



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 07:26 PM GMT

PDB ID : 4RV6
Title : Human ARTD1 (PARP1) catalytic domain in complex with inhibitor Rucaparib
Authors : Karlberg, T.; Thorsell, A.G.; Schuler, H.
Deposited on : 2014-11-25
Resolution : 3.19 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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-20026982

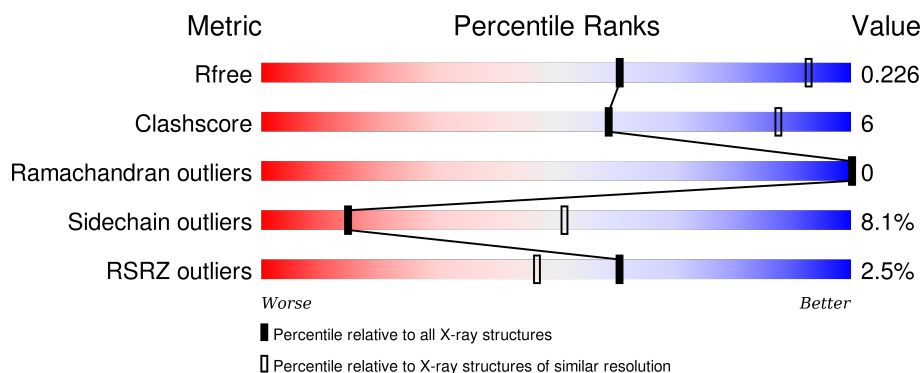
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.19 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}	91344	1124 (3.24-3.16)
Clashscore	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)
RSRZ outliers	91569	1129 (3.24-3.16)

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	352	<div> <div>%</div> <div> <div></div> <div>76%</div> <div>21%</div> <div>..</div> </div> </div>
1	B	352	<div> <div>%</div> <div> <div></div> <div>78%</div> <div>20%</div> <div>..</div> </div> </div>
1	C	352	<div> <div>4%</div> <div> <div></div> <div>79%</div> <div>19%</div> <div>..</div> </div> </div>
1	D	352	<div> <div>4%</div> <div> <div></div> <div>80%</div> <div>18%</div> <div>..</div> </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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	RPB	A	1103	-	-	-	X
3	RPB	B	1103	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11027 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 Poly [ADP-ribose] polymerase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	347	Total	C	N	O	S	0	0	0
			2730	1739	462	518	11			
1	B	349	Total	C	N	O	S	0	0	0
			2744	1746	464	523	11			
1	C	347	Total	C	N	O	S	0	0	0
			2728	1738	462	517	11			
1	D	349	Total	C	N	O	S	0	0	0
			2744	1746	464	523	11			

There are 12 discrepancies between the modelled and reference sequences:

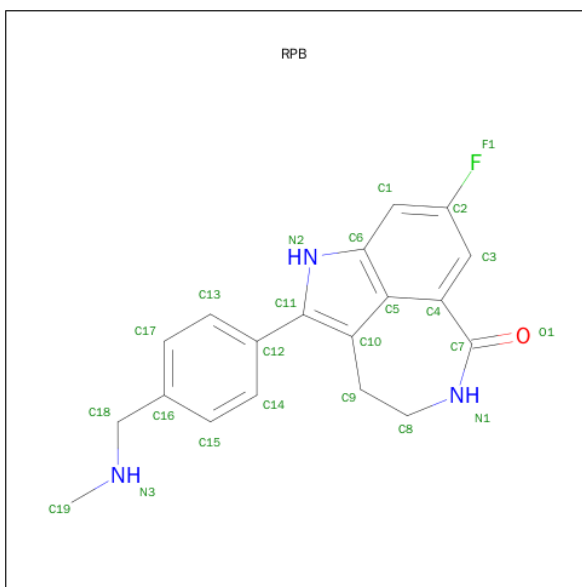
Chain	Residue	Modelled	Actual	Comment	Reference
A	660	SER	-	EXPRESSION TAG	UNP P09874
A	661	MET	-	EXPRESSION TAG	UNP P09874
A	762	ALA	VAL	variant	UNP P09874
B	660	SER	-	EXPRESSION TAG	UNP P09874
B	661	MET	-	EXPRESSION TAG	UNP P09874
B	762	ALA	VAL	variant	UNP P09874
C	660	SER	-	EXPRESSION TAG	UNP P09874
C	661	MET	-	EXPRESSION TAG	UNP P09874
C	762	ALA	VAL	variant	UNP P09874
D	660	SER	-	EXPRESSION TAG	UNP P09874
D	661	MET	-	EXPRESSION TAG	UNP P09874
D	762	ALA	VAL	variant	UNP P09874

