



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 09:39 am BST

PDB ID : 6DAQ
Title : PhdJ bound to substrate intermediate
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Deposited on : 2018-05-01
Resolution : 2.00 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 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 : 2.11
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) : 2.11

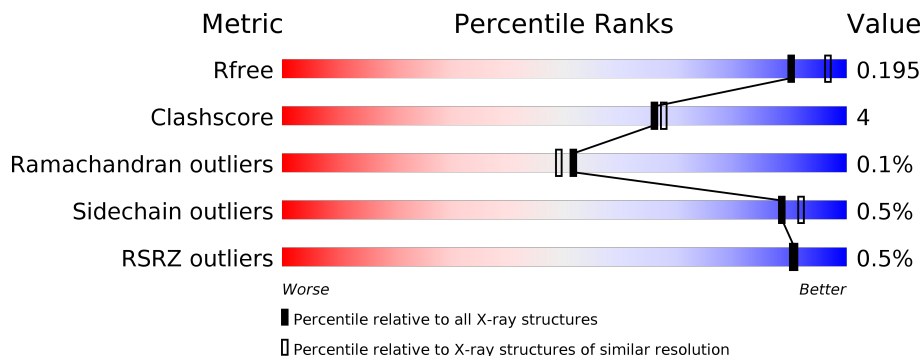
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 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 $\leq 5\%$. 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	334	 90% 7%
1	C	334	 92% 5%
1	D	334	 87% 8%
2	B	334	 88% 8%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11044 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 PhdJ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	323	2466	1569	419	469	9	0	0	0
1	C	323	2466	1569	419	469	9	0	0	0
1	D	323	2466	1569	419	469	9	0	0	0

- Molecule 2 is a protein called PhdJ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	323	2456	1561	419	467	9	0	0	0

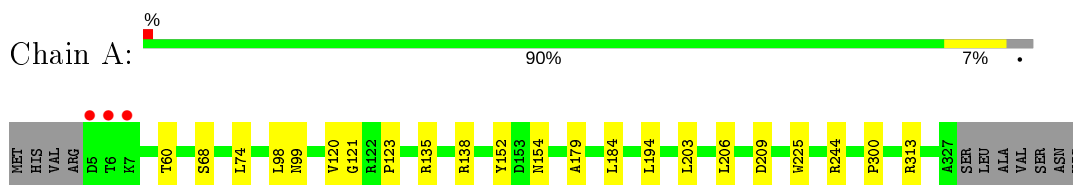
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	302	Total 302	O 302	0	0
3	B	306	Total 306	O 306	0	0
3	C	289	Total 289	O 289	0	0
3	D	293	Total 293	O 293	0	0

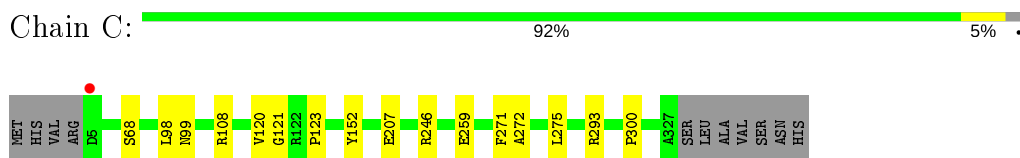
3 Residue-property plots [i](#)

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.

