



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 3, 2017 – 04:56 AM EDT

PDB ID : 5IBZ
Title : Crystal structure of a novel cyclase (pfam04199).
Authors : Nocek, B.; Skarina, T.; Brown, G.; Joachimiak, A.; Savchenko, A.; Yakunin, A.
Deposited on : unknown
Resolution : 1.61 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

<http://wwpdb.org/validation/2016/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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029824
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20029824

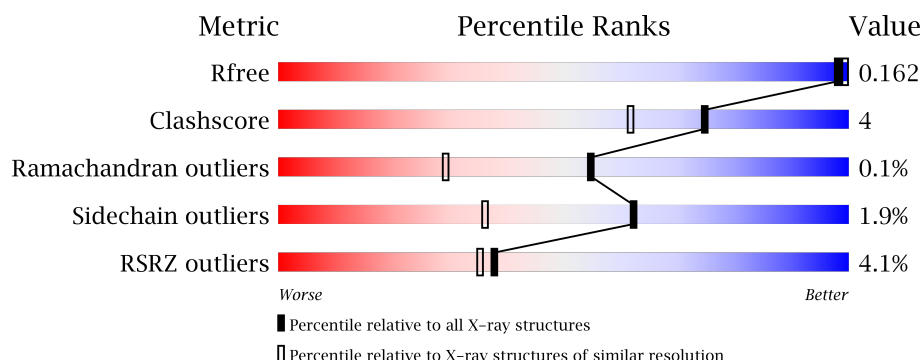
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.61 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}	100719	3539 (1.64-1.60)
Clashscore	112137	3855 (1.64-1.60)
Ramachandran outliers	110173	3764 (1.64-1.60)
Sidechain outliers	110143	3763 (1.64-1.60)
RSRZ outliers	101464	3562 (1.64-1.60)

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. 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	323	<div> <div>4%</div> <div>89%</div> <div>5% • 5%</div> </div>
1	B	323	<div> <div>6%</div> <div>90%</div> <div>5% • 5%</div> </div>
1	C	323	<div> <div>3%</div> <div>89%</div> <div>• 7%</div> </div>
1	D	323	<div> <div>3%</div> <div>85%</div> <div>8% 7%</div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	PGE	C	402	-	-	-	X

2 Entry composition [i](#)

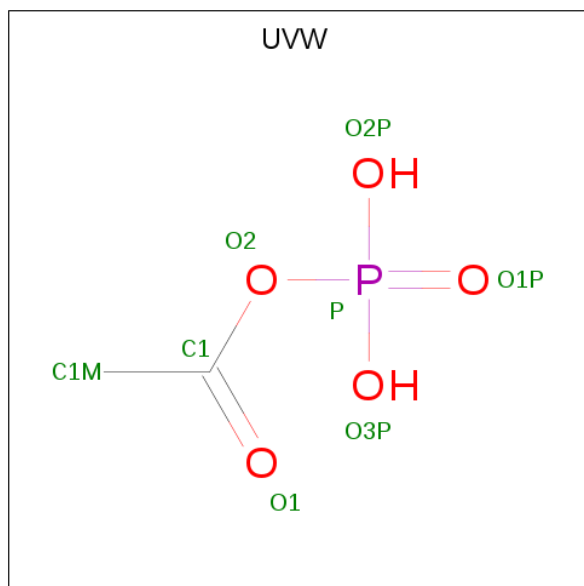
There are 4 unique types of molecules in this entry. The entry contains 20359 atoms, of which 9282 are hydrogens and 0 are deuteriums.

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 Uncharacterized protein.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
1	A	306	Total	C	H	N	O	S	Se	0	3	0
			4678	1514	2320	405	435	1	3			
1	B	308	Total	C	H	N	O	S	Se	0	11	0
			4754	1540	2353	412	445	1	3			
1	C	301	Total	C	H	N	O	S	Se	0	3	0
			4583	1483	2270	398	428	1	3			
1	D	300	Total	C	H	N	O	S	Se	0	13	0
			4660	1508	2313	398	437	1	3			

- Molecule 2 is ACETYLPHOSPHATE (three-letter code: UVW) (formula: C₂H₅O₅P).



