



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 10:46 am BST

PDB ID : 4XR5
Title : X-ray structure of the unliganded thymidine phosphorylase from Salmonella typhimurium at 2.05 Å resolution
Authors : Balaev, V.V.; Lashkov, A.A.; Gabdulkhakov, A.G.; Betzel, C.; Mikhailov, A.M.
Deposited on : 2015-01-20
Resolution : 2.05 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

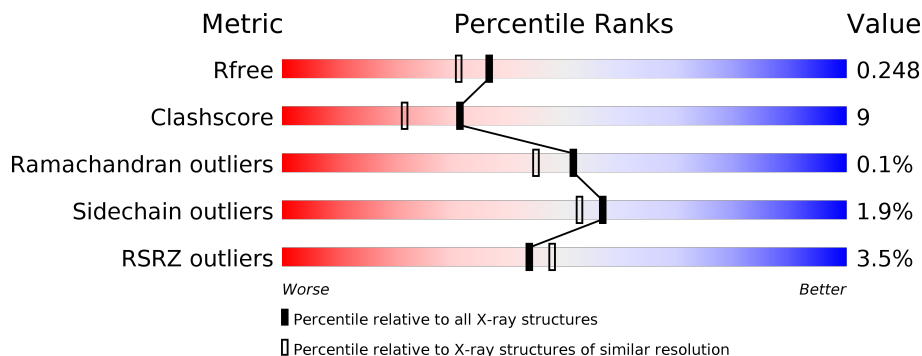
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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}	130704	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

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 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	440	 3% 83% 17%
1	B	440	 4% 75% 24%

2 Entry composition [i](#)

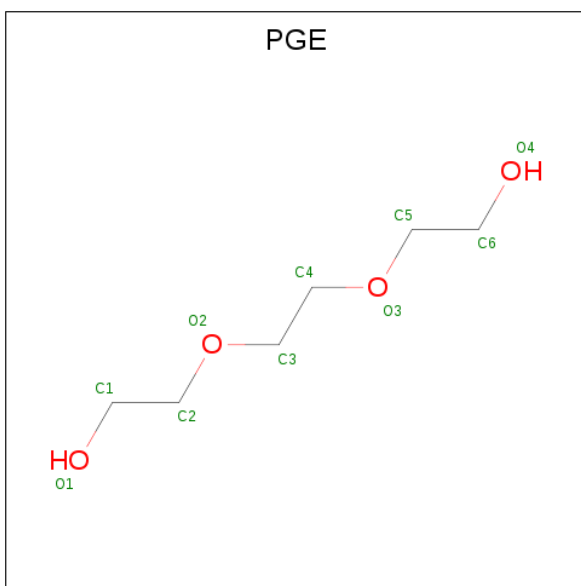
There are 5 unique types of molecules in this entry. The entry contains 6717 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Thymidine phosphorylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	440	Total 3291	C 2062	N 566	O 641	S 22	0	0	0
1	B	440	Total 3299	C 2066	N 567	O 644	S 22	0	1	0

- Molecule 2 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C₆H₁₄O₄).

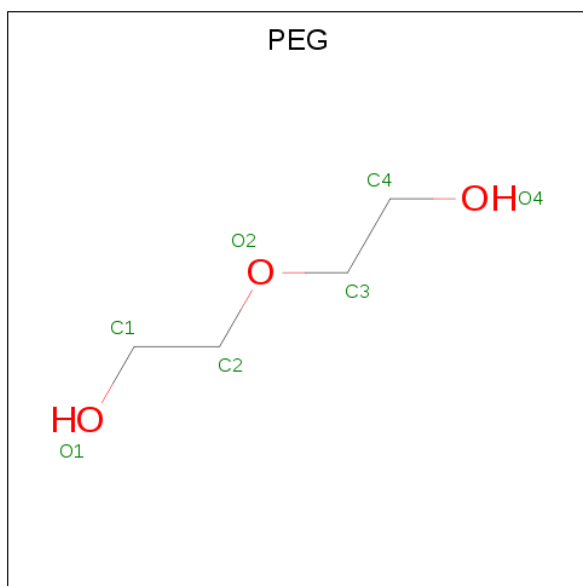


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	Total 10	C 6	O 4	0	0
2	A	1	Total 10	C 6	O 4	0	0
2	B	1	Total 10	C 6	O 4	0	0