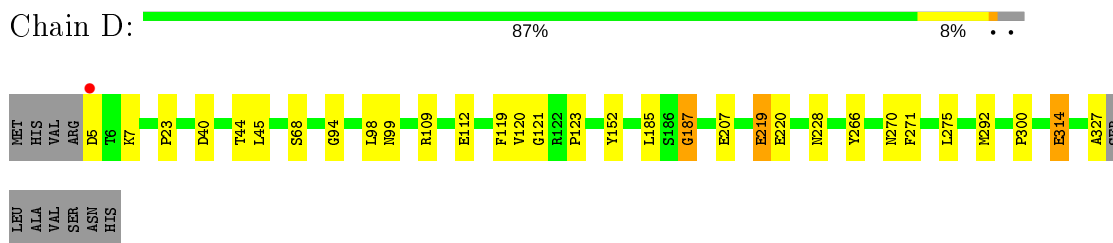
- Molecule 1: PhdJ



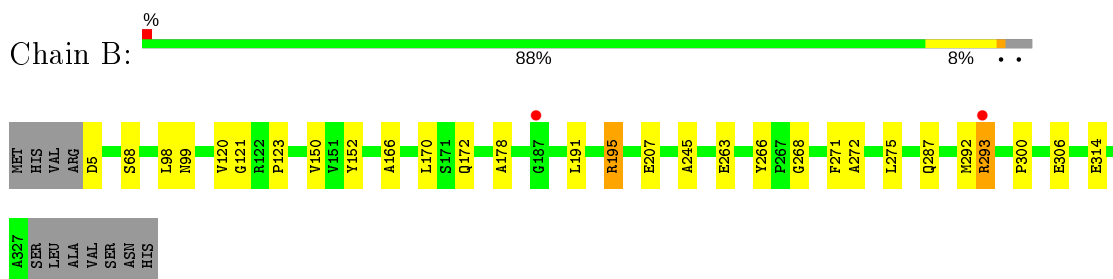
- Molecule 1: PhdJ



- Molecule 1: PhdJ



- Molecule 2: PhdJ



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	52.53Å 137.20Å 88.27Å 90.00° 98.91° 90.00°	Depositor
Resolution (Å)	45.29 – 2.00 45.29 – 2.00	Depositor EDS
% Data completeness (in resolution range)	93.4 (45.29-2.00) 93.3 (45.29-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.91 (at 2.00Å)	Xtrriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.160 , 0.194 0.160 , 0.195	Depositor DCC
R_{free} test set	2000 reflections (2.59%)	wwPDB-VP
Wilson B-factor (Å ²)	20.5	Xtrriage
Anisotropy	0.470	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 44.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11044	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.70% 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, 0.333 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: KPI, 9KP

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	0.30	0/2498	0.48	0/3411
1	C	0.31	0/2498	0.52	2/3411 (0.1%)
1	D	0.31	0/2498	0.49	0/3411
2	B	0.31	0/2498	0.55	3/3411 (0.1%)
All	All	0.31	0/9992	0.51	5/13644 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	195	ARG	NE-CZ-NH2	-11.68	114.46	120.30
1	C	293	ARG	NE-CZ-NH1	10.92	125.76	120.30
1	C	293	ARG	NE-CZ-NH2	-7.84	116.38	120.30
2	B	195	ARG	NE-CZ-NH1	6.06	123.33	120.30
2	B	293	ARG	NE-CZ-NH2	5.81	123.21	120.30

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	2466	0	2393	17	0
1	C	2466	0	2393	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2466	0	2393	27	0
2	B	2456	0	2406	26	0
3	A	302	0	0	3	0
3	B	306	0	0	7	2
3	C	289	0	0	5	2
3	D	293	0	0	9	0
All	All	11044	0	9585	77	2

