wwPDB X-ray Structure Validation Summary Report (i)

Nov 6, 2023 – 07:35 PM EST

PDB ID	:	8F88
Title	:	Crystal structure of PTP1B D181A/Q262A/C215A phosphatase domain with monophosphorylated JAK2 activation loop phosphopeptide
Authors	:	Morris, R.; Kershaw, N.J.; Babon, J.J.
Deposited on	:	2022-11-21
Resolution	:	3.10 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (1)) were used in the production of this report:

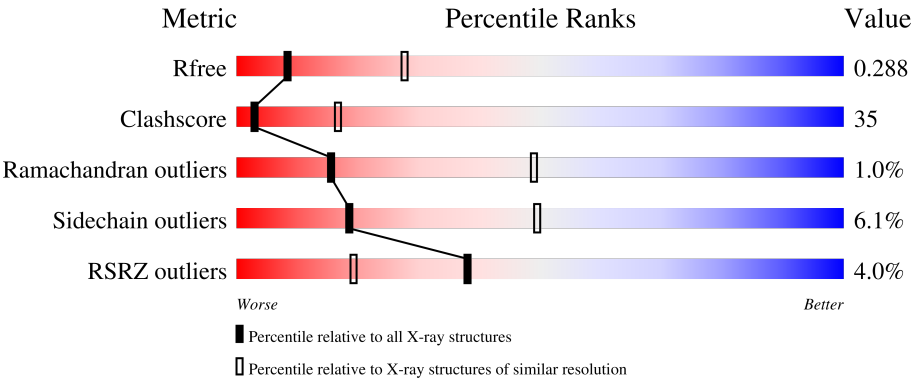
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 3.10 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	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)
RSRZ outliers	127900	1067 (3.10-3.10)

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	A	321	% 55% 33% · 8%
1	B	321	% 54% 36% · 8%
1	C	321	9% 42% 44% 5% 9%
2	E	16	31% 6% 62%
2	F	16	38% 6% 6% 50%

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Mol	Chain	Length	Quality of chain
2	G	16	6% 12% 6% 81%

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	TRS	A	401	-	X	-	-
3	TRS	A	402	-	X	-	-
3	TRS	C	401	-	X	-	-

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 7020 atoms, of which 36 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 Tyrosine-protein phosphatase non-receptor type 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	296	Total	C	N	O	S	0	0	0
			2280	1446	396	423	15			
1	B	296	Total	C	N	O	S	0	0	0
			2313	1475	400	424	14			
1	C	293	Total	C	N	O	S	0	0	0
			2185	1391	373	410	11			

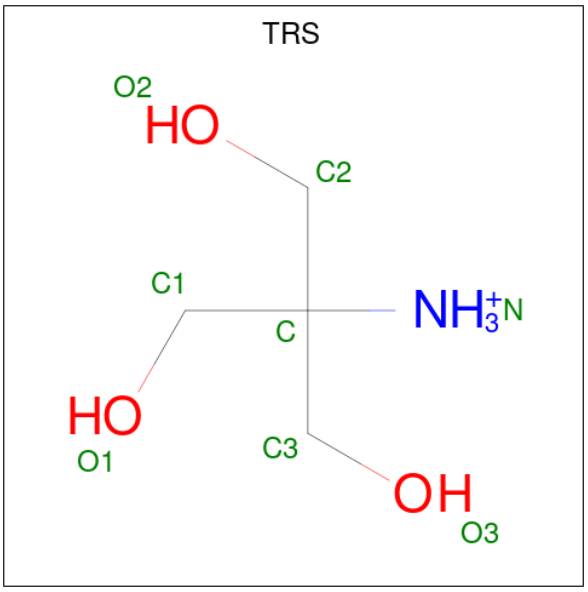
There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	181	ALA	ASP	engineered mutation	UNP P18031
A	215	ALA	CYS	engineered mutation	UNP P18031
A	262	ALA	GLN	engineered mutation	UNP P18031
B	181	ALA	ASP	engineered mutation	UNP P18031
B	215	ALA	CYS	engineered mutation	UNP P18031
B	262	ALA	GLN	engineered mutation	UNP P18031
C	181	ALA	ASP	engineered mutation	UNP P18031
C	215	ALA	CYS	engineered mutation	UNP P18031
C	262	ALA	GLN	engineered mutation	UNP P18031

- Molecule 2 is a protein called Tyrosine-protein kinase JAK2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	6	Total	C	N	O	P	0	0	0
			49	31	6	11	1			
2	F	8	Total	C	N	O	P	0	0	0
			63	39	8	15	1			
2	G	3	Total	C	N	O	P	0	0	0
			33	21	3	8	1			

- Molecule 3 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	0	0
			20	4	12	1	3		
3	A	1	Total	C	H	N	O	0	0
			20	4	12	1	3		
3	C	1	Total	C	H	N	O	0	0
			20	4	12	1	3		

- Molecule 4 is water.

