wwPDB X-ray Structure Validation Summary Report ⓘ

Nov 11, 2023 – 09:21 am GMT

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PDB ID	:	5NKM
Title	:	SMG8-SMG9 complex
Authors	:	Li, L.; Basquin, J.; Conti, E.
Deposited on	:	2017-03-31
Resolution	:	2.49 Å(reported)

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 2.49 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.
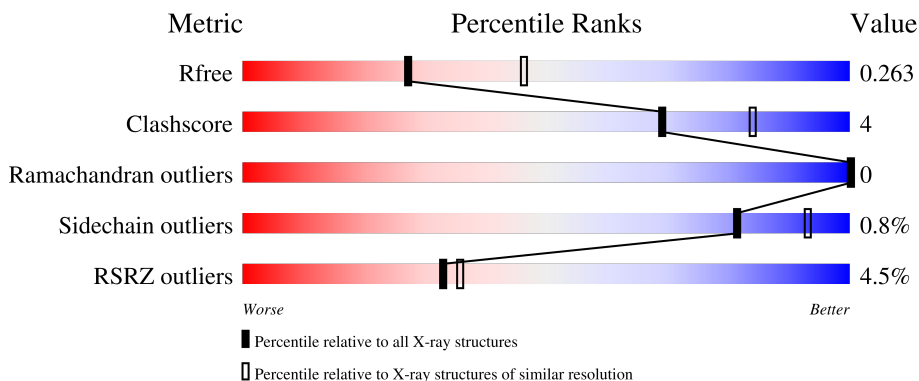
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 <=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	429	
1	C	429	
2	B	317	
3	D	305	
3	F	305	

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Mol	Chain	Length	Quality of chain
4	E	419	2% 76% 8% 16%

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
5	EDO	D	401	-	-	-	X
7	PEG	C	505	-	-	-	X

2 Entry composition ⓘ

There are 8 unique types of molecules in this entry. The entry contains 27997 atoms, of which 13638 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 Protein smg-8.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
1	A	345	Total 5217	C 1671	H 2550	N 462	O 514	S 11	Se 9	0	0	0
1	C	347	Total 5160	C 1661	H 2504	N 464	O 511	S 11	Se 9	0	0	0

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	initiating methionine	UNP O62301
A	424	HIS	-	expression tag	UNP O62301
A	425	HIS	-	expression tag	UNP O62301
A	426	HIS	-	expression tag	UNP O62301
A	427	HIS	-	expression tag	UNP O62301
A	428	HIS	-	expression tag	UNP O62301
A	429	HIS	-	expression tag	UNP O62301
C	1	MSE	-	initiating methionine	UNP O62301
C	424	HIS	-	expression tag	UNP O62301
C	425	HIS	-	expression tag	UNP O62301
C	426	HIS	-	expression tag	UNP O62301
C	427	HIS	-	expression tag	UNP O62301
C	428	HIS	-	expression tag	UNP O62301
C	429	HIS	-	expression tag	UNP O62301

- Molecule 2 is a protein called Protein smg-9.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
2	B	260	Total 4035	C 1299	H 1996	N 345	O 386	S 2	Se 7	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	59	MSE	-	initiating methionine	UNP Q9XWJ1

- Molecule 3 is a protein called Protein smg-9.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
3	D	263	Total 3999	C 1295	H 1960	N 341	O 393	S 2	Se 8	0	0	0
3	F	258	Total 3931	C 1276	H 1928	N 342	O 376	S 2	Se 7	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	59	MSE	-	initiating methionine	UNP Q9XWJ1
F	59	MSE	-	initiating methionine	UNP Q9XWJ1

- Molecule 4 is a protein called Protein smg-8.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
4	E	351	Total 5301	C 1704	H 2590	N 473	O 514	S 11	Se 9	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	1	MSE	-	initiating methionine	UNP O62301

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).

