



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 09:55 PM UTC

PDB ID : 6OIV / pdb_00006oiv
Title : XFEL structure of Escherichia coli dGTPase
Authors : Barnes, C.O.; Wu, Y.; Calero, G.
Deposited on : 2019-04-09
Resolution : 3.06 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:

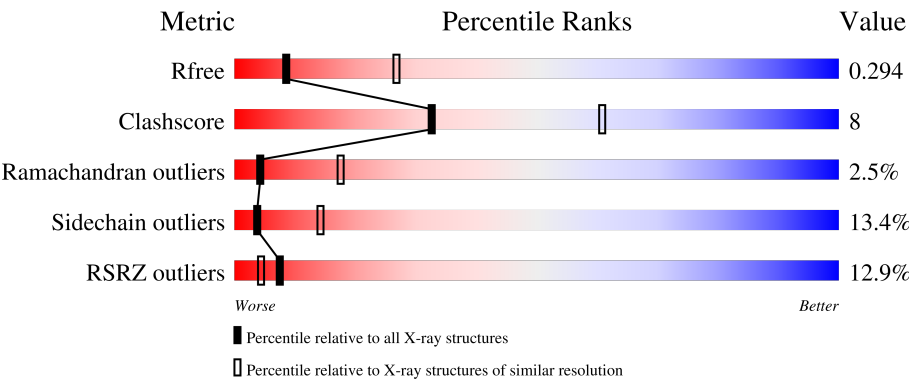
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.06 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	2469 (3.10-3.02)
Clashscore	190562	2569 (3.10-3.02)
Ramachandran outliers	187476	2424 (3.10-3.02)
Sidechain outliers	187428	2423 (3.10-3.02)
RSRZ outliers	180081	2469 (3.10-3.02)

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	505	
1	B	505	
1	C	505	
1	D	505	
1	E	505	

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Mol	Chain	Length	Quality of chain
1	F	505	 4% 71% 22% 5% **

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SO4	C	602	-	-	X	-
3	SO4	D	602	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 24778 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 Deoxyguanosinetriphosphate triphosphohydrolase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	498	Total	C	N	O	S	Se	0	0	0
			4135	2639	737	743	6	10			
1	B	488	Total	C	N	O	S	Se	0	0	0
			4042	2586	720	720	6	10			
1	C	497	Total	C	N	O	S	Se	0	0	0
			4125	2633	736	740	6	10			
1	D	495	Total	C	N	O	S	Se	0	0	0
			4118	2630	734	738	6	10			
1	E	497	Total	C	N	O	S	Se	0	0	0
			4130	2636	736	742	6	10			
1	F	502	Total	C	N	O	S	Se	0	0	0
			4169	2658	745	750	6	10			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	initiating methionine	UNP P15723
B	1	MSE	-	initiating methionine	UNP P15723
C	1	MSE	-	initiating methionine	UNP P15723
D	1	MSE	-	initiating methionine	UNP P15723
E	1	MSE	-	initiating methionine	UNP P15723
F	1	MSE	-	initiating methionine	UNP P15723

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mn	0	0
			1	1		
2	B	1	Total	Mn	0	0
			1	1		
2	C	1	Total	Mn	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	1	Total	Mn	0	0
			1	1		
2	E	1	Total	Mn	0	0
			1	1		
2	F	1	Total	Mn	0	0
			1	1		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	F	1	Total	O	S	0	0
			5	4	1		