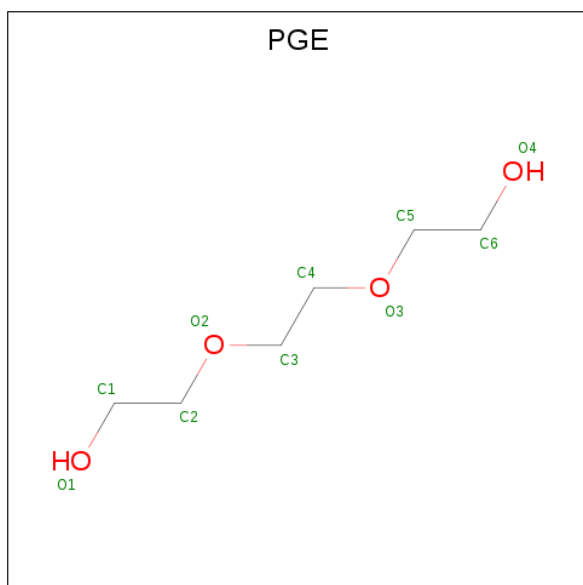
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	H	O	P	0	0
			11	2	3	5	1		
2	B	1	Total	C	H	O	P	0	0
			11	2	3	5	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	H	O	P	0	0
			11	2	3	5	1		
2	D	1	Total	C	H	O	P	0	0
			11	2	3	5	1		

- Molecule 3 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total	C	H	O	0	0
			24	6	14	4		

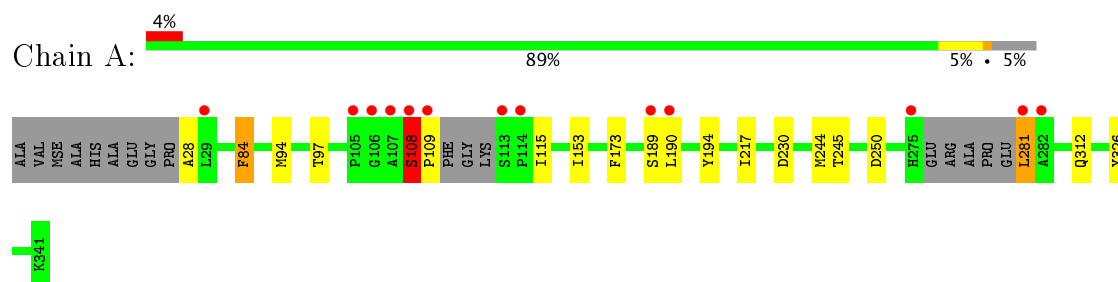
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	403	Total	O	0	0
			403	403		
4	B	403	Total	O	0	0
			403	403		
4	C	421	Total	O	0	0
			421	421		
4	D	389	Total	O	0	0
			389	389		

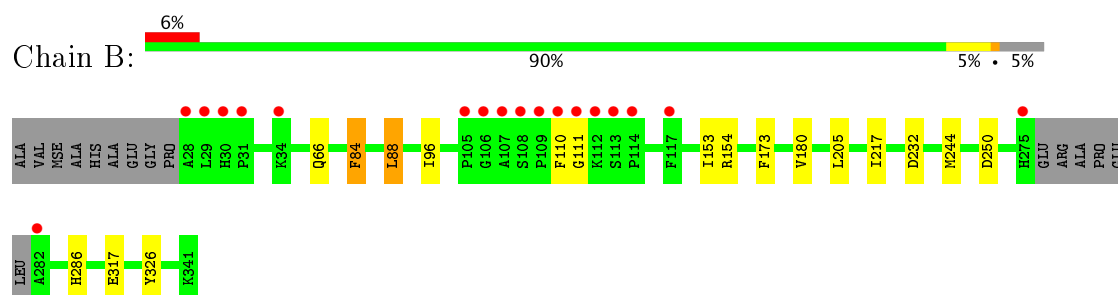
3 Residue-property plots [i](#)

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.

