Full wwPDB X-ray Structure Validation Report ⓘ

Nov 5, 2023 – 12:07 PM EST

PDB ID	:	5JR6
Title	:	The Xray Crystal Structure of P. falciparum Aminopeptidase P in Complex With Apstatin
Authors	:	Drinkwater, N.; McGowan, S.
Deposited on	:	2016-05-05
Resolution	:	2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references ⓘ) were used in the production of this report:

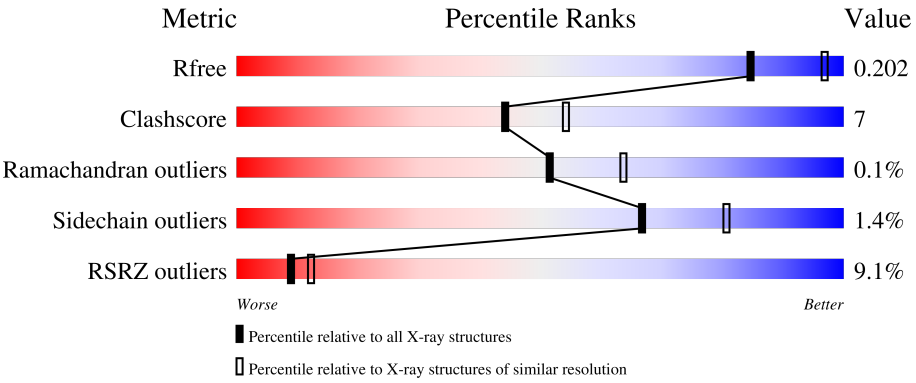
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, 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 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	664	9% 78% 16% 5%
1	B	664	7% 71% 11% 18%
2	F	5	20% 40% 60%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 9214 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 Peptidase, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	628	Total	C	N	O	S	0	0	0
			4841	3120	794	911	16			
1	B	545	Total	C	N	O	S	0	0	0
			4138	2659	687	778	14			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	120	MET	-	initiating methionine	UNP Q8IKT5
A	778	HIS	-	expression tag	UNP Q8IKT5
A	779	HIS	-	expression tag	UNP Q8IKT5
A	780	HIS	-	expression tag	UNP Q8IKT5
A	781	HIS	-	expression tag	UNP Q8IKT5
A	782	HIS	-	expression tag	UNP Q8IKT5
A	783	HIS	-	expression tag	UNP Q8IKT5
B	120	MET	-	initiating methionine	UNP Q8IKT5
B	778	HIS	-	expression tag	UNP Q8IKT5
B	779	HIS	-	expression tag	UNP Q8IKT5
B	780	HIS	-	expression tag	UNP Q8IKT5
B	781	HIS	-	expression tag	UNP Q8IKT5
B	782	HIS	-	expression tag	UNP Q8IKT5
B	783	HIS	-	expression tag	UNP Q8IKT5

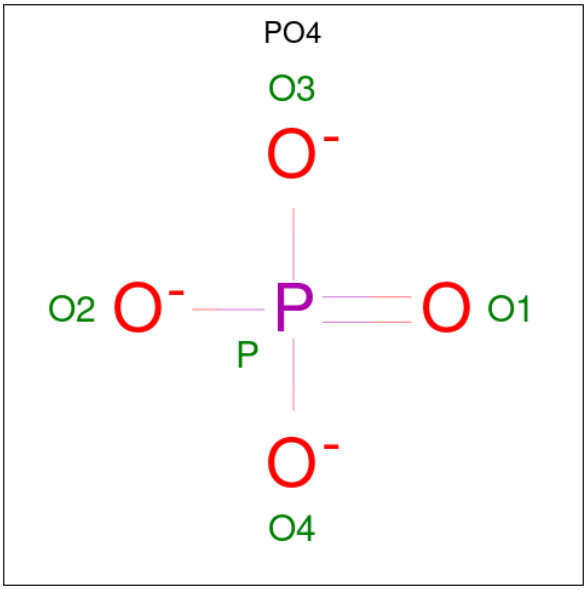
- Molecule 2 is a protein called Apstatin.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	F	5	Total	C	N	O	0	0	1
			33	23	5	5			

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

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

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



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

- Molecule 5 is water.

