



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:20 PM UTC

PDB ID : 4E7K / pdb_00004e7k
Title : PFV integrase Target Capture Complex (TCC-Mn), freeze-trapped prior to strand transfer, at 3.0 Å resolution
Authors : Maertens, G.N.; Cherepanov, P.
Deposited on : 2012-03-17
Resolution : 3.02 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

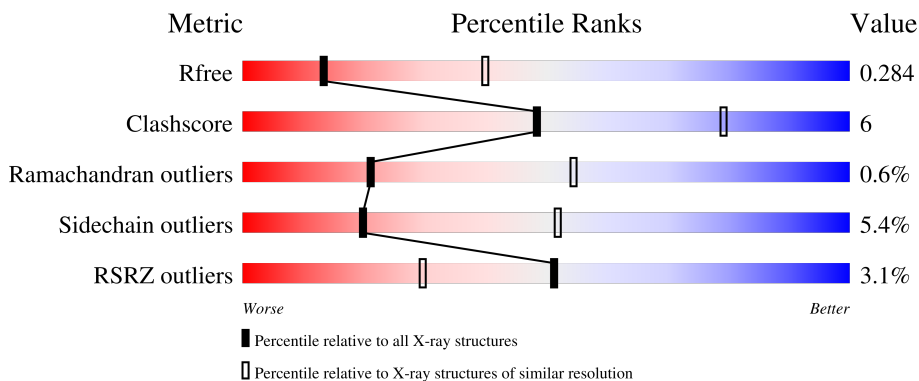
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.02 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	3131 (3.04-3.00)
Clashscore	190562	3444 (3.04-3.00)
Ramachandran outliers	187476	3319 (3.04-3.00)
Sidechain outliers	187428	3322 (3.04-3.00)
RSRZ outliers	180081	3130 (3.04-3.00)

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	395	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 74%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 18%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 7%; height: 10px; background-color: grey;"></div> </div> <p style="font-size: small; margin-top: 5px;">2% 74% 18% • 7%</p>
1	B	395	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 32%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 9%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 59%; height: 10px; background-color: grey;"></div> </div> <p style="font-size: small; margin-top: 5px;">2% 32% 9% • 59%</p>
2	C	19	<div style="display: flex; align-items: center;"> <div style="width: 5%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 68%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 32%; height: 10px; background-color: yellow;"></div> </div> <p style="font-size: small; margin-top: 5px;">5% 68% 32%</p>
3	D	17	<div style="display: flex; align-items: center;"> <div style="width: 82%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 18%; height: 10px; background-color: yellow;"></div> </div> <p style="font-size: small; margin-top: 5px;">82% 18%</p>
4	T	30	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 50%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 10%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 40%; height: 10px; background-color: grey;"></div> </div> <p style="font-size: small; margin-top: 5px;">3% 50% 10% 40%</p>

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 5307 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 Pro-Pol polyprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	366	Total	C	N	O	S	0	0	0
			2905	1863	511	527	4			
1	B	163	Total	C	N	O	S	0	0	0
			1270	833	204	232	1			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P14350
A	-1	PRO	-	expression tag	UNP P14350
A	0	GLY	-	expression tag	UNP P14350
A	217	SER	GLY	variant	UNP P14350
A	218	GLY	SER	variant	UNP P14350
B	-2	GLY	-	expression tag	UNP P14350
B	-1	PRO	-	expression tag	UNP P14350
B	0	GLY	-	expression tag	UNP P14350
B	217	SER	GLY	variant	UNP P14350
B	218	GLY	SER	variant	UNP P14350

- Molecule 2 is a DNA chain called DNA (5'-D(*AP*TP*TP*GP*TP*CP*AP*TP*GP*GP*AP*AP*TP*TP*TP*CP*GP*CP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	C	19	Total	C	N	O	P	0	0	0
			387	187	68	114	18			

- Molecule 3 is a DNA chain called DNA (5'-D(*TP*GP*CP*GP*AP*AP*AP*TP*TP*CP*CP*AP*TP*GP*AP*CP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	D	17	Total	C	N	O	P	0	0	0
			345	166	65	98	16			

- Molecule 4 is a DNA chain called DNA (5'-D(*CP*CP*CP*GP*AP*GP*GP*CP*AP*CP*GP*TP*GP*CP*TP*AP*GP*CP*AP*CP*GP*TP*GP*CP*CP*TP*CP*GP*GP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
4	T	18	369	174	69	108	18	0	0	0