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Cl 1 1	0	0

- Molecule 4 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C O 7 4 3	0	0

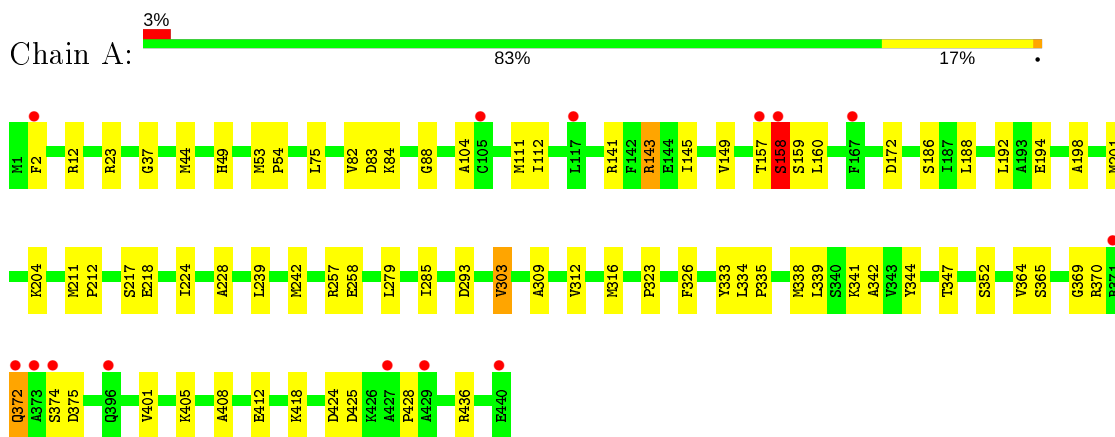
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	37	Total O 37 37	0	0
5	B	52	Total O 52 52	0	0

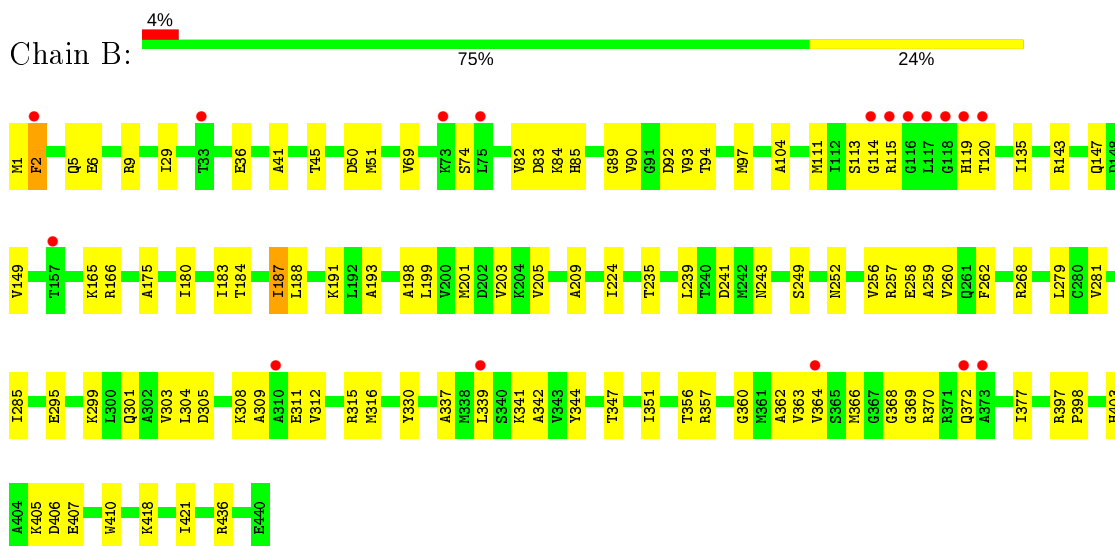
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Thymidine phosphorylase



