



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:34 PM UTC

PDB ID : 4DI4 / pdb_00004di4
Title : Crystal structure of a 3:1 complex of Treponema pallidum TatP(T) (Tp0957) bound to TatT (Tp0956)
Authors : Brautigam, C.A.; Deka, R.K.; Norgard, M.V.
Deposited on : 2012-01-30
Resolution : 2.71 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

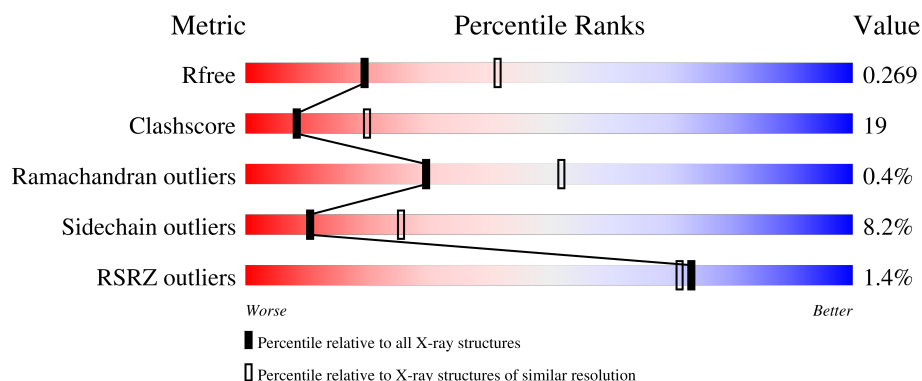
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.71 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	4348 (2.74-2.70)
Clashscore	190562	4665 (2.74-2.70)
Ramachandran outliers	187476	4584 (2.74-2.70)
Sidechain outliers	187428	4585 (2.74-2.70)
RSRZ outliers	180081	4348 (2.74-2.70)

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	301	
2	B	324	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4572 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 TatT (Tp0956).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	273	Total	C	N	O	S	0	0	0
			2166	1386	374	398	8			

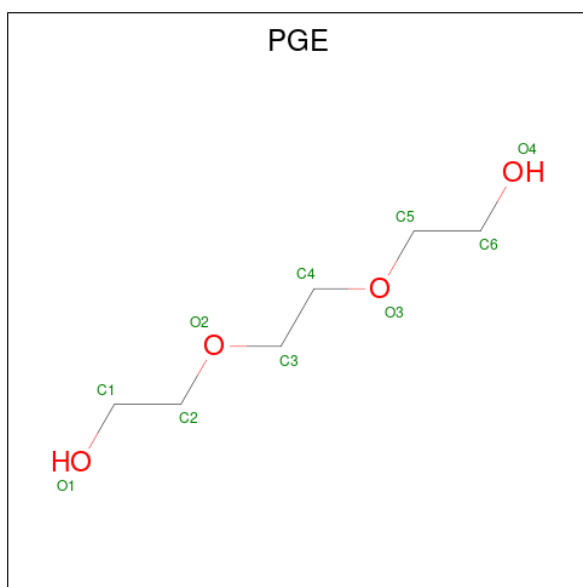
- Molecule 2 is a protein called TatP(T) (Tp0957).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	300	Total	C	N	O	S	0	0	0
			2373	1524	413	425	11			

There are 3 discrepancies between the modelled and reference sequences:

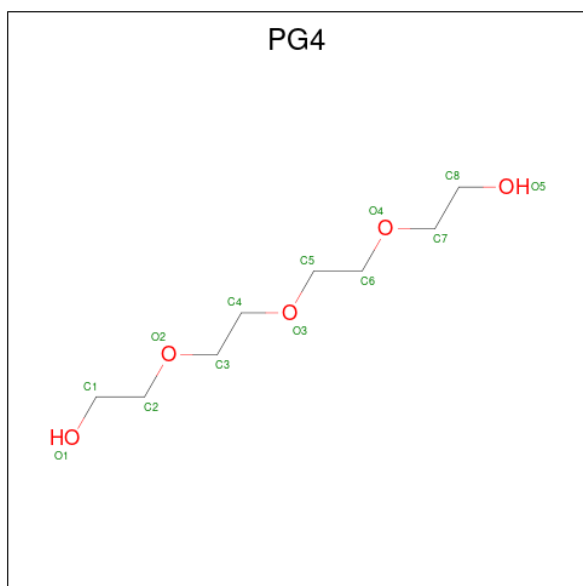
Chain	Residue	Modelled	Actual	Comment	Reference
B	5	GLY	-	expression tag	UNP O83923
B	6	ARG	-	expression tag	UNP O83923
B	185	VAL	ALA	engineered mutation	UNP O83923

