



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 09:13 AM EDT

PDB ID : 6T5A
Title : Crystal structure of herpes simplex virus 1 pUL7:pUL51 complex
Authors : Butt, B.G.; Graham, S.C.
Deposited on : 2019-10-15
Resolution : 1.83 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 : 2.10.1
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
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) : 2.10.1

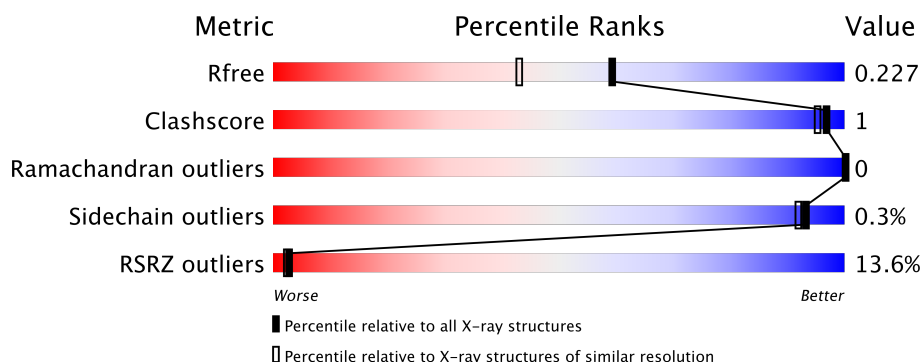
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.83 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}	111664	3313 (1.86-1.82)
Clashscore	122126	3530 (1.86-1.82)
Ramachandran outliers	120053	3495 (1.86-1.82)
Sidechain outliers	120020	3496 (1.86-1.82)
RSRZ outliers	108989	3265 (1.86-1.82)

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	136	<div> <div>4%</div> <div>69%</div> <div>29%</div> </div>
1	B	136	<div> <div>6%</div> <div>70%</div> <div>29%</div> </div>
1	C	136	<div> <div>6%</div> <div>70%</div> <div>29%</div> </div>
1	D	136	<div> <div>7%</div> <div>70%</div> <div>29%</div> </div>
2	E	304	<div> <div>16%</div> <div>87%</div> <div>9%</div> </div>

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Mol	Chain	Length	Quality of chain
2	F	304	<div><div></div><div>16%</div><div>88%</div><div>9%</div></div>
2	G	304	<div><div></div><div>11%</div><div>88%</div><div>9%</div></div>
2	H	304	<div><div></div><div>13%</div><div>87%</div><div>9%</div></div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12669 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 Tegument protein UL51.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	96	Total	C	N	O	S	0	1	0
			738	457	138	136	7			
1	B	96	Total	C	N	O	S	0	1	0
			738	457	138	136	7			
1	C	96	Total	C	N	O	S	0	2	0
			742	461	138	136	7			
1	D	96	Total	C	N	O	S	0	1	0
			738	457	138	136	7			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	7	MET	-	initiating methionine	UNP D3YPL0
A	9	SER	CYS	engineered mutation	UNP D3YPL0
B	7	MET	-	initiating methionine	UNP D3YPL0
B	9	SER	CYS	engineered mutation	UNP D3YPL0
C	7	MET	-	initiating methionine	UNP D3YPL0
C	9	SER	CYS	engineered mutation	UNP D3YPL0
D	7	MET	-	initiating methionine	UNP D3YPL0
D	9	SER	CYS	engineered mutation	UNP D3YPL0