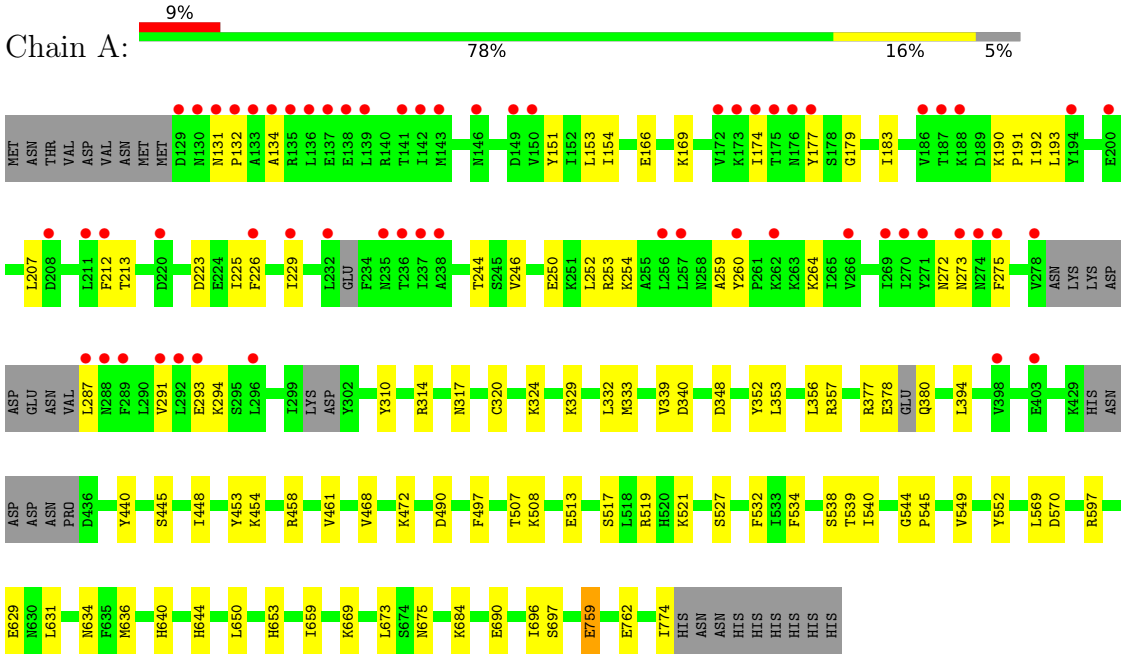
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	105	Total O 105 105	0	0
5	B	88	Total O 88 88	0	0

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.