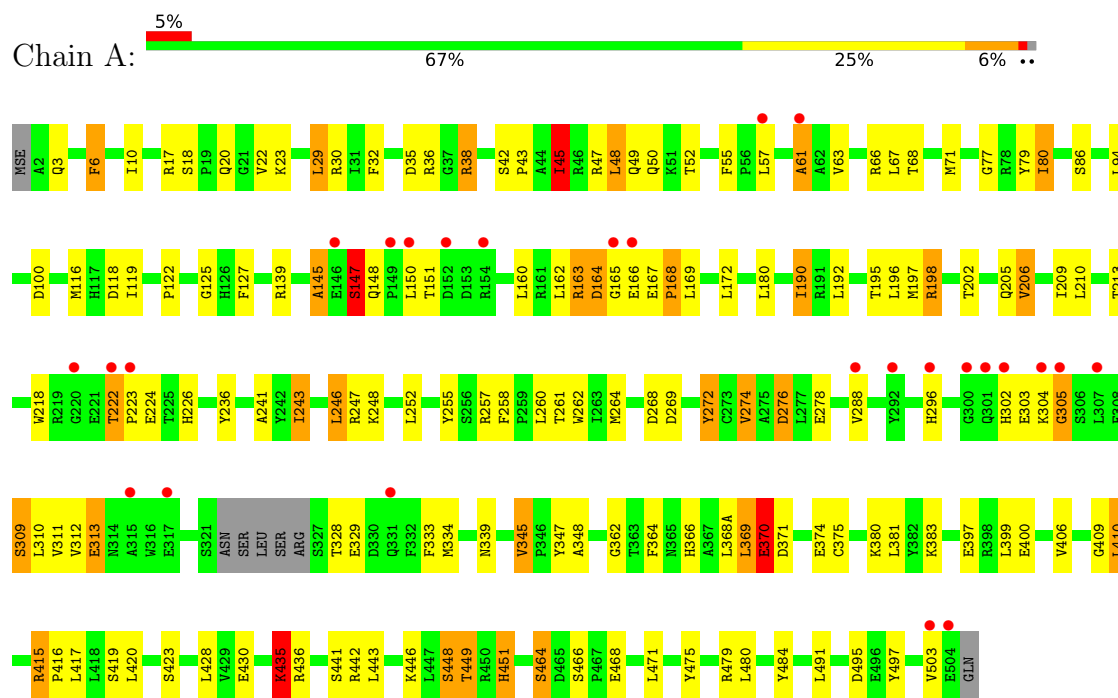
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	3	Total	O	0	0
			3	3		
4	E	3	Total	O	0	0
			3	3		
4	F	2	Total	O	0	0
			2	2		

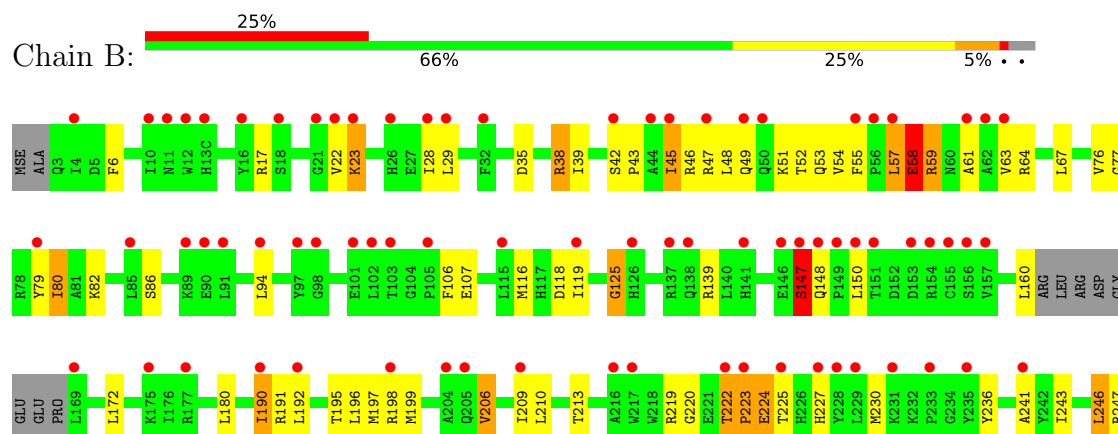
3 Residue-property plots [i](#)

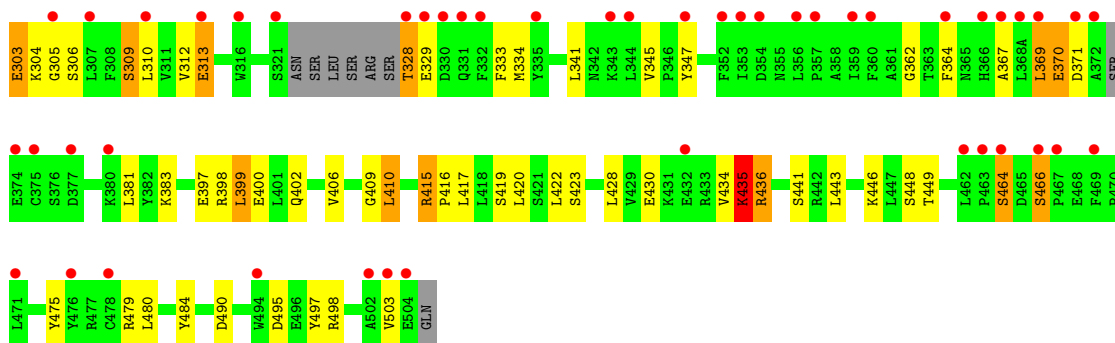
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Deoxyguanosinetriphosphate triphosphohydrolase

