



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 1, 2026 – 09:44 AM UTC

PDB ID : 1L8I / pdb_0000118i
Title : Dna Protection and Binding by E. Coli DPS Protein
Authors : Luo, J.; Liu, D.; White, M.A.; Fox, R.O.
Deposited on : 2002-03-20
Resolution : 3.00 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

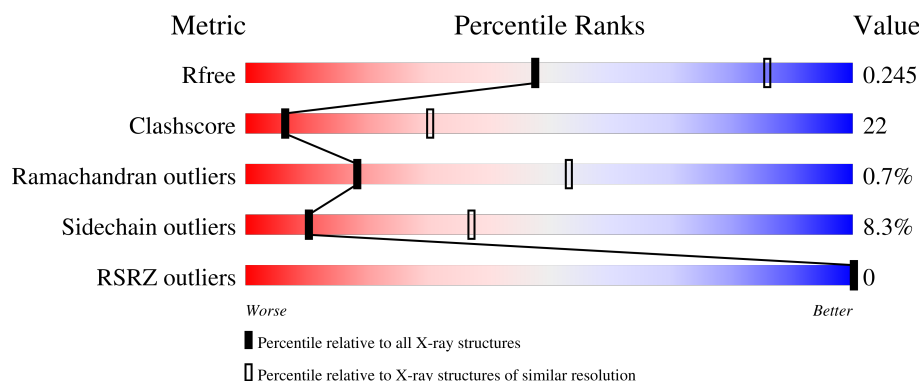
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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 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	167	51% 40% • 7%
1	B	167	53% 35% 5% 7%
1	C	167	56% 34% • 7%
1	D	167	54% 35% • 7%
1	E	167	54% 35% • 7%

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Mol	Chain	Length	Quality of chain
1	F	167	 58%31%5%6%
1	G	167	 56%34%•7%
1	H	167	 58%32%•7%
1	I	167	 54%35%•7%
1	J	167	 54%36%•6%
1	K	167	 52%37%••7%
1	L	167	 56%33%•7%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14968 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 DNA PROTECTION DURING STARVATION PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	156	Total	C	N	O	S	0	0	0
			1231	774	216	237	4			
1	B	156	Total	C	N	O	S	0	0	0
			1231	774	216	237	4			
1	C	155	Total	C	N	O	S	0	0	0
			1224	770	215	235	4			
1	D	156	Total	C	N	O	S	0	0	0
			1231	774	216	237	4			
1	E	156	Total	C	N	O	S	0	0	0
			1231	774	216	237	4			
1	F	157	Total	C	N	O	S	0	0	0
			1236	777	217	238	4			
1	G	155	Total	C	N	O	S	0	0	0
			1224	770	215	235	4			
1	H	155	Total	C	N	O	S	0	0	0
			1224	770	215	235	4			
1	I	156	Total	C	N	O	S	0	0	0
			1231	774	216	237	4			
1	J	157	Total	C	N	O	S	0	0	0
			1236	777	217	238	4			
1	K	156	Total	C	N	O	S	0	0	0
			1231	774	216	237	4			
1	L	156	Total	C	N	O	S	0	0	0
			1231	774	216	237	4			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	75	CYS	ASP	engineered mutation	UNP P0ABT2
A	78	ALA	ASP	engineered mutation	UNP P0ABT2
B	75	CYS	ASP	engineered mutation	UNP P0ABT2
B	78	ALA	ASP	engineered mutation	UNP P0ABT2
C	75	CYS	ASP	engineered mutation	UNP P0ABT2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	78	ALA	ASP	engineered mutation	UNP P0ABT2
D	75	CYS	ASP	engineered mutation	UNP P0ABT2
D	78	ALA	ASP	engineered mutation	UNP P0ABT2
E	75	CYS	ASP	engineered mutation	UNP P0ABT2
E	78	ALA	ASP	engineered mutation	UNP P0ABT2
F	75	CYS	ASP	engineered mutation	UNP P0ABT2
F	78	ALA	ASP	engineered mutation	UNP P0ABT2
G	75	CYS	ASP	engineered mutation	UNP P0ABT2
G	78	ALA	ASP	engineered mutation	UNP P0ABT2
H	75	CYS	ASP	engineered mutation	UNP P0ABT2
H	78	ALA	ASP	engineered mutation	UNP P0ABT2
I	75	CYS	ASP	engineered mutation	UNP P0ABT2
I	78	ALA	ASP	engineered mutation	UNP P0ABT2
J	75	CYS	ASP	engineered mutation	UNP P0ABT2
J	78	ALA	ASP	engineered mutation	UNP P0ABT2
K	75	CYS	ASP	engineered mutation	UNP P0ABT2
K	78	ALA	ASP	engineered mutation	UNP P0ABT2
L	75	CYS	ASP	engineered mutation	UNP P0ABT2
L	78	ALA	ASP	engineered mutation	UNP P0ABT2

- Molecule 2 is POTASSIUM ION (CCD ID: K) (formula: K).