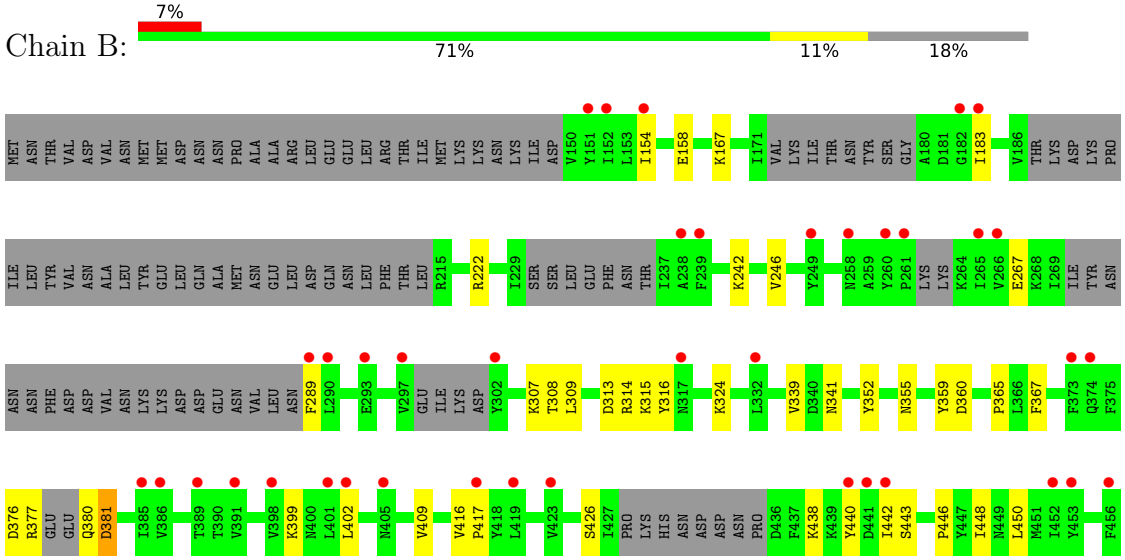
• Molecule 1: Peptidase, putative

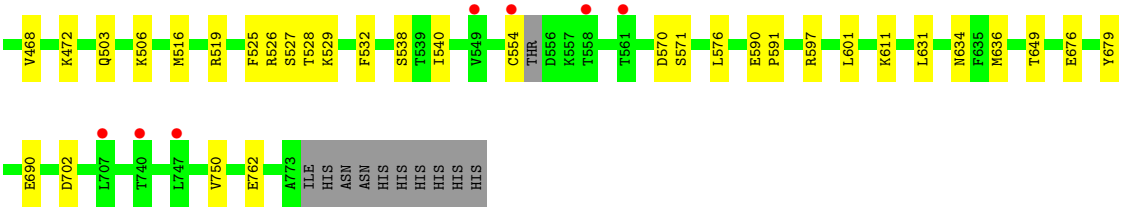
Chain A:



• Molecule 1: Peptidase, putative

Chain B:





• Molecule 2: Apstatin

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	147.22Å 99.93Å 105.17Å 90.00° 105.21° 90.00°	Depositor
Resolution (Å)	42.79 – 2.30 59.43 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.5 (42.79-2.30) 99.5 (59.43-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.70 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.197 , 0.241 0.202 , 0.202	Depositor DCC
R_{free} test set	3232 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å^2)	42.2	Xtriage
Anisotropy	0.499	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 62.3	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	9214	wwPDB-VP
Average B, all atoms (Å^2)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.21% 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: NH2, MN, 01B, PO4

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	0.42	0/4938	0.58	0/6706
1	B	0.40	0/4218	0.56	0/5724
2	F	2.74	2/20 (10.0%)	1.20	0/27
All	All	0.43	2/9176 (0.0%)	0.57	0/12457

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	2	PRO	C-N	8.24	1.50	1.34
2	F	3	PRO	C-N	7.11	1.50	1.34

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	4841	0	4574	69	0
1	B	4138	0	3835	46	0
2	F	33	0	24	3	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	B	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	105	0	0	2	0
5	B	88	0	0	2	0
All	All	9214	0	8433	115	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 7.