- Molecule 3 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



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

- Molecule 4 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			13	8	5		

- Molecule 5 is water.

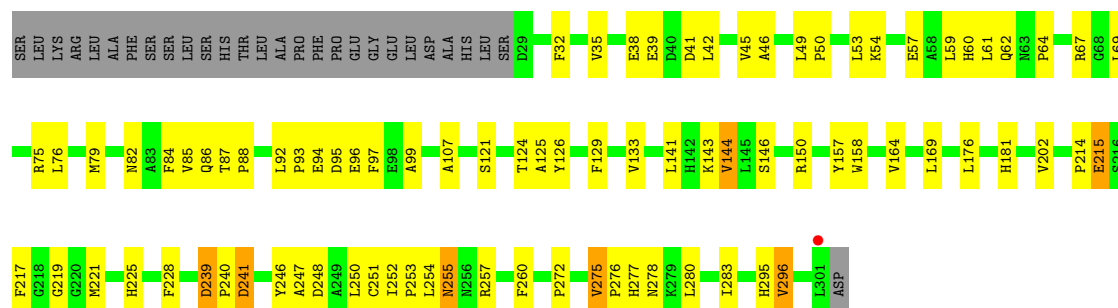
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	8	Total 8	O 8	0	0
5	B	2	Total 2	O 2	0	0

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.

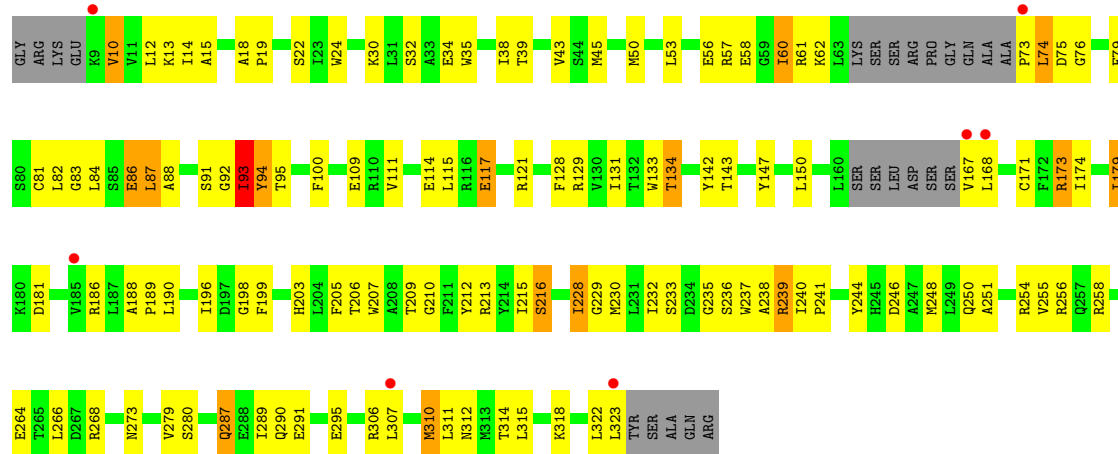
• Molecule 1: TatT (Tp0956)

Chain A: 



• Molecule 2: TatP(T) (Tp0957)

