



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:33 PM UTC

PDB ID : 4PV8 / pdb_00004pv8
Title : Crystal Structure of H2Kb-Q600F complex
Authors : Twist, K.-A.; Rossjohn, J.; Gras, S.
Deposited on : 2014-03-15
Resolution : 2.31 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

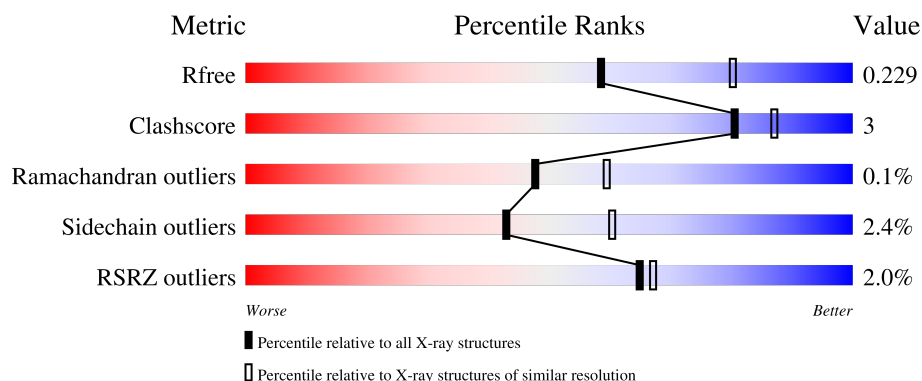
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.31 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}	180053	7754 (2.34-2.30)
Clashscore	190562	8383 (2.34-2.30)
Ramachandran outliers	187476	8303 (2.34-2.30)
Sidechain outliers	187428	8303 (2.34-2.30)
RSRZ outliers	180081	7760 (2.34-2.30)

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 $\leq 5\%$. 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	278	3% 90% 10%
1	C	278	2% 87% 13%
2	B	99	2% 88% 12%
2	D	99	% 98% .
3	E	8	100%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	F	8	 A horizontal bar chart showing the quality of chain F. The bar is divided into three segments: green (62%), yellow (25%), and orange (12%).

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6850 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 H-2 class I histocompatibility antigen, K-B alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	278	Total	C	N	O	S	7	8	0
			2312	1461	406	436	9			
1	C	278	Total	C	N	O	S	5	11	0
			2323	1470	408	435	10			

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	99	Total	C	N	O	S	0	1	0
			829	530	139	153	7			
2	D	99	Total	C	N	O	S	0	1	0
			829	530	139	153	7			

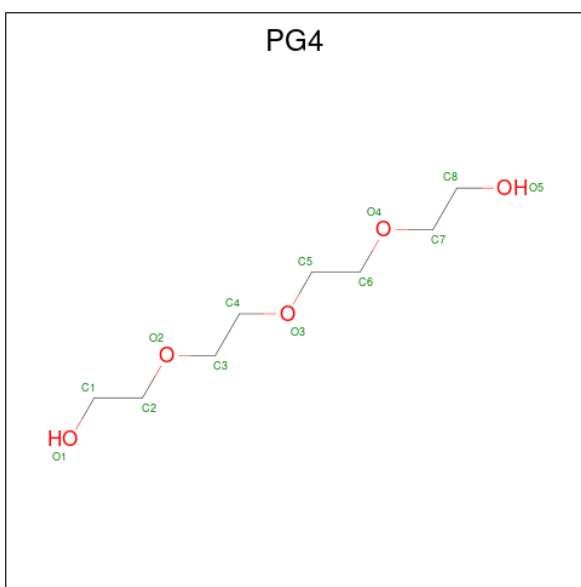
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	85	ASP	ALA	conflict	UNP P01887
D	85	ASP	ALA	conflict	UNP P01887

- Molecule 3 is a protein called S598 peptide modified Q600F.