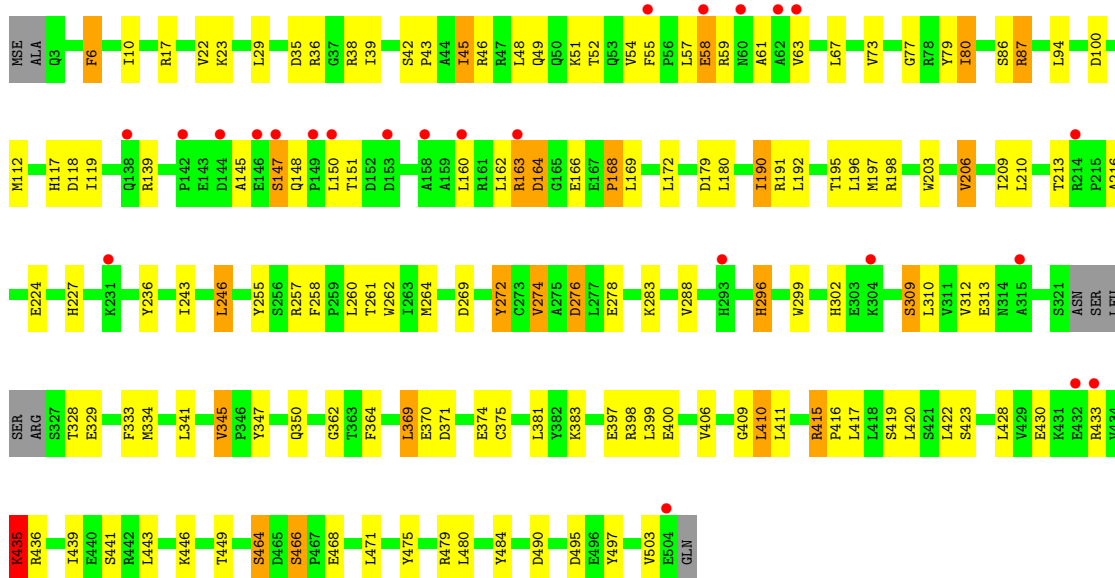


- Molecule 1: Deoxyguanosinetriphosphate triphosphohydrolase

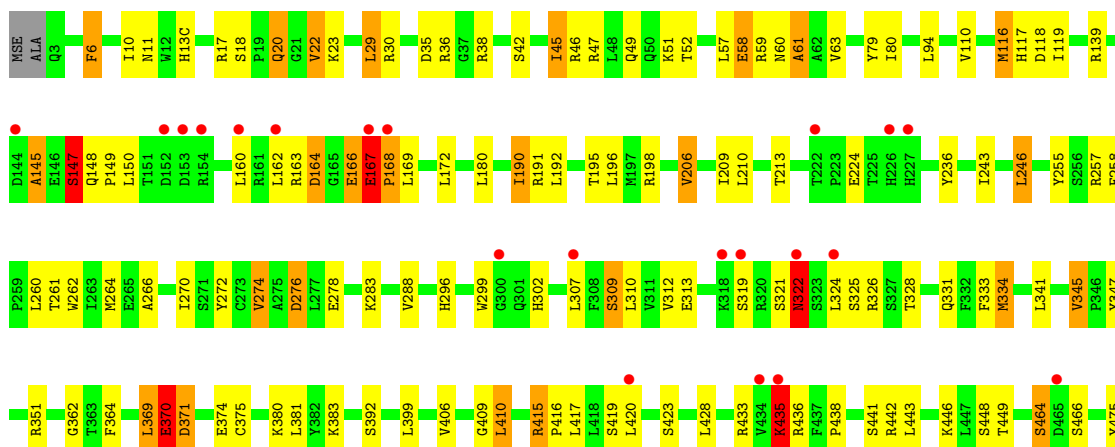


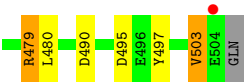


• Molecule 1: Deoxyguanosinetriphosphate triphosphohydrolase



• Molecule 1: Deoxyguanosinetriphosphate triphosphohydrolase





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	192.59Å 192.59Å 291.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	61.20 – 3.06 61.20 – 3.06	Depositor EDS
% Data completeness (in resolution range)	92.3 (61.20-3.06) 92.3 (61.20-3.06)	Depositor EDS
R_{merge}	0.66	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.39 (at 3.07Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, R_{free}	0.225 , 0.248 (Not available) , 0.294	Depositor DCC
R_{free} test set	2922 reflections (2.83%)	wwPDB-VP
Wilson B-factor (Å ²)	29.1	Xtriage
Anisotropy	0.268	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 91.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.31$, $\langle L^2 \rangle = 0.15$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.73	EDS
Total number of atoms	24778	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.84% 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: SO4, MN

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	0.90	1/4227 (0.0%)	1.55	44/5699 (0.8%)
1	B	0.81	1/4132 (0.0%)	1.58	40/5572 (0.7%)
1	C	0.90	4/4217 (0.1%)	1.59	47/5685 (0.8%)
1	D	0.80	1/4209 (0.0%)	1.57	54/5673 (1.0%)
1	E	0.88	1/4222 (0.0%)	1.56	45/5692 (0.8%)
1	F	0.89	2/4262 (0.0%)	1.56	36/5747 (0.6%)
All	All	0.87	10/25269 (0.0%)	1.57	266/34068 (0.8%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	334	MSE	SE-CE	8.58	2.21	1.95