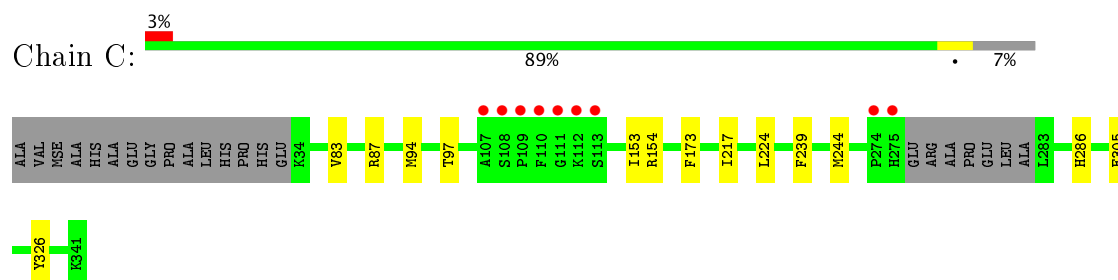
- Molecule 1: Uncharacterized protein



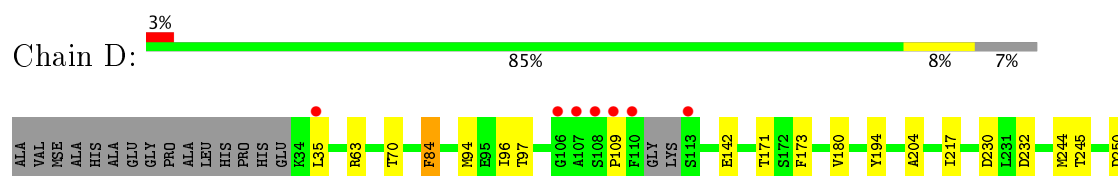
- Molecule 1: Uncharacterized protein

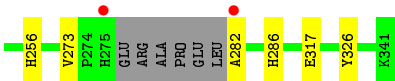


- Molecule 1: Uncharacterized protein



- Molecule 1: Uncharacterized protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	64.51Å 120.62Å 148.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.81 – 1.61 42.24 – 1.61	Depositor EDS
% Data completeness (in resolution range)	99.8 (38.81-1.61) 99.8 (42.24-1.61)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.40 (at 1.61Å)	Xtriage
Refinement program	PHENIX dev_1888	Depositor
R, R_{free}	0.133 , 0.161 0.133 , 0.162	Depositor DCC
R_{free} test set	7523 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	13.3	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 53.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	20359	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 70.89 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 2.8883e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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, UVW

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	0.45	0/2429	0.65	0/3310
1	B	0.45	0/2496	0.65	0/3402
1	C	0.45	0/2382	0.64	0/3244
1	D	0.44	0/2445	0.64	0/3333
All	All	0.45	0/9752	0.64	0/13289

There are no bond length outliers.

There are no bond angle outliers.

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	2358	2320	2318	20	0
1	B	2401	2353	2334	16	0
1	C	2313	2270	2273	15	0
1	D	2347	2313	2303	25	0
2	A	8	3	3	1	0
2	B	8	3	3	1	0
2	C	8	3	3	1	0
2	D	8	3	3	1	0
3	C	10	14	14	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	403	0	0	11	1
4	B	403	0	0	3	6
4	C	421	0	0	1	4
4	D	389	0	0	8	3
All	All	11077	9282	9254	70	7