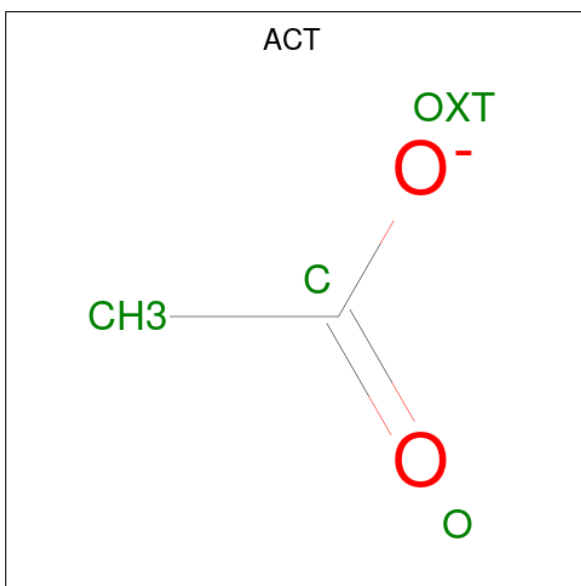
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	E	8	Total	C	N	O	0	0	0
			69	47	12	10			
3	F	8	Total	C	N	O	0	0	0
			69	47	12	10			

- Molecule 4 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: C₈H₁₈O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			13	8	5		

- Molecule 5 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2^-$).



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

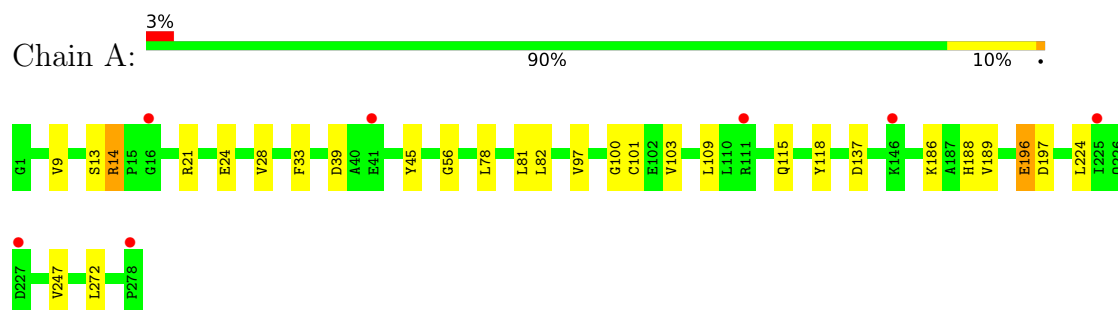
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	150	Total 150	O 150	0	0
6	B	58	Total 58	O 58	0	0
6	C	123	Total 123	O 123	0	0
6	D	61	Total 61	O 61	0	0
6	E	7	Total 7	O 7	0	0
6	F	3	Total 3	O 3	0	0

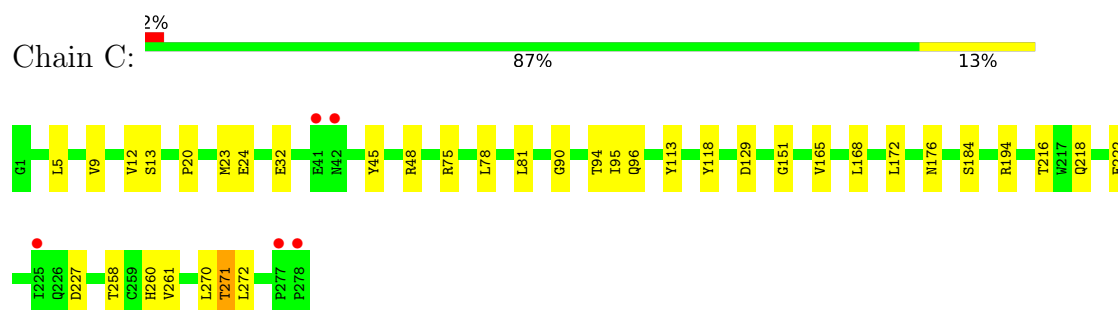
3 Residue-property plots [i](#)

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.