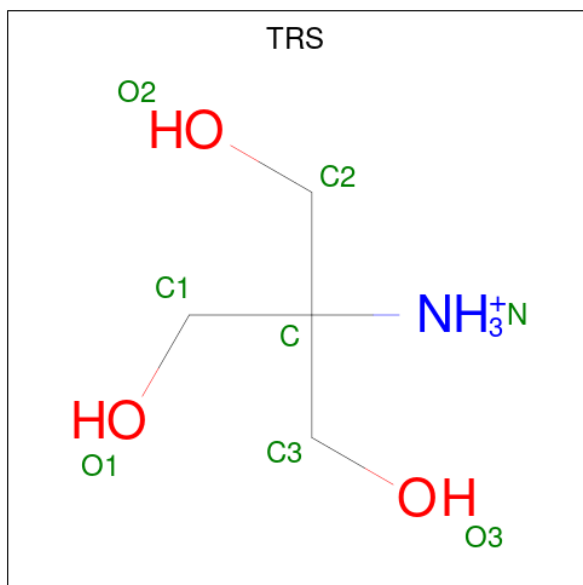
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total K 1 1	0	0
2	B	1	Total K 1 1	0	0
2	C	1	Total K 1 1	0	0
2	D	1	Total K 1 1	0	0
2	E	2	Total K 2 2	0	0
2	F	1	Total K 1 1	0	0
2	G	1	Total K 1 1	0	0
2	H	1	Total K 1 1	0	0
2	I	1	Total K 1 1	0	0
2	J	1	Total K 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	L	1	Total	K	0	0
			1	1		

- Molecule 3 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: C₄H₁₂NO₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			8	4	1	3		
3	A	1	Total	C	N	O	0	0
			8	4	1	3		
3	A	1	Total	C	N	O	0	0
			8	4	1	3		
3	B	1	Total	C	N	O	0	0
			8	4	1	3		
3	B	1	Total	C	N	O	0	0
			8	4	1	3		
3	C	1	Total	C	N	O	0	0
			8	4	1	3		
3	C	1	Total	C	N	O	0	0
			8	4	1	3		
3	C	1	Total	C	N	O	0	0
			8	4	1	3		
3	D	1	Total	C	N	O	0	0
			8	4	1	3		
3	D	1	Total	C	N	O	0	0
			8	4	1	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	G	1	Total	C	N	O	0	0
			8	4	1	3		
3	L	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	8	Total	O	0	0
			8	8		
4	B	10	Total	O	0	0
			10	10		
4	C	6	Total	O	0	0
			6	6		
4	D	10	Total	O	0	0
			10	10		
4	E	8	Total	O	0	0
			8	8		
4	F	3	Total	O	0	0
			3	3		
4	G	14	Total	O	0	0
			14	14		
4	H	10	Total	O	0	0
			10	10		
4	I	6	Total	O	0	0
			6	6		
4	J	9	Total	O	0	0
			9	9		
4	K	7	Total	O	0	0
			7	7		
4	L	8	Total	O	0	0
			8	8		

- Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN





• Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN

Chain E: 54% 35% 7%



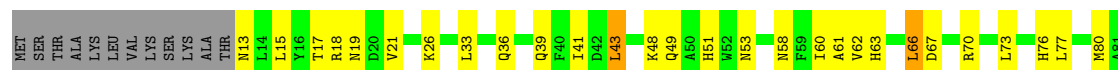
• Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN