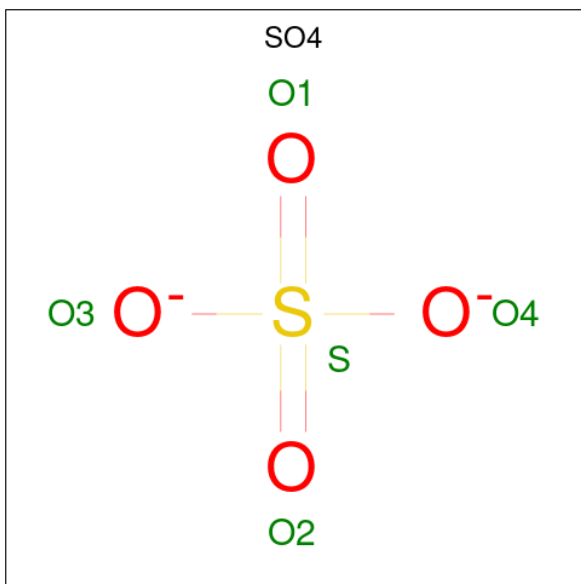
- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
5	A	1	1	1	0	0

- Molecule 6 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mn		
6	A	2	2	2	0	0
6	B	1	1	1	0	0

- Molecule 7 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
7	B	1	5	4	1	0	0

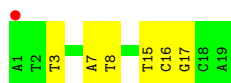
- Molecule 8 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



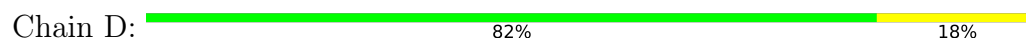
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	T	1	Total	C	O	0	0
			10	6	4		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	7	Total	O	0	0
			7	7		
9	B	5	Total	O	0	0
			5	5		



- Molecule 3: DNA (5'-D(*TP*GP*CP*GP*AP*AP*AP*TP*TP*CP*CP*AP*TP*GP*AP*CP*A)-3')



- Molecule 4: DNA (5'-D(*CP*CP*CP*GP*AP*GP*GP*CP*AP*CP*GP*TP*GP*CP*TP*AP*GP*CP*AP*CP*GP*TP*GP*CP*CP*TP*CP*GP*GP*G)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	163.91Å 163.91Å 129.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.01 – 3.02 38.01 – 3.02	Depositor EDS
% Data completeness (in resolution range)	99.5 (38.01-3.02) 99.5 (38.01-3.02)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 3.01Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.245 , 0.289 0.242 , 0.284	Depositor DCC
R_{free} test set	1739 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	71.2	Xtrriage
Anisotropy	0.320	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 32.3	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	0.91	EDS
Total number of atoms	5307	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.13% 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 i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: PGE, SO4, ZN, MN

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	0.53	0/2984	0.92	14/4074 (0.3%)
1	B	0.56	0/1310	0.96	2/1795 (0.1%)
2	C	0.37	0/433	1.09	2/667 (0.3%)
3	D	0.31	0/387	1.05	2/595 (0.3%)
4	T	0.31	0/413	0.93	2/635 (0.3%)
All	All	0.50	0/5527	0.96	22/7766 (0.3%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	15	DT	P-O5'-C5'	-7.42	108.87	120.00
1	A	178	ILE	CA-C-N	6.74	126.91	120.31
1	A	178	ILE	C-N-CA	6.74	126.91	120.31
1	A	307	THR	CA-C-N	6.59	127.16	120.38
1	A	307	THR	C-N-CA	6.59	127.16	120.38
3	D	17	DA	O4'-C1'-N9	6.50	118.15	108.40
3	D	3	DC	P-O5'-C5'	-6.49	110.26	120.00
4	T	4	DC	N1-C1'-C2'	-6.01	104.48	113.50
1	A	151	GLY	N-CA-C	-5.70	107.23	115.27
1	B	238	ARG	CA-C-N	5.55	125.64	119.87
1	B	238	ARG	C-N-CA	5.55	125.64	119.87
1	A	49	PRO	CA-C-N	5.51	125.51	119.89
1	A	49	PRO	C-N-CA	5.51	125.51	119.89
1	A	348	ASN	CA-C-N	-5.34	113.16	119.84
1	A	348	ASN	C-N-CA	-5.34	113.16	119.84
1	A	358	LEU	N-CA-C	5.32	119.77	113.12
4	T	5	DA	N9-C1'-C2'	-5.22	105.66	113.50
1	A	82	TRP	N-CA-C	5.13	112.64	108.07
2	C	15	DT	O5'-C5'-C4'	5.13	118.49	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	112	ILE	N-CA-C	5.10	116.81	109.51
1	A	176	ILE	CB-CA-C	-5.10	105.34	112.02
1	A	49	PRO	O-C-N	5.05	123.52	121.15

