:A:77:ILE:HG12	1.87	0.73
1:C:99:LYS:HD2	1:C:99:LYS:H	1.52	0.73
1:C:55:ARG:HD2	1:C:236:ASP:OD2	1.87	0.73
1:A:22:PRO:HG3	1:A:45:MET:SD	2.30	0.71
1:C:88:VAL:HB	1:C:224:MET:HE3	1.73	0.70
1:B:55:ARG:HD2	1:B:236:ASP:OD2	1.91	0.70
1:C:22:PRO:HG3	1:C:45:MET:SD	2.31	0.70
1:A:141:ALA:HA	1:A:227:ILE:HD12	1.73	0.69
1:C:211:GLN:H	1:C:211:GLN:CD	2.01	0.68
1:C:146:ILE:HD13	1:C:223:ASP:HB3	1.75	0.67
1:A:14:ASP:OD1	1:A:15:VAL:HG23	1.95	0.66
1:C:94:VAL:CG1	1:C:171:GLU:HB2	2.26	0.66
1:B:24:ARG:O	1:B:28:ILE:HG12	1.97	0.65
1:A:20:GLY:H	1:A:24:ARG:NH2	1.95	0.65
1:A:24:ARG:O	1:A:28:ILE:HG12	1.97	0.65
1:A:144:LEU:HD12	1:A:227:ILE:HD13	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:165:GLU:H	1:B:165:GLU:CD	2.05	0.64
1:A:141:ALA:HA	1:A:227:ILE:CD1	2.30	0.62
1:C:26:LYS:HA	1:C:48:PHE:CE2	2.35	0.62
1:C:99:LYS:HG2	1:C:102:ASP:OD1	2.00	0.62
1:B:73:THR:HG22	1:B:85:ILE:HD13	1.82	0.62
1:C:74:LYS:HD3	1:C:74:LYS:C	2.26	0.60
1:B:229:LEU:O	1:B:232:VAL:HG13	2.02	0.60
1:A:26:LYS:O	1:A:30:GLU:HG2	2.02	0.59
1:B:6:ASN:H	1:B:40:ASN:ND2	1.99	0.59
1:C:100:LEU:HD12	1:C:100:LEU:H	1.67	0.59
1:B:14:ASP:OD2	1:B:15:VAL:HG23	2.03	0.58
1:A:6:ASN:H	1:A:40:ASN:ND2	2.01	0.58
1:C:99:LYS:HD2	1:C:99:LYS:N	2.17	0.58
1:C:14:ASP:OD1	1:C:15:VAL:HG13	2.03	0.58
1:B:146:ILE:HD13	1:B:223:ASP:HB3	1.85	0.58
1:B:95:LEU:HD12	1:B:176:LEU:HD23	1.87	0.57
1:A:123:HIS:HD2	1:C:114:LYS:NZ	2.04	0.56
1:A:210:GLU:HG2	1:A:211:GLN:N	2.19	0.56
1:A:21:ASP:H	1:A:24:ARG:NH1	2.03	0.56
1:B:226:LYS:NZ	1:B:226:LYS:HB3	2.21	0.56
1:C:205:HIS:HB2	1:C:208:THR:OG1	2.06	0.56
1:A:74:LYS:C	1:A:74:LYS:HD3	2.31	0.56
1:B:101:ARG:NH1	1:B:216:GLU:HB3	2.21	0.54
1:C:76:LEU:HB3	1:C:82:VAL:HG21	1.87	0.54
1:B:197:LEU:HD21	1:B:199:ILE:HD11	1.88	0.54
1:C:90:SER:HB2	1:C:203:SER:OG	2.07	0.54
1:C:62:HIS:CE1	1:C:181:GLU:HG2	2.43	0.54
1:B:105:ILE:HG12	1:B:199:ILE:HD12	1.91	0.53
1:B:146:ILE:CD1	1:B:223:ASP:HB3	2.39	0.53