Chain F: 58% 31% 5% 6%



• Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN

Chain G: 56% 34% 7%



• Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN

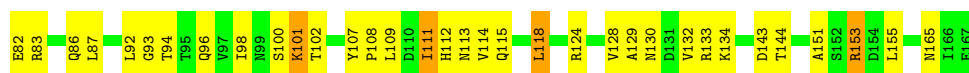
Chain H: 58% 32% 7%



• Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN

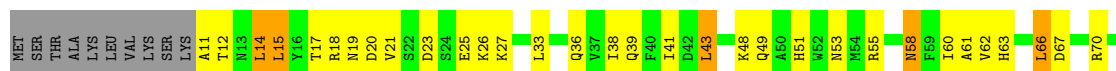
Chain I: 54% 35% 7%





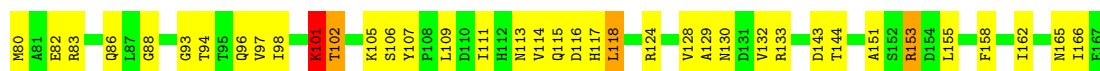
• Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN

Chain J: 54% 36% 6%



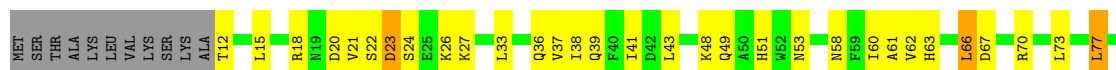
• Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN

Chain K: 52% 37% 7%



• Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN

Chain L: 56% 33% 7%



4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	90.54Å 90.54Å 226.08Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.86 – 3.00 29.86 – 3.00	Depositor EDS
% Data completeness (in resolution range)	98.0 (29.86-3.00) 98.0 (29.86-3.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.25 (at 3.00Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.202 , 0.250 0.198 , 0.245	Depositor DCC
R_{free} test set	2074 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	61.8	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 63.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.016 for -h,-k,l 0.045 for h,-h-k,-l 0.029 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14968	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: TRS, K

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	0.55	0/1249	0.98	5/1691 (0.3%)
1	B	0.57	0/1249	0.91	2/1691 (0.1%)
1	C	0.52	0/1242	0.95	5/1681 (0.3%)
1	D	0.51	0/1249	0.94	6/1691 (0.4%)
1	E	0.49	0/1249	0.93	4/1691 (0.2%)
1	F	0.49	0/1254	0.90	3/1698 (0.2%)
1	G	0.57	0/1242	0.92	1/1681 (0.1%)
1	H	0.53	0/1242	0.93	4/1681 (0.2%)
1	I	0.49	0/1249	0.91	2/1691 (0.1%)
1	J	0.48	0/1254	0.89	2/1698 (0.1%)
1	K	0.51	0/1249	0.91	2/1691 (0.1%)
1	L	0.49	0/1249	0.91	2/1691 (0.1%)
All	All	0.52	0/14977	0.92	38/20276 (0.2%)

There are no bond length outliers.

The worst 5 of 38 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	53	ASN	N-CA-C	8.55	122.50	112.93
1	G	53	ASN	N-CA-C	8.20	122.12	112.93
1	K	53	ASN	N-CA-C	7.84	121.71	112.93
1	I	53	ASN	N-CA-C	7.80	121.67	112.93
1	C	70	ARG	NE-CZ-NH2	7.74	126.16	119.20

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	1231	0	1234	73	0
1	B	1231	0	1234	67	0
1	C	1224	0	1227	68	0
1	D	1231	0	1234	67	0
1	E	1231	0	1234	66	0
1	F	1236	0	1239	57	0
1	G	1224	0	1227	54	0
1	H	1224	0	1227	49	0
1	I	1231	0	1234	57	0
1	J	1236	0	1239	67	0
1	K	1231	0	1234	64	0
1	L	1231	0	1234	59	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	2	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
2	L	1	0	0	0	0
3	A	24	0	36	3	0
3	B	16	0	24	0	0
3	C	24	0	36	3	0
3	D	16	0	24	2	0
3	G	8	0	12	0	0
3	L	8	0	12	4	0
4	A	8	0	0	0	0
4	B	10	0	0	1	0
4	C	6	0	0	0	0
4	D	10	0	0	0	0
4	E	8	0	0	0	0
4	F	3	0	0	0	0
4	G	14	0	0	1	0
4	H	10	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	I	6	0	0	0	0
4	J	9	0	0	1	0
4	K	7	0	0	0	0
4	L	8	0	0	0	0
All	All	14968	0	14941	650	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 22.

