Full wwPDB X-ray Structure Validation Report (i)

Aug 28, 2023 – 02:29 PM EDT

PDB ID	:	3KD5
Title	:	Closed ternary complex of an RB69 gp43 fingers domain mutant complexed with an acyclic GMP terminated primer template pair and phosphonoformic acid.
Authors	:	Zahn, K.E.; Doublie, S.
Deposited on	:	2009-10-22
Resolution	:	2.69 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 (i) 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 (i)) were used in the production of this report:

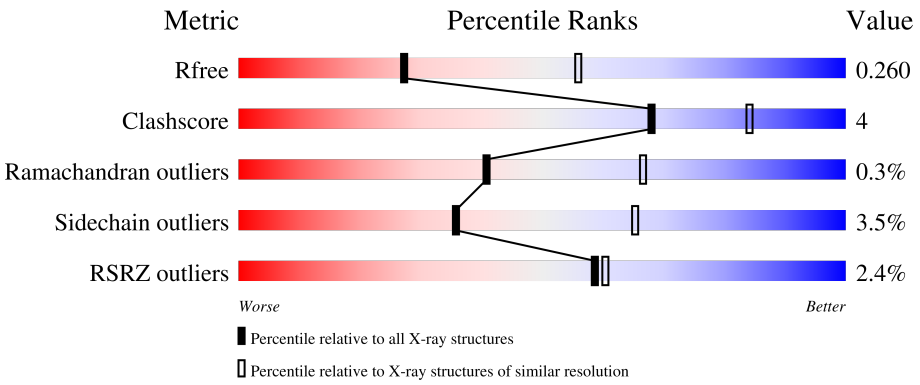
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 2.69 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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 <=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	T	18	83% 11% 6%
2	P	14	57% 36% 7%
3	E	913	2% 87% 12% ••

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 8283 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 DNA chain called DNA (5'-D(*CP*GP*TP*CP*TP*TP*AP*TP*GP*AP*CP*AP*GP*CP*CP*GP*CP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	T	18	364	174	66	107	17	0	0	0

- Molecule 2 is a DNA chain called DNA (5'-D(*GP*CP*GP*GP*CP*TP*GP*TP*CP*AP*TP*AP*AP*(4DG))-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	P	14	284	135	55	81	13	0	0	0

- Molecule 3 is a protein called DNA polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	E	905	7379	4737	1224	1383	35	0	0	0

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	222	ALA	ASP	engineered mutation	UNP Q38087
E	478	TRP	VAL	engineered mutation	UNP Q38087
E	479	VAL	PHE	engineered mutation	UNP Q38087
E	480	SER	ASN	engineered mutation	UNP Q38087
E	557	MET	ILE	engineered mutation	UNP Q38087
E	558	ALA	ASN	engineered mutation	UNP Q38087
E	559	LEU	ARG	engineered mutation	UNP Q38087
E	561	VAL	LEU	engineered mutation	UNP Q38087
E	562	THR	LEU	engineered mutation	UNP Q38087
E	563	CYS	ILE	engineered mutation	UNP Q38087
E	904	SER	-	expression tag	UNP Q38087
E	905	ALA	-	expression tag	UNP Q38087

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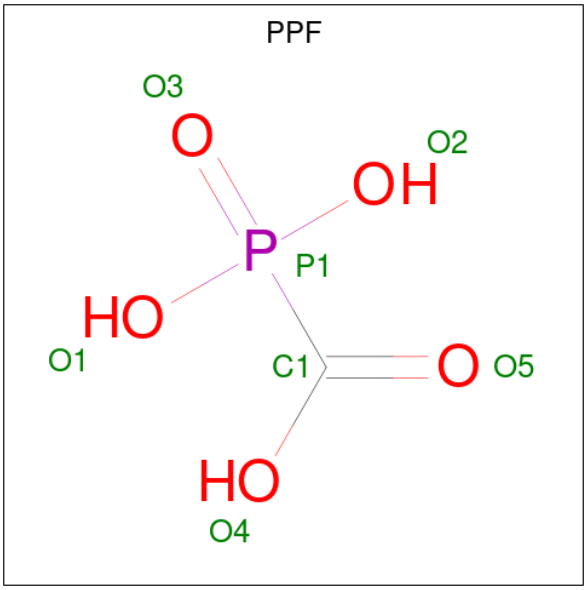
Chain	Residue	Modelled	Actual	Comment	Reference
E	906	TRP	-	expression tag	UNP Q38087
E	907	SER	-	expression tag	UNP Q38087
E	908	HIS	-	expression tag	UNP Q38087
E	909	PRO	-	expression tag	UNP Q38087
E	910	GLN	-	expression tag	UNP Q38087
E	911	PHE	-	expression tag	UNP Q38087
E	912	GLU	-	expression tag	UNP Q38087
E	913	LYS	-	expression tag	UNP Q38087