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 (70) 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:94:MSE:HE2	1:A:97[B]:THR:HG23	1.38	1.03
1:D:94:MSE:HE2	1:D:97[B]:THR:HG23	1.41	1.00
1:A:97[B]:THR:HG22	1:C:153:ILE:HG22	1.54	0.89
1:C:224:LEU:HD21	1:C:244:MSE:HE2	1.57	0.85
1:A:230:ASP:OD2	4:A:501:HOH:O	1.96	0.84
1:D:63:ARG:NH1	4:D:505:HOH:O	2.10	0.83
1:D:142[A]:GLU:OE1	4:D:501:HOH:O	1.96	0.83
1:D:230:ASP:OD2	4:D:502:HOH:O	1.98	0.82
1:A:28:ALA:N	4:A:502:HOH:O	2.14	0.81
1:B:153:ILE:HG22	1:D:97[B]:THR:HG22	1.63	0.80
1:D:171:THR:O	4:D:503:HOH:O	1.99	0.79
1:B:317[A]:GLU:OE1	4:B:501:HOH:O	1.99	0.78
1:D:317[A]:GLU:OE1	4:D:504:HOH:O	2.01	0.78
1:D:282:ALA:N	4:D:507:HOH:O	2.16	0.77
4:A:643:HOH:O	1:C:97:THR:HB	1.86	0.75
1:D:245[B]:THR:HG23	4:D:688:HOH:O	1.87	0.75
1:D:84:PHE:CE2	1:D:244:MSE:HE1	2.22	0.74
1:A:194:TYR:H	1:A:245[B]:THR:HG22	1.51	0.74
1:A:281:LEU:N	4:A:503:HOH:O	2.20	0.73
1:D:244:MSE:HE3	1:D:273[B]:VAL:HG21	1.74	0.69
1:B:84:PHE:CE2	1:B:244:MSE:HE1	2.28	0.69
1:A:312:GLN:NE2	4:A:505:HOH:O	2.25	0.66
1:D:35:LEU:HD11	1:D:256:HIS:CD2	2.31	0.65
1:C:305:GLU:OE1	4:C:501:HOH:O	2.13	0.65
1:A:194:TYR:N	1:A:245[B]:THR:HG22	2.12	0.64
1:D:244:MSE:HE3	1:D:273[B]:VAL:CG2	2.29	0.63
1:A:84:PHE:CE2	1:A:244:MSE:HE1	2.34	0.62
1:B:154:ARG:O	1:D:97[B]:THR:HG21	1.99	0.61
1:A:94:MSE:CE	1:A:97[B]:THR:HG23	2.24	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:194:TYR:H	1:D:245[B]:THR:HG22	1.65	0.61
1:A:84:PHE:CZ	1:A:244:MSE:HE1	2.36	0.60
1:A:245[B]:THR:HG23	4:A:692:HOH:O	2.02	0.59
1:C:224:LEU:CD2	1:C:244:MSE:HE2	2.30	0.58
1:A:94:MSE:HE2	1:A:97[B]:THR:CG2	2.23	0.58
1:B:88:LEU:HD12	4:B:753:HOH:O	2.03	0.58
1:B:110:PHE:N	1:B:111:GLY:HA2	2.19	0.57
4:A:643:HOH:O	1:C:94:MSE:HE2	2.07	0.54
1:B:84:PHE:HE2	1:B:244:MSE:HE1	1.72	0.54
1:C:173:PHE:HA	1:C:217:ILE:HD11	1.91	0.53
4:A:643:HOH:O	1:C:97:THR:CB	2.48	0.53
1:A:97[B]:THR:CG2	1:C:153:ILE:HG22	2.34	0.52
4:A:504:HOH:O	1:B:88:LEU:CD1	2.57	0.52
1:C:239:PHE:CE2	1:C:244:MSE:HE3	2.45	0.51
1:B:180[B]:VAL:HG22	1:B:205:LEU:CD2	2.41	0.50
1:A:84:PHE:CD2	2:A:401:UVW:C1M	2.95	0.50
1:A:153:ILE:HG22	4:A:643:HOH:O	2.12	0.49
1:D:194:TYR:N	1:D:245[B]:THR:HG22	2.27	0.49
1:D:96:ILE:HD13	1:D:109:PRO:HG3	1.94	0.49
1:D:96:ILE:HD13	1:D:109:PRO:HB3	1.96	0.48
1:A:173:PHE:HA	1:A:217:ILE:HD11	1.95	0.48
1:D:173:PHE:HA	1:D:217:ILE:HD11	1.94	0.48
1:D:180[A]:VAL:CG2	1:D:204:ALA:HB3	2.44	0.48
1:B:66[A]:GLN:OE1	4:B:502:HOH:O	2.20	0.47
1:B:180[B]:VAL:HG22	1:B:205:LEU:HD21	1.97	0.47
1:A:97[B]:THR:HG21	1:C:154:ARG:O	2.14	0.46
1:B:180[B]:VAL:CG2	1:B:205:LEU:HD21	2.46	0.46
1:D:84:PHE:CZ	1:D:244:MSE:HE1	2.50	0.46
1:C:239:PHE:CE2	1:C:244:MSE:CE	2.99	0.46
1:C:224:LEU:HD21	1:C:244:MSE:CE	2.38	0.46
1:C:83:VAL:HG22	1:C:87:ARG:HB2	1.97	0.45
4:A:504:HOH:O	1:B:88:LEU:HD11	2.13	0.45
1:D:70:THR:OG1	4:D:506:HOH:O	2.12	0.45
1:D:96:ILE:HD13	1:D:109:PRO:CG	2.46	0.44
1:C:286:HIS:NE2	2:C:401:UVW:O1P	2.53	0.42
1:B:154:ARG:C	1:D:97[B]:THR:HG21	2.39	0.42
1:A:108:SER:CB	1:A:109:PRO:CD	2.99	0.41
1:A:190:LEU:N	1:A:190:LEU:HD12	2.36	0.41
1:D:286:HIS:NE2	2:D:401:UVW:O3P	2.53	0.40
1:B:286:HIS:NE2	2:B:401:UVW:O2P	2.54	0.40
1:B:173:PHE:HA	1:B:217:ILE:HD11	2.03	0.40