1:C:138:VAL:O	1:C:142:LYS:HG2	2.08	0.53
1:C:212:THR:HB	1:C:217:ARG:CD	2.39	0.53
1:C:101:ARG:HD3	1:C:216:GLU:HB3	1.91	0.53
1:B:74:LYS:HD3	1:B:74:LYS:C	2.34	0.53
1:C:88:VAL:HG23	1:C:88:VAL:O	2.10	0.52
1:A:141:ALA:CA	1:A:227:ILE:HD12	2.39	0.52
1:C:167:PHE:CE2	1:C:206:ILE:HD13	2.45	0.52
1:A:221:PHE:O	1:A:225:ILE:HG12	2.11	0.51
1:B:212:THR:HG23	1:B:217:ARG:HD2	1.93	0.51
1:C:221:PHE:O	1:C:225:ILE:HG12	2.11	0.50
1:A:100:LEU:H	1:A:100:LEU:HD12	1.76	0.50
1:A:212:THR:OG1	1:A:217:ARG:HD3	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:88:VAL:HG23	1:B:88:VAL:O	2.12	0.49
1:C:100:LEU:HD12	1:C:100:LEU:N	2.26	0.49
1:A:100:LEU:HD12	1:A:100:LEU:N	2.27	0.49
1:B:93:ALA:HB2	1:B:202:VAL:HG11	1.94	0.49
1:B:6:ASN:H	1:B:40:ASN:HD22	1.61	0.49
1:B:114:LYS:O	1:B:118:ILE:HG12	2.12	0.48
1:A:21:ASP:OD2	1:A:23:LEU:HB2	2.13	0.48
1:B:93:ALA:HB3	1:B:205:HIS:HA	1.95	0.48
1:C:230:GLU:O	1:C:234:LEU:HD13	2.14	0.48
1:B:165:GLU:CD	1:B:165:GLU:N	2.72	0.48
1:A:210:GLU:CG	1:A:211:GLN:H	2.17	0.48
1:A:95:LEU:HD23	1:A:171:GLU:HG3	1.96	0.48
1:B:226:LYS:HB3	1:B:226:LYS:HZ3	1.78	0.47
1:C:214:ALA:HB3	1:C:216:GLU:OE1	2.14	0.47
1:B:27:TYR:HD1	1:B:28:ILE:HD13	1.78	0.47
1:B:213:THR:O	1:B:214:ALA:HB2	2.15	0.47
1:A:95:LEU:HD13	1:A:97:HIS:CE1	2.50	0.47
1:A:88:VAL:HB	1:A:224:MET:HE3	1.97	0.46
1:C:218:GLN:HG2	1:C:222:ASN:ND2	2.31	0.46
1:B:222:ASN:O	1:B:226:LYS:HG3	2.16	0.46
1:B:100:LEU:N	1:B:100:LEU:HD22	2.30	0.46
1:C:228:ALA:O	1:C:232:VAL:HG23	2.16	0.46
1:A:2:THR:HB	1:A:3:PRO:HD2	1.97	0.46
1:C:116:ASN:ND2	1:C:119:ARG:HH11	2.14	0.46
1:A:45:MET:HE3	1:A:72:TYR:CE2	2.51	0.46
1:C:214:ALA:HB3	1:C:216:GLU:CD	2.41	0.46
1:A:100:LEU:CD2	1:A:210:GLU:HB3	2.46	0.46
1:C:223:ASP:O	1:C:227:ILE:HG13	2.15	0.46
1:C:138:VAL:HG13	1:C:142:LYS:NZ	2.32	0.45
1:C:21:ASP:HA	1:C:22:PRO:HD3	1.87	0.45
1:C:26:LYS:O	1:C:30:GLU:HG3	2.16	0.45
1:C:34:GLU:O	1:C:35:ASP:HB2	2.17	0.45
1:A:2:THR:HB	1:A:3:PRO:CD	2.47	0.45