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is RUCAPARIB (three-letter code: RPB) (formula: C₁₉H₁₈FN₃O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			24	19	1	3	1		
3	B	1	Total	C	F	N	O	0	0
			24	19	1	3	1		

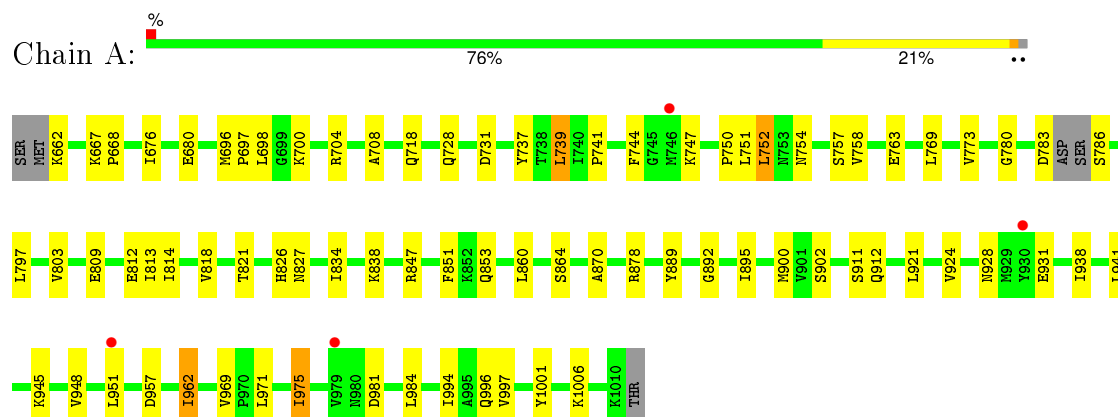
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	O	0	0
			1	1		
4	C	1	Total	O	0	0
			1	1		
4	D	1	Total	O	0	0
			1	1		

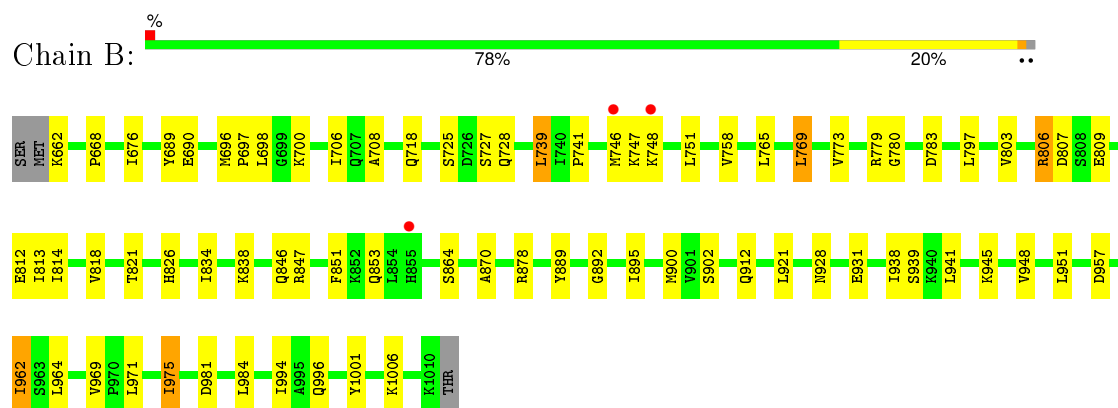
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