- Molecule 1: Thymidine phosphorylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	114.94Å 171.21Å 45.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.72 – 2.05 47.72 – 2.00	Depositor EDS
% Data completeness (in resolution range)	95.5 (47.72-2.05) 95.3 (47.72-2.00)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.36 (at 2.00Å)	Xtrriage
Refinement program	PHENIX 1.9-1692	Depositor
R, R_{free}	0.221 , 0.247 0.227 , 0.248	Depositor DCC
R_{free} test set	2946 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	37.7	Xtrriage
Anisotropy	0.464	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 54.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6717	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.63% 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: PEG, PGE, CL

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	0.44	0/3339	0.63	0/4514
1	B	0.42	0/3347	0.65	0/4525
All	All	0.43	0/6686	0.64	0/9039

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	2	PHE	Peptide
1	B	2	PHE	Peptide

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	3291	0	3306	50	0
1	B	3299	0	3309	69	0
2	A	20	0	28	5	0
2	B	10	0	14	1	0
3	B	1	0	0	0	0
4	B	7	0	10	1	0
5	A	37	0	0	2	0
5	B	52	0	0	1	1
All	All	6717	0	6667	119	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 9.

All (119) 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:157:THR:O	1:A:158:SER:HB3	1.30	1.04
1:A:157:THR:OG1	1:A:159:SER:OG	1.88	0.89
1:A:372:GLN:HG2	1:A:375:ASP:N	1.89	0.88
1:A:372:GLN:HG2	1:A:375:ASP:H	1.34	0.88
1:A:157:THR:O	1:A:158:SER:CB	2.16	0.86
1:A:75:LEU:HD11	1:A:143:ARG:HG2	1.58	0.85
1:A:201:MET:HG3	1:A:224:ILE:HG21	1.59	0.84
1:B:201:MET:HG3	1:B:224:ILE:HG21	1.63	0.81
1:B:364:VAL:HA	1:B:369:GLY:H	1.51	0.75
1:A:370:ARG:HB3	1:A:375:ASP:HB2	1.69	0.74
1:A:257:ARG:NH1	2:A:502:PGE:O1	2.20	0.74
1:A:12:ARG:NH1	1:B:175:ALA:O	2.25	0.69
1:A:239:LEU:HD12	1:A:436:ARG:HG2	1.74	0.69
1:B:303:VAL:HG13	1:B:308:LYS:HB2	1.75	0.69
1:A:352:SER:HA	1:A:428:PRO:HG3	1.76	0.67
1:A:23:ARG:NH1	5:A:601:HOH:O	2.22	0.66
1:A:339:LEU:HD11	1:A:341:LYS:HD2	1.77	0.65
1:B:342:ALA:HB1	1:B:398:PRO:HB3	1.78	0.65
1:A:157:THR:HG1	1:A:159:SER:HG	1.20	0.64