The worst 5 of 650 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:C:130:ASN:ND2	1:C:133:ARG:HH22	1.64	0.95
1:L:108:PRO:HG2	1:L:111:ILE:HD11	1.49	0.95
1:J:130:ASN:ND2	1:J:133:ARG:HH22	1.66	0.93
1:J:130:ASN:HD22	1:J:133:ARG:HH22	1.16	0.93
1:G:49:GLN:HE21	1:H:93:GLY:H	1.17	0.91

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	154/167 (92%)	140 (91%)	12 (8%)	2 (1%)	9	38
1	B	154/167 (92%)	145 (94%)	9 (6%)	0	100	100
1	C	153/167 (92%)	141 (92%)	12 (8%)	0	100	100
1	D	154/167 (92%)	141 (92%)	11 (7%)	2 (1%)	9	38
1	E	154/167 (92%)	147 (96%)	6 (4%)	1 (1%)	21	56
1	F	155/167 (93%)	144 (93%)	9 (6%)	2 (1%)	9	38

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	G	153/167 (92%)	142 (93%)	11 (7%)	0	100	100
1	H	153/167 (92%)	145 (95%)	8 (5%)	0	100	100
1	I	154/167 (92%)	143 (93%)	9 (6%)	2 (1%)	9	38
1	J	155/167 (93%)	144 (93%)	9 (6%)	2 (1%)	9	38
1	K	154/167 (92%)	139 (90%)	13 (8%)	2 (1%)	9	38
1	L	154/167 (92%)	141 (92%)	13 (8%)	0	100	100
All	All	1847/2004 (92%)	1712 (93%)	122 (7%)	13 (1%)	18	53

5 of 13 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	12	THR
1	J	14	LEU
1	K	101	LYS
1	A	14	LEU
1	I	101	LYS

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	133/142 (94%)	123 (92%)	10 (8%)	12	41
1	B	133/142 (94%)	119 (90%)	14 (10%)	6	27
1	C	132/142 (93%)	123 (93%)	9 (7%)	14	45
1	D	133/142 (94%)	122 (92%)	11 (8%)	10	37
1	E	133/142 (94%)	123 (92%)	10 (8%)	12	41
1	F	133/142 (94%)	122 (92%)	11 (8%)	10	37
1	G	132/142 (93%)	121 (92%)	11 (8%)	10	37
1	H	132/142 (93%)	121 (92%)	11 (8%)	10	37
1	I	133/142 (94%)	122 (92%)	11 (8%)	10	37
1	J	133/142 (94%)	123 (92%)	10 (8%)	12	41

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	K	133/142 (94%)	121 (91%)	12 (9%)	9	34
1	L	133/142 (94%)	120 (90%)	13 (10%)	7	30
All	All	1593/1704 (94%)	1460 (92%)	133 (8%)	10	37

5 of 133 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	K	101	LYS
1	K	155	LEU
1	L	118	LEU
1	E	67	ASP
1	E	66	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 125 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	63	HIS
1	K	63	HIS
1	G	113	ASN
1	K	58	ASN
1	L	51	HIS

5.3.3 RNA [i](#)

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 24 ligands modelled in this entry, 12 are monoatomic - leaving 12 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	TRS	L	309	-	7,7,7	0.27	0	9,9,9	1.12	1 (11%)
3	TRS	B	310	-	7,7,7	0.39	0	9,9,9	1.09	0
3	TRS	C	302	-	7,7,7	0.56	0	9,9,9	1.07	0
3	TRS	D	311	-	7,7,7	0.56	0	9,9,9	1.06	0
3	TRS	B	308	-	7,7,7	0.40	0	9,9,9	1.13	0
3	TRS	A	306	-	7,7,7	0.22	0	9,9,9	1.09	1 (11%)
3	TRS	A	305	-	7,7,7	0.55	0	9,9,9	1.06	0
3	TRS	G	307	-	7,7,7	0.48	0	9,9,9	1.12	0
3	TRS	C	303	-	7,7,7	0.50	0	9,9,9	1.10	0
3	TRS	D	312	-	7,7,7	0.32	0	9,9,9	1.07	0
3	TRS	C	304	-	7,7,7	0.66	0	9,9,9	1.26	1 (11%)
3	TRS	A	301	-	7,7,7	0.56	0	9,9,9	1.16	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	TRS	L	309	-	-	1/9/9/9	-
3	TRS	B	310	-	-	3/9/9/9	-
3	TRS	C	302	-	-	0/9/9/9	-
3	TRS	D	311	-	-	0/9/9/9	-
3	TRS	B	308	-	-	0/9/9/9	-
3	TRS	A	306	-	-	1/9/9/9	-
3	TRS	A	305	-	-	0/9/9/9	-
3	TRS	G	307	-	-	3/9/9/9	-
3	TRS	C	303	-	-	1/9/9/9	-
3	TRS	D	312	-	-	1/9/9/9	-
3	TRS	C	304	-	-	3/9/9/9	-
3	TRS	A	301	-	-	3/9/9/9	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	L	309	TRS	C1-C-N	-2.17	102.64	108.17
3	A	306	TRS	C1-C-N	-2.07	102.90	108.17
3	C	304	TRS	C3-C-N	2.04	113.38	108.17