All (7) 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)
4:B:790:HOH:O	4:D:760:HOH:O[1_455]	1.88	0.32
4:B:827:HOH:O	4:C:835:HOH:O[3_746]	2.07	0.13
4:B:648:HOH:O	4:C:732:HOH:O[3_746]	2.08	0.12
4:A:614:HOH:O	4:D:679:HOH:O[3_746]	2.14	0.06
4:B:765:HOH:O	4:C:772:HOH:O[3_746]	2.18	0.02
4:B:758:HOH:O	4:D:558:HOH:O[1_455]	2.19	0.01
4:B:765:HOH:O	4:C:787:HOH:O[3_746]	2.19	0.01

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	303/323 (94%)	292 (96%)	10 (3%)	1 (0%)	44	22
1	B	315/323 (98%)	304 (96%)	11 (4%)	0	100	100
1	C	300/323 (93%)	293 (98%)	7 (2%)	0	100	100
1	D	307/323 (95%)	298 (97%)	9 (3%)	0	100	100
All	All	1225/1292 (95%)	1187 (97%)	37 (3%)	1 (0%)	55	29

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	108	SER

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	243/249 (98%)	236 (97%)	7 (3%)	48	19
1	B	247/249 (99%)	241 (98%)	6 (2%)	54	25
1	C	238/249 (96%)	237 (100%)	1 (0%)	93	87
1	D	244/249 (98%)	240 (98%)	4 (2%)	68	44
All	All	972/996 (98%)	954 (98%)	18 (2%)	62	36

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	PHE
1	A	108	SER
1	A	115	ILE
1	A	189	SER
1	A	250	ASP
1	A	281	LEU
1	A	326	TYR
1	B	84	PHE
1	B	88	LEU
1	B	96	ILE
1	B	232	ASP
1	B	250	ASP
1	B	326	TYR
1	C	326	TYR
1	D	84	PHE
1	D	232	ASP
1	D	250	ASP
1	D	326	TYR

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

Mol	Chain	Res	Type
1	A	131	GLN
1	B	131	GLN

5.3.3 RNA ⓘ

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](#)