- Molecule 4 is PHOSPHONOFORMIC ACID (three-letter code: PPF) (formula: CH_3O_5P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	P		
4	E	1	7	1	5	1	0	0

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
5	E	3	3	3	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	T	30	30	30	0	0

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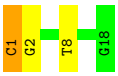
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	P	16	16	16	0	0
			Total	O		
6	E	200	200	200	0	0

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.

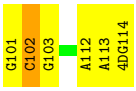
• Molecule 1: DNA (5'-D(*CP*GP*TP*CP*TP*TP*AP*TP*GP*AP*CP*AP*GP*CP*CP*GP*CP*G)-3')

Chain T:



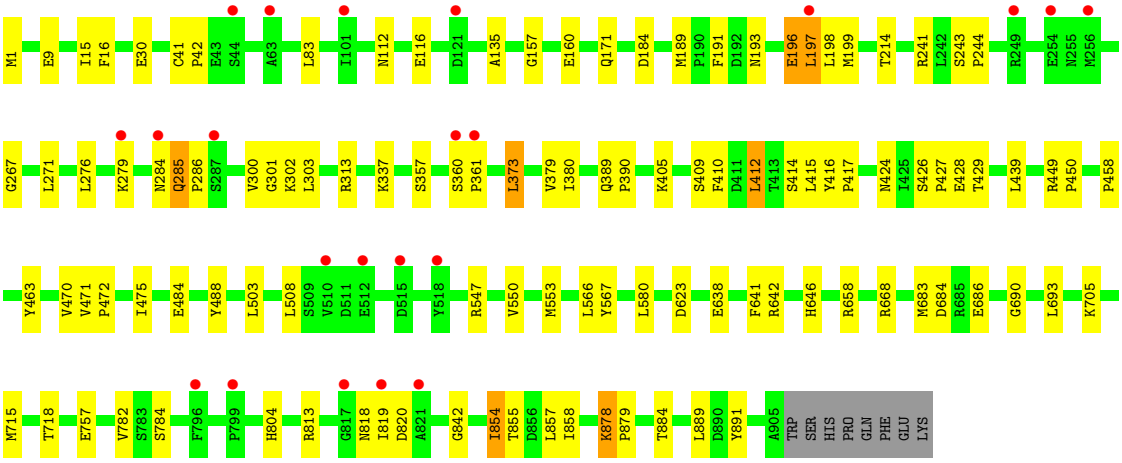
• Molecule 2: DNA (5'-D(*GP*CP*GP*GP*CP*TP*GP*TP*CP*AP*TP*AP*AP*(4DG))-3')

Chain P:



• Molecule 3: DNA polymerase

Chain E:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	77.86Å 122.48Å 133.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.69 39.04 – 2.69	Depositor EDS
% Data completeness (in resolution range)	99.2 (30.00-2.69) 99.2 (39.04-2.69)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.81 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.6.0041, CNS	Depositor
R, R_{free}	0.215 , 0.263 0.218 , 0.260	Depositor DCC
R_{free} test set	3471 reflections (9.64%)	wwPDB-VP
Wilson B-factor (Å^2)	71.8	Xtriage
Anisotropy	0.509	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 36.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	8283	wwPDB-VP
Average B, all atoms (Å^2)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: MG, 4DG, PPF

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	T	0.26	0/407	0.81	1/626 (0.2%)
2	P	0.32	0/297	0.89	2/457 (0.4%)
3	E	0.35	0/7561	0.49	0/10219
All	All	0.35	0/8265	0.53	3/11302 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	P	102	DC	P-O3'-C3'	6.51	127.51	119.70
1	T	1	DC	P-O3'-C3'	6.21	127.16	119.70
2	P	101	DG	P-O3'-C3'	5.37	126.15	119.70

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	T	364	0	204	4	0
2	P	284	0	157	4	0
3	E	7379	0	7259	54	0
4	E	7	0	0	0	0
5	E	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	E	200	0	0	4	0
6	P	16	0	0	0	0
6	T	30	0	0	1	0
All	All	8283	0	7620	60	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 4.