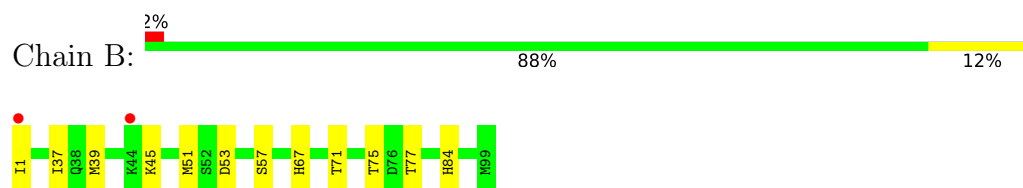
- Molecule 1: H-2 class I histocompatibility antigen, K-B alpha chain



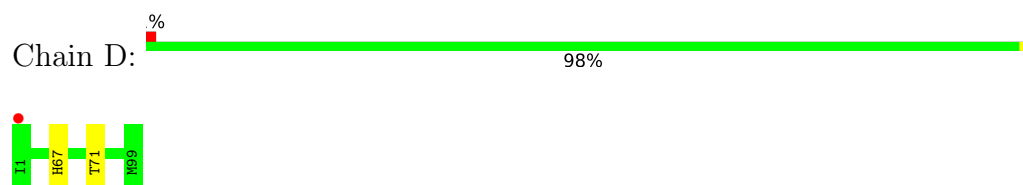
- Molecule 1: H-2 class I histocompatibility antigen, K-B alpha chain



- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin



- Molecule 3: S598 peptide modified Q600F

Chain E:  100%

There are no outlier residues recorded for this chain.

- Molecule 3: S598 peptide modified Q600F

Chain F:  62% 25% 12%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.46Å 90.99Å 89.61Å 90.00° 111.65° 90.00°	Depositor
Resolution (Å)	33.82 – 2.31 33.82 – 2.31	Depositor EDS
% Data completeness (in resolution range)	97.5 (33.82-2.31) 97.4 (33.82-2.31)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 2.32Å)	Xtriage
Refinement program	BUSTER-TNT, BUSTER 2.10.0	Depositor
R, R_{free}	0.170 , 0.216 (Not available) , 0.229	Depositor DCC
R_{free} test set	2140 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	37.6	Xtriage
Anisotropy	0.476	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 54.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.046 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6850	wwPDB-VP
Average B, all atoms (Å ²)	44.0	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 41.04 % 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 2.5028e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ABA, ACT, PG4

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	0.86	0/2391	1.25	7/3251 (0.2%)
1	C	0.82	1/2411 (0.0%)	1.24	7/3275 (0.2%)
2	B	0.82	0/855	1.20	5/1159 (0.4%)
2	D	0.85	0/855	1.16	1/1159 (0.1%)
3	E	0.75	0/63	0.95	0/80
3	F	0.81	0/63	1.13	1/80 (1.2%)
All	All	0.83	1/6638 (0.0%)	1.23	21/9004 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	194	ARG	CA-C	5.30	1.57	1.52

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	75	THR	N-CA-C	6.02	120.90	113.50
1	A	100	GLY	N-CA-C	5.75	118.85	110.63
1	C	24	GLU	CB-CG-CD	5.66	122.23	112.60
1	C	90	GLY	N-CA-C	5.65	121.67	114.66
1	A	82	LEU	CA-C-N	5.53	125.97	119.94
1	A	82	LEU	C-N-CA	5.53	125.97	119.94
1	C	95	ILE	N-CA-C	-5.50	100.20	108.12
1	C	151	GLY	CA-C-N	5.50	127.92	120.38
1	C	151	GLY	C-N-CA	5.50	127.92	120.38
2	B	71	THR	CB-CA-C	5.43	117.40	110.34
1	A	137	ASP	CA-C-N	5.34	127.43	120.28
1	A	137	ASP	C-N-CA	5.34	127.43	120.28
1	C	129	ASP	CA-CB-CG	5.32	117.92	112.60
2	D	71	THR	CB-CA-C	5.29	117.21	110.34