5 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	$\# Z > 2$
2	UVW	A	401	-	6,7,7	1.56	1 (16%)	7,10,10	1.41	2 (28%)
2	UVW	B	401	-	6,7,7	1.38	1 (16%)	7,10,10	0.91	0
2	UVW	C	401	-	6,7,7	1.41	1 (16%)	7,10,10	1.13	0
3	PGE	C	402	-	9,9,9	0.32	0	8,8,8	0.67	0
2	UVW	D	401	-	6,7,7	1.32	1 (16%)	7,10,10	0.82	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	UVW	A	401	-	-	0/3/5/5	0/0/0/0
2	UVW	B	401	-	-	0/3/5/5	0/0/0/0
2	UVW	C	401	-	-	0/3/5/5	0/0/0/0
3	PGE	C	402	-	-	0/7/7/7	0/0/0/0
2	UVW	D	401	-	-	0/3/5/5	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	401	UVW	P-O2	2.89	1.64	1.59
2	B	401	UVW	P-O2	2.94	1.64	1.59
2	C	401	UVW	P-O2	3.00	1.65	1.59
2	A	401	UVW	P-O2	3.50	1.65	1.59

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	UVW	O2-P-O1P	-2.08	99.75	108.44
2	A	401	UVW	O2P-P-O2	2.17	112.84	105.57

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	UVW	1	0
2	B	401	UVW	1	0
2	C	401	UVW	1	0
2	D	401	UVW	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, 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	303/323 (93%)	-0.40	13 (4%) 36 33	8, 12, 28, 61	0
1	B	305/323 (94%)	-0.32	18 (5%) 23 21	8, 12, 38, 55	0
1	C	298/323 (92%)	-0.42	9 (3%) 51 49	8, 12, 30, 50	0
1	D	297/323 (91%)	-0.47	9 (3%) 51 49	7, 12, 28, 50	0
All	All	1203/1292 (93%)	-0.40	49 (4%) 38 35	7, 12, 31, 61	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	110	PHE	8.6
1	B	109	PRO	8.5
1	C	110	PHE	7.1
1	A	113	SER	7.0
1	A	107	ALA	6.3
1	C	275	HIS	6.1
1	A	109	PRO	6.1
1	B	113	SER	5.9
1	A	106	GLY	5.9
1	B	29	LEU	5.9
1	A	281	LEU	5.5
1	A	108	SER	5.5
1	D	107	ALA	5.4
1	D	282	ALA	5.3
1	B	108	SER	5.2
1	B	106	GLY	5.0
1	B	282	ALA	5.0
1	C	111	GLY	5.0
1	C	109	PRO	4.9
1	B	28	ALA	4.8
1	A	189	SER	4.7

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Mol	Chain	Res	Type	RSRZ
1	C	107	ALA	4.3
1	A	282	ALA	4.1
1	A	114	PRO	4.0
1	B	275	HIS	3.9
1	B	111	GLY	3.8
1	B	107	ALA	3.8
1	B	34	LYS	3.8
1	D	113	SER	3.7
1	B	30	HIS	3.7
1	C	113	SER	3.6
1	A	29	LEU	3.6
1	D	110	PHE	3.6
1	D	109	PRO	3.3
1	B	31	PRO	3.2
1	B	114	PRO	3.2
1	D	106	GLY	3.2
1	C	112	LYS	3.1
1	A	105	PRO	2.8
1	B	112	LYS	2.8
1	D	35	LEU	2.7
1	C	274	PRO	2.7
1	D	108	SER	2.7
1	A	190	LEU	2.5
1	B	117	PHE	2.4
1	C	108	SER	2.4
1	A	275	HIS	2.3
1	B	105	PRO	2.3
1	D	275	HIS	2.1

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. LLDF column lists the quality of electron

density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	PGE	C	402	10/10	0.87	0.33	13.59	66,88,113,118	0
2	UVW	C	401	8/8	0.97	0.09	0.11	23,32,51,55	0
2	UVW	D	401	8/8	0.96	0.11	0.09	20,29,57,57	0
2	UVW	A	401	8/8	0.97	0.09	0.02	17,26,58,58	0
2	UVW	B	401	8/8	0.98	0.09	-0.00	20,27,51,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.