EDO

O2

OH

C1

C2

HO

O1

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total 10	C 2	H 6	O 2	0	0
5	A	1	Total 10	C 2	H 6	O 2	0	0
5	C	1	Total 10	C 2	H 6	O 2	0	0
5	C	1	Total 10	C 2	H 6	O 2	0	0
5	D	1	Total 10	C 2	H 6	O 2	0	0
5	D	1	Total 10	C 2	H 6	O 2	0	0
5	E	1	Total 10	C 2	H 6	O 2	0	0
5	E	1	Total 10	C 2	H 6	O 2	0	0
5	E	1	Total 10	C 2	H 6	O 2	0	0
5	E	1	Total 10	C 2	H 6	O 2	0	0
5	E	1	Total 10	C 2	H 6	O 2	0	0

- Molecule 6 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: $C_4H_{12}NO_3$).

TRS

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	C	1	Total	C	H	N	O	0	0
			20	4	12	1	3		
6	E	1	Total	C	H	N	O	0	0
			20	4	12	1	3		

- Molecule 7 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).
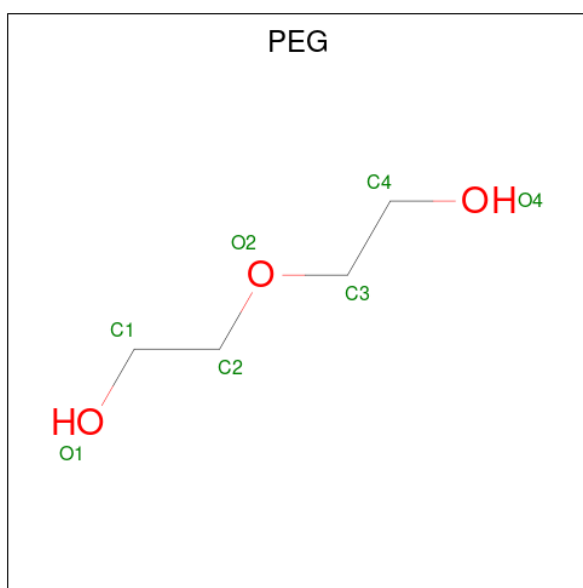
PEG

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	C	1	Total	C	H	O	0	0
			17	4	10	3		
7	C	1	Total	C	H	O	0	0
			17	4	10	3		

- Molecule 8 is water.

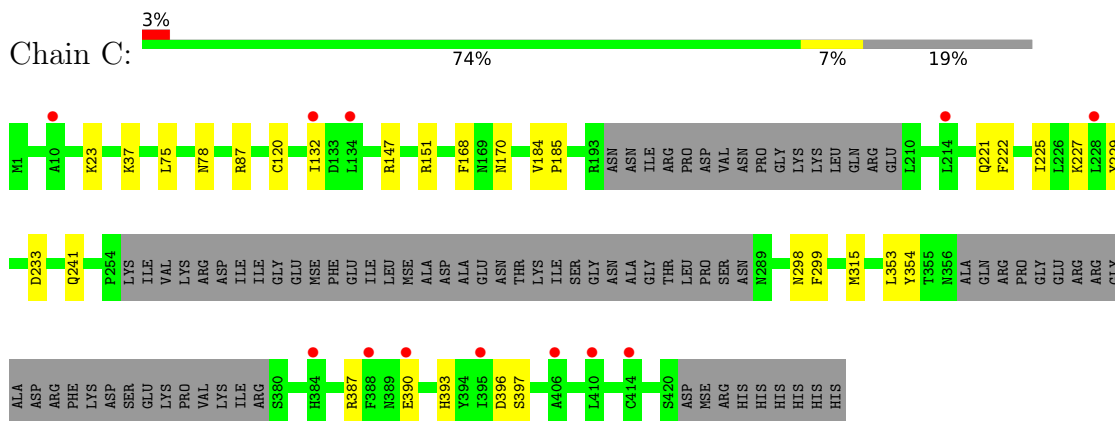| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	38	Total 38	O 38	0	0
8	B	27	Total 27	O 27	0	0
8	C	34	Total 34	O 34	0	0
8	D	22	Total 22	O 22	0	0
8	E	34	Total 34	O 34	0	0
8	F	15	Total 15	O 15	0	0