- Molecule 2 is a protein called Cytoplasmic envelopment protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	276	Total	C	N	O	S	0	7	0
			2242	1441	388	396	17			
2	F	276	Total	C	N	O	S	0	7	0
			2242	1441	388	396	17			
2	G	276	Total	C	N	O	S	0	7	0
			2242	1441	388	396	17			
2	H	276	Total	C	N	O	S	0	7	0
			2242	1441	388	396	17			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	297	GLY	-	expression tag	UNP A0A110B4Q7
E	298	SER	-	expression tag	UNP A0A110B4Q7
E	299	LEU	-	expression tag	UNP A0A110B4Q7
E	300	GLU	-	expression tag	UNP A0A110B4Q7
E	301	VAL	-	expression tag	UNP A0A110B4Q7
E	302	LEU	-	expression tag	UNP A0A110B4Q7
E	303	PHE	-	expression tag	UNP A0A110B4Q7
E	304	GLN	-	expression tag	UNP A0A110B4Q7
F	297	GLY	-	expression tag	UNP A0A110B4Q7
F	298	SER	-	expression tag	UNP A0A110B4Q7
F	299	LEU	-	expression tag	UNP A0A110B4Q7
F	300	GLU	-	expression tag	UNP A0A110B4Q7
F	301	VAL	-	expression tag	UNP A0A110B4Q7
F	302	LEU	-	expression tag	UNP A0A110B4Q7
F	303	PHE	-	expression tag	UNP A0A110B4Q7
F	304	GLN	-	expression tag	UNP A0A110B4Q7
G	297	GLY	-	expression tag	UNP A0A110B4Q7
G	298	SER	-	expression tag	UNP A0A110B4Q7
G	299	LEU	-	expression tag	UNP A0A110B4Q7
G	300	GLU	-	expression tag	UNP A0A110B4Q7
G	301	VAL	-	expression tag	UNP A0A110B4Q7
G	302	LEU	-	expression tag	UNP A0A110B4Q7
G	303	PHE	-	expression tag	UNP A0A110B4Q7
G	304	GLN	-	expression tag	UNP A0A110B4Q7
H	297	GLY	-	expression tag	UNP A0A110B4Q7
H	298	SER	-	expression tag	UNP A0A110B4Q7
H	299	LEU	-	expression tag	UNP A0A110B4Q7
H	300	GLU	-	expression tag	UNP A0A110B4Q7
H	301	VAL	-	expression tag	UNP A0A110B4Q7
H	302	LEU	-	expression tag	UNP A0A110B4Q7
H	303	PHE	-	expression tag	UNP A0A110B4Q7
H	304	GLN	-	expression tag	UNP A0A110B4Q7

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	1	Total	Cl	0	0
			1	1		
4	G	1	Total	Cl	0	0
			1	1		
4	F	1	Total	Cl	0	0
			1	1		
4	E	1	Total	Cl	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	56	Total	O	0	0
			56	56		

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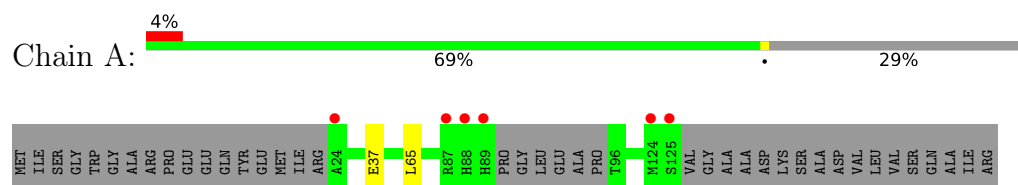
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	48	Total 48	O 48	0	0
5	C	49	Total 49	O 49	0	0
5	D	56	Total 56	O 56	0	0
5	E	126	Total 126	O 126	0	0
5	F	111	Total 111	O 111	0	0
5	G	139	Total 139	O 139	0	0
5	H	132	Total 132	O 132	0	0

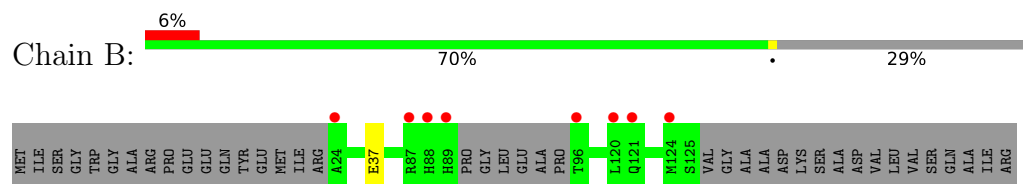
3 Residue-property plots [i](#)

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.