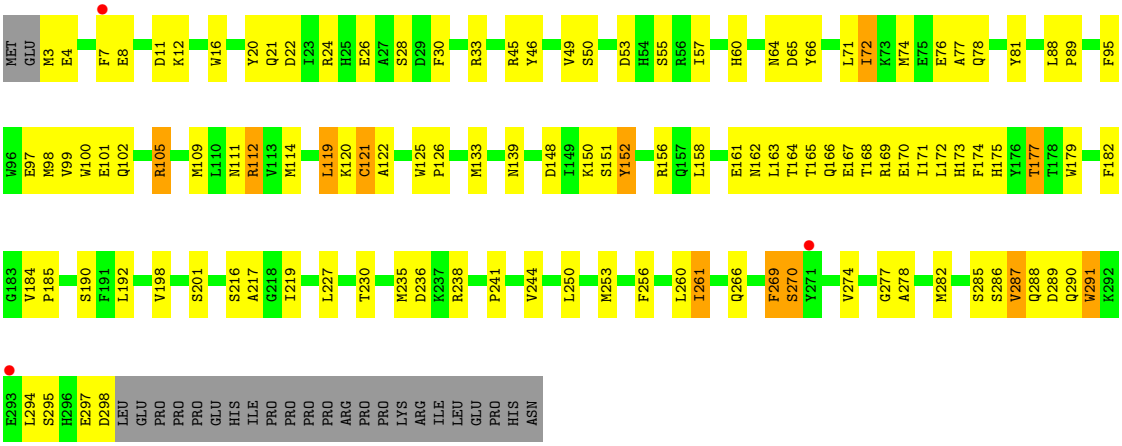
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	11	Total	O	0	0
			11	11		
4	B	16	Total	O	0	0
			16	16		
4	C	10	Total	O	0	0
			10	10		

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.

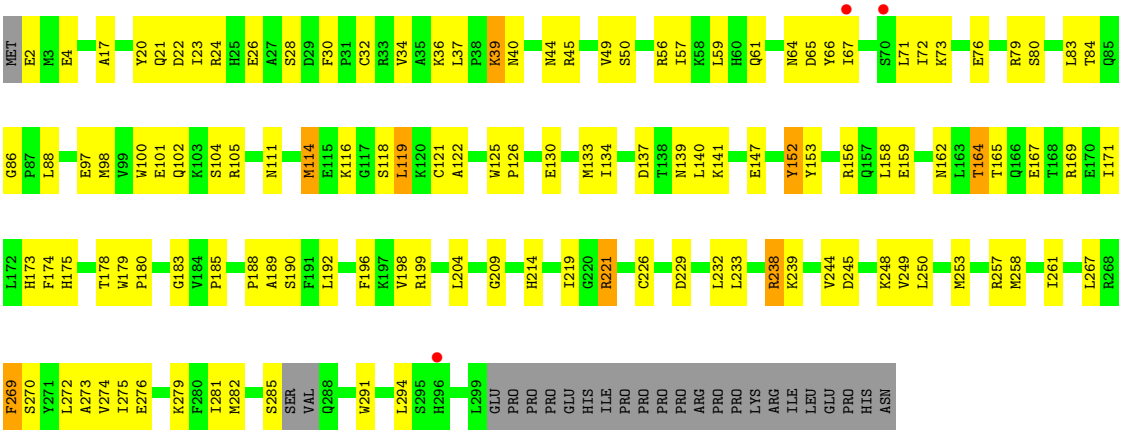
• Molecule 1: Tyrosine-protein phosphatase non-receptor type 1

Chain A:



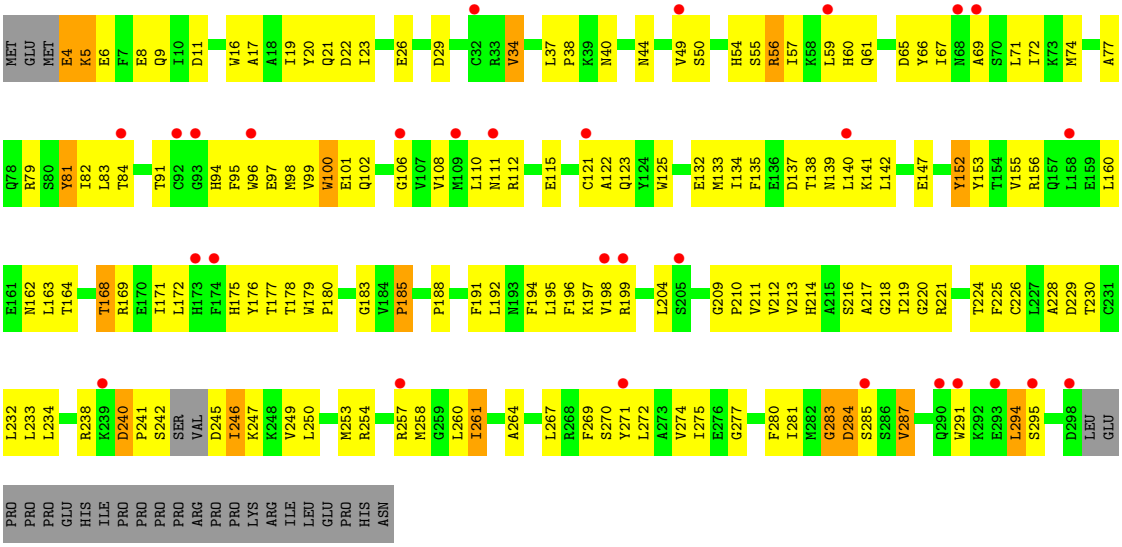
• Molecule 1: Tyrosine-protein phosphatase non-receptor type 1