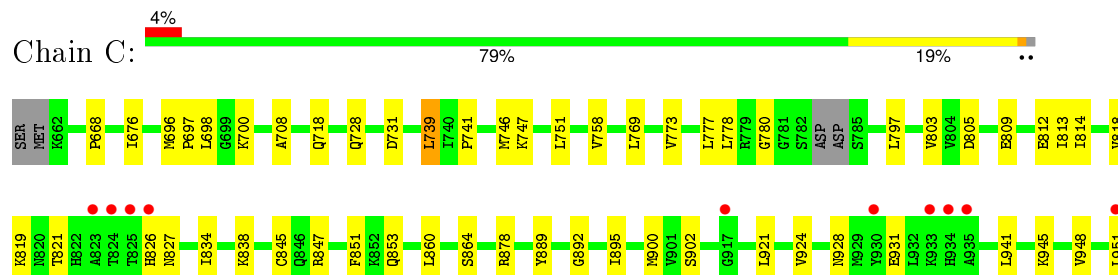
- Molecule 1: Poly [ADP-ribose] polymerase 1

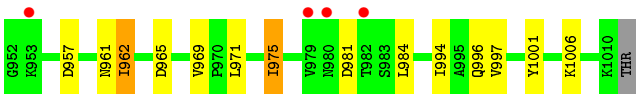


- Molecule 1: Poly [ADP-ribose] polymerase 1

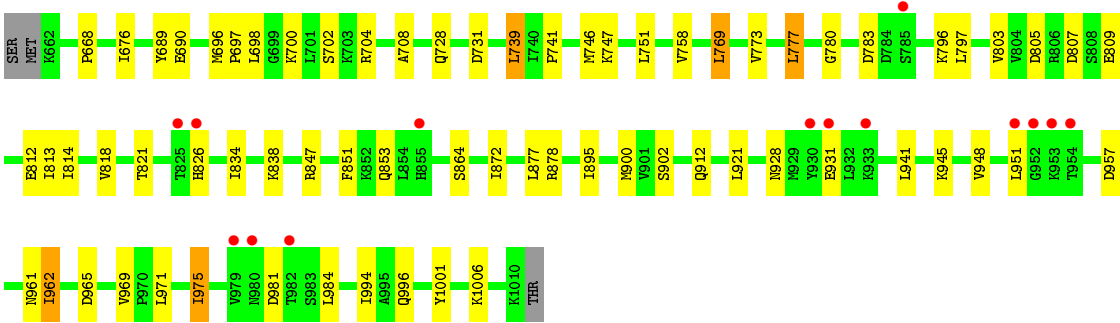
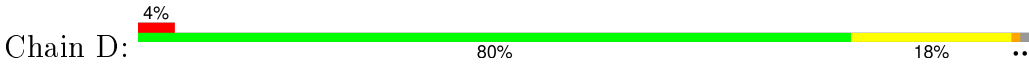


- Molecule 1: Poly [ADP-ribose] polymerase 1





● Molecule 1: Poly [ADP-ribose] polymerase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	93.63Å 85.10Å 101.94Å 90.00° 100.97° 90.00°	Depositor
Resolution (Å)	37.94 – 3.19 47.95 – 3.19	Depositor EDS
% Data completeness (in resolution range)	99.9 (37.94-3.19) 99.8 (47.95-3.19)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 3.19Å)	Xtriage
Refinement program	BUSTER 2.11.5	Depositor
R, R_{free}	0.181 , 0.212 0.202 , 0.226	Depositor DCC
R_{free} test set	1322 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	65.4	Xtriage
Anisotropy	0.562	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 77.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 26498 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11027	wwPDB-VP
Average B, all atoms (Å ²)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.63% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: RPB, SO4

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	0/2781	0.70	0/3751
1	B	0.51	0/2796	0.70	0/3773
1	C	0.49	0/2779	0.69	0/3748
1	D	0.48	0/2796	0.67	0/3773
All	All	0.50	0/11152	0.69	0/15045