1:A:364:VAL:HG22	1:A:369:GLY:HA3	1.78	0.64
1:B:363:VAL:HG12	1:B:368:GLY:HA3	1.80	0.63
1:B:249:SER:OG	1:B:258:GLU:OE1	2.12	0.63
1:A:424:ASP:OD1	1:A:425:ASP:N	2.27	0.62
1:B:308:LYS:O	1:B:312:VAL:HG23	2.00	0.62
1:B:252:ASN:O	1:B:256:VAL:HG23	1.99	0.61
1:A:84:LYS:NZ	5:A:607:HOH:O	2.30	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:113:SER:OG	1:B:114:GLY:N	2.34	0.60
1:B:119:HIS:CG	1:B:357:ARG:HG2	2.37	0.60
1:B:104:ALA:HB2	1:B:312:VAL:HG21	1.82	0.59
1:B:311:GLU:HG2	1:B:315:ARG:HH11	1.67	0.58
1:B:119:HIS:O	1:B:360:GLY:HA3	2.03	0.58
1:B:241:ASP:OD2	1:B:243:ASN:ND2	2.32	0.58
1:B:93:VAL:HG21	1:B:259:ALA:HB2	1.86	0.58
1:A:211:MET:HG3	1:A:217:SER:HA	1.85	0.58
1:B:203:VAL:HB	1:B:239:LEU:HD23	1.87	0.57
1:A:344:TYR:O	1:A:418:LYS:HE3	2.04	0.57
1:B:143:ARG:O	1:B:147:GLN:HG3	2.05	0.56
1:A:303:VAL:HG13	1:A:309:ALA:HB2	1.88	0.55
1:B:360:GLY:O	1:B:364:VAL:HG23	2.08	0.54
1:A:323:PRO:HB2	1:A:326:PHE:HB2	1.90	0.54
1:B:341:LYS:HG3	1:B:410:TRP:CE2	2.43	0.53
1:A:204:LYS:HB3	1:A:242:MET:HB2	1.90	0.53
2:A:501:PGE:H32	2:A:501:PGE:O4	2.09	0.53
1:A:218:GLU:OE2	1:A:436:ARG:NH1	2.38	0.53
1:B:312:VAL:HA	1:B:315:ARG:HD3	1.90	0.52
1:B:239:LEU:HD12	1:B:436:ARG:HG2	1.91	0.52
1:A:112:ILE:HD12	1:A:194:GLU:HG2	1.91	0.52
1:A:104:ALA:HB2	1:A:312:VAL:HG21	1.92	0.52
1:B:362:ALA:O	1:B:366:MET:HG3	2.10	0.52
1:B:1:MET:N	5:B:645:HOH:O	2.40	0.52
1:B:84:LYS:HD2	1:B:279:LEU:HD13	1.91	0.52
1:A:342:ALA:HA	1:A:401:VAL:HA	1.91	0.51
1:A:84:LYS:HD2	1:A:279:LEU:HD13	1.91	0.51
1:B:268:ARG:CZ	1:B:304:LEU:HG	2.41	0.51
1:B:90:VAL:HG21	1:B:356:THR:O	2.11	0.51
1:B:205:VAL:HA	1:B:209:ALA:HB2	1.93	0.50
1:A:37:GLY:N	1:B:36:GLU:OE2	2.34	0.50
1:A:338:MET:HE3	1:A:405:LYS:O	2.11	0.50
1:B:1:MET:HG3	1:B:2:PHE:H	1.76	0.50
1:B:82:VAL:HG13	1:B:198:ALA:HB3	1.94	0.50
1:B:303:VAL:HG12	1:B:309:ALA:HB2	1.93	0.49
1:B:370:ARG:HB3	1:B:372:GLN:O	2.12	0.49
1:B:29:ILE:O	1:B:166:ARG:HD3	2.12	0.49
1:B:83:ASP:OD1	1:B:111:MET:HG3	2.13	0.49
1:A:83:ASP:OD1	1:A:111:MET:HG3	2.13	0.49
1:B:85:HIS:NE2	1:B:187:ILE:HG12	2.28	0.49
1:B:149:VAL:HG11	1:B:316:MET:HB2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:ARG:O	1:A:49:HIS:NE2	2.44	0.48
1:B:97:MET:HG2	1:B:262:PHE:CD1	2.49	0.48
1:B:92:ASP:OD1	1:B:94:THR:HG23	2.14	0.47