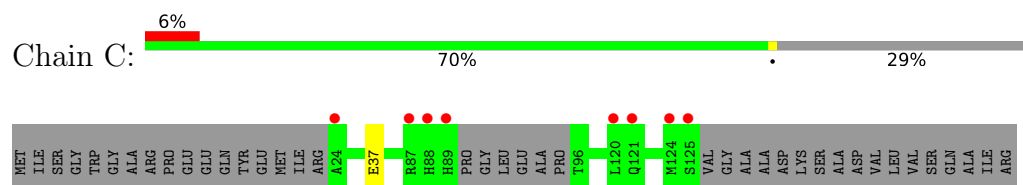
- Molecule 1: Tegument protein UL51



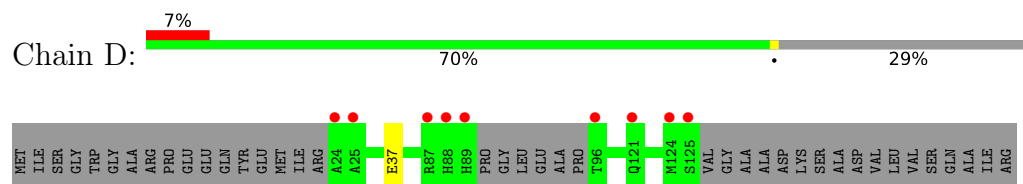
- Molecule 1: Tegument protein UL51



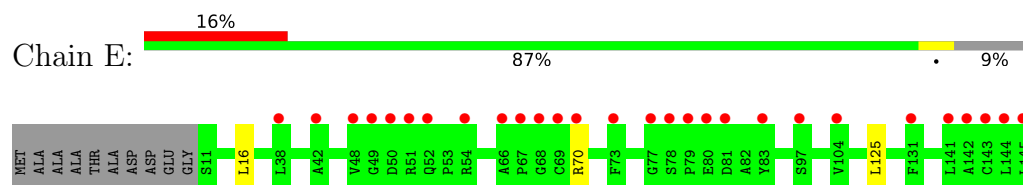
- Molecule 1: Tegument protein UL51

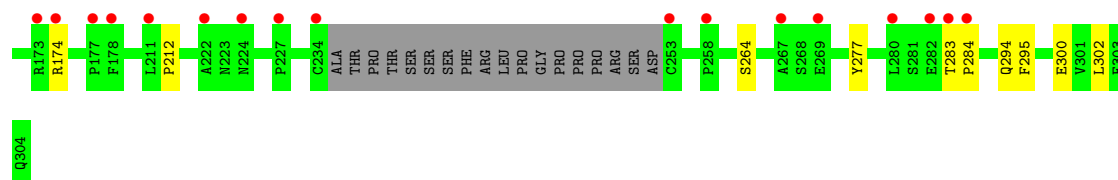


- Molecule 1: Tegument protein UL51

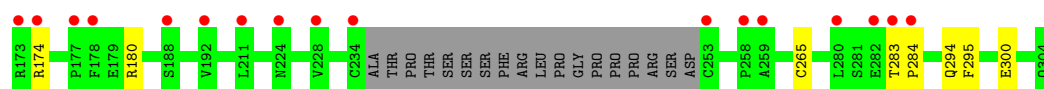
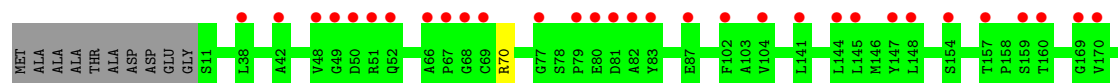
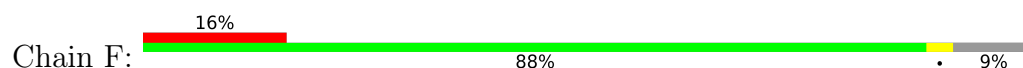


- Molecule 2: Cytoplasmic envelopment protein 1

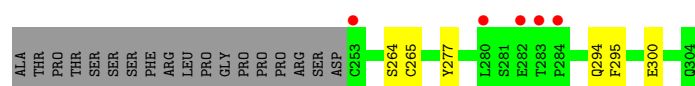
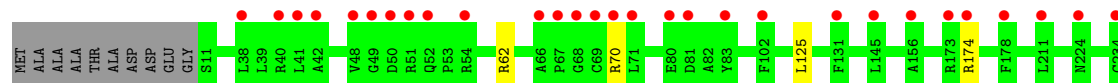
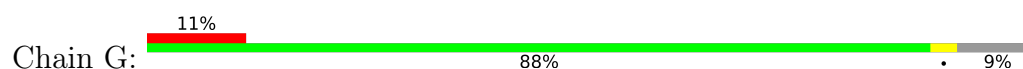




