



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 23, 2017 – 10:30 AM EST

PDB ID : 5JBS
Title : Conformational changes during monomer-to-dimer transition of Brucella suis VirB8
Authors : Arya, T.; Sharifahmadian, M.; Sygusch, J.; Baron, B.
Deposited on : 2016-04-13
Resolution : 1.95 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

<http://wwpdb.org/validation/2016/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	:	1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20028442
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20028442

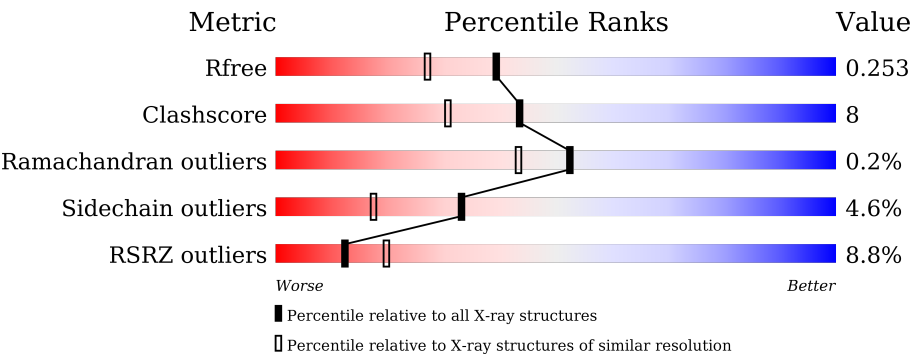
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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}	91344	1833 (1.96-1.96)
Clashscore	102246	1953 (1.96-1.96)
Ramachandran outliers	100387	1936 (1.96-1.96)
Sidechain outliers	100360	1936 (1.96-1.96)
RSRZ outliers	91569	1835 (1.96-1.96)

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. 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	142	<div><div>11%</div><div><div></div><div>84%</div><div>13%</div><div>..</div></div></div>
1	B	142	<div><div>4%</div><div><div></div><div>75%</div><div>12%</div><div>..</div><div>11%</div></div></div>
1	C	142	<div><div>7%</div><div><div></div><div>72%</div><div>13%</div><div>.</div><div>15%</div></div></div>
1	D	142	<div><div>11%</div><div><div></div><div>82%</div><div>13%</div><div>.</div></div></div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	CL	A	301	-	-	X	X
3	GOL	A	302	-	-	-	X
4	PEG	A	303[A]	-	-	X	X
4	PEG	A	303[B]	-	-	X	X
4	PEG	B	303	-	-	X	X
4	PEG	C	302	-	-	-	X
4	PEG	C	303	-	-	-	X
4	PEG	D	301	-	-	X	X
5	EPE	B	301	-	-	-	X
5	EPE	C	301	-	-	-	X
5	EPE	C	304	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4580 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 Type IV secretion system protein virB8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	139	Total	C	N	O	S	0	1	0
			1110	699	188	220	3			
1	B	126	Total	C	N	O	S	0	0	0
			1006	634	168	202	2			
1	C	121	Total	C	N	O	S	0	0	0
			965	612	159	192	2			
1	D	136	Total	C	N	O	S	0	0	0
			1083	681	181	218	3			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	102	ARG	MET	engineered mutation	UNP Q7CEG3
B	102	ARG	MET	engineered mutation	UNP Q7CEG3
C	102	ARG	MET	engineered mutation	UNP Q7CEG3
D	102	ARG	MET	engineered mutation	UNP Q7CEG3

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- 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		

- Molecule 4 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



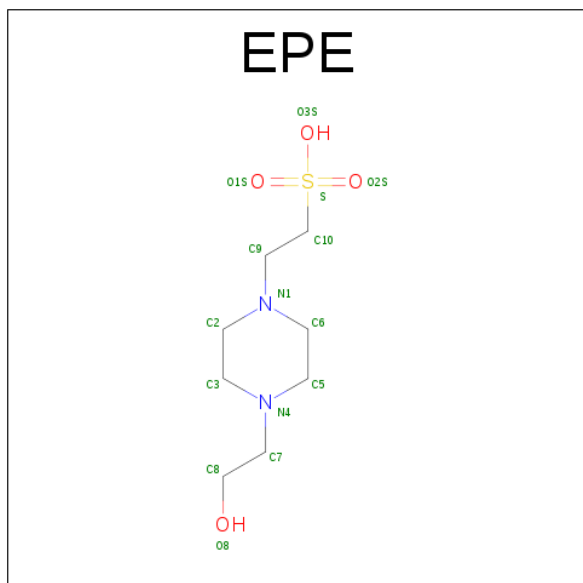
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	1
			14	8	6		
4	B	1	Total	C	O	0	0
			7	4	3		
4	C	1	Total	C	O	0	0
			7	4	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			7	4	3		
4	D	1	Total	C	O	0	0
			7	4	3		