3 Residue-property plots ⓘ

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.
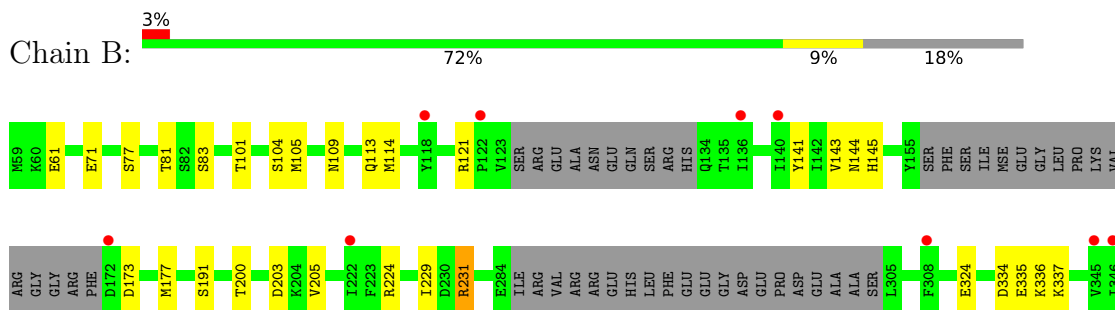
- Molecule 1: Protein smg-8
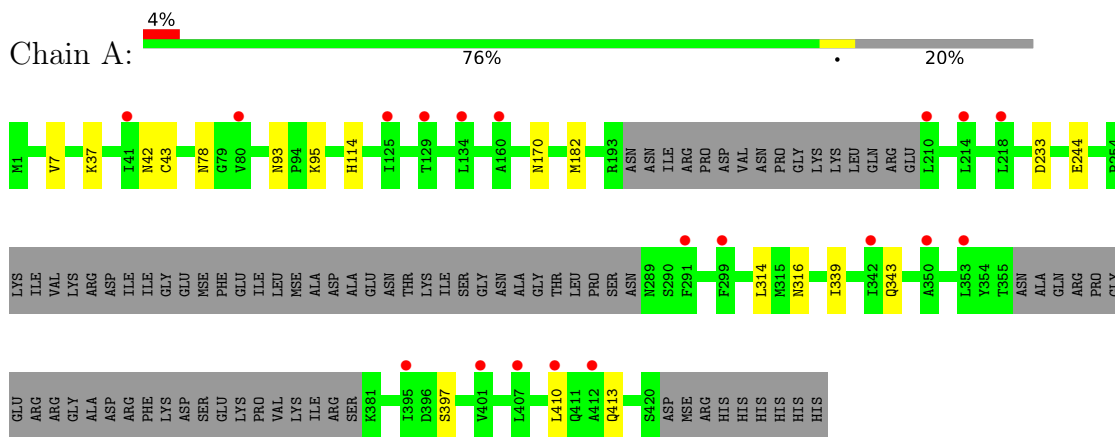
Chain A:

- Molecule 1: Protein smg-8
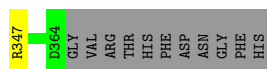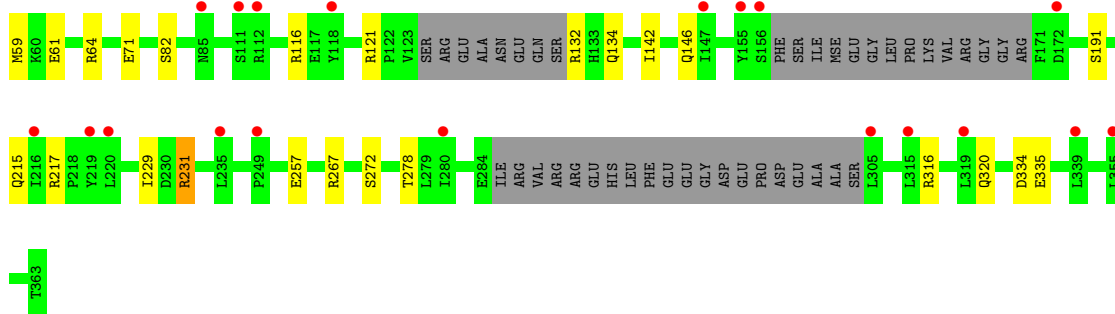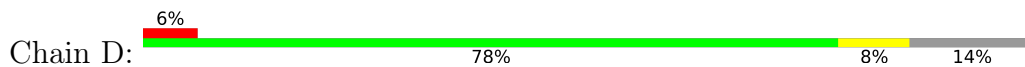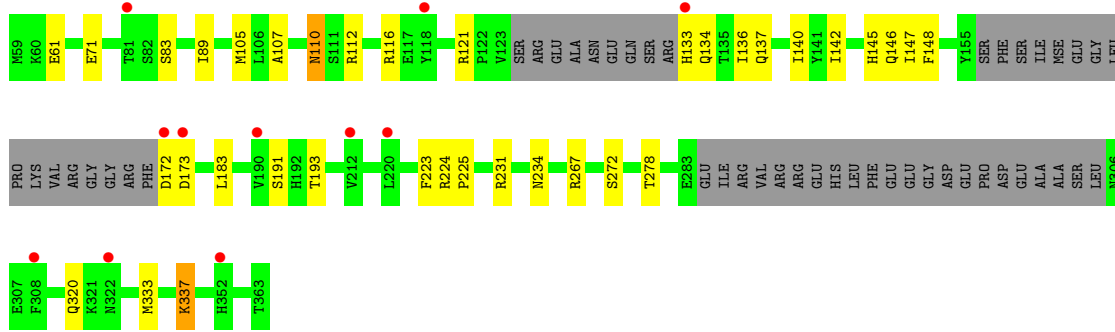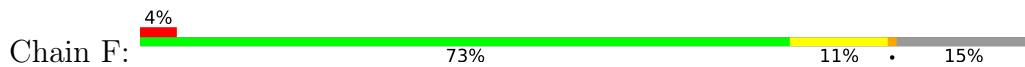
Chain C:

- Molecule 2: Protein smg-9

Chain B:

• Molecule 3: Protein smg-9

Chain D:

6%   78%   8%   14%

• Molecule 3: Protein smg-9

Chain F:

4%   73%   11%   15%

• Molecule 4: Protein smg-8

Chain E:

2%   76%   8%   16%
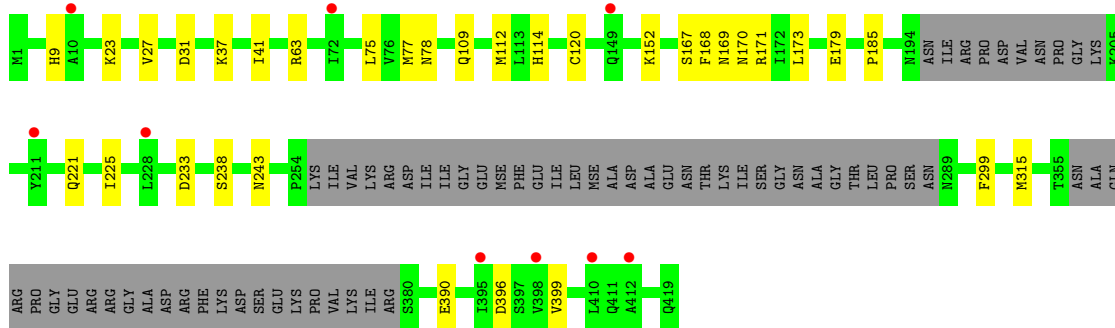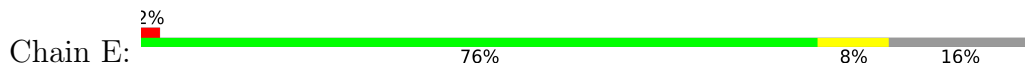
4   Data and refinement statistics   ⓘ

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α, β, γ	111.08Å 111.08Å 374.47Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	62.41 – 2.49 96.20 – 2.49	Depositor EDS
% Data completeness (in resolution range)	99.8 (62.41-2.49) 99.8 (96.20-2.49)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.01 (at 2.48Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.225 , 0.260 0.226 , 0.263	Depositor DCC
R_{free} test set	4711 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	63.3	Xtriage
Anisotropy	0.635	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 60.4	EDS
L-test for twinning[2]	$< \lvert L \rvert > = 0.48, < L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.024 for -h,-k,l	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	27997	wwPDB-VP
Average B, all atoms (Å^2)	96.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.06% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< \lvert L \rvert >$, $< L^2 >$ 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: EDO, PEG, TRS