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	172.88Å 172.88Å 150.72Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.44 – 2.71 43.44 – 2.71	Depositor EDS
% Data completeness (in resolution range)	99.7 (43.44-2.71) 99.7 (43.44-2.71)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.56 (at 2.73Å)	Xtriage
Refinement program	PHENIX 1.7.1 _743	Depositor
R, R_{free}	0.194 , 0.272 0.197 , 0.269	Depositor DCC
R_{free} test set	1207 reflections (5.14%)	wwPDB-VP
Wilson B-factor (Å ²)	70.0	Xtriage
Anisotropy	0.423	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 75.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4572	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.89% 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: PG4, PGE

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	0.60	0/2224	0.94	4/3029 (0.1%)
2	B	0.51	0/2422	0.94	4/3275 (0.1%)
All	All	0.56	0/4646	0.94	8/6304 (0.1%)

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
2	B	0	1

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	B	18	ALA	CA-C-N	9.96	130.04	119.89
2	B	18	ALA	C-N-CA	9.96	130.04	119.89
1	A	275	VAL	CA-C-N	7.82	127.38	118.85
1	A	275	VAL	C-N-CA	7.82	127.38	118.85
1	A	239	ASP	CA-C-N	6.54	126.12	119.19
1	A	239	ASP	C-N-CA	6.54	126.12	119.19
2	B	10	VAL	N-CA-C	5.41	116.02	108.89
2	B	230	MET	N-CA-C	5.22	117.23	109.25

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	228	ILE	Peptide

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	2166	0	2100	55	0
2	B	2373	0	2421	120	0
3	A	10	0	14	3	0
4	A	13	0	18	5	0
5	A	8	0	0	0	0
5	B	2	0	0	0	0
All	All	4572	0	4553	174	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 19.