Chain B:



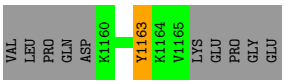
• Molecule 1: Tyrosine-protein phosphatase non-receptor type 1

Chain C:



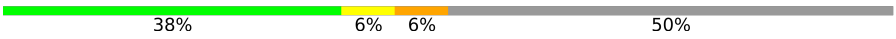
• Molecule 2: Tyrosine-protein kinase JAK2

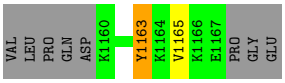
Chain E:



• Molecule 2: Tyrosine-protein kinase JAK2

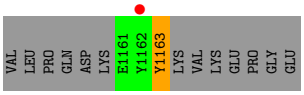
Chain F:





• Molecule 2: Tyrosine-protein kinase JAK2

Chain G:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	88.23Å 176.23Å 43.80Å 90.00° 90.29° 90.00°	Depositor
Resolution (Å)	42.79 – 3.10 48.90 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.1 (42.79-3.10) 93.1 (48.90-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.60 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, R_{free}	0.262 , 0.291 0.257 , 0.288	Depositor DCC
R_{free} test set	2013 reflections (8.33%)	wwPDB-VP
Wilson B-factor (Å^2)	78.5	Xtriage
Anisotropy	0.461	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 58.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.439 for -h,-k,l	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	7020	wwPDB-VP
Average B, all atoms (Å^2)	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.83% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PTR, TRS

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	0.43	1/2333 (0.0%)	0.64	0/3168
1	B	0.41	0/2367	0.63	0/3205
1	C	0.43	0/2236	0.67	0/3052
2	E	0.38	0/32	0.39	0/41
2	F	0.44	0/46	0.36	0/60
2	G	0.52	0/17	0.59	0/22
All	All	0.42	1/7031 (0.0%)	0.64	0/9548

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	C	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	121	CYS	CB-SG	-5.54	1.72	1.81

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	8	GLU	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	2280	0	2134	116	0
1	B	2313	0	2210	124	0
1	C	2185	0	1979	221	0
2	E	49	0	25	4	0
2	F	63	0	33	2	0
2	G	33	0	17	1	0
3	A	16	24	24	3	0
3	C	8	12	12	2	0
4	A	11	0	0	2	0
4	B	16	0	0	2	0
4	C	10	0	0	0	0
All	All	6984	36	6434	463	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 35.

The worst 5 of 463 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:C:134:ILE:HD11	1:C:141:LYS:HG3	1.25	1.14
1:C:242:SER:HG	1:C:245:ASP:N	1.45	1.14
1:C:160:LEU:HD22	1:C:171:ILE:HG13	1.26	1.10
1:C:57:ILE:HG13	1:C:98:MET:HE3	1.32	1.09
1:C:112:ARG:NE	1:C:177:THR:O	1.93	1.01

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	294/321 (92%)	274 (93%)	17 (6%)	3 (1%)	15 49
1	B	292/321 (91%)	274 (94%)	17 (6%)	1 (0%)	41 73
1	C	289/321 (90%)	273 (94%)	12 (4%)	4 (1%)	11 40
2	E	3/16 (19%)	3 (100%)	0	0	100 100
2	F	5/16 (31%)	3 (60%)	1 (20%)	1 (20%)	0 0
2	G	1/16 (6%)	1 (100%)	0	0	100 100
All	All	884/1011 (87%)	828 (94%)	47 (5%)	9 (1%)	15 49

5 of 9 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	5	LYS
1	A	285	SER
1	A	287	VAL
1	B	261	ILE
1	C	283	GLY