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	2730	0	2774	41	0
1	B	2744	0	2784	32	0
1	C	2728	0	2775	27	0
1	D	2744	0	2784	28	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	1	0
3	A	24	0	18	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	24	0	18	3	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
All	All	11027	0	11153	129	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 (129) 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:B:1103:RPB:H92C	3:B:1103:RPB:H14	1.50	0.94
1:C:860:LEU:HD12	1:C:924:VAL:HG11	1.58	0.86
3:A:1103:RPB:H92C	3:A:1103:RPB:H14	1.59	0.84
1:A:860:LEU:HD12	1:A:924:VAL:HG11	1.60	0.83
1:B:895:ILE:HD11	1:B:994:ILE:HG22	1.62	0.81
1:A:895:ILE:HD11	1:A:994:ILE:HG22	1.63	0.80
1:C:895:ILE:HD11	1:C:994:ILE:HG22	1.66	0.76
1:B:826:HIS:ND1	1:B:902:SER:HB3	2.02	0.74
1:A:826:HIS:ND1	1:A:902:SER:HB3	2.03	0.73
1:D:826:HIS:ND1	1:D:902:SER:HB3	2.03	0.73
1:C:826:HIS:ND1	1:C:902:SER:HB3	2.03	0.73
1:D:895:ILE:HD11	1:D:994:ILE:HG22	1.72	0.72
3:B:1103:RPB:C9	3:B:1103:RPB:H14	2.22	0.69
3:B:1103:RPB:H92C	3:B:1103:RPB:C14	2.17	0.65
1:A:962:ILE:HD13	1:A:971:LEU:HD11	1.78	0.65
1:B:962:ILE:HD13	1:B:971:LEU:HD11	1.78	0.63
1:C:962:ILE:HD13	1:C:971:LEU:HD11	1.80	0.62
1:D:962:ILE:HD13	1:D:971:LEU:HD11	1.80	0.62
1:D:872:ILE:HG23	1:D:877:LEU:HD12	1.83	0.60
1:C:931:GLU:HB3	1:C:951:LEU:HD11	1.88	0.56
3:A:1103:RPB:H14	3:A:1103:RPB:C9	2.32	0.56
1:D:931:GLU:HB3	1:D:951:LEU:HD11	1.87	0.56
1:B:806:ARG:CZ	1:B:807:ASP:OD2	2.55	0.55
3:A:1103:RPB:H92C	3:A:1103:RPB:C14	2.24	0.55
1:A:931:GLU:HB3	1:A:951:LEU:HD11	1.89	0.54
1:C:860:LEU:HD12	1:C:924:VAL:CG1	2.32	0.54
1:A:780:GLY:HA2	1:D:780:GLY:HA2	1.90	0.54
1:A:851:PHE:CD2	1:A:996:GLN:HG2	2.41	0.54
1:D:872:ILE:HG23	1:D:877:LEU:CD1	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:676:ILE:HD11	1:B:797:LEU:HD11	1.90	0.53
1:A:860:LEU:HD12	1:A:924:VAL:CG1	2.34	0.52
1:C:696:MET:HB2	1:C:741:PRO:HG2	1.91	0.52
1:C:809:GLU:O	1:C:813:ILE:HG12	2.10	0.52
1:D:809:GLU:O	1:D:813:ILE:HG12	2.10	0.51
1:C:676:ILE:HD11	1:C:797:LEU:HD11	1.92	0.51
1:C:838:LYS:HG2	1:C:1001:TYR:CE1	2.45	0.51
1:B:931:GLU:HB3	1:B:951:LEU:HD11	1.93	0.51
1:B:809:GLU:O	1:B:813:ILE:HG12	2.10	0.51
1:A:769:LEU:O	1:A:773:VAL:HG23	2.10	0.51
1:A:827:ASN:OD1	1:B:725:SER:HB3	2.11	0.51
1:D:696:MET:HB2	1:D:741:PRO:HG2	1.92	0.51
1:A:676:ILE:HD11	1:A:797:LEU:HD11	1.92	0.51
1:D:769:LEU:O	1:D:773:VAL:HG23	2.11	0.51
1:A:737:TYR:CE2	1:A:752:LEU:HD11	2.46	0.50
1:D:838:LYS:HG2	1:D:1001:TYR:CE1	2.46	0.50
1:A:838:LYS:HG2	1:A:1001:TYR:CE1	2.46	0.50
1:A:696:MET:HB2	1:A:741:PRO:HG2	1.92	0.50
1:B:780:GLY:HA2	1:C:780:GLY:HA2	1.94	0.50
1:D:676:ILE:HD11	1:D:797:LEU:HD11	1.92	0.50
1:B:821:THR:HB	1:B:900:MET:HA	1.94	0.50
1:B:696:MET:HB2	1:B:741:PRO:HG2	1.93	0.50
1:A:809:GLU:O	1:A:813:ILE:HG12	2.12	0.49
1:A:834:ILE:HD11	1:A:1006:LYS:HB2	1.94	0.49
1:C:769:LEU:O	1:C:773:VAL:HG23	2.13	0.49
1:A:737:TYR:CZ	1:A:752:LEU:HD11	2.48	0.49
1:C:928:ASN:HB3	1:C:945:LYS:HA	1.95	0.49
1:B:921:LEU:HB2	1:B:1001:TYR:HB2	1.94	0.49
1:C:921:LEU:HB2	1:C:1001:TYR:HB2	1.94	0.48
1:A:847:ARG:NH1	1:A:994:ILE:HD12	2.28	0.48
1:A:821:THR:HB	1:A:900:MET:HA	1.95	0.48
1:B:697:PRO:HD2	1:B:700:LYS:HB2	1.95	0.48
1:C:957:ASP:HA	1:C:975:ILE:HG12	1.96	0.48
1:A:752:LEU:HD12	1:A:752:LEU:N	2.29	0.48
1:D:957:ASP:HA	1:D:975:ILE:HG12	1.96	0.48
1:A:957:ASP:HA	1:A:975:ILE:HG12	1.94	0.48