All (115) 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:B:377:ARG:O	1:B:380:GLN:CB	2.18	0.91
1:A:669:LYS:HD2	1:A:697:SER:HB2	1.58	0.85
1:B:468:VAL:O	5:B:901:HOH:O	1.98	0.82
1:A:132:PRO:HB3	1:A:174:ILE:O	1.81	0.80
1:B:380:GLN:O	1:B:381:ASP:HB2	1.80	0.80
1:A:154:ILE:HG12	1:A:183:ILE:HG23	1.67	0.77
1:A:272:ASN:HA	1:A:294:LYS:HE2	1.69	0.75
1:A:348:ASP:OD2	1:A:472:LYS:NZ	2.20	0.74
1:A:273:ASN:N	1:A:294:LYS:HZ1	1.88	0.71
1:A:378:GLU:O	1:A:380:GLN:N	2.23	0.71
1:A:291:VAL:HG21	1:A:454:LYS:HG3	1.74	0.69
1:A:225:ILE:O	1:A:229:ILE:HG12	1.93	0.69
1:B:472:LYS:N	5:B:901:HOH:O	2.04	0.68
1:B:519:ARG:NH2	1:B:538:SER:OG	2.26	0.67
1:B:376:ASP:O	1:B:377:ARG:C	2.32	0.67
1:A:629:GLU:OE2	5:A:901:HOH:O	2.13	0.66
1:B:154:ILE:HG23	1:B:183:ILE:HG12	1.81	0.63
1:A:519:ARG:NH2	1:A:538:SER:OG	2.31	0.62
1:B:516:MET:SD	1:B:554:CYS:HB3	2.41	0.61
1:A:549:VAL:HB	1:A:552:TYR:HB2	1.83	0.61
1:A:453:TYR:CE1	1:A:461:VAL:HG21	2.38	0.59
1:B:360:ASP:HB2	1:B:367:PHE:HA	1.84	0.58
1:B:324:LYS:NZ	1:B:355:ASN:OD1	2.38	0.56
1:A:273:ASN:H	1:A:294:LYS:HZ1	1.53	0.55
1:A:759:GLU:HA	1:A:762:GLU:HG2	1.89	0.55
1:A:320:CYS:SG	5:A:1004:HOH:O	2.58	0.54
1:B:313:ASP:OD1	1:B:314:ARG:N	2.40	0.54
1:A:207:LEU:HD11	1:A:212:PHE:HB2	1.90	0.54
1:B:324:LYS:HE2	1:B:352:TYR:CZ	2.43	0.53
1:A:631:LEU:HD23	1:A:636:MET:HE3	1.90	0.52
1:B:365:PRO:HG3	1:B:649:THR:HG22	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:540:ILE:HB	1:B:570:ASP:HB3	1.91	0.52
1:A:329:LYS:NZ	1:A:380:GLN:O	2.39	0.52
1:A:453:TYR:HE1	1:A:461:VAL:HG21	1.75	0.52
1:A:659:ILE:HG13	1:A:673:LEU:HD22	1.91	0.51
1:A:640:HIS:CG	2:F:2:PRO:HB3	2.46	0.51
1:A:250:GLU:HB3	1:A:254:LYS:HZ3	1.76	0.51
1:B:527:SER:HA	1:B:532:PHE:CD2	2.45	0.50
1:A:597:ARG:NH2	1:A:634:ASN:OD1	2.45	0.50
1:B:246:VAL:HG13	1:B:448:ILE:O	2.11	0.50
1:A:329:LYS:HG2	1:A:333:MET:HE2	1.94	0.50
1:B:601:LEU:HD13	1:B:636:MET:HE1	1.94	0.50
1:A:356:LEU:O	1:A:357:ARG:HD3	2.12	0.50
1:B:402:LEU:HD13	1:B:409:VAL:HG21	1.95	0.49
1:A:332:LEU:O	1:A:377:ARG:NH1	2.35	0.49
1:B:158:GLU:HA	1:B:242:LYS:O	2.13	0.49
1:A:513:GLU:OE1	1:A:517:SER:OG	2.28	0.49
1:B:438:LYS:HE3	1:B:440:TYR:CE1	2.48	0.48
1:A:527:SER:HA	1:A:532:PHE:CD2	2.49	0.48
1:A:250:GLU:OE1	1:A:458:ARG:NH2	2.46	0.48
1:B:315:LYS:HE3	1:B:316:TYR:CZ	2.48	0.47
1:A:250:GLU:OE2	1:A:254:LYS:NZ	2.45	0.47
1:A:507:THR:O	1:A:508:LYS:HB2	2.15	0.47
1:A:253:ARG:HH21	1:A:254:LYS:HZ2	1.62	0.47
1:A:273:ASN:HD22	1:A:275:PHE:HE2	1.63	0.47
1:A:229:ILE:CD1	1:A:252:LEU:HD13	2.44	0.47
1:A:183:ILE:O	1:A:193:LEU:HD12	2.15	0.47
1:B:339:VAL:HG22	1:B:341:ASN:H	1.80	0.46
1:B:503:GLN:HA	1:B:506:LYS:HE2	1.96	0.46
1:A:540:ILE:HB	1:A:570:ASP:HB3	1.98	0.46
1:B:315:LYS:HE3	1:B:316:TYR:CE2	2.50	0.46