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	232/291 (80%)	216 (93%)	16 (7%)	15 45
1	B	240/291 (82%)	229 (95%)	11 (5%)	27 59
1	C	213/291 (73%)	198 (93%)	15 (7%)	15 45
2	E	1/14 (7%)	1 (100%)	0	100 100
2	F	2/14 (14%)	2 (100%)	0	100 100
2	G	1/14 (7%)	1 (100%)	0	100 100
All	All	689/915 (75%)	647 (94%)	42 (6%)	18 49

5 of 42 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	34	VAL
1	C	185	PRO
1	C	56	ARG
1	C	100	TRP
1	C	240	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	61	GLN
1	B	139	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)

3 non-standard protein/DNA/RNA residues 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	PTR	E	1163	2	15,16,17	1.19	1 (6%)	19,22,24	0.54	0
2	PTR	G	1163	2	15,16,17	1.40	2 (13%)	19,22,24	0.81	0
2	PTR	F	1163	2	15,16,17	1.52	2 (13%)	19,22,24	0.83	1 (5%)

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	PTR	E	1163	2	-	0/10/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PTR	G	1163	2	-	1/10/11/13	0/1/1/1
2	PTR	F	1163	2	-	0/10/11/13	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	1163	PTR	OH-CZ	-4.25	1.31	1.40
2	E	1163	PTR	OH-CZ	-4.13	1.31	1.40
2	G	1163	PTR	OH-CZ	-3.79	1.32	1.40
2	G	1163	PTR	P-OH	3.28	1.64	1.59
2	F	1163	PTR	P-OH	2.23	1.62	1.59

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	F	1163	PTR	O3P-P-O2P	2.24	116.21	107.64

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	G	1163	PTR	CZ-OH-P-O3P

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	1163	PTR	4	0
2	G	1163	PTR	1	0
2	F	1163	PTR	2	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

3 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	TRS	A	402	-	7,7,7	2.16	2 (28%)	9,9,9	2.89	5 (55%)
3	TRS	C	401	-	7,7,7	1.64	1 (14%)	9,9,9	2.22	4 (44%)
3	TRS	A	401	-	7,7,7	2.00	1 (14%)	9,9,9	2.83	4 (44%)

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	TRS	A	402	-	-	7/9/9/9	-
3	TRS	C	401	-	-	9/9/9/9	-
3	TRS	A	401	-	-	7/9/9/9	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	TRS	C-N	4.70	1.65	1.49
3	A	401	TRS	C-N	4.50	1.64	1.49
3	C	401	TRS	C-N	3.63	1.61	1.49
3	A	402	TRS	C2-C	2.05	1.59	1.53

The worst 5 of 13 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	402	TRS	C2-C-N	5.73	125.08	107.98
3	A	401	TRS	C3-C-N	5.08	123.15	107.98
3	A	401	TRS	C3-C-C1	-4.25	97.64	110.81
3	C	401	TRS	C3-C-N	4.21	120.55	107.98
3	A	402	TRS	C2-C-C1	-4.12	98.05	110.81

There are no chirality outliers.

5 of 23 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	401	TRS	C3-C-C2-O2
3	A	401	TRS	N-C-C2-O2
3	A	402	TRS	C1-C-C2-O2
3	A	402	TRS	C1-C-C3-O3
3	A	402	TRS	C2-C-C3-O3

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	402	TRS	2	0
3	C	401	TRS	2	0
3	A	401	TRS	1	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, 95^{th} 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	296/321 (92%)	0.07	3 (1%) 82 67	44, 66, 114, 128	0
1	B	296/321 (92%)	0.02	3 (1%) 82 67	45, 65, 113, 124	0
1	C	293/321 (91%)	0.68	29 (9%) 7 2	80, 111, 157, 172	0
2	E	5/16 (31%)	-0.35	0 100 100	76, 79, 85, 99	0
2	F	7/16 (43%)	-0.56	0 100 100	80, 87, 99, 100	0
2	G	2/16 (12%)	1.67	1 (50%) 0 0	103, 103, 103, 109	0
All	All	899/1011 (88%)	0.25	36 (4%) 38 19	44, 80, 135, 172	0

The worst 5 of 36 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	291	TRP	5.3
1	C	295	SER	4.7
1	C	121	CYS	4.6
1	C	298	ASP	4.2
1	B	296	HIS	4.1

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

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, 95^{th} 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	PTR	G	1163	16/17	0.87	0.27	110,114,121,124	0
2	PTR	F	1163	16/17	0.95	0.17	54,59,73,73	0
2	PTR	E	1163	16/17	0.96	0.25	55,58,73,73	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, 95^{th} 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
3	TRS	C	401	8/8	0.76	0.23	92,116,129,129	0
3	TRS	A	401	8/8	0.77	0.32	71,101,121,121	0
3	TRS	A	402	8/8	0.82	0.26	66,88,106,106	0

6.5 Other polymers (i)

There are no such residues in this entry.