All (60) 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)
3:E:424:ASN:O	3:E:429:THR:HG21	1.78	0.82
3:E:279:LYS:HE2	6:E:1025:HOH:O	1.91	0.71
3:E:415:LEU:HD22	3:E:623:ASP:HB3	1.74	0.70
3:E:813:ARG:HH21	3:E:842:GLY:HA3	1.58	0.69
2:P:112:DA:H2"	2:P:113:DA:H5"	1.75	0.68
3:E:409:SER:OG	3:E:686:GLU:HB2	1.95	0.65
3:E:112:ASN:HB3	3:E:214:THR:HG23	1.79	0.63
3:E:414:SER:HB3	3:E:417:PRO:HG2	1.81	0.63
3:E:271:LEU:HB3	3:E:276:LEU:HD21	1.82	0.61
3:E:412:LEU:HB2	3:E:623:ASP:HB2	1.82	0.61
1:T:1:DC:H5"	3:E:357:SER:HA	1.83	0.60
3:E:171:GLN:HG3	3:E:303:LEU:HD21	1.84	0.59
3:E:410:PHE:HB3	3:E:683:MET:HG2	1.86	0.57
3:E:471:VAL:HG13	3:E:566:LEU:HD21	1.87	0.56
3:E:475:ILE:HD12	3:E:566:LEU:HD23	1.87	0.56
3:E:116:GLU:HB2	3:E:135:ALA:HB3	1.88	0.55
3:E:715:MET:O	3:E:718:THR:HG22	2.06	0.55
1:T:1:DC:H2"	1:T:2:DG:C8	2.43	0.54
3:E:757:GLU:HB2	3:E:889:LEU:HD22	1.90	0.52
3:E:157:GLY:O	3:E:313:ARG:NH2	2.44	0.51
3:E:373:LEU:HD23	3:E:380:ILE:HG22	1.91	0.51
3:E:337:LYS:HE2	3:E:550:VAL:HG11	1.92	0.51
3:E:684:ASP:HB2	6:E:1116:HOH:O	2.10	0.50
3:E:286:PRO:HB3	3:E:782:VAL:HG21	1.94	0.49
1:T:8:DT:H5"	3:E:705:LYS:HD3	1.94	0.49
3:E:267:GLY:HA2	6:E:933:HOH:O	2.12	0.48
3:E:193:ASN:ND2	3:E:196:GLU:H	2.12	0.47
3:E:41:CYS:HB2	3:E:42:PRO:HD2	1.97	0.47
3:E:300:VAL:HG23	3:E:301:GLY:H	1.80	0.47
3:E:300:VAL:HG23	3:E:301:GLY:N	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:102:DC:H2"	2:P:103:DG:O5'	2.15	0.46
3:E:429:THR:HG23	3:E:463:TYR:HD1	1.81	0.46
3:E:285:GLN:HE21	3:E:285:GLN:HA	1.80	0.45
3:E:428:GLU:OE2	3:E:470:VAL:HG23	2.17	0.45
1:T:1:DC:H5'	6:T:90:HOH:O	2.18	0.44
3:E:416:TYR:OH	6:E:1099:HOH:O	2.15	0.44
3:E:878:LYS:HB3	3:E:879:PRO:HD3	2.00	0.44
3:E:191:PHE:CD2	3:E:197:LEU:HD12	2.53	0.43
3:E:484:GLU:HG2	3:E:488:TYR:CE2	2.54	0.43
3:E:642:ARG:HG2	3:E:646:HIS:ND1	2.33	0.43
3:E:855:THR:HG22	3:E:857:LEU:H	1.83	0.43
2:P:102:DC:H2"	2:P:103:DG:C8	2.54	0.43
3:E:389:GLN:HA	3:E:390:PRO:HD3	1.91	0.43
3:E:373:LEU:HD11	3:E:470:VAL:HG21	2.00	0.42
3:E:471:VAL:HB	3:E:472:PRO:HD3	1.99	0.42
3:E:784:SER:O	3:E:804:HIS:CD2	2.73	0.42
3:E:9:GLU:OE1	3:E:267:GLY:N	2.52	0.42
3:E:412:LEU:HD23	3:E:415:LEU:HD13	2.02	0.42
3:E:16:PHE:CE1	3:E:30:GLU:HG3	2.55	0.41
3:E:449:ARG:HA	3:E:450:PRO:HD3	1.90	0.41
3:E:83:LEU:HB3	3:E:379:VAL:HG12	2.03	0.41
3:E:405:LYS:O	3:E:690:GLY:HA2	2.21	0.41
3:E:854:ILE:H	3:E:854:ILE:HG13	1.72	0.41
3:E:884:THR:HG21	3:E:891:TYR:HB3	2.02	0.41
3:E:426:SER:OG	3:E:427:PRO:HD2	2.21	0.41
2:P:102:DC:C2'	2:P:103:DG:C8	3.05	0.40
3:E:360:SER:HA	3:E:361:PRO:HD2	1.90	0.40
3:E:243:SER:HA	3:E:244:PRO:HD3	1.96	0.40
3:E:638:GLU:HA	3:E:641:PHE:HD2	1.85	0.40
3:E:302:LYS:HG2	3:E:303:LEU:N	2.36	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	
3	E	903/913 (99%)	855 (95%)	45 (5%)	3 (0%)	41	66