1:A:314:ARG:HA	1:A:317:ASN:OD1	2.16	0.46
1:A:352:TYR:CD2	1:A:472:LYS:HG2	2.51	0.46
1:A:653:HIS:NE2	2:F:1:01B:O	2.39	0.46
1:A:519:ARG:HG3	1:A:539:THR:HB	1.98	0.45
1:A:177:TYR:CZ	1:A:179:GLY:HA3	2.52	0.45
1:A:339:VAL:CG1	1:A:440:TYR:HB2	2.47	0.45
1:B:246:VAL:HG12	1:B:450:LEU:HD23	1.97	0.45
1:A:151:TYR:CE2	1:A:153:LEU:HG	2.51	0.45
1:A:246:VAL:HG13	1:A:448:ILE:O	2.16	0.45
1:B:416:VAL:HB	1:B:417:PRO:HD3	1.98	0.45
1:A:223:ASP:HA	1:A:226:PHE:HD2	1.81	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:497:PHE:CE2	1:A:569:LEU:HD22	2.52	0.44
1:A:259:ALA:C	1:A:260:TYR:HD1	2.20	0.44
1:B:636:MET:HB3	1:B:679:TYR:CE1	2.52	0.44
1:A:310:TYR:CZ	1:A:357:ARG:HB2	2.51	0.44
1:A:675:ASN:O	1:A:690:GLU:HA	2.18	0.44
1:B:267:GLU:HA	1:B:289:PHE:O	2.18	0.43
1:B:246:VAL:HG22	1:B:446:PRO:O	2.18	0.43
1:B:308:THR:HG22	1:B:309:LEU:O	2.18	0.43
1:B:676:GLU:HB3	1:B:690:GLU:HG3	1.99	0.43
1:A:190:LYS:HA	1:A:191:PRO:HD2	1.84	0.43
1:B:529:LYS:HA	1:B:529:LYS:HD3	1.83	0.43
1:A:468:VAL:O	1:A:472:LYS:HG3	2.18	0.43
1:B:526:ARG:HH12	1:B:571:SER:HB2	1.83	0.43
1:B:380:GLN:O	1:B:381:ASP:CB	2.55	0.43
1:B:519:ARG:HH11	1:B:554:CYS:HB2	1.84	0.43
1:A:644:HIS:HB3	2:F:3:PRO:HB3	2.00	0.43
1:A:759:GLU:HA	1:A:762:GLU:CG	2.49	0.43
1:A:166:GLU:HA	1:A:169:LYS:HG3	2.01	0.43
1:B:525:PHE:O	1:B:528:THR:HG22	2.19	0.43
1:A:684:LYS:HB3	1:A:684:LYS:HE3	1.90	0.42
1:B:631:LEU:HD23	1:B:636:MET:HE3	2.01	0.42
1:A:544:GLY:HA3	1:A:545:PRO:HD3	1.88	0.42
1:A:192:ILE:HG13	1:A:213:THR:HB	2.02	0.42
1:A:264:LYS:O	1:A:287:LEU:N	2.53	0.42
1:A:636:MET:HE2	1:A:636:MET:HB2	1.79	0.42
1:B:601:LEU:HD13	1:B:636:MET:CE	2.49	0.41
1:A:352:TYR:HB2	1:A:650:LEU:HD11	2.01	0.41
1:B:590:GLU:HA	1:B:591:PRO:HD3	1.90	0.41
1:B:750:VAL:HG12	1:B:762:GLU:HG2	2.01	0.41
1:A:340:ASP:OD1	1:A:377:ARG:NE	2.50	0.41
1:B:442:ILE:HG13	1:B:443:SER:N	2.35	0.41
1:B:611:LYS:HD3	1:B:611:LYS:HA	1.86	0.41
1:A:131:ASN:O	1:A:134:ALA:HB3	2.21	0.41
1:B:316:TYR:CE2	1:B:576:LEU:HD13	2.55	0.41
1:A:352:TYR:CG	1:A:472:LYS:HG2	2.55	0.41
1:B:597:ARG:NH2	1:B:634:ASN:OD1	2.52	0.41
1:B:307:LYS:HE2	1:B:359:TYR:HB3	2.03	0.40
1:B:399:LYS:HA	1:B:402:LEU:HD12	2.03	0.40
1:A:253:ARG:NH2	1:A:254:LYS:HZ2	2.19	0.40
1:A:272:ASN:HA	1:A:294:LYS:CE	2.44	0.40
1:A:273:ASN:H	1:A:294:LYS:NZ	2.19	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:324:LYS:HB3	1:A:353:LEU:HD12	2.04	0.40
1:A:394:LEU:HD12	1:A:394:LEU:HA	1.92	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	616/664 (93%)	602 (98%)	14 (2%)	0	100 100
1	B	525/664 (79%)	511 (97%)	13 (2%)	1 (0%)	47 58
2	F	3/5 (60%)	1 (33%)	2 (67%)	0	100 100
All	All	1144/1333 (86%)	1114 (97%)	29 (2%)	1 (0%)	51 64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	381	ASP