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	2905	0	2933	33	0
1	B	1270	0	1252	20	0
2	C	387	0	218	3	0
3	D	345	0	193	1	0
4	T	369	0	202	1	0
5	A	1	0	0	0	0
6	A	2	0	0	0	0
6	B	1	0	0	0	0
7	B	5	0	0	0	0
8	T	10	0	14	0	0
9	A	7	0	0	1	0
9	B	5	0	0	1	0
All	All	5307	0	4812	56	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 6.

All (56) 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:275:ASN:HB3	1:B:178:ILE:HG22	1.69	0.74
1:A:67:THR:HB	1:A:71:ALA:HB3	1.75	0.69
1:B:137:GLN:NE2	1:B:244:ASP:OD2	2.30	0.64
1:B:134:PRO:HG2	1:B:235:LEU:HD11	1.84	0.60
1:A:358:LEU:N	1:A:359:GLY:HA2	2.19	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:136:SER:HB2	1:B:242:TRP:CD1	2.40	0.56
1:A:127:ILE:HG22	1:A:145:VAL:HG13	1.88	0.56
1:A:324:GLN:HG2	1:A:340:PRO:HA	1.91	0.53
1:A:358:LEU:H	1:A:359:GLY:HA2	1.73	0.53
1:A:280:ASN:ND2	9:A:505:HOH:O	2.42	0.53
1:B:167:VAL:HG13	1:B:198:TRP:CD1	2.45	0.52
1:A:229:ARG:HG2	1:A:233:LYS:HE3	1.90	0.52
1:A:182:ILE:HG13	1:A:204:ILE:HG21	1.91	0.51
1:B:174:THR:HA	1:B:177:ALA:O	2.11	0.51
1:B:235:LEU:O	1:B:239:PRO:HB3	2.11	0.51
1:A:150:THR:HG21	1:A:269:LEU:HD13	1.94	0.50
1:B:266:HIS:ND1	9:B:505:HOH:O	2.32	0.50
1:B:134:PRO:O	1:B:136:SER:N	2.46	0.48
1:A:110:GLY:O	1:A:350:ARG:NH1	2.46	0.48
1:A:137:GLN:H	1:A:241:LYS:HZ2	1.61	0.48
1:B:136:SER:OG	1:B:242:TRP:N	2.48	0.47
1:A:73:LEU:HD22	1:A:86:ARG:CZ	2.45	0.47
1:A:231:LEU:HD23	1:A:231:LEU:HA	1.79	0.47
1:B:153:THR:HG21	1:B:227:ILE:HD11	1.96	0.46
2:C:16:DC:H2''	2:C:17:DG:C8	2.50	0.46
1:B:222:ARG:O	1:B:225:SER:HB3	2.16	0.46
1:A:56:LYS:O	1:A:60:GLN:HG3	2.16	0.45
1:A:257:TYR:O	1:A:259:PRO:HD3	2.15	0.45
1:B:182:ILE:HG13	1:B:204:ILE:HG21	1.98	0.45
1:B:230:LEU:HD12	1:B:230:LEU:HA	1.83	0.45
1:A:54:ARG:O	1:A:58:VAL:HG23	2.16	0.45
1:A:230:LEU:HD23	1:A:249:VAL:HG13	1.99	0.45
1:A:271:GLY:HA2	1:A:295:LEU:HD13	2.00	0.44
1:A:115:PRO:HG3	2:C:3:DT:C2	2.52	0.44
1:A:133:LEU:HD23	1:A:133:LEU:HA	1.81	0.44
1:B:127:ILE:HA	1:B:144:VAL:O	2.17	0.44
3:D:1:DT:H6	3:D:1:DT:H2'	1.63	0.43
1:A:246:LEU:HD23	1:A:246:LEU:HA	1.88	0.43
1:A:9:ASP:HA	1:A:12:LEU:HB2	2.00	0.43
1:A:97:GLN:O	1:A:101:ILE:HG23	2.19	0.43
1:B:235:LEU:HD12	1:B:235:LEU:HA	1.82	0.43
2:C:7:DA:H2''	2:C:8:DT:O5'	2.18	0.43
4:T:9:DG:H8	4:T:9:DG:H5''	1.83	0.43
1:A:163:THR:HG23	1:A:195:PHE:HB2	2.00	0.43
1:A:304:HIS:ND1	1:A:305:PRO:HD2	2.34	0.42
1:A:348:ASN:C	1:A:350:ARG:H	2.27	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:141:TYR:CE2	1:B:161:PRO:HD3	2.54	0.42
1:B:170:LEU:HD23	1:B:170:LEU:HA	1.88	0.42
1:B:234:LEU:HD13	1:B:245:LEU:HD13	2.02	0.42
1:A:322:LEU:HD23	1:A:322:LEU:HA	1.80	0.42
1:A:58:VAL:HG11	1:A:89:VAL:HA	2.01	0.41
1:A:245:LEU:O	1:A:249:VAL:HG23	2.20	0.41
1:A:51:GLN:HA	1:A:54:ARG:HD2	2.03	0.41
1:A:315:TRP:CE2	1:A:371:PRO:HD3	2.54	0.41
1:A:358:LEU:HD13	1:A:358:LEU:HA	1.86	0.41
1:B:258:SER:O	1:B:261:LEU:O	2.40	0.40