1	F	116	MSE	SE-CE	7.63	2.18	1.95
1	D	231	LYS	C-N	-6.26	1.20	1.33
1	A	116	MSE	SE-CE	5.61	2.12	1.95
1	E	345	VAL	CA-CB	5.54	1.56	1.54

The worst 5 of 266 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	163	ARG	CA-C-N	8.92	138.58	121.54
1	A	163	ARG	C-N-CA	8.92	138.58	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	231	LYS	O-C-N	-8.86	112.87	122.09
1	F	322	ASN	CA-CB-CG	8.81	121.41	112.60
1	C	150	LEU	N-CA-C	8.53	122.78	112.38

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	47	ARG	Sidechain
1	A	66	ARG	Sidechain

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	4135	0	4091	72	0
1	B	4042	0	3996	76	0
1	C	4125	0	4080	72	0
1	D	4118	0	4075	76	0
1	E	4130	0	4086	50	0
1	F	4169	0	4127	64	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	1	0	0	0	0
2	F	1	0	0	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
3	C	20	0	0	3	0
3	D	5	0	0	2	0
3	E	5	0	0	0	0
3	F	5	0	0	0	0
4	C	3	0	0	0	0
4	E	3	0	0	1	0
4	F	2	0	0	0	0
All	All	24778	0	24455	387	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 8.

The worst 5 of 387 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:F:116:MSE:SE	1:F:116:MSE:CE	2.18	1.40
1:C:334:MSE:SE	1:C:334:MSE:CE	2.21	1.38
1:C:66:ARG:NH1	1:C:118:ASP:OD1	1.58	1.33
1:A:17:ARG:NH2	1:F:326:ARG:HH22	1.54	1.05
1:A:17:ARG:HH21	1:F:326:ARG:NH2	1.62	0.98

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	494/505 (98%)	434 (88%)	47 (10%)	13 (3%)	4	17
1	B	482/505 (95%)	421 (87%)	46 (10%)	15 (3%)	3	14
1	C	493/505 (98%)	439 (89%)	41 (8%)	13 (3%)	4	17
1	D	489/505 (97%)	432 (88%)	45 (9%)	12 (2%)	4	17
1	E	493/505 (98%)	441 (90%)	40 (8%)	12 (2%)	4	18
1	F	500/505 (99%)	443 (89%)	47 (9%)	10 (2%)	6	21
All	All	2951/3030 (97%)	2610 (88%)	266 (9%)	75 (2%)	4	17