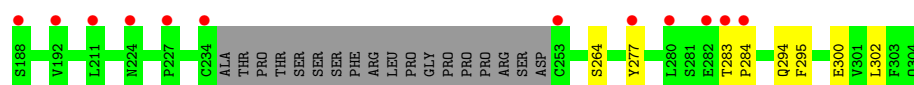
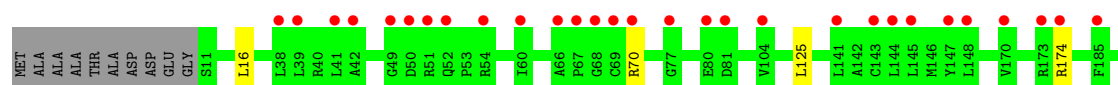
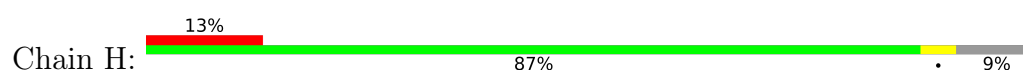
• Molecule 2: Cytoplasmic envelopment protein 1



• Molecule 2: Cytoplasmic envelopment protein 1



• Molecule 2: Cytoplasmic envelopment protein 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	79.51Å 106.30Å 106.01Å 90.00° 92.05° 90.00°	Depositor
Resolution (Å)	26.57 – 1.83 26.58 – 1.83	Depositor EDS
% Data completeness (in resolution range)	100.0 (26.57-1.83) 100.0 (26.58-1.83)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 1.83Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.194 , 0.220 0.199 , 0.227	Depositor DCC
R_{free} test set	8093 reflections (5.23%)	wwPDB-VP
Wilson B-factor (Å ²)	33.1	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 48.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.247 for -h,-l,-k 0.032 for -h,l,k 0.086 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12669	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.24% 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: GOL, CL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.45	0/750	0.54	0/1017
1	B	0.45	0/750	0.53	0/1017
1	C	0.47	0/757	0.53	0/1027
1	D	0.46	0/750	0.54	0/1017
2	E	0.45	0/2306	0.57	0/3139
2	F	0.45	0/2306	0.57	0/3139
2	G	0.46	0/2306	0.57	0/3139
2	H	0.47	0/2306	0.57	0/3139
All	All	0.46	0/12231	0.56	0/16634

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	738	0	756	2	0
1	B	738	0	756	1	0
1	C	742	0	765	1	0
1	D	738	0	756	1	0
2	E	2242	0	2219	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	2242	0	2219	5	0
2	G	2242	0	2219	6	0
2	H	2242	0	2219	7	0
3	A	6	0	8	0	0
3	B	6	0	8	0	0
3	C	6	0	8	0	0
3	D	6	0	8	0	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
4	G	1	0	0	0	0
4	H	1	0	0	0	0
5	A	56	0	0	0	0
5	B	48	0	0	0	0
5	C	49	0	0	0	0
5	D	56	0	0	0	0
5	E	126	0	0	0	0
5	F	111	0	0	1	0
5	G	139	0	0	2	0
5	H	132	0	0	0	0
All	All	12669	0	11941	24	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 1.