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	0.25	0/2694	0.41	0/3628
1	C	0.25	0/2682	0.42	0/3614
2	B	0.25	0/2063	0.47	0/2784
3	D	0.25	0/2062	0.44	0/2788
3	F	0.24	0/2026	0.45	0/2738
4	E	0.25	0/2739	0.42	0/3687
All	All	0.25	0/14266	0.43	0/19239

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
3	F	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	F	224	ARG	Peptide

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	2667	2550	2550	11	0
1	C	2656	2504	2509	24	0
2	B	2039	1996	2001	20	0
3	D	2039	1960	1965	18	0
3	F	2003	1928	1947	22	0
4	E	2711	2590	2591	19	0
5	A	8	12	12	1	0
5	C	8	12	12	0	0
5	D	8	12	12	2	0
5	E	20	30	30	2	0
6	C	8	12	12	0	0
6	E	8	12	12	1	0
7	C	14	20	20	0	0
8	A	38	0	0	0	0
8	B	27	0	0	1	0
8	C	34	0	0	6	0
8	D	22	0	0	1	0
8	E	34	0	0	0	0
8	F	15	0	0	0	0
All	All	14359	13638	13673	103	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 4.

The worst 5 of 103 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:23:LYS:NZ	1:C:299:PHE:O	2.03	0.92
4:E:233:ASP:O	4:E:238:SER:OG	1.88	0.90
4:E:23:LYS:NZ	4:E:299:PHE:O	2.05	0.89
1:C:185:PRO:O	8:C:601:HOH:O	1.92	0.85
3:F:61:GLU:OE1	3:F:121:ARG:NH2	2.09	0.84

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	337/429 (79%)	324 (96%)	13 (4%)	0	100	100
1	C	339/429 (79%)	324 (96%)	15 (4%)	0	100	100
2	B	252/317 (80%)	245 (97%)	7 (3%)	0	100	100
3	D	255/305 (84%)	246 (96%)	9 (4%)	0	100	100
3	F	250/305 (82%)	241 (96%)	9 (4%)	0	100	100
4	E	343/419 (82%)	332 (97%)	11 (3%)	0	100	100
All	All	1776/2204 (81%)	1712 (96%)	64 (4%)	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	288/378 (76%)	287 (100%)	1 (0%)	92	97
1	C	281/378 (74%)	280 (100%)	1 (0%)	91	97
2	B	219/279 (78%)	217 (99%)	2 (1%)	78	92
3	D	216/269 (80%)	215 (100%)	1 (0%)	88	96
3	F	211/269 (78%)	208 (99%)	3 (1%)	67	86
4	E	287/369 (78%)	283 (99%)	4 (1%)	67	86
All	All	1502/1942 (77%)	1490 (99%)	12 (1%)	81	93

5 of 12 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
4	E	114	HIS
4	E	168	PHE
3	F	337	LYS
3	F	105	MSE
1	C	168	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	114	HIS
3	D	145	HIS

5.3.3   RNA   ⓘ

There are no RNA molecules in this entry.

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

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

5.5   Carbohydrates   ⓘ

There are no monosaccharides in this entry.