1:A:89:GLY:O	1:A:200:CYS:HA	2.17	0.45
1:B:51:THR:HA	1:B:55:ARG:O	2.16	0.45
1:A:161:SER:HA	1:A:162:PRO:HD3	1.82	0.45
1:C:26:LYS:NZ	1:C:30:GLU:OE2	2.51	0.44
1:B:14:ASP:OD2	1:B:55:ARG:NH2	2.40	0.44
1:A:90:SER:HB2	1:A:203:SER:HB3	2.00	0.44
1:C:115:VAL:HG23	1:C:116:ASN:N	2.33	0.44
1:A:20:GLY:H	1:A:24:ARG:HH22	1.61	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:77:ILE:HD13	1:B:194:ALA:HB3	1.99	0.44
1:A:17:LEU:O	1:A:59:VAL:HA	2.18	0.44
1:C:211:GLN:CD	1:C:211:GLN:N	2.74	0.43
1:B:213:THR:HA	1:B:217:ARG:HD3	1.99	0.43
1:C:51:THR:HA	1:C:55:ARG:O	2.19	0.43
1:A:135:ARG:HD2	4:A:255:HOH:O	2.18	0.43
1:C:216:GLU:OE1	1:C:216:GLU:N	2.51	0.43
1:B:94:VAL:HG22	1:B:175:ILE:O	2.19	0.42
1:C:98:VAL:O	1:C:205:HIS:CE1	2.72	0.42
1:C:96:PRO:HD3	1:C:171:GLU:OE2	2.20	0.42
1:A:206:ILE:N	1:A:206:ILE:HD12	2.35	0.42
1:C:24:ARG:HH11	1:C:24:ARG:HG3	1.85	0.42
1:A:163:ASP:OD1	1:A:163:ASP:C	2.62	0.42
1:B:95:LEU:HD13	1:B:97:HIS:NE2	2.35	0.42
1:B:228:ALA:O	1:B:232:VAL:HG12	2.19	0.42
1:B:212:THR:OG1	1:B:217:ARG:HB2	2.19	0.42
1:C:13:ALA:HB2	1:C:56:LYS:HG2	2.02	0.42
1:A:48:PHE:O	1:A:58:SER:HA	2.20	0.42
1:C:218:GLN:HG2	1:C:222:ASN:HD21	1.84	0.41
1:C:208:THR:O	1:C:209:HIS:HB2	2.19	0.41
1:A:71:ILE:O	1:A:75:GLU:HG3	2.21	0.41
1:C:20:GLY:HA2	1:C:62:HIS:CE1	2.55	0.41
1:A:82:VAL:HG11	1:A:85:ILE:HD11	2.02	0.41
1:A:201:THR:HG23	1:A:224:MET:SD	2.60	0.41
1:A:213:THR:OG1	1:A:216:GLU:HG3	2.21	0.41
1:B:93:ALA:HB2	1:B:202:VAL:CG1	2.51	0.41
1:A:13:ALA:HB2	1:A:56:LYS:HG2	2.01	0.41
1:A:21:ASP:HA	1:A:22:PRO:HD3	1.90	0.41
1:A:210:GLU:O	1:A:211:GLN:HG2	2.20	0.41
1:B:206:ILE:O	1:B:206:ILE:HG22	2.20	0.41
1:C:100:LEU:O	1:C:101:ARG:HB2	2.21	0.41
1:A:66:ILE:HB	1:A:67:PRO:HD3	2.02	0.41
1:C:19:PRO:O	1:C:61:GLY:HA2	2.21	0.41
1:C:66:ILE:HB	1:C:67:PRO:HD3	2.03	0.41
1:A:76:LEU:HB3	1:A:82:VAL:HG21	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:107:MET:CE	1:C:107:MET:CE[12_555]	2.15	0.05