1:B:957:ASP:HA	1:B:975:ILE:HG12	1.95	0.48
1:B:834:ILE:HD11	1:B:1006:LYS:HB2	1.95	0.47
1:C:851:PHE:CD1	1:C:996:GLN:HG2	2.49	0.47
1:D:834:ILE:HD11	1:D:1006:LYS:HB2	1.95	0.47
1:B:847:ARG:NH1	1:B:994:ILE:HD12	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:921:LEU:HB2	1:D:1001:TYR:HB2	1.95	0.47
1:B:851:PHE:CD1	1:B:996:GLN:HG2	2.50	0.47
1:A:928:ASN:HB3	1:A:945:LYS:HA	1.97	0.47
1:C:821:THR:HB	1:C:900:MET:HA	1.97	0.47
1:D:928:ASN:HB3	1:D:945:LYS:HA	1.95	0.47
1:C:834:ILE:HD11	1:C:1006:LYS:HB2	1.95	0.47
1:D:962:ILE:HG22	1:D:969:VAL:HB	1.97	0.47
1:C:941:LEU:HD11	1:C:948:VAL:HG23	1.97	0.47
1:B:769:LEU:O	1:B:773:VAL:HG23	2.15	0.47
1:D:851:PHE:CD1	1:D:996:GLN:HG2	2.49	0.47
1:A:921:LEU:HB2	1:A:1001:TYR:HB2	1.95	0.47
1:B:928:ASN:HB3	1:B:945:LYS:HA	1.97	0.47
1:D:941:LEU:HD11	1:D:948:VAL:HG23	1.96	0.46
1:A:962:ILE:HG22	1:A:969:VAL:HB	1.98	0.46
1:B:962:ILE:HG22	1:B:969:VAL:HB	1.98	0.46
1:A:941:LEU:HD11	1:A:948:VAL:HG23	1.97	0.46
1:A:911:SER:HB2	1:B:748:LYS:NZ	2.30	0.46
1:D:697:PRO:HD2	1:D:700:LYS:HB2	1.98	0.45
1:B:938:ILE:HG21	1:B:941:LEU:HD12	1.98	0.45
1:D:702:SER:HB2	2:D:1101:SO4:O3	2.15	0.45
1:D:821:THR:HB	1:D:900:MET:HA	1.98	0.45
1:B:838:LYS:HG2	1:B:1001:TYR:CE2	2.52	0.45
1:C:962:ILE:HG22	1:C:969:VAL:HB	1.98	0.44
1:B:668:PRO:HG2	1:B:803:VAL:HG21	1.99	0.44
1:C:697:PRO:HD2	1:C:700:LYS:HB2	1.98	0.44
1:A:697:PRO:HD2	1:A:700:LYS:HB2	2.00	0.44
1:C:847:ARG:NH1	1:C:994:ILE:HD12	2.32	0.44
1:B:941:LEU:HD11	1:B:948:VAL:HG23	2.00	0.44
1:A:752:LEU:N	1:A:752:LEU:CD1	2.81	0.44
1:A:668:PRO:HG2	1:A:803:VAL:HG21	2.00	0.43
1:D:668:PRO:HG2	1:D:803:VAL:HG21	2.01	0.43
1:B:676:ILE:HD12	1:B:870:ALA:HA	2.02	0.42
1:D:847:ARG:NH1	1:D:994:ILE:HD12	2.34	0.42
1:A:814:ILE:O	1:A:818:VAL:HG23	2.20	0.42
1:C:668:PRO:HG2	1:C:803:VAL:HG21	2.02	0.42
1:A:763:GLU:HG3	3:A:1103:RPB:H191	2.02	0.42
1:C:814:ILE:O	1:C:818:VAL:HG23	2.19	0.42
1:B:689:TYR:O	1:B:690:GLU:HB2	2.20	0.41
1:D:814:ILE:O	1:D:818:VAL:HG23	2.19	0.41
1:A:924:VAL:HG23	1:A:997:VAL:HG22	2.01	0.41
1:A:827:ASN:ND2	1:B:727:SER:HB3	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:889:TYR:HB3	1:A:892:GLY:O	2.20	0.41
1:A:708:ALA:HB3	1:A:739:LEU:HD21	2.02	0.41
1:B:706:ILE:HG23	1:B:765:LEU:HD22	2.02	0.41
1:C:924:VAL:HG23	1:C:997:VAL:HG22	2.03	0.41
1:A:737:TYR:OH	1:A:750:PRO:HG2	2.21	0.41
1:B:889:TYR:HB3	1:B:892:GLY:O	2.21	0.41
1:A:744:PHE:CZ	1:A:750:PRO:HD2	2.56	0.41
1:D:777:LEU:HD22	1:D:796:LYS:HB3	2.03	0.41
1:A:938:ILE:HG21	1:A:941:LEU:HD12	2.01	0.41
1:D:708:ALA:HB3	1:D:739:LEU:HD21	2.03	0.41
1:B:708:ALA:HB3	1:B:739:LEU:HD21	2.03	0.41
1:D:689:TYR:O	1:D:690:GLU:HB2	2.20	0.41
1:C:889:TYR:HB3	1:C:892:GLY:O	2.21	0.40
1:B:814:ILE:O	1:B:818:VAL:HG23	2.21	0.40
1:A:754:ASN:ND2	1:A:757:SER:H	2.19	0.40
1:A:737:TYR:CE2	1:A:752:LEU:CD1	3.04	0.40
1:A:676:ILE:HD12	1:A:870:ALA:HA	2.02	0.40
1:C:708:ALA:HB3	1:C:739:LEU:HD21	2.03	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	343/352 (97%)	330 (96%)	13 (4%)	0	100	100
1	B	347/352 (99%)	332 (96%)	15 (4%)	0	100	100
1	C	343/352 (97%)	329 (96%)	14 (4%)	0	100	100
1	D	347/352 (99%)	332 (96%)	15 (4%)	0	100	100
All	All	1380/1408 (98%)	1323 (96%)	57 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	304/309 (98%)	280 (92%)	24 (8%)	15	53
1	B	306/309 (99%)	281 (92%)	25 (8%)	14	50
1	C	304/309 (98%)	279 (92%)	25 (8%)	14	50
1	D	306/309 (99%)	281 (92%)	25 (8%)	14	50
All	All	1220/1236 (99%)	1121 (92%)	99 (8%)	15	51