All (174) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:117:GLU:H	2:B:117:GLU:CD	1.61	1.08
2:B:306:ARG:O	2:B:307:LEU:HG	1.72	0.90
1:A:54:LYS:HA	1:A:54:LYS:HE2	1.56	0.85
2:B:117:GLU:OE1	2:B:117:GLU:N	2.16	0.79
1:A:53:LEU:CD2	1:A:76:LEU:HB3	2.14	0.78
2:B:50:MET:HG3	2:B:53:LEU:HD11	1.66	0.77
2:B:167:VAL:HG23	2:B:168:LEU:H	1.48	0.77
2:B:142:TYR:OH	2:B:203:HIS:ND1	2.17	0.76
2:B:287:GLN:O	2:B:291:GLU:HG3	1.86	0.76
1:A:252:ILE:HB	1:A:253:PRO:HD3	1.68	0.76
2:B:142:TYR:HH	2:B:203:HIS:HD1	1.33	0.75
2:B:62:LYS:O	2:B:73:PRO:HG2	1.89	0.73
2:B:117:GLU:CD	2:B:117:GLU:N	2.42	0.71
2:B:93:ILE:HG22	2:B:94:TYR:H	1.55	0.70
2:B:143:THR:HG21	2:B:147:TYR:HE2	1.57	0.69
2:B:15:ALA:HB1	2:B:50:MET:HE2	1.75	0.69
2:B:88:ALA:O	2:B:91:SER:N	2.23	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:15:ALA:CB	2:B:50:MET:HE2	2.24	0.68
2:B:114:GLU:C	2:B:115:LEU:HD23	2.18	0.68
2:B:30:LYS:HD2	2:B:258:ARG:HH12	1.58	0.68
2:B:22:SER:HB2	2:B:24:TRP:HE3	1.59	0.67
2:B:10:VAL:HB	2:B:43:VAL:HG22	1.75	0.67
2:B:310:MET:O	2:B:311:LEU:HD23	1.96	0.66
2:B:131:ILE:HD11	2:B:232:ILE:CG2	2.26	0.66
2:B:79:PHE:HB2	2:B:84:LEU:HD13	1.79	0.65
2:B:210:GLY:HA2	2:B:212:TYR:CE2	2.32	0.65
1:A:247:ALA:O	1:A:251:CYS:HB2	1.96	0.64
2:B:79:PHE:O	2:B:229:GLY:HA3	1.97	0.64
1:A:221:MET:HE3	1:A:225:HIS:CE1	2.32	0.64
2:B:167:VAL:HG23	2:B:168:LEU:HD12	1.79	0.63
2:B:22:SER:HB2	2:B:24:TRP:CE3	2.32	0.63
1:A:255:ASN:O	1:A:255:ASN:ND2	2.31	0.63
2:B:100:PHE:HE1	2:B:289:ILE:HG23	1.63	0.63
2:B:196:ILE:HG22	2:B:198:GLY:H	1.64	0.62
1:A:39:GLU:HG2	1:A:45:VAL:HG23	1.82	0.62
2:B:61:ARG:HB3	2:B:61:ARG:NH1	2.14	0.61
1:A:214:PRO:HD2	1:A:217:PHE:HB2	1.82	0.61
2:B:60:ILE:HD11	2:B:86:GLU:HB3	1.83	0.61
1:A:82:ASN:HD22	4:A:402:PG4:H32	1.66	0.61
2:B:237:TRP:CE3	2:B:240:ILE:HD12	2.36	0.61
2:B:131:ILE:HD11	2:B:232:ILE:HG22	1.81	0.60
2:B:244:TYR:HB3	2:B:248:MET:CE	2.32	0.60
1:A:53:LEU:HD23	1:A:76:LEU:HB3	1.83	0.59
2:B:60:ILE:HG12	2:B:87:LEU:HD12	1.83	0.59
2:B:244:TYR:HB3	2:B:248:MET:HE1	1.83	0.59
2:B:93:ILE:CG2	2:B:94:TYR:N	2.66	0.58
1:A:221:MET:HE2	1:A:254:LEU:HD11	1.85	0.58
1:A:272:PRO:O	1:A:275:VAL:HG22	2.04	0.58
2:B:93:ILE:HG22	2:B:94:TYR:N	2.18	0.58
2:B:196:ILE:CG2	2:B:198:GLY:H	2.17	0.57
2:B:30:LYS:O	2:B:30:LYS:HD3	2.04	0.57
2:B:100:PHE:CE1	2:B:289:ILE:HG23	2.40	0.56
1:A:46:ALA:HB2	1:A:84:PHE:CD1	2.40	0.56
2:B:76:GLY:HA2	2:B:232:ILE:HA	1.88	0.56
1:A:75:ARG:HG3	1:A:158:TRP:CE2	2.42	0.55
2:B:32:SER:HB2	2:B:45:MET:HE3	1.89	0.55
2:B:84:LEU:HD21	2:B:133:TRP:CZ3	2.42	0.55
2:B:87:LEU:H	2:B:87:LEU:HD13	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:150:LEU:HD23	2:B:295:GLU:CD	2.31	0.55