5.6   Ligand geometry   ⓘ

15 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	$\#|Z| > 2$
5	EDO	C	502	-	3,3,3	0.43	0	2,2,2	0.36	0
5	EDO	E	503	-	3,3,3	0.46	0	2,2,2	0.18	0
5	EDO	E	504	-	3,3,3	0.47	0	2,2,2	0.30	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
6	TRS	E	506	-	7,7,7	0.31	0	9,9,9	0.34	0
5	EDO	A	501	-	3,3,3	0.45	0	2,2,2	0.24	0
7	PEG	C	504	-	6,6,6	0.48	0	5,5,5	0.46	0
5	EDO	C	501	-	3,3,3	0.46	0	2,2,2	0.28	0
5	EDO	E	505	-	3,3,3	0.45	0	2,2,2	0.28	0
6	TRS	C	503	-	7,7,7	0.33	0	9,9,9	0.27	0
5	EDO	E	502	-	3,3,3	0.47	0	2,2,2	0.30	0
7	PEG	C	505	-	6,6,6	0.48	0	5,5,5	0.30	0
5	EDO	A	502	-	3,3,3	0.47	0	2,2,2	0.27	0
5	EDO	D	402	-	3,3,3	0.39	0	2,2,2	0.44	0
5	EDO	E	501	-	3,3,3	0.47	0	2,2,2	0.30	0
5	EDO	D	401	-	3,3,3	0.43	0	2,2,2	0.15	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
5	EDO	C	502	-	-	1/1/1/1	-
5	EDO	E	503	-	-	0/1/1/1	-
5	EDO	E	504	-	-	0/1/1/1	-
6	TRS	E	506	-	-	<mark>7/9/9/9</mark>	-
5	EDO	A	501	-	-	0/1/1/1	-
7	PEG	C	504	-	-	1/4/4/4	-
5	EDO	C	501	-	-	1/1/1/1	-
5	EDO	E	505	-	-	0/1/1/1	-
6	TRS	C	503	-	-	0/9/9/9	-
5	EDO	E	502	-	-	0/1/1/1	-
7	PEG	C	505	-	-	<mark>2/4/4/4</mark>	-
5	EDO	A	502	-	-	1/1/1/1	-
5	EDO	D	402	-	-	1/1/1/1	-
5	EDO	E	501	-	-	0/1/1/1	-
5	EDO	D	401	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 14 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	E	506	TRS	C1-C-C3-O3
6	E	506	TRS	C2-C-C3-O3
6	E	506	TRS	N-C-C3-O3
7	C	505	PEG	O1-C1-C2-O2
5	C	502	EDO	O1-C1-C2-O2

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	503	EDO	2	0
6	E	506	TRS	1	0
5	A	501	EDO	1	0
5	D	402	EDO	1	0
5	D	401	EDO	1	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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(Å²)	Q<0.9
1	A	335/429 (78%)	0.47	19 (5%)	23	25	57, 83, 132, 163	0
1	C	337/429 (78%)	0.53	12 (3%)	42	46	56, 83, 149, 197	0
2	B	252/317 (79%)	0.49	9 (3%)	42	46	53, 83, 136, 166	0
3	D	255/305 (83%)	0.70	19 (7%)	14	14	56, 90, 143, 195	0
3	F	250/305 (81%)	0.61	11 (4%)	34	37	59, 89, 132, 150	0
4	E	341/419 (81%)	0.41	9 (2%)	56	59	54, 79, 141, 171	0
All	All	1770/2204 (80%)	0.53	79 (4%)	33	36	53, 84, 141, 197	0

The worst 5 of 79 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	155	TYR	7.0
3	F	133	HIS	4.9
1	C	10	ALA	4.7
3	D	156	SER	4.5
3	F	173	ASP	4.3

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

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

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

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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
5	EDO	D	401	4/4	0.17	1.07	107,129,136,136	0
5	EDO	A	501	4/4	0.46	0.28	124,148,149,150	0
5	EDO	E	501	4/4	0.60	0.30	83,99,101,102	0
5	EDO	A	502	4/4	0.69	0.26	121,145,146,146	0
5	EDO	D	402	4/4	0.72	0.36	98,118,118,119	0
5	EDO	E	502	4/4	0.78	0.17	99,119,120,120	0
6	TRS	E	506	8/8	0.78	0.26	98,119,121,121	0
7	PEG	C	504	7/7	0.79	0.16	110,132,133,133	0
7	PEG	C	505	7/7	0.79	0.49	133,160,160,160	0
5	EDO	C	502	4/4	0.81	0.45	92,111,111,111	0
5	EDO	E	503	4/4	0.81	0.79	112,134,135,135	0
5	EDO	E	504	4/4	0.81	0.27	78,94,96,97	0
5	EDO	E	505	4/4	0.89	0.50	89,107,108,108	0
5	EDO	C	501	4/4	0.90	0.33	85,103,105,105	0
6	TRS	C	503	8/8	0.92	0.33	88,108,111,112	0

6.5 Other polymers ⓘ

There are no such residues in this entry.