All (99) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	662	LYS
1	A	667	LYS
1	A	680	GLU
1	A	698	LEU
1	A	704	ARG
1	A	718	GLN
1	A	728	GLN
1	A	731	ASP
1	A	739	LEU
1	A	747	LYS
1	A	751	LEU
1	A	752	LEU
1	A	758	VAL
1	A	783	ASP
1	A	786	SER
1	A	812	GLU
1	A	853	GLN
1	A	864	SER
1	A	878	ARG
1	A	912	GLN
1	A	962	ILE
1	A	975	ILE
1	A	981	ASP
1	A	984	LEU

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Mol	Chain	Res	Type
1	B	662	LYS
1	B	698	LEU
1	B	718	GLN
1	B	728	GLN
1	B	739	LEU
1	B	746	MET
1	B	747	LYS
1	B	751	LEU
1	B	758	VAL
1	B	769	LEU
1	B	779	ARG
1	B	783	ASP
1	B	806	ARG
1	B	812	GLU
1	B	846	GLN
1	B	853	GLN
1	B	864	SER
1	B	878	ARG
1	B	912	GLN
1	B	939	SER
1	B	962	ILE
1	B	964	LEU
1	B	975	ILE
1	B	981	ASP
1	B	984	LEU
1	C	698	LEU
1	C	718	GLN
1	C	728	GLN
1	C	731	ASP
1	C	739	LEU
1	C	746	MET
1	C	747	LYS
1	C	751	LEU
1	C	758	VAL
1	C	777	LEU
1	C	778	LEU
1	C	805	ASP
1	C	812	GLU
1	C	819	LYS
1	C	827	ASN
1	C	845	CYS
1	C	853	GLN