1:A:42:LEU:HD23	1:A:280:LEU:HD11	1.87	0.55
2:B:93:ILE:C	2:B:95:THR:N	2.62	0.54
2:B:196:ILE:HG22	2:B:198:GLY:N	2.22	0.54
2:B:83:GLY:HA2	2:B:86:GLU:HB2	1.90	0.54
2:B:237:TRP:CZ3	2:B:240:ILE:HD12	2.43	0.54
2:B:129:ARG:HB2	2:B:129:ARG:NH1	2.24	0.53
2:B:60:ILE:CD1	2:B:86:GLU:HB3	2.38	0.53
2:B:114:GLU:OE2	2:B:322:LEU:HD11	2.09	0.53
2:B:30:LYS:O	2:B:34:GLU:HB2	2.09	0.53
2:B:75:ASP:HB2	2:B:236:SER:OG	2.10	0.52
2:B:81:CYS:SG	2:B:93:ILE:HG23	2.50	0.52
2:B:115:LEU:HD23	2:B:115:LEU:N	2.23	0.51
1:A:214:PRO:HD2	1:A:217:PHE:CB	2.40	0.51
2:B:93:ILE:O	2:B:95:THR:N	2.43	0.51
2:B:39:THR:HG21	2:B:43:VAL:HB	1.93	0.51
1:A:79:MET:HG2	4:A:402:PG4:H11	1.93	0.50
2:B:75:ASP:CB	2:B:236:SER:OG	2.59	0.50
2:B:188:ALA:HB3	2:B:189:PRO:HD3	1.94	0.50
2:B:312:ASN:HB3	2:B:315:LEU:HB3	1.93	0.50
1:A:141:LEU:O	1:A:144:VAL:HG12	2.11	0.50
2:B:199:PHE:HE1	2:B:206:THR:HG22	1.77	0.50
2:B:128:PHE:CD1	2:B:233:SER:HA	2.47	0.49
2:B:235:GLY:O	2:B:238:ALA:HB3	2.12	0.49
2:B:241:PRO:HG2	2:B:244:TYR:CD2	2.47	0.49
1:A:126:TYR:O	1:A:129:PHE:HB2	2.12	0.49
2:B:39:THR:HG22	2:B:248:MET:HE1	1.94	0.49
2:B:83:GLY:O	2:B:87:LEU:HD13	2.12	0.49
2:B:173:ARG:HD2	2:B:179:ILE:HB	1.95	0.49
1:A:181:HIS:HA	3:A:401:PGE:H5	1.95	0.49
2:B:186:ARG:O	2:B:190:LEU:HG	2.13	0.49
2:B:134:THR:N	2:B:228:ILE:O	2.39	0.48
2:B:93:ILE:C	2:B:95:THR:H	2.22	0.48
1:A:86:GLN:OE1	1:A:169:LEU:HD11	2.12	0.48
1:A:82:ASN:ND2	4:A:402:PG4:H32	2.28	0.48
2:B:129:ARG:HD2	2:B:237:TRP:CD1	2.49	0.48
1:A:143:LYS:O	1:A:146:SER:OG	2.26	0.48
2:B:10:VAL:HG21	2:B:239:ARG:HB3	1.96	0.48
2:B:186:ARG:C	2:B:189:PRO:HD2	2.39	0.48
1:A:85:VAL:HG12	1:A:107:ALA:HB2	1.95	0.48
2:B:50:MET:SD	2:B:56:GLU:HA	2.53	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:57:ARG:HG2	2:B:61:ARG:NH2	2.28	0.48
2:B:94:TYR:CE1	2:B:168:LEU:HG	2.49	0.47
1:A:79:MET:HG2	4:A:402:PG4:C1	2.44	0.47
1:A:96:GLU:HG3	1:A:99:ALA:HB3	1.96	0.47
2:B:13:LYS:C	2:B:14:ILE:HD13	2.40	0.47
2:B:88:ALA:O	2:B:91:SER:CB	2.62	0.47
2:B:134:THR:HG22	2:B:228:ILE:HD11	1.96	0.47
2:B:143:THR:HG21	2:B:147:TYR:CE2	2.45	0.47
1:A:62:GLN:C	1:A:64:PRO:HD3	2.39	0.47
2:B:171:CYS:O	2:B:174:ILE:HB	2.14	0.47
2:B:15:ALA:HB3	2:B:50:MET:HE2	1.97	0.46
2:B:92:GLY:HA2	2:B:312:ASN:HB2	1.97	0.46
2:B:199:PHE:CE1	2:B:206:THR:HG22	2.51	0.46
1:A:54:LYS:HA	1:A:54:LYS:CE	2.37	0.46
1:A:67:ARG:HE	1:A:125:ALA:HB2	1.79	0.46
1:A:221:MET:HE1	1:A:250:LEU:HD22	1.97	0.46
1:A:217:PHE:O	3:A:401:PGE:H4	2.16	0.45
2:B:205:PHE:O	2:B:209:THR:HG23	2.16	0.45
1:A:38:GLU:OE2	1:A:278:ASN:N	2.41	0.45