1:B:337:ALA:CB	1:B:403:HIS:HB3	2.44	0.47
1:B:344:TYR:O	1:B:418:LYS:HE2	2.14	0.47
2:A:502:PGE:H32	2:A:502:PGE:H5	1.29	0.47
1:B:41:ALA:O	1:B:45:THR:HG23	2.16	0.46
1:B:69:VAL:HG13	1:B:193:ALA:HA	1.98	0.46
1:A:82:VAL:HG13	1:A:198:ALA:HB3	1.96	0.46
1:A:211:MET:HA	1:A:212:PRO:HD3	1.74	0.45
2:B:503:PGE:H22	2:B:503:PGE:H4	1.46	0.45
1:A:188:LEU:HD21	1:A:228:ALA:HB2	1.99	0.45
1:B:115:ARG:HH21	1:B:135:ILE:HD11	1.80	0.45
1:B:180:ILE:HA	1:B:183:ILE:HD12	1.99	0.45
1:B:256:VAL:O	1:B:260:VAL:HG23	2.17	0.45
1:B:50:ASP:OD1	1:B:51:MET:N	2.50	0.45
1:B:191:LYS:HA	1:B:191:LYS:HD2	1.75	0.44
1:B:364:VAL:HA	1:B:369:GLY:N	2.24	0.44
1:B:281:VAL:O	1:B:285:ILE:HG13	2.18	0.44
1:A:53:MET:HB3	1:A:54:PRO:HD3	2.00	0.44
1:B:6:GLU:OE1	1:B:9:ARG:NH1	2.51	0.43
1:A:285:ILE:HD11	1:A:293:ASP:N	2.33	0.43
1:B:199:LEU:O	1:B:235:THR:HA	2.18	0.43
1:B:257:ARG:HG3	1:B:330:TYR:CE1	2.54	0.43
1:A:149:VAL:HG11	1:A:316:MET:HB2	2.00	0.42
1:A:364:VAL:HG13	1:A:369:GLY:C	2.39	0.42
1:B:119:HIS:CD2	1:B:120:THR:HG23	2.54	0.42
1:B:370:ARG:NE	1:B:377:ILE:HG13	2.34	0.42
1:A:188:LEU:O	1:A:192:LEU:HG	2.19	0.42
1:A:372:GLN:NE2	1:A:374:SER:OG	2.53	0.42
1:B:397:ARG:HA	1:B:398:PRO:HD3	1.69	0.42
1:B:405:LYS:HE3	1:B:405:LYS:HB3	1.69	0.42
1:A:88:GLY:O	1:A:204:LYS:HG3	2.20	0.41
1:A:258:GLU:HB2	2:A:502:PGE:C6	2.50	0.41
1:B:165:LYS:HE2	1:B:165:LYS:HB2	1.71	0.41
1:A:372:GLN:HG3	1:A:374:SER:H	1.84	0.41
1:B:89:GLY:HA2	1:B:120:THR:HG22	2.02	0.41
1:B:312:VAL:HA	1:B:315:ARG:CD	2.51	0.41
1:A:334:LEU:HA	1:A:335:PRO:HD3	1.99	0.41
1:B:119:HIS:ND1	1:B:357:ARG:HG2	2.36	0.41
1:B:339:LEU:HD21	1:B:341:LYS:HE2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:301:GLN:HE21	1:B:305[A]:ASP:CG	2.25	0.41
1:A:172:ASP:O	1:B:5:GLN:NE2	2.53	0.41
1:A:408:ALA:O	1:A:412:GLU:HG2	2.21	0.41
1:B:74:SER:HB3	4:B:502:PEG:H12	2.02	0.41
1:A:333:TYR:CZ	2:A:501:PGE:H6	2.56	0.40
1:B:184:THR:O	1:B:188:LEU:HB2	2.20	0.40
1:B:337:ALA:HB2	1:B:403:HIS:HB3	2.01	0.40
1:A:141:ARG:O	1:A:145:ILE:HG13	2.21	0.40
1:A:159:SER:HG	1:A:159:SER:H	1.25	0.40
1:B:351:ILE:HG23	1:B:421:ILE:HG23	2.03	0.40
1:B:295:GLU:OE2	1:B:299:LYS:NZ	2.55	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)
5:B:601:HOH:O	5:B:601:HOH:O[2_655]	2.16	0.04