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Mol	Chain	Res	Type
1	C	864	SER
1	C	878	ARG
1	C	961	ASN
1	C	962	ILE
1	C	965	ASP
1	C	975	ILE
1	C	981	ASP
1	C	984	LEU
1	D	698	LEU
1	D	704	ARG
1	D	728	GLN
1	D	731	ASP
1	D	739	LEU
1	D	746	MET
1	D	747	LYS
1	D	751	LEU
1	D	758	VAL
1	D	769	LEU
1	D	777	LEU
1	D	783	ASP
1	D	805	ASP
1	D	807	ASP
1	D	812	GLU
1	D	853	GLN
1	D	864	SER
1	D	878	ARG
1	D	912	GLN
1	D	961	ASN
1	D	962	ILE
1	D	965	ASP
1	D	975	ILE
1	D	981	ASP
1	D	984	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	705	GLN
1	A	906	ASN
1	B	705	GLN
1	C	705	GLN
1	D	705	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

8 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	SO4	A	1101	-	4,4,4	0.42	0	6,6,6	1.08	1 (16%)
2	SO4	A	1102	-	4,4,4	0.10	0	6,6,6	0.32	0
3	RPB	A	1103	-	25,27,27	5.92	10 (40%)	26,39,39	2.09	5 (19%)
2	SO4	B	1101	-	4,4,4	0.50	0	6,6,6	0.37	0
2	SO4	B	1102	-	4,4,4	0.34	0	6,6,6	0.30	0
3	RPB	B	1103	-	25,27,27	6.36	10 (40%)	26,39,39	1.99	5 (19%)
2	SO4	C	1101	-	4,4,4	0.54	0	6,6,6	0.34	0
2	SO4	D	1101	-	4,4,4	0.57	0	6,6,6	0.45	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	SO4	A	1101	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SO4	A	1102	-	-	0/0/0/0	0/0/0/0
3	RPB	A	1103	-	-	0/6/17/17	0/3/4/4
2	SO4	B	1101	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1102	-	-	0/0/0/0	0/0/0/0
3	RPB	B	1103	-	-	0/6/17/17	0/3/4/4
2	SO4	C	1101	-	-	0/0/0/0	0/0/0/0
2	SO4	D	1101	-	-	0/0/0/0	0/0/0/0

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1103	RPB	C9-C10	-16.60	1.36	1.51
3	B	1103	RPB	F1-C2	-15.92	0.98	1.36
3	A	1103	RPB	C9-C10	-15.82	1.37	1.51
3	B	1103	RPB	C12-C11	-14.56	1.32	1.49
3	A	1103	RPB	C12-C11	-13.74	1.33	1.49
3	A	1103	RPB	F1-C2	-13.32	1.04	1.36
3	A	1103	RPB	C4-C7	-11.57	1.37	1.49
3	B	1103	RPB	C4-C7	-11.24	1.38	1.49
3	B	1103	RPB	C1-C6	-7.17	1.30	1.41
3	A	1103	RPB	C1-C6	-6.42	1.31	1.41
3	B	1103	RPB	C18-C16	-5.43	1.39	1.51
3	A	1103	RPB	C18-C16	-5.36	1.39	1.51
3	B	1103	RPB	C10-C5	-4.70	1.32	1.41
3	B	1103	RPB	C4-C5	-4.32	1.32	1.41
3	A	1103	RPB	C10-C5	-4.32	1.33	1.41
3	A	1103	RPB	C4-C5	-4.14	1.32	1.41
3	B	1103	RPB	C11-C10	-3.38	1.32	1.40
3	A	1103	RPB	C11-C10	-3.29	1.33	1.40
3	B	1103	RPB	C6-N2	-2.34	1.31	1.39
3	A	1103	RPB	C6-N2	-2.24	1.31	1.39