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	213/238 (90%)	199 (93%)	12 (6%)	2 (1%)	14	27
1	B	235/238 (99%)	220 (94%)	13 (6%)	2 (1%)	14	27
1	C	235/238 (99%)	218 (93%)	17 (7%)	0	100	100
All	All	683/714 (96%)	637 (93%)	42 (6%)	4 (1%)	21	38

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	207	ARG
1	A	210	GLU
1	B	211	GLN
1	B	214	ALA

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	162/189 (86%)	159 (98%)	3 (2%)	50	76
1	B	184/189 (97%)	176 (96%)	8 (4%)	26	51
1	C	103/189 (54%)	100 (97%)	3 (3%)	37	65
All	All	449/567 (79%)	435 (97%)	14 (3%)	35	62

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	LEU
1	A	41	ASN
1	A	142	LYS
1	B	24	ARG
1	B	55	ARG
1	B	94	VAL
1	B	112	ASP
1	B	152	ASN
1	B	179	GLU
1	B	211	GLN
1	B	232	VAL
1	C	51	THR
1	C	55	ARG
1	C	179	GLU

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

Mol	Chain	Res	Type
1	A	40	ASN
1	A	41	ASN
1	A	97	HIS
1	A	116	ASN
1	A	123	HIS
1	A	136	ASN
1	A	222	ASN
1	B	40	ASN
1	B	152	ASN
1	B	218	GLN
1	B	222	ASN
1	C	205	HIS
1	C	222	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	248	-	4,4,4	1.78	2 (50%)	6,6,6	0.47	0
2	PO4	B	249	-	4,4,4	2.08	3 (75%)	6,6,6	0.47	0
3	6MP	C	245	-	10,11,11	3.68	8 (80%)	10,15,15	3.88	7 (70%)
3	6MP	A	247	-	10,11,11	3.79	8 (80%)	10,15,15	3.77	7 (70%)
3	6MP	B	246	-	10,11,11	3.87	7 (70%)	10,15,15	3.98	7 (70%)
2	PO4	C	250	-	4,4,4	1.85	2 (50%)	6,6,6	0.47	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	6MP	B	246	-	-	-	0/2/2/2
3	6MP	A	247	-	-	-	0/2/2/2
3	6MP	C	245	-	-	-	0/2/2/2

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	246	6MP	C4-N9	8.72	1.45	1.38
3	A	247	6MP	C4-N9	8.48	1.44	1.38
3	C	245	6MP	C4-N9	8.35	1.44	1.38
3	B	246	6MP	C6-N1	4.55	1.41	1.33
3	A	247	6MP	C6-N1	4.49	1.41	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	245	6MP	C6-N1	4.22	1.41	1.33
3	B	246	6MP	C4-N3	3.73	1.43	1.35
3	A	247	6MP	C4-N3	3.54	1.43	1.35
3	C	245	6MP	C4-N3	3.43	1.42	1.35
3	B	246	6MP	C2-N3	3.17	1.39	1.33
3	B	246	6MP	C8-N9	3.06	1.38	1.32
3	C	245	6MP	C8-N9	3.03	1.38	1.32
3	A	247	6MP	C8-N9	2.94	1.38	1.32
3	C	245	6MP	C2-N3	2.92	1.39	1.33
3	A	247	6MP	C2-N3	2.91	1.39	1.33
3	B	246	6MP	C7-C6	2.84	1.54	1.50
3	A	247	6MP	C7-C6	2.78	1.54	1.50
3	A	247	6MP	C2-N1	2.75	1.38	1.33
3	B	246	6MP	C2-N1	2.64	1.38	1.33
3	C	245	6MP	C2-N1	2.62	1.38	1.33
3	C	245	6MP	C7-C6	2.52	1.54	1.50
2	B	249	PO4	P-O4	-2.40	1.47	1.54
3	A	247	6MP	C5-C4	2.36	1.44	1.40
2	B	249	PO4	P-O3	-2.29	1.48	1.54
2	C	250	PO4	P-O2	-2.20	1.48	1.54
2	B	249	PO4	P-O2	-2.18	1.48	1.54
3	C	245	6MP	C5-C4	2.10	1.43	1.40
2	A	248	PO4	P-O2	-2.05	1.48	1.54
2	A	248	PO4	P-O4	-2.02	1.48	1.54
2	C	250	PO4	P-O4	-2.02	1.48	1.54