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	438/440 (100%)	425 (97%)	12 (3%)	1 (0%)	47	39
1	B	439/440 (100%)	433 (99%)	6 (1%)	0	100	100
All	All	877/880 (100%)	858 (98%)	18 (2%)	1 (0%)	51	45

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	158	SER

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	344/344 (100%)	335 (97%)	9 (3%)	46	39
1	B	345/344 (100%)	341 (99%)	4 (1%)	71	70
All	All	689/688 (100%)	676 (98%)	13 (2%)	57	53

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	MET
1	A	143	ARG
1	A	158	SER
1	A	160	LEU
1	A	186	SER
1	A	303	VAL
1	A	347	THR
1	A	365	SER
1	A	372	GLN
1	B	187	ILE
1	B	347	THR
1	B	406	ASP
1	B	407	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	119	HIS
1	A	372	GLN
1	B	301	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 for Mogul analysis.

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	PGE	B	503	-	9,9,9	0.28	0	8,8,8	0.37	0
4	PEG	B	502	-	6,6,6	0.58	0	5,5,5	0.55	0
2	PGE	A	502	-	9,9,9	0.27	0	8,8,8	0.84	0
2	PGE	A	501	-	9,9,9	0.37	0	8,8,8	0.25	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	PGE	B	503	-	-	5/7/7/7	-
4	PEG	B	502	-	-	2/4/4/4	-
2	PGE	A	502	-	-	4/7/7/7	-
2	PGE	A	501	-	-	5/7/7/7	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	502	PGE	C3-C4-O3-C5
2	A	501	PGE	C4-C3-O2-C2
2	B	503	PGE	C4-C3-O2-C2
2	A	501	PGE	O1-C1-C2-O2
2	A	501	PGE	O3-C5-C6-O4
2	A	501	PGE	O2-C3-C4-O3
2	B	503	PGE	O1-C1-C2-O2
2	B	503	PGE	O3-C5-C6-O4
4	B	502	PEG	O1-C1-C2-O2
4	B	502	PEG	O2-C3-C4-O4
2	A	502	PGE	O1-C1-C2-O2
2	A	501	PGE	C3-C4-O3-C5
2	B	503	PGE	C1-C2-O2-C3
2	A	502	PGE	O3-C5-C6-O4
2	A	502	PGE	C4-C3-O2-C2
2	B	503	PGE	O2-C3-C4-O3

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	503	PGE	1	0
4	B	502	PEG	1	0
2	A	502	PGE	3	0
2	A	501	PGE	2	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	440/440 (100%)	0.22	14 (3%) 47 52	36, 49, 68, 85	0
1	B	440/440 (100%)	0.47	17 (3%) 39 42	27, 51, 77, 91	0
All	All	880/880 (100%)	0.35	31 (3%) 44 48	27, 50, 73, 91	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	117	LEU	9.0
1	B	116	GLY	7.9
1	B	118	GLY	6.6
1	A	372	GLN	5.1
1	A	117	LEU	4.6
1	B	119	HIS	4.5
1	B	115	ARG	4.5
1	A	429	ALA	4.5
1	A	2	PHE	4.3
1	A	157	THR	4.1
1	B	339	LEU	4.0
1	A	374	SER	3.5
1	B	373	ALA	3.3
1	A	158	SER	3.3
1	A	373	ALA	3.1
1	B	372	GLN	3.1
1	B	157	THR	2.7
1	A	440	GLU	2.5
1	B	364	VAL	2.4
1	A	371	ARG	2.3
1	B	33	THR	2.3
1	B	73	LYS	2.3
1	B	2	PHE	2.3
1	B	310	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	105	CYS	2.2
1	B	120	THR	2.2
1	A	427	ALA	2.1
1	B	75	LEU	2.1
1	B	114	GLY	2.1
1	A	167	PHE	2.1
1	A	396	GLN	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. 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	PGE	A	501	10/10	0.69	0.21	52,63,67,69	0
4	PEG	B	502	7/7	0.72	0.23	55,57,67,69	0
2	PGE	A	502	10/10	0.83	0.16	51,53,55,56	0
2	PGE	B	503	10/10	0.89	0.14	62,66,73,74	0
3	CL	B	501	1/1	0.91	0.13	60,60,60,60	0

6.5 Other polymers [i](#)

There are no such residues in this entry.