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 (77) 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:D:187:GLY:O	3:D:401:HOH:O	1.90	0.88
2:B:5:ASP:N	3:B:405:HOH:O	2.05	0.88
1:C:246:ARG:NH1	3:C:401:HOH:O	2.08	0.87
2:B:314:GLU:OE1	3:B:401:HOH:O	1.94	0.85
2:B:263:GLU:OE2	3:B:402:HOH:O	1.98	0.82
1:C:272:ALA:HA	1:C:275:LEU:HD23	1.63	0.80
2:B:266:TYR:OH	3:B:403:HOH:O	1.98	0.79
1:C:271:PHE:O	1:C:275:LEU:HD22	1.84	0.75
1:D:314:GLU:OE2	3:D:402:HOH:O	2.04	0.75
1:D:112:GLU:OE1	3:D:403:HOH:O	2.07	0.71
1:C:272:ALA:O	3:C:402:HOH:O	2.09	0.70
1:D:270:ASN:OD1	3:D:405:HOH:O	2.12	0.67
2:B:191:LEU:O	2:B:195:ARG:NH1	2.29	0.65
1:C:300:PRO:HG3	1:D:123:PRO:HG2	1.80	0.64
1:C:259:GLU:OE1	3:C:403:HOH:O	2.15	0.64
1:D:266:TYR:OH	3:D:404:HOH:O	2.09	0.62
1:A:300:PRO:HG3	2:B:123:PRO:HG2	1.82	0.61
2:B:293:ARG:HA	2:B:293:ARG:CZ	2.31	0.61
1:D:23:PRO:HG3	1:D:292:MET:HE1	1.85	0.58
2:B:292:MET:O	2:B:293:ARG:NH1	2.36	0.58
1:A:123:PRO:HG2	2:B:300:PRO:HG3	1.84	0.57
1:C:123:PRO:HG2	1:D:300:PRO:HG3	1.86	0.57
1:D:327:ALA:O	3:D:406:HOH:O	2.16	0.57
2:B:287:GLN:HG2	2:B:293:ARG:NH2	2.21	0.56
1:C:120:VAL:HG12	1:C:121:GLY:O	2.06	0.56
1:D:185:LEU:HG	3:D:533:HOH:O	2.06	0.55
2:B:287:GLN:HG2	2:B:293:ARG:HH22	1.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:120:VAL:HG12	1:D:121:GLY:O	2.07	0.54
1:D:109:ARG:NH1	3:D:403:HOH:O	2.34	0.53
1:A:184:LEU:O	1:D:187:GLY:HA3	2.07	0.53
1:A:313:ARG:NE	3:A:402:HOH:O	2.26	0.53
1:C:99:ASN:HB3	1:D:68:SER:HA	1.91	0.53
1:A:99:ASN:HB3	2:B:68:SER:HA	1.91	0.51
1:D:45:LEU:HD22	1:D:292:MET:HE3	1.93	0.51
1:A:120:VAL:HG12	1:A:121:GLY:O	2.12	0.50
1:A:60:THR:HG21	1:A:74:LEU:HD12	1.93	0.50
1:C:68:SER:HA	1:D:99:ASN:HB3	1.94	0.50
1:D:219:GLU:H	1:D:219:GLU:CD	2.14	0.50
2:B:172:GLN:NE2	3:B:406:HOH:O	2.30	0.49
2:B:272:ALA:HA	2:B:275:LEU:HD23	1.95	0.49
2:B:120:VAL:HG12	2:B:121:GLY:O	2.13	0.49
2:B:98:LEU:O	2:B:123:PRO:HB3	2.12	0.49
2:B:245:ALA:HB1	3:B:474:HOH:O	2.11	0.48
1:D:271:PHE:CE2	1:D:275:LEU:HD21	2.48	0.48
1:A:206:LEU:HG	1:A:225:TRP:CZ2	2.50	0.47
1:A:135:ARG:HH22	1:A:138:ARG:HD2	1.79	0.47
1:D:5:ASP:N	3:D:415:HOH:O	2.47	0.47
1:A:179:ALA:HB3	1:A:203:LEU:HD23	1.98	0.46
2:B:271:PHE:CE2	2:B:275:LEU:HD21	2.50	0.46
1:A:244:ARG:NE	3:A:401:HOH:O	2.18	0.46
2:B:150:VAL:HG22	2:B:178:ALA:HB3	1.97	0.46
2:B:293:ARG:NE	2:B:293:ARG:HA	2.28	0.46
1:A:209:ASP:HA	3:A:584:HOH:O	2.16	0.46
1:A:98:LEU:O	1:A:123:PRO:HB3	2.16	0.46
1:A:135:ARG:NH2	1:A:138:ARG:HD2	2.30	0.45
1:A:194:LEU:HA	1:A:203:LEU:HD12	1.98	0.45
1:A:68:SER:HA	2:B:99:ASN:HB3	1.97	0.45
2:B:268:GLY:N	3:B:404:HOH:O	2.02	0.45
1:C:275:LEU:H	1:C:275:LEU:HD22	1.81	0.45
1:D:219:GLU:HG2	1:D:220:GLU:HG3	1.98	0.45
2:B:287:GLN:HA	2:B:293:ARG:HH22	1.82	0.44
1:C:272:ALA:HA	1:C:275:LEU:CD2	2.43	0.44
1:D:98:LEU:O	1:D:123:PRO:HB3	2.16	0.44
1:D:45:LEU:HD22	1:D:292:MET:CE	2.47	0.44
1:C:246:ARG:NH2	3:C:404:HOH:O	2.16	0.44
1:D:121:GLY:HA2	1:D:152:TYR:HB3	2.00	0.43
2:B:306:GLU:H	2:B:306:GLU:CD	2.21	0.43
1:C:108:ARG:NH1	3:C:417:HOH:O	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:94:GLY:HA2	1:D:119:PHE:O	2.19	0.43
1:C:121:GLY:HA2	1:C:152:TYR:HB3	2.01	0.42
1:D:207:GLU:CD	1:D:228:ASN:HD22	2.21	0.42
1:D:207:GLU:OE2	1:D:228:ASN:ND2	2.49	0.42
1:D:40:ASP:O	1:D:44:THR:HG23	2.19	0.42
1:C:98:LEU:O	1:C:123:PRO:HB3	2.19	0.42
2:B:166:ALA:O	2:B:170:LEU:HG	2.20	0.41
1:A:152:TYR:CE1	1:A:154:ASN:HB2	2.55	0.41
2:B:121:GLY:HA2	2:B:152:TYR:HB3	2.03	0.41