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	246	6MP	C4-C5-N7	8.93	111.28	105.28
3	C	245	6MP	C4-C5-N7	8.67	111.10	105.28
3	A	247	6MP	C4-C5-N7	8.27	110.83	105.28
3	B	246	6MP	C5-C4-N9	-6.23	105.50	110.46
3	C	245	6MP	C5-C4-N9	-6.08	105.62	110.46
3	A	247	6MP	C5-C4-N9	-6.01	105.68	110.46
3	B	246	6MP	C5-N7-C8	-3.83	103.06	106.27
3	A	247	6MP	C5-N7-C8	-3.69	103.18	106.27
3	C	245	6MP	C5-N7-C8	-3.66	103.20	106.27
3	A	247	6MP	N7-C8-N9	2.70	117.99	113.87
3	C	245	6MP	N7-C8-N9	2.65	117.91	113.87
3	C	245	6MP	C4-N9-C8	-2.63	102.16	103.84
3	B	246	6MP	N7-C8-N9	2.63	117.89	113.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	246	6MP	C4-N9-C8	-2.46	102.28	103.84
3	A	247	6MP	C4-N9-C8	-2.40	102.31	103.84
3	B	246	6MP	N1-C2-N3	-2.36	125.01	128.58
3	C	245	6MP	N1-C2-N3	-2.25	125.17	128.58
3	A	247	6MP	N1-C2-N3	-2.16	125.31	128.58
3	B	246	6MP	N3-C4-N9	2.13	129.31	125.78
3	A	247	6MP	N3-C4-N9	2.06	129.18	125.78
3	C	245	6MP	N3-C4-N9	2.05	129.17	125.78

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	237/238 (99%)	0.30	21 (8%) 15 13	13, 32, 73, 119	0
1	B	237/238 (99%)	0.14	18 (7%) 20 17	13, 26, 65, 133	0
1	C	237/238 (99%)	0.21	19 (8%) 18 16	17, 32, 66, 114	0
All	All	711/714 (99%)	0.22	58 (8%) 17 15	13, 29, 71, 133	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	212	THR	11.0
1	B	214	ALA	10.2
1	B	213	THR	8.7
1	B	206	ILE	7.4
1	C	212	THR	7.2
1	B	211	GLN	6.2
1	C	213	THR	6.1
1	A	209	HIS	5.8
1	C	214	ALA	5.6
1	A	208	THR	5.4
1	B	205	HIS	5.3
1	B	204	ASP	5.3
1	B	208	THR	5.0
1	C	211	GLN	4.9
1	A	206	ILE	4.8
1	B	209	HIS	4.5
1	A	207	ARG	4.4
1	B	207	ARG	4.3
1	B	163	ASP	4.3
1	C	215	ALA	4.3
1	C	217	ARG	4.0
1	B	147	ASP	4.0
1	A	212	THR	3.8

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Mol	Chain	Res	Type	RSRZ
1	C	210	GLU	3.8
1	A	147	ASP	3.6
1	B	97	HIS	3.5
1	B	101	ARG	3.5
1	A	210	GLU	3.5
1	C	142	LYS	3.5
1	B	210	GLU	3.3
1	C	208	THR	3.2
1	A	21	ASP	3.2
1	B	37	ARG	3.1
1	A	97	HIS	3.0
1	A	211	GLN	3.0
1	A	24	ARG	2.9
1	C	37	ARG	2.9
1	A	203	SER	2.7
1	A	217	ARG	2.7
1	B	215	ALA	2.6
1	C	206	ILE	2.5
1	A	205	HIS	2.5
1	A	214	ALA	2.5
1	A	213	THR	2.4
1	A	44	GLY	2.4
1	C	207	ARG	2.3
1	A	164	GLY	2.3
1	C	209	HIS	2.2
1	A	5	ILE	2.2
1	C	216	GLU	2.2
1	C	135	ARG	2.2
1	A	163	ASP	2.2
1	C	101	ARG	2.2
1	C	147	ASP	2.1
1	C	205	HIS	2.1
1	B	34	GLU	2.0
1	C	15	VAL	2.0
1	A	41	ASN	2.0

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

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($\text{\AA}^2$)	Q<0.9
3	6MP	A	247	10/10	0.80	0.21	57,60,62,62	0
3	6MP	B	246	10/10	0.84	0.16	55,56,58,61	0
3	6MP	C	245	10/10	0.90	0.12	37,41,43,44	0
2	PO4	A	248	5/5	0.92	0.15	58,58,60,60	0
2	PO4	C	250	5/5	0.97	0.07	36,36,37,37	0
2	PO4	B	249	5/5	0.99	0.06	27,28,29,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.