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	491/613 (80%)	482 (98%)	9 (2%)	59 75
1	B	409/613 (67%)	405 (99%)	4 (1%)	76 87

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	F	2/2 (100%)	2 (100%)	0	100 100
All	All	902/1228 (74%)	889 (99%)	13 (1%)	67 81

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	244	THR
1	A	293	GLU
1	A	445	SER
1	A	490	ASP
1	A	521	LYS
1	A	534	PHE
1	A	696	ILE
1	A	759	GLU
1	A	774	ILE
1	B	167	LYS
1	B	222	ARG
1	B	426	SER
1	B	702	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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	01B	F	1	2,3	11,13,14	3.66	4 (36%)	11,16,18	0.68	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	01B	F	1	2,3	-	8/9/10/12	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	1	01B	C12-C7	7.23	1.54	1.38
2	F	1	01B	C9-C8	7.09	1.53	1.38
2	F	1	01B	C11-C10	5.37	1.52	1.38
2	F	1	01B	C6-C7	2.58	1.57	1.51

There are no bond angle outliers.

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	F	1	01B	O2-CA-CB-N
2	F	1	01B	C-CA-CB-N
2	F	1	01B	O2-CA-CB-C6
2	F	1	01B	C-CA-CB-C6
2	F	1	01B	CB-C6-C7-C8
2	F	1	01B	CB-C6-C7-C12
2	F	1	01B	C7-C6-CB-N
2	F	1	01B	C7-C6-CB-CA

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	1	01B	1	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 5 ligands modelled in this entry, 4 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	PO4	B	803	3	4,4,4	0.82	0	6,6,6	0.64	0

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

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	628/664 (94%)	0.49	60 (9%)	8	10	26, 60, 124, 163	3 (0%)
1	B	545/664 (82%)	0.65	46 (8%)	11	15	31, 69, 124, 168	1 (0%)
2	F	3/5 (60%)	2.10	1 (33%)	0	0	82, 82, 101, 106	0
All	All	1176/1333 (88%)	0.57	107 (9%)	9	12	26, 65, 124, 168	4 (0%)