All (24) 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)
2:F:283[B]:THR:HB	2:F:284[B]:PRO:HD3	1.67	0.76
2:E:70:ARG:HH22	2:E:174:ARG:HH12	1.42	0.68
2:G:70:ARG:HH22	2:G:174:ARG:HH12	1.42	0.68
2:H:70:ARG:HH22	2:H:174:ARG:HH12	1.42	0.67
2:F:70:ARG:HH22	2:F:174:ARG:HH12	1.41	0.65
2:H:283[A]:THR:HB	2:H:284[A]:PRO:HD3	1.80	0.62
1:C:37:GLU:HG2	2:G:300:GLU:HG2	1.84	0.60
2:H:125:LEU:HD11	2:H:277:TYR:HD2	1.67	0.59
1:D:37:GLU:HG2	2:H:300:GLU:HG2	1.82	0.59
2:G:125:LEU:HD11	2:G:277:TYR:HD2	1.67	0.59
1:B:37:GLU:HG2	2:F:300:GLU:HG2	1.85	0.58
1:A:37:GLU:HG2	2:E:300:GLU:HG2	1.90	0.54
2:F:265:CYS:HB3	5:F:508:HOH:O	2.09	0.52
2:E:283[B]:THR:HB	2:E:284[B]:PRO:HD3	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:294:GLN:HG2	2:H:295:PHE:N	2.30	0.46
2:E:294:GLN:HG2	2:E:295:PHE:N	2.31	0.46
2:G:294:GLN:HG2	2:G:295:PHE:N	2.30	0.45
2:F:294:GLN:HG2	2:F:295:PHE:N	2.30	0.45
2:E:125:LEU:HD11	2:E:277:TYR:HD2	1.83	0.43
1:A:65:LEU:HD13	2:E:212:PRO:HB2	2.00	0.42
2:G:62:ARG:HG3	5:G:503:HOH:O	2.19	0.41
2:E:16:LEU:HB2	2:H:302:LEU:HD21	2.02	0.41
2:G:265:CYS:HB3	5:G:527:HOH:O	2.21	0.40
2:E:302:LEU:HD21	2:H:16:LEU:HB2	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	93/136 (68%)	92 (99%)	1 (1%)	0	100	100
1	B	93/136 (68%)	92 (99%)	1 (1%)	0	100	100
1	C	94/136 (69%)	93 (99%)	1 (1%)	0	100	100
1	D	93/136 (68%)	92 (99%)	1 (1%)	0	100	100
2	E	279/304 (92%)	272 (98%)	7 (2%)	0	100	100
2	F	279/304 (92%)	270 (97%)	9 (3%)	0	100	100
2	G	279/304 (92%)	274 (98%)	5 (2%)	0	100	100
2	H	279/304 (92%)	276 (99%)	3 (1%)	0	100	100
All	All	1489/1760 (85%)	1461 (98%)	28 (2%)	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	78/107 (73%)	78 (100%)	0	100	100
1	B	78/107 (73%)	78 (100%)	0	100	100
1	C	79/107 (74%)	79 (100%)	0	100	100
1	D	78/107 (73%)	78 (100%)	0	100	100
2	E	244/258 (95%)	243 (100%)	1 (0%)	92	90
2	F	244/258 (95%)	243 (100%)	1 (0%)	92	90
2	G	244/258 (95%)	243 (100%)	1 (0%)	92	90
2	H	244/258 (95%)	243 (100%)	1 (0%)	92	90
All	All	1289/1460 (88%)	1285 (100%)	4 (0%)	93	92

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	E	264	SER
2	F	180	ARG
2	G	264	SER
2	H	264	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
2	E	224	ASN
2	F	224	ASN
2	G	224	ASN
2	H	224	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	$\# Z > 2$
3	GOL	B	201	-	5,5,5	0.05	0	5,5,5	0.13	0
3	GOL	A	201	-	5,5,5	0.08	0	5,5,5	0.16	0
3	GOL	D	201	-	5,5,5	0.05	0	5,5,5	0.12	0
3	GOL	C	201	-	5,5,5	0.06	0	5,5,5	0.13	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	GOL	B	201	-	-	0/4/4/4	-
3	GOL	A	201	-	-	2/4/4/4	-
3	GOL	D	201	-	-	2/4/4/4	-
3	GOL	C	201	-	-	1/4/4/4	-

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
3	A	201	GOL	C1-C2-C3-O3
3	D	201	GOL	C1-C2-C3-O3
3	D	201	GOL	O2-C2-C3-O3
3	C	201	GOL	C1-C2-C3-O3
3	A	201	GOL	O2-C2-C3-O3

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, 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	96/136 (70%)	0.39	6 (6%)	20 17	26, 37, 65, 108	0
1	B	96/136 (70%)	0.37	8 (8%)	11 10	26, 37, 67, 109	0
1	C	96/136 (70%)	0.37	8 (8%)	11 10	25, 36, 66, 111	0
1	D	96/136 (70%)	0.43	9 (9%)	8 7	26, 37, 68, 107	0
2	E	276/304 (90%)	0.81	48 (17%)	1 1	25, 43, 77, 97	0
2	F	276/304 (90%)	0.90	48 (17%)	1 1	27, 45, 85, 112	0
2	G	276/304 (90%)	0.63	34 (12%)	4 3	26, 39, 71, 89	0
2	H	276/304 (90%)	0.63	41 (14%)	2 2	26, 40, 71, 93	0
All	All	1488/1760 (84%)	0.65	202 (13%)	3 2	25, 40, 75, 112	0