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	53	ASP	N-CA-C	-5.22	102.56	110.28
1	A	186	LYS	N-CA-C	-5.17	100.74	109.07
1	C	222	GLU	N-CA-C	5.16	118.16	110.52
1	A	56	GLY	N-CA-C	5.13	118.41	112.10
3	F	3	PHE	CA-CB-CG	5.12	118.92	113.80
2	B	84	HIS	CA-C-N	5.08	127.79	120.38
2	B	84	HIS	C-N-CA	5.08	127.79	120.38

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	2312	0	2216	11	0
1	C	2323	0	2241	14	0
2	B	829	0	806	4	0
2	D	829	0	806	1	0
3	E	69	0	72	0	0
3	F	69	0	72	5	0
4	C	13	0	18	3	0
5	D	4	0	3	0	0
6	A	150	0	0	1	0
6	B	58	0	0	1	0
6	C	123	0	0	0	0
6	D	61	0	0	1	0
6	E	7	0	0	0	0
6	F	3	0	0	0	0
All	All	6850	0	6234	33	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 3.

All (33) 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:188:HIS:HD2	6:A:410:HOH:O	1.82	0.61
4:C:301:PG4:H32	3:F:7:ASN:ND2	2.18	0.58
4:C:301:PG4:H32	3:F:7:ASN:HD21	1.69	0.57
1:C:81:LEU:HD11	3:F:8:ILE:HD11	1.86	0.57
1:C:32:GLU:OE2	1:C:48:ARG:HD2	2.05	0.56
1:A:13:SER:HB3	1:A:78:LEU:HD13	1.88	0.56
1:C:12:VAL:HG22	1:C:94[B]:THR:HG22	1.88	0.56
2:B:39:MET:H	2:B:45:LYS:NZ	2.03	0.56
2:B:67:HIS:HD2	6:B:157:HOH:O	1.90	0.53
4:C:301:PG4:H22	3:F:7:ASN:HD21	1.74	0.53
1:C:81:LEU:HD11	3:F:8:ILE:CD1	2.39	0.53
2:B:39:MET:H	2:B:45:LYS:HZ3	1.62	0.48
2:B:37:ILE:HD12	2:B:51:MET:HE1	1.96	0.48
1:C:9:VAL:O	1:C:96:GLN:HA	2.15	0.47
1:C:218:GLN:HB2	1:C:258[B]:THR:HG23	1.97	0.47
1:C:261:VAL:HB	1:C:270:LEU:HB2	1.96	0.47
1:A:28:VAL:HG23	1:A:33:PHE:CD1	2.50	0.46
1:C:258[A]:THR:CG2	1:C:271:THR:HG23	2.45	0.46
1:A:14:ARG:NH2	1:A:21[B]:ARG:HH11	2.13	0.46
1:A:81:LEU:HD13	1:A:118:TYR:CD1	2.51	0.46
1:A:224:LEU:HD13	1:A:247[B]:VAL:HG21	1.98	0.46
1:C:13:SER:HB3	1:C:78:LEU:HD13	1.98	0.45
1:A:189:VAL:HG23	1:A:272:LEU:HD12	1.97	0.45
2:D:67:HIS:HD2	6:D:237:HOH:O	1.99	0.45
1:A:196:GLU:HB3	1:A:197:ASP:H	1.59	0.44
1:C:216:THR:OG1	1:C:260:HIS:HB2	2.18	0.44
1:C:172:LEU:O	1:C:176:ASN:HB2	2.18	0.43
1:A:9:VAL:HG22	1:A:24:GLU:HG3	2.01	0.43
1:A:97:VAL:HA	1:A:115:GLN:O	2.19	0.42
1:C:81:LEU:HD13	1:C:118:TYR:CD1	2.54	0.42
1:C:5:LEU:HB2	1:C:168:LEU:HD13	2.01	0.42
1:C:20:PRO:HG2	1:C:75:ARG:HD3	2.01	0.42
1:A:101:CYS:HB2	1:A:109:LEU:CD1	2.52	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	284/278 (102%)	277 (98%)	6 (2%)	1 (0%)	30	37
1	C	286/278 (103%)	275 (96%)	11 (4%)	0	100	100
2	B	98/99 (99%)	97 (99%)	1 (1%)	0	100	100
2	D	98/99 (99%)	96 (98%)	2 (2%)	0	100	100
3	E	5/8 (62%)	5 (100%)	0	0	100	100
3	F	5/8 (62%)	5 (100%)	0	0	100	100
All	All	776/770 (101%)	755 (97%)	20 (3%)	1 (0%)	48	59