There are no symmetry-related clashes.

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	364/395 (92%)	338 (93%)	24 (7%)	2 (0%)	24	59
1	B	161/395 (41%)	152 (94%)	8 (5%)	1 (1%)	21	54
All	All	525/790 (66%)	490 (93%)	32 (6%)	3 (1%)	21	54

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	282	ASP
1	A	356	ASP
1	B	135	PRO

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	325/354 (92%)	308 (95%)	17 (5%)	21	53
1	B	139/354 (39%)	131 (94%)	8 (6%)	18	49
All	All	464/708 (66%)	439 (95%)	25 (5%)	20	52

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	TYR
1	A	113	LEU
1	A	149	MET
1	A	178	ILE
1	A	184	SER
1	A	210	THR
1	A	219	LYS
1	A	236	VAL
1	A	256	THR
1	A	307	THR
1	A	311	SER
1	A	312	SER
1	A	316	SER
1	A	329	ARG
1	A	342	THR
1	A	352	VAL
1	A	358	LEU
1	B	142	VAL
1	B	206	LEU
1	B	224	ASN
1	B	230	LEU
1	B	240	THR
1	B	248	VAL
1	B	250	GLN
1	B	258	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	GLN
1	A	78	ASN
1	A	137	GLN
1	A	296	GLN
1	A	321	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	$\# Z > 2$
7	SO4	B	402	-	4,4,4	0.22	0	6,6,6	0.14	0
8	PGE	T	101	-	9,9,9	0.72	0	8,8,8	0.59	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
8	PGE	T	101	-	-	3/7/7/7	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	T	101	PGE	C3-C4-O3-C5
8	T	101	PGE	C4-C3-O2-C2
8	T	101	PGE	O2-C3-C4-O3

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 [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	366/395 (92%)	0.05	8 (2%) 62 39	39, 58, 89, 117	0
1	B	163/395 (41%)	0.42	8 (4%) 35 18	44, 61, 88, 100	0
2	C	19/19 (100%)	0.23	1 (5%) 32 16	47, 65, 80, 90	0
3	D	17/17 (100%)	0.04	0 100 100	46, 58, 81, 92	0
4	T	18/30 (60%)	0.74	1 (5%) 30 15	59, 72, 133, 134	0
All	All	583/856 (68%)	0.18	18 (3%) 51 30	39, 59, 90, 134	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	9	ASP	4.7
1	B	237	GLY	4.1
4	T	10	DC	4.1
2	C	1	DA	3.3
1	B	212	TYR	3.0
1	A	19	HIS	3.0
1	A	29	THR	2.8
1	B	262	LYS	2.8
1	A	28	TYR	2.6
1	B	234	LEU	2.5
1	B	215	GLN	2.3
1	B	261	LEU	2.3
1	B	240	THR	2.3
1	B	239	PRO	2.1
1	A	374	HIS	2.1
1	A	282	ASP	2.1
1	A	10	ALA	2.1
1	A	211	PRO	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(Å ²)	Q<0.9
7	SO4	B	402	5/5	0.85	0.13	88,88,93,100	0
8	PGE	T	101	10/10	0.87	0.20	65,69,69,69	10
6	MN	A	403	1/1	0.98	0.04	81,81,81,81	0
6	MN	A	402	1/1	0.99	0.02	47,47,47,47	0
6	MN	B	401	1/1	0.99	0.05	52,52,52,52	0
5	ZN	A	401	1/1	1.00	0.02	53,53,53,53	0

6.5 Other polymers [i](#)

There are no such residues in this entry.