All (202) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	124	MET	7.8
2	E	67	PRO	7.7
2	G	67	PRO	7.5
2	F	67	PRO	7.5
1	D	24	ALA	7.4
1	D	124	MET	7.2
2	F	81	ASP	6.9
2	G	50	ASP	6.7
1	B	124	MET	6.7
1	A	124	MET	6.7
2	F	69	CYS	6.4
1	B	24	ALA	6.2
2	H	50	ASP	6.1
2	F	68	GLY	6.1
2	H	67	PRO	6.0
2	F	284[A]	PRO	6.0

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Mol	Chain	Res	Type	RSRZ
2	G	253	CYS	6.0
2	F	80	GLU	5.9
1	C	24	ALA	5.7
2	E	69	CYS	5.7
1	D	88	HIS	5.6
2	F	50	ASP	5.6
2	E	81	ASP	5.6
2	F	79	PRO	5.5
1	A	88	HIS	5.4
2	E	284[A]	PRO	5.3
2	E	68	GLY	5.2
2	F	224	ASN	5.2
1	D	89	HIS	5.1
2	H	283[A]	THR	5.1
1	A	24	ALA	5.0
1	C	88	HIS	4.8
2	G	68	GLY	4.8
2	G	283[A]	THR	4.8
2	F	253	CYS	4.8
2	H	224	ASN	4.7
1	B	89	HIS	4.7
2	E	224	ASN	4.6
1	A	125	SER	4.6
2	G	224	ASN	4.5
2	F	283[A]	THR	4.5
1	B	87	ARG	4.4
2	E	283[A]	THR	4.4
2	H	145	LEU	4.4
2	H	284[A]	PRO	4.3
2	F	52	GLN	4.3
2	G	52	GLN	4.3
2	E	66	ALA	4.3
2	E	50	ASP	4.2
2	F	66	ALA	4.2
1	C	89	HIS	4.1
2	E	253	CYS	4.1
2	G	48	VAL	4.1
2	E	145	LEU	4.1
1	D	87	ARG	4.0
2	F	77	GLY	4.0
1	A	87	ARG	4.0
2	G	80	GLU	4.0

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Mol	Chain	Res	Type	RSRZ
1	B	88	HIS	3.9
2	E	173	ARG	3.9
2	E	222	ALA	3.9
2	F	48	VAL	3.8
2	F	174	ARG	3.8
2	E	49	GLY	3.8
2	F	49	GLY	3.8
2	E	51	ARG	3.8
2	E	80	GLU	3.8
2	F	42	ALA	3.7
2	G	173	ARG	3.7
2	F	51	ARG	3.7
2	G	145	LEU	3.7
2	F	145	LEU	3.7
2	E	178	PHE	3.6
2	F	178	PHE	3.6
2	G	42	ALA	3.6
1	C	125	SER	3.6
2	H	280[A]	LEU	3.6
2	H	69	CYS	3.6
2	G	51	ARG	3.6
2	F	282[A]	GLU	3.6
2	H	80	GLU	3.6
2	G	280[A]	LEU	3.6
1	C	87	ARG	3.5
2	E	144	LEU	3.5
2	H	144	LEU	3.5
2	F	160	THR	3.4
2	E	282[A]	GLU	3.4
2	E	227	PRO	3.4
2	F	82	ALA	3.4
2	E	83	TYR	3.4
2	H	68	GLY	3.4
2	F	234	CYS	3.4
2	G	81	ASP	3.3
2	H	253	CYS	3.3
2	H	173	ARG	3.3
2	F	173	ARG	3.3
2	G	284[A]	PRO	3.2
1	D	125	SER	3.2
2	G	83	TYR	3.2
2	E	234	CYS	3.2