All (2) 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)
3:B:508:HOH:O	3:C:618:HOH:O[2_453]	2.07	0.13
3:B:619:HOH:O	3:C:677:HOH:O[2_453]	2.14	0.06

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	320/334 (96%)	317 (99%)	3 (1%)	0	100	100
1	C	320/334 (96%)	316 (99%)	4 (1%)	0	100	100
1	D	320/334 (96%)	314 (98%)	5 (2%)	1 (0%)	41	37
2	B	320/334 (96%)	315 (98%)	5 (2%)	0	100	100
All	All	1280/1336 (96%)	1262 (99%)	17 (1%)	1 (0%)	51	49

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	187	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	250/260 (96%)	250 (100%)	0	100	100
1	C	250/260 (96%)	249 (100%)	1 (0%)	91	93
1	D	250/260 (96%)	247 (99%)	3 (1%)	71	76
2	B	250/260 (96%)	249 (100%)	1 (0%)	91	93
All	All	1000/1040 (96%)	995 (100%)	5 (0%)	88	92

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	207	GLU
1	C	207	GLU
1	D	7	LYS
1	D	219	GLU
1	D	314	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	324	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

4 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
1	9KP	A	180	1	19,24,25	3.13	4 (21%)	15,30,32	3.11	3 (20%)
2	KPI	B	180	2	10,13,14	2.40	2 (20%)	6,15,17	4.23	3 (50%)
1	9KP	C	180	1	19,24,25	3.09	4 (21%)	15,30,32	3.22	3 (20%)
1	9KP	D	180	1	19,24,25	3.10	4 (21%)	15,30,32	2.98	3 (20%)

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
1	9KP	A	180	1	-	4/13/23/25	0/1/1/1
2	KPI	B	180	2	-	1/9/14/16	-
1	9KP	C	180	1	-	4/13/23/25	0/1/1/1
1	9KP	D	180	1	-	4/13/23/25	0/1/1/1