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	196	GLU

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	244/236 (103%)	240 (98%)	4 (2%)	55	72
1	C	246/236 (104%)	237 (96%)	9 (4%)	30	44
2	B	95/94 (101%)	92 (97%)	3 (3%)	34	50
2	D	95/94 (101%)	95 (100%)	0	100	100
3	E	6/6 (100%)	6 (100%)	0	100	100
3	F	6/6 (100%)	5 (83%)	1 (17%)	2	2
All	All	692/672 (103%)	675 (98%)	17 (2%)	43	59

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	ARG
1	A	39	ASP
1	A	45	TYR
1	A	103	VAL
2	B	1	ILE
2	B	57	SER
2	B	77	THR
1	C	23	MET
1	C	45	TYR
1	C	113	TYR
1	C	165	VAL
1	C	184	SER
1	C	227	ASP
1	C	271	THR
1	C	272[A]	LEU
1	C	272[B]	LEU
3	F	8	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	96	GLN
1	A	188	HIS
2	B	13	HIS
2	B	29	GLN
1	C	54	GLN
1	C	255	GLN
2	D	67	HIS
3	F	7	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 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	$\# Z > 2$
3	ABA	F	2	3	4,5,6	1.15	0	1,5,7	1.41	0
3	ABA	E	2	3	4,5,6	0.84	0	1,5,7	0.64	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
3	ABA	F	2	3	-	0/3/4/6	-
3	ABA	E	2	3	-	0/3/4/6	-

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.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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
4	PG4	C	301	-	12,12,12	0.22	0	11,11,11	0.15	0
5	ACT	D	101	-	3,3,3	1.17	0	3,3,3	0.98	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
4	PG4	C	301	-	-	5/10/10/10	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	301	PG4	O2-C3-C4-O3
4	C	301	PG4	C4-C3-O2-C2
4	C	301	PG4	C8-C7-O4-C6
4	C	301	PG4	C6-C5-O3-C4
4	C	301	PG4	C1-C2-O2-C3

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	301	PG4	3	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	278/278 (100%)	0.06	7 (2%) 58 61	15, 40, 65, 84	31 (11%)
1	C	278/278 (100%)	0.07	5 (1%) 67 70	16, 41, 64, 91	34 (12%)
2	B	99/99 (100%)	-0.00	2 (2%) 65 67	20, 42, 60, 74	13 (13%)
2	D	99/99 (100%)	-0.12	1 (1%) 79 81	18, 41, 58, 75	12 (12%)
3	E	7/8 (87%)	-0.07	0 100 100	32, 33, 56, 61	0
3	F	7/8 (87%)	0.03	0 100 100	33, 39, 55, 59	1 (14%)
All	All	768/770 (99%)	0.03	15 (1%) 65 67	15, 41, 64, 91	91 (11%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	225	ILE	6.0
2	B	1	ILE	5.6
1	C	277	PRO	5.0
2	D	1	ILE	4.5
1	C	278	PRO	4.2
1	A	41	GLU	3.4
1	A	146	LYS	3.3
1	C	41	GLU	2.8
1	A	278	PRO	2.8
2	B	44	LYS	2.7
1	A	111	ARG	2.3
1	A	225	ILE	2.3
1	A	227	ASP	2.1
1	C	42	ASN	2.1
1	A	16	GLY	2.1

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($\text{\AA}^2$)	Q<0.9
3	ABA	F	2	6/7	0.95	0.07	27,30,31,32	0
3	ABA	E	2	6/7	0.98	0.04	19,25,27,29	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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, 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($\text{\AA}^2$)	Q<0.9
4	PG4	C	301	13/13	0.67	0.16	89,97,101,101	0
5	ACT	D	101	4/4	0.76	0.21	71,73,74,75	0

6.5 Other polymers [i](#)

There are no such residues in this entry.