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Mol	Chain	Res	Type	RSRZ
2	G	234	CYS	3.1
1	A	89	HIS	3.1
2	H	38	LEU	3.1
2	G	66	ALA	3.1
2	H	81	ASP	3.0
2	E	141	LEU	2.9
2	E	52	GLN	2.9
2	H	51	ARG	2.9
2	E	48	VAL	2.9
1	D	25	ALA	2.9
2	G	174	ARG	2.9
2	F	177	PRO	2.9
2	G	38	LEU	2.9
2	H	39	LEU	2.8
2	F	38	LEU	2.8
1	B	96	THR	2.8
2	H	52	GLN	2.8
2	E	143	CYS	2.8
2	F	280[A]	LEU	2.8
2	H	143	CYS	2.8
2	F	170	VAL	2.8
2	F	87	GLU	2.8
2	H	42	ALA	2.7
2	H	148	LEU	2.7
2	H	174	ARG	2.7
2	G	49	GLY	2.7
2	F	141	LEU	2.7
2	H	192	VAL	2.7
2	E	104	VAL	2.7
2	F	192	VAL	2.7
2	E	280[A]	LEU	2.7
2	F	258	PRO	2.6
2	H	66	ALA	2.6
2	E	269	GLU	2.6
2	H	104	VAL	2.6
2	H	277	TYR	2.6
2	E	79	PRO	2.6
2	F	157	THR	2.6
2	F	104	VAL	2.6
2	E	77	GLY	2.5
2	G	54	ARG	2.5
2	F	154	SER	2.5

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Mol	Chain	Res	Type	RSRZ
2	E	78	SER	2.5
2	E	70	ARG	2.5
2	G	71	LEU	2.5
2	F	159	SER	2.5
2	F	83	TYR	2.4
2	F	147	TYR	2.4
2	E	142	ALA	2.4
2	E	54	ARG	2.4
2	F	102	PHE	2.4
2	G	131	PHE	2.4
2	G	69	CYS	2.4
2	H	41	LEU	2.3
2	E	42	ALA	2.3
2	H	185	PHE	2.3
2	E	148	LEU	2.3
2	F	169	GLY	2.3
1	D	96	THR	2.3
2	G	178	PHE	2.3
2	E	174	ARG	2.3
2	E	267	ALA	2.3
2	F	148	LEU	2.3
2	G	282[A]	GLU	2.2
1	D	121	GLN	2.2
1	B	120	LEU	2.2
2	E	211	LEU	2.2
2	F	144	LEU	2.2
2	H	141	LEU	2.2
2	H	211	LEU	2.2
2	F	188	SER	2.2
2	G	70	ARG	2.2
2	H	227	PRO	2.2
2	G	211	LEU	2.2
2	F	259	ALA	2.2
2	H	147	TYR	2.2
2	G	102	PHE	2.2
2	G	156	ALA	2.1
2	F	211	LEU	2.1
2	H	282[A]	GLU	2.1
2	H	188	SER	2.1
1	C	120	LEU	2.1
2	E	97	SER	2.1
2	E	258	PRO	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	121	GLN	2.1
2	H	170	VAL	2.1
2	H	54	ARG	2.1
2	H	234	CYS	2.1
2	E	147	TYR	2.1
2	H	60	ILE	2.1
2	H	77	GLY	2.1
2	E	73	PHE	2.0
2	E	131	PHE	2.0
2	F	228	VAL	2.0
2	E	38	LEU	2.0
2	G	41	LEU	2.0
1	B	121	GLN	2.0
2	H	70	ARG	2.0
2	E	177	PRO	2.0
2	H	49	GLY	2.0
2	E	160	THR	2.0
2	G	40	ARG	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. 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
3	GOL	A	201	6/6	0.78	0.14	66,67,71,71	0
3	GOL	C	201	6/6	0.87	0.18	63,64,68,69	0
3	GOL	D	201	6/6	0.88	0.12	66,67,72,73	0
3	GOL	B	201	6/6	0.88	0.14	67,69,72,73	0
4	CL	E	401	1/1	0.94	0.04	49,49,49,49	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	CL	H	401	1/1	0.97	0.04	45,45,45,45	0
4	CL	G	401	1/1	0.97	0.07	47,47,47,47	0
4	CL	F	401	1/1	0.97	0.04	47,47,47,47	0

6.5 Other polymers [i](#)

There are no such residues in this entry.