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	E	819	ILE
3	E	858	ILE
3	E	458	PRO

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	
3	E	801/809 (99%)	773 (96%)	28 (4%)	36	65

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

Mol	Chain	Res	Type
3	E	1	MET
3	E	15	ILE
3	E	160	GLU
3	E	184	ASP
3	E	189	MET
3	E	196	GLU
3	E	197	LEU
3	E	198	LEU
3	E	199	MET
3	E	241	ARG
3	E	284	ASN
3	E	285	GLN
3	E	373	LEU
3	E	412	LEU
3	E	439	LEU
3	E	503	LEU

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Mol	Chain	Res	Type
3	E	508	LEU
3	E	547	ARG
3	E	553	MET
3	E	567	TYR
3	E	580	LEU
3	E	658	ARG
3	E	668	ARG
3	E	693	LEU
3	E	818	ASN
3	E	820	ASP
3	E	854	ILE
3	E	878	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
3	E	193	ASN
3	E	207	GLN
3	E	245	HIS
3	E	284	ASN
3	E	285	GLN
3	E	444	ASN
3	E	678	GLN
3	E	775	ASN
3	E	786	ASN
3	E	787	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	4DG	P	114	2,5	12,20,21	1.25	2 (16%)	11,27,30	0.98	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
2	4DG	P	114	2,5	-	3/4/8/9	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	P	114	4DG	C5-C6	-2.51	1.42	1.47
2	P	114	4DG	C8-N7	-2.42	1.30	1.35

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	P	114	4DG	O6-C6-C5	2.13	128.53	124.37

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	P	114	4DG	C5'-C4'-O4'-C1'
2	P	114	4DG	N9-C1'-O4'-C4'
2	P	114	4DG	O4'-C4'-C5'-O5'

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 4 ligands modelled in this entry, 3 are monoatomic - leaving 1 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	$\#\|Z\| > 2$
4	PPF	E	914	5	4,6,6	1.12	0	3,9,9	1.18	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
4	PPF	E	914	5	-	0/0/6/6	-

There are no bond length outliers.

There are no bond angle outliers.

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 (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	T	18/18 (100%)	-0.31	0	100	100	43, 60, 86, 86	0
2	P	13/14 (92%)	0.07	0	100	100	41, 60, 96, 98	0
3	E	905/913 (99%)	0.28	22 (2%)	59	60	25, 36, 47, 63	0
All	All	936/945 (99%)	0.27	22 (2%)	59	60	25, 36, 50, 98	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	E	515	ASP	3.5
3	E	510	VAL	3.2
3	E	256	MET	2.9
3	E	284	ASN	2.9
3	E	197	LEU	2.8
3	E	254	GLU	2.8
3	E	512	GLU	2.8
3	E	63	ALA	2.7
3	E	518	TYR	2.7
3	E	819	ILE	2.6
3	E	796	PHE	2.5
3	E	121	ASP	2.5
3	E	817	GLY	2.3
3	E	279	LYS	2.3
3	E	101	ILE	2.3
3	E	799	PRO	2.2
3	E	821	ALA	2.1
3	E	44	SER	2.1
3	E	287	SER	2.1
3	E	360	SER	2.0
3	E	249	ARG	2.0
3	E	361	PRO	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (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	4DG	P	114	19/20	0.97	0.24	41,42,43,44	0

6.3 Carbohydrates (i)

There are no monosaccharides 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
5	MG	E	917	1/1	0.84	0.07	58,58,58,58	0
5	MG	E	915	1/1	0.92	0.09	43,43,43,43	1
5	MG	E	916	1/1	0.94	0.17	42,42,42,42	0
4	PPF	E	914	7/7	0.97	0.14	41,43,44,45	0

6.5 Other polymers (i)

There are no such residues in this entry.