All (107) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	212	PHE	8.1
1	A	296	LEU	7.4
1	B	401	LEU	7.3
1	A	134	ALA	6.7
1	A	174	ILE	6.6
1	A	260	TYR	5.9
1	B	238	ALA	5.7
1	B	405	ASN	5.5
1	A	136	LEU	5.5
1	A	236	THR	5.3
1	A	269	ILE	5.1
1	A	289	PHE	5.0
1	A	131	ASN	5.0
1	A	270	ILE	4.8
1	A	237	ILE	4.5
1	B	260	TYR	4.5
1	A	137	GLU	4.4
1	B	398	VAL	4.3
1	A	208	ASP	4.2
1	B	419	LEU	4.2
1	A	229	ILE	4.2
2	F	4	ALA	4.2
1	A	211	LEU	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	232	LEU	4.1
1	B	456	PHE	4.1
1	A	150	VAL	4.1
1	A	262	LYS	4.0
1	A	235	ASN	4.0
1	A	403	GLU	4.0
1	A	288	ASN	4.0
1	A	146	ASN	4.0
1	A	275	PHE	3.8
1	B	554	CYS	3.8
1	A	176	ASN	3.6
1	B	402	LEU	3.6
1	B	239	PHE	3.6
1	A	186	VAL	3.5
1	B	258	ASN	3.5
1	B	249	TYR	3.5
1	A	129	ASP	3.5
1	B	385	ILE	3.5
1	A	238	ALA	3.4
1	A	135	ARG	3.3
1	A	143	MET	3.3
1	B	266	VAL	3.3
1	B	152	ILE	3.1
1	B	417	PRO	3.1
1	A	138	GLU	3.1
1	A	271	TYR	3.1
1	A	141	THR	3.1
1	B	452	ILE	3.1
1	A	130	ASN	3.0
1	A	266	VAL	2.9
1	B	441	ASP	2.9
1	B	302	TYR	2.9
1	A	291	VAL	2.8
1	A	139	LEU	2.8
1	B	154	ILE	2.8
1	A	175	THR	2.7
1	B	558	THR	2.7
1	A	177	TYR	2.7
1	A	149	ASP	2.6
1	B	423	VAL	2.6
1	A	132	PRO	2.6
1	B	151	TYR	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	374	GLN	2.5
1	A	256	LEU	2.5
1	A	293	GLU	2.5
1	B	293	GLU	2.5
1	A	220	ASP	2.5
1	A	287	LEU	2.5
1	B	391	VAL	2.5
1	A	273	ASN	2.5
1	B	453	TYR	2.4
1	A	172	VAL	2.4
1	A	226	PHE	2.4
1	B	317	ASN	2.3
1	B	707	LEU	2.3
1	A	187	THR	2.3
1	A	173	LYS	2.3
1	B	265	ILE	2.3
1	B	386	VAL	2.3
1	B	261	PRO	2.3
1	A	194	TYR	2.3
1	B	183	ILE	2.3
1	A	278	VAL	2.3
1	A	274	ASN	2.3
1	B	389	THR	2.2
1	B	549	VAL	2.2
1	A	142	ILE	2.2
1	A	292	LEU	2.2
1	B	440	TYR	2.2
1	B	289	PHE	2.2
1	A	133	ALA	2.2
1	A	188	LYS	2.1
1	B	290	LEU	2.1
1	A	200	GLU	2.1
1	B	297	VAL	2.1
1	B	561	THR	2.1
1	B	182	GLY	2.1
1	A	398	VAL	2.1
1	B	442	ILE	2.1
1	B	373	PHE	2.0
1	A	257	LEU	2.0
1	B	332	LEU	2.0
1	B	747	LEU	2.0
1	B	740	THR	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, 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	01B	F	1	13/14	0.91	0.28	35,72,82,82	6

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	MN	B	801	1/1	0.95	0.06	74,74,74,74	0
3	MN	A	801	1/1	0.98	0.09	62,62,62,62	0
3	MN	B	802	1/1	0.98	0.07	57,57,57,57	0
4	PO4	B	803	5/5	0.98	0.10	66,73,79,82	0
3	MN	A	802	1/1	0.99	0.13	42,42,42,42	0

6.5 Other polymers (i)

There are no such residues in this entry.