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1103	RPB	O1-C7-N1	-6.25	113.26	121.80
3	A	1103	RPB	O1-C7-N1	-6.17	113.37	121.80
3	A	1103	RPB	C13-C12-C11	-3.32	115.42	120.60
3	B	1103	RPB	C13-C12-C11	-3.12	115.73	120.60
3	A	1103	RPB	C19-N3-C18	-2.22	107.64	111.97
2	A	1101	SO4	O2-S-O1	-2.07	102.68	109.59
3	B	1103	RPB	C19-N3-C18	-2.03	108.01	111.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1103	RPB	C14-C12-C11	2.45	124.41	120.60
3	A	1103	RPB	C14-C12-C11	3.47	126.00	120.60
3	A	1103	RPB	O1-C7-C4	4.22	123.23	119.32
3	B	1103	RPB	O1-C7-C4	4.54	123.53	119.32

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1103	RPB	4	0
3	B	1103	RPB	3	0
2	D	1101	SO4	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	347/352 (98%)	-0.21	4 (1%) 81 69	31, 64, 106, 144	0
1	B	349/352 (99%)	-0.18	3 (0%) 85 78	33, 71, 113, 147	0
1	C	347/352 (98%)	-0.00	14 (4%) 42 27	42, 85, 147, 169	0
1	D	349/352 (99%)	0.09	14 (4%) 42 27	40, 93, 158, 179	0
All	All	1392/1408 (98%)	-0.08	35 (2%) 61 47	31, 77, 144, 179	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	951	LEU	6.0
1	D	980	ASN	5.1
1	D	933	LYS	5.0
1	C	825	THR	4.4
1	C	826	HIS	4.2
1	D	826	HIS	3.5
1	D	979	VAL	3.4
1	D	785	SER	3.2
1	C	951	LEU	3.0
1	D	954	THR	3.0
1	B	746	MET	2.9
1	A	930	TYR	2.9
1	C	933	LYS	2.8
1	D	952	GLY	2.7
1	D	953	LYS	2.6
1	C	979	VAL	2.6
1	C	980	ASN	2.6
1	B	855	HIS	2.3
1	C	935	ALA	2.3
1	C	917	GLY	2.3
1	C	953	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	855	HIS	2.3
1	C	982	THR	2.2
1	D	931	GLU	2.2
1	A	979	VAL	2.2
1	D	982	THR	2.2
1	C	823	ALA	2.2
1	C	934	HIS	2.1
1	D	825	THR	2.1
1	A	746	MET	2.1
1	C	824	THR	2.1
1	C	930	TYR	2.1
1	D	930	TYR	2.1
1	A	951	LEU	2.1
1	B	748	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 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	RPB	B	1103	24/24	0.88	0.29	2.52	67,84,87,88	0
3	RPB	A	1103	24/24	0.90	0.26	2.01	57,66,74,74	0
2	SO4	D	1101	5/5	0.99	0.21	1.59	39,42,48,50	0
2	SO4	C	1101	5/5	0.99	0.21	1.03	51,52,54,56	0
2	SO4	B	1101	5/5	0.98	0.20	0.02	54,57,62,62	0
2	SO4	A	1101	5/5	0.97	0.19	0.01	50,51,55,62	0
2	SO4	B	1102	5/5	0.87	0.20	-0.40	95,97,100,103	0
2	SO4	A	1102	5/5	0.92	0.17	-0.55	97,99,99,99	0

6.5 Other polymers [i](#)

There are no such residues in this entry.