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	180	9KP	C14-C15	8.69	1.55	1.33
1	D	180	9KP	C14-C15	8.61	1.55	1.33
1	C	180	9KP	C14-C15	8.57	1.55	1.33
1	A	180	9KP	C17-C22	7.04	1.54	1.47
1	C	180	9KP	C17-C22	6.95	1.54	1.47
1	D	180	9KP	C17-C22	6.83	1.54	1.47
1	C	180	9KP	C16-C15	5.74	1.56	1.47
1	D	180	9KP	C16-C15	5.70	1.56	1.47
1	A	180	9KP	C16-C15	5.67	1.56	1.47
2	B	180	KPI	CX1-NZ	5.61	1.45	1.29
2	B	180	KPI	C1-CX1	-4.89	1.40	1.50
1	A	180	9KP	C14-C10	4.77	1.54	1.46
1	D	180	9KP	C14-C10	4.69	1.54	1.46
1	C	180	9KP	C14-C10	4.47	1.53	1.46

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	180	9KP	C16-C15-C14	-9.44	108.17	126.93
1	C	180	9KP	C16-C15-C14	-8.71	109.62	126.93
1	D	180	9KP	C16-C15-C14	-8.43	110.18	126.93
1	C	180	9KP	C15-C14-C10	-8.15	115.16	125.43
1	D	180	9KP	C15-C14-C10	-7.20	116.36	125.43
2	B	180	KPI	C1-CX1-CX2	-7.09	110.07	117.92
1	A	180	9KP	C15-C14-C10	-6.86	116.79	125.43
2	B	180	KPI	C1-CX1-NZ	-5.95	108.27	123.12
2	B	180	KPI	CE-NZ-CX1	-4.50	109.66	121.77
1	C	180	9KP	CD-CE-NZ	-2.66	105.81	110.66
1	A	180	9KP	CD-CE-NZ	-2.26	106.55	110.66
1	D	180	9KP	CD-CE-NZ	-2.17	106.72	110.66

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	180	9KP	C11-C10-C14-C15
1	A	180	9KP	NZ-C10-C14-C15
2	B	180	KPI	C1-CX1-NZ-CE
1	C	180	9KP	C11-C10-C14-C15
1	C	180	9KP	NZ-C10-C14-C15
1	A	180	9KP	C14-C15-C16-C21
1	C	180	9KP	C14-C15-C16-C17
1	A	180	9KP	C14-C15-C16-C17
1	C	180	9KP	C14-C15-C16-C21
1	D	180	9KP	C14-C15-C16-C17
1	D	180	9KP	NZ-C10-C14-C15
1	D	180	9KP	C11-C10-C14-C15
1	D	180	9KP	C14-C15-C16-C21

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	322/334 (96%)	-0.51	3 (0%) 84 83	13, 19, 32, 53	0
1	C	322/334 (96%)	-0.51	1 (0%) 94 93	15, 20, 34, 55	0
1	D	322/334 (96%)	-0.56	1 (0%) 94 93	15, 20, 33, 55	0
2	B	322/334 (96%)	-0.61	2 (0%) 89 88	14, 19, 33, 45	0
All	All	1288/1336 (96%)	-0.55	7 (0%) 91 90	13, 20, 33, 55	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	5	ASP	4.7
1	A	5	ASP	3.9
1	D	5	ASP	3.6
2	B	187	GLY	2.7
2	B	293	ARG	2.2
1	A	6	THR	2.1
1	A	7	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	9KP	A	180	24/25	0.91	0.13	14,21,33,35	0
1	9KP	C	180	24/25	0.92	0.13	15,24,33,36	0
2	KPI	B	180	14/15	0.93	0.10	14,18,24,32	0
1	9KP	D	180	24/25	0.93	0.10	15,24,32,35	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.