1:A:239:ASP:OD1	1:A:241:ASP:HB2	2.17	0.45
2:B:35:TRP:CD1	2:B:248:MET:HG3	2.51	0.45
2:B:61:ARG:HB3	2:B:61:ARG:HH11	1.80	0.45
2:B:251:ALA:HA	2:B:254:ARG:CZ	2.46	0.45
2:B:266:LEU:HD12	2:B:266:LEU:O	2.17	0.45
2:B:173:ARG:CD	2:B:179:ILE:HB	2.47	0.45
2:B:142:TYR:HH	2:B:203:HIS:CE1	2.34	0.45
1:A:57:GLU:O	1:A:61:LEU:HG	2.16	0.45
2:B:57:ARG:O	2:B:60:ILE:HB	2.17	0.45
2:B:58:GLU:CD	2:B:61:ARG:HE	2.25	0.44
1:A:277:HIS:O	1:A:278:ASN:HB2	2.18	0.44
1:A:32:PHE:O	1:A:35:VAL:HG22	2.18	0.44
1:A:157:TYR:CD1	1:A:202:VAL:HG22	2.52	0.44
2:B:251:ALA:HB2	2:B:254:ARG:NH2	2.32	0.44
1:A:164:VAL:HG12	4:A:402:PG4:H82	2.00	0.44
3:A:401:PGE:H22	2:B:213:ARG:HH22	1.82	0.44
1:A:257:ARG:O	1:A:260:PHE:HB3	2.17	0.44
2:B:12:LEU:HD13	2:B:232:ILE:HD11	1.98	0.44
1:A:59:LEU:HD23	1:A:59:LEU:HA	1.65	0.44
1:A:62:GLN:O	1:A:64:PRO:HD3	2.18	0.43
2:B:117:GLU:O	2:B:121:ARG:HG3	2.18	0.43
1:A:39:GLU:HG3	1:A:41:ASP:H	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:167:VAL:HG23	2:B:168:LEU:N	2.25	0.43
1:A:221:MET:HE2	1:A:254:LEU:CD1	2.48	0.43
2:B:87:LEU:HD13	2:B:87:LEU:N	2.33	0.43
2:B:323:LEU:C	2:B:323:LEU:HD13	2.44	0.43
2:B:79:PHE:CB	2:B:84:LEU:HD13	2.46	0.43
2:B:188:ALA:O	2:B:189:PRO:C	2.61	0.43
2:B:15:ALA:N	2:B:74:LEU:HD11	2.33	0.43
2:B:111:VAL:O	2:B:115:LEU:HB2	2.18	0.43
2:B:205:PHE:C	2:B:205:PHE:CD2	2.96	0.43
1:A:275:VAL:HA	1:A:276:PRO:HD3	1.70	0.43
1:A:49:LEU:O	1:A:50:PRO:C	2.62	0.42
1:A:97:PHE:HZ	2:B:30:LYS:HG2	1.83	0.42
1:A:239:ASP:HA	1:A:240:PRO:HD3	1.78	0.42
2:B:248:MET:HE3	2:B:248:MET:HB2	1.86	0.42
2:B:216:SER:C	2:B:279:VAL:HG13	2.44	0.42
1:A:60:HIS:HA	1:A:69:LEU:HD23	2.00	0.42
2:B:84:LEU:HD21	2:B:133:TRP:HZ3	1.81	0.42
2:B:129:ARG:HB2	2:B:129:ARG:HH11	1.85	0.42
1:A:92:LEU:HA	1:A:93:PRO:HD3	1.89	0.42
2:B:19:PRO:HB2	2:B:22:SER:HB3	2.01	0.42
2:B:94:TYR:HE1	2:B:168:LEU:HG	1.84	0.42
1:A:75:ARG:HG3	1:A:158:TRP:CZ2	2.54	0.42
1:A:295:HIS:O	1:A:296:VAL:C	2.62	0.42
2:B:207:TRP:CH2	2:B:273:ASN:HB3	2.54	0.41
2:B:82:LEU:CD1	2:B:167:VAL:HG11	2.51	0.41
1:A:215:GLU:HA	1:A:219:GLY:O	2.21	0.41
2:B:268:ARG:HA	2:B:268:ARG:NE	2.35	0.41
1:A:228:PHE:CZ	1:A:246:TYR:HB2	2.56	0.41
2:B:60:ILE:HG22	2:B:61:ARG:N	2.36	0.41
2:B:109:GLU:HA	2:B:109:GLU:OE1	2.21	0.41
1:A:59:LEU:HB3	1:A:69:LEU:HD21	2.02	0.40
2:B:10:VAL:O	2:B:43:VAL:HA	2.20	0.40
2:B:134:THR:HG22	2:B:228:ILE:CD1	2.51	0.40
1:A:87:THR:HB	1:A:88:PRO:HD3	2.04	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	271/301 (90%)	264 (97%)	7 (3%)	0	100	100
2	B	294/324 (91%)	276 (94%)	16 (5%)	2 (1%)	18	39
All	All	565/625 (90%)	540 (96%)	23 (4%)	2 (0%)	30	52