There are no chirality outliers.

5 of 16 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	301	TRS	N-C-C2-O2
3	B	310	TRS	N-C-C2-O2
3	C	304	TRS	N-C-C2-O2
3	G	307	TRS	N-C-C2-O2
3	A	301	TRS	C3-C-C2-O2

There are no ring outliers.

4 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	L	309	TRS	4	0
3	A	306	TRS	3	0
3	C	303	TRS	3	0
3	D	312	TRS	2	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 [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	156/167 (93%)	-0.58	0 100 100	27, 48, 65, 86	0
1	B	156/167 (93%)	-0.59	0 100 100	25, 48, 69, 85	0
1	C	155/167 (92%)	-0.53	0 100 100	36, 54, 76, 88	0
1	D	156/167 (93%)	-0.44	0 100 100	36, 57, 70, 80	0
1	E	156/167 (93%)	-0.55	0 100 100	31, 50, 78, 94	0
1	F	157/167 (94%)	-0.39	0 100 100	39, 59, 76, 100	0
1	G	155/167 (92%)	-0.56	0 100 100	26, 48, 74, 86	0
1	H	155/167 (92%)	-0.52	0 100 100	33, 52, 68, 76	0
1	I	156/167 (93%)	-0.44	0 100 100	34, 59, 75, 89	0
1	J	157/167 (94%)	-0.39	0 100 100	40, 58, 77, 108	0
1	K	156/167 (93%)	-0.56	0 100 100	32, 51, 72, 91	0
1	L	156/167 (93%)	-0.57	0 100 100	33, 52, 73, 87	0
All	All	1871/2004 (93%)	-0.51	0 100 100	25, 53, 74, 108	0

There are no RSRZ outliers to report.

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 oligosaccharides in this entry.

6.4 Ligands ⓘ

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	TRS	A	305	8/8	0.65	0.21	99,100,100,102	0
3	TRS	B	308	8/8	0.69	0.20	95,95,96,99	0
3	TRS	D	311	8/8	0.69	0.18	106,107,107,108	0
3	TRS	C	302	8/8	0.73	0.19	95,97,99,101	0
3	TRS	D	312	8/8	0.80	0.11	69,70,71,72	0
2	K	E	207	1/1	0.86	0.11	67,67,67,67	0
3	TRS	C	303	8/8	0.87	0.09	61,66,67,69	0
3	TRS	L	309	8/8	0.87	0.09	47,49,51,52	0
3	TRS	A	306	8/8	0.88	0.09	51,52,53,54	0
2	K	G	201	1/1	0.88	0.10	79,79,79,79	0
3	TRS	B	310	8/8	0.88	0.14	64,65,66,67	0
2	K	E	208	1/1	0.88	0.07	55,55,55,55	0
2	K	J	202	1/1	0.89	0.07	64,64,64,64	0
3	TRS	A	301	8/8	0.89	0.13	54,56,57,60	0
3	TRS	C	304	8/8	0.89	0.10	52,54,54,55	0
2	K	H	205	1/1	0.90	0.07	52,52,52,52	0
2	K	I	204	1/1	0.91	0.08	67,67,67,67	0
2	K	A	209	1/1	0.93	0.08	61,61,61,61	0
2	K	F	211	1/1	0.94	0.05	49,49,49,49	0
3	TRS	G	307	8/8	0.94	0.07	30,31,33,34	0
2	K	D	203	1/1	0.94	0.14	76,76,76,76	0
2	K	L	210	1/1	0.95	0.07	60,60,60,60	0
2	K	B	206	1/1	0.96	0.09	62,62,62,62	0
2	K	C	212	1/1	0.97	0.03	52,52,52,52	0

6.5 Other polymers ⓘ

There are no such residues in this entry.