5 of 75 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	164	ASP
1	A	371	ASP
1	B	58	GLU

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Mol	Chain	Res	Type
1	B	125	GLY
1	B	222	THR

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	443/439 (101%)	383 (86%)	60 (14%)	4	14
1	B	431/439 (98%)	374 (87%)	57 (13%)	4	14
1	C	442/439 (101%)	381 (86%)	61 (14%)	3	13
1	D	441/439 (100%)	380 (86%)	61 (14%)	3	13
1	E	443/439 (101%)	387 (87%)	56 (13%)	4	16
1	F	448/439 (102%)	388 (87%)	60 (13%)	4	14
All	All	2648/2634 (100%)	2293 (87%)	355 (13%)	4	14

5 of 355 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	446	LYS
1	E	436	ARG
1	D	503	VAL
1	E	224	GLU
1	F	59	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 51 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	69	HIS
1	E	26	HIS
1	F	322	ASN
1	D	126	HIS
1	D	186	ASN

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 15 ligands modelled in this entry, 6 are monoatomic - leaving 9 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	SO4	C	604	-	4,4,4	0.22	0	6,6,6	0.33	0
3	SO4	F	602	-	4,4,4	0.32	0	6,6,6	0.12	0
3	SO4	B	602	-	4,4,4	0.29	0	6,6,6	0.14	0
3	SO4	E	602	-	4,4,4	0.29	0	6,6,6	0.14	0
3	SO4	C	605	-	4,4,4	0.45	0	6,6,6	0.35	0
3	SO4	C	602	-	4,4,4	0.37	0	6,6,6	0.32	0
3	SO4	C	603	-	4,4,4	0.35	0	6,6,6	0.32	0
3	SO4	A	602	-	4,4,4	0.37	0	6,6,6	0.21	0
3	SO4	D	602	-	4,4,4	0.54	0	6,6,6	0.14	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.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	605	SO4	1	0
3	C	602	SO4	2	0
3	D	602	SO4	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	488/505 (96%)	0.38	26 (5%) 32 16	4, 43, 93, 126	0
1	B	478/505 (94%)	1.46	127 (26%) 1 1	22, 95, 126, 141	0
1	C	487/505 (96%)	0.36	26 (5%) 32 16	5, 44, 87, 122	0
1	D	485/505 (96%)	1.57	150 (30%) 1 1	34, 94, 125, 145	0
1	E	487/505 (96%)	0.42	24 (4%) 35 17	4, 50, 95, 125	0
1	F	492/505 (97%)	0.51	22 (4%) 38 20	5, 49, 95, 126	0
All	All	2917/3030 (96%)	0.78	375 (12%) 7 4	4, 60, 116, 145	0

The worst 5 of 375 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	60	ASN	5.2
1	D	150	LEU	5.0
1	B	155	CYS	4.8
1	D	74	GLN	4.8
1	B	22	VAL	4.8

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 [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	SO4	D	602	5/5	0.58	0.15	143,144,144,144	0
3	SO4	B	602	5/5	0.66	0.14	140,141,142,143	0
3	SO4	E	602	5/5	0.76	0.15	147,147,147,147	0
3	SO4	A	602	5/5	0.86	0.22	98,98,100,100	0
3	SO4	F	602	5/5	0.87	0.12	113,113,114,114	0
3	SO4	C	603	5/5	0.90	0.10	70,72,72,74	0
3	SO4	C	602	5/5	0.90	0.11	61,63,63,64	0
2	MN	D	601	1/1	0.93	0.08	112,112,112,112	0
3	SO4	C	605	5/5	0.95	0.11	54,58,60,60	0
2	MN	A	601	1/1	0.95	0.06	69,69,69,69	0
2	MN	E	601	1/1	0.97	0.04	43,43,43,43	1
2	MN	F	601	1/1	0.97	0.04	26,26,26,26	1
2	MN	B	601	1/1	0.97	0.05	43,43,43,43	1
3	SO4	C	604	5/5	0.97	0.08	49,50,50,53	0
2	MN	C	601	1/1	0.98	0.03	39,39,39,39	1

6.5 Other polymers [i](#)

There are no such residues in this entry.