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	93	ILE
2	B	94	TYR

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	223/247 (90%)	209 (94%)	14 (6%)	16	37
2	B	253/272 (93%)	228 (90%)	25 (10%)	7	18
All	All	476/519 (92%)	437 (92%)	39 (8%)	10	25

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	94	GLU
1	A	95	ASP
1	A	121	SER
1	A	124	THR

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Mol	Chain	Res	Type
1	A	133	VAL
1	A	144	VAL
1	A	150	ARG
1	A	176	LEU
1	A	215	GLU
1	A	241	ASP
1	A	248	ASP
1	A	255	ASN
1	A	283	ILE
1	A	296	VAL
2	B	38	ILE
2	B	60	ILE
2	B	74	LEU
2	B	86	GLU
2	B	87	LEU
2	B	93	ILE
2	B	117	GLU
2	B	134	THR
2	B	173	ARG
2	B	179	ILE
2	B	181	ASP
2	B	215	ILE
2	B	216	SER
2	B	239	ARG
2	B	246	ASP
2	B	250	GLN
2	B	255	VAL
2	B	256	ARG
2	B	264	GLU
2	B	280	SER
2	B	287	GLN
2	B	290	GLN
2	B	310	MET
2	B	314	THR
2	B	318	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	278	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 ⓘ

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	PGE	A	401	-	9,9,9	0.62	0	8,8,8	1.56	0
4	PG4	A	402	-	12,12,12	0.60	0	11,11,11	1.68	1 (9%)

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	PGE	A	401	-	-	4/7/7/7	-
4	PG4	A	402	-	-	8/10/10/10	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
4	A	402	PG4	O3-C4-C3	2.13	120.04	110.35

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	402	PG4	O2-C3-C4-O3
4	A	402	PG4	O3-C5-C6-O4
3	A	401	PGE	O1-C1-C2-O2
4	A	402	PG4	O1-C1-C2-O2
4	A	402	PG4	C5-C6-O4-C7
3	A	401	PGE	C4-C3-O2-C2
4	A	402	PG4	C4-C3-O2-C2
3	A	401	PGE	C1-C2-O2-C3
4	A	402	PG4	C6-C5-O3-C4
4	A	402	PG4	O4-C7-C8-O5
4	A	402	PG4	C1-C2-O2-C3
3	A	401	PGE	O3-C5-C6-O4

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	PGE	3	0
4	A	402	PG4	5	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	273/301 (90%)	-0.55	1 (0%) 88 88	41, 65, 96, 120	0
2	B	300/324 (92%)	-0.09	7 (2%) 61 58	56, 91, 136, 154	0
All	All	573/625 (91%)	-0.31	8 (1%) 73 71	41, 77, 127, 154	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	301	LEU	3.3
2	B	307	LEU	3.1
2	B	167	VAL	3.1
2	B	323	LEU	2.7
2	B	73	PRO	2.6
2	B	185	VAL	2.3
2	B	168	LEU	2.3
2	B	9	LYS	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 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	PGE	A	401	10/10	0.92	0.12	67,85,94,96	0
4	PG4	A	402	13/13	0.93	0.15	80,90,104,109	0

6.5 Other polymers [i](#)

There are no such residues in this entry.