- Molecule 5 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C₈H₁₈N₂O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
5	C	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
5	C	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

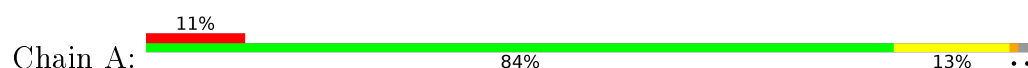
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	84	Total	O	0	0
			84	84		
6	B	86	Total	O	0	0
			86	86		
6	C	67	Total	O	0	0
			67	67		
6	D	84	Total	O	0	0
			84	84		

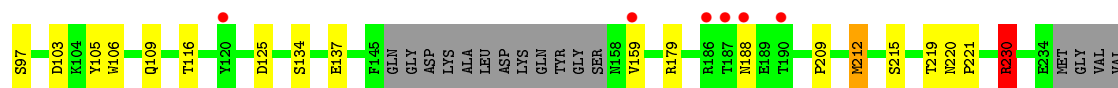
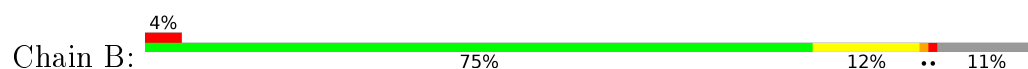
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

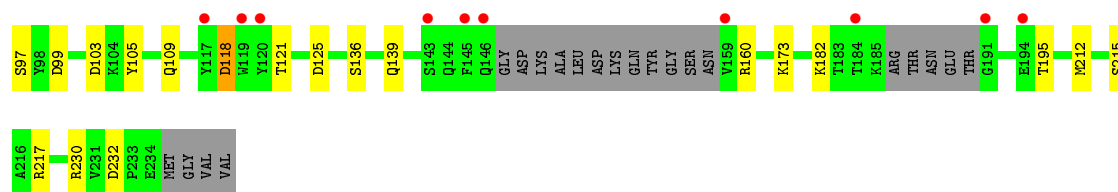
- Molecule 1: Type IV secretion system protein virB8



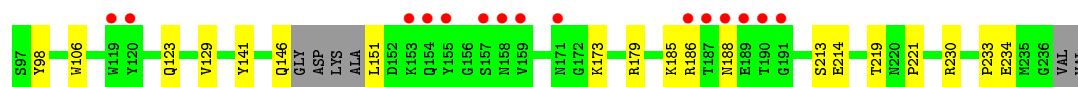
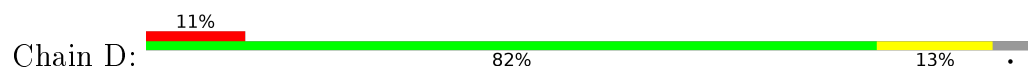
- Molecule 1: Type IV secretion system protein virB8



- Molecule 1: Type IV secretion system protein virB8



- Molecule 1: Type IV secretion system protein virB8



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.50 Å 78.72 Å 70.84 Å 90.00° 111.56° 90.00°	Depositor
Resolution (Å)	42.20 – 1.95 38.87 – 1.95	Depositor EDS
% Data completeness (in resolution range)	96.9 (42.20-1.95) 96.9 (38.87-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.86 (at 1.95 Å)	Xtriage
Refinement program	REFMAC 5.8.0123	Depositor
R, R_{free}	0.191 , 0.241 0.206 , 0.253	Depositor DCC
R_{free} test set	2497 reflections (5.39%)	DCC
Wilson B-factor (Å ²)	36.0	Xtriage
Anisotropy	0.150	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 58.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.023 for l,-k,h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4580	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.45% 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: PEG, GOL, EPE, 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	0.98	0/1134	0.97	1/1539 (0.1%)
1	B	0.93	0/1029	0.95	4/1401 (0.3%)
1	C	0.95	0/987	0.99	6/1342 (0.4%)
1	D	0.91	0/1107	0.95	1/1504 (0.1%)
All	All	0.94	0/4257	0.96	12/5786 (0.2%)

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	D	0	1

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	103	ASP	CB-CG-OD2	-9.94	109.36	118.30
1	B	103	ASP	CB-CG-OD2	-7.84	111.25	118.30
1	B	103	ASP	CB-CG-OD1	7.05	124.65	118.30
1	C	103	ASP	CB-CG-OD1	6.38	124.05	118.30
1	A	103	ASP	CB-CG-OD1	5.65	123.38	118.30
1	C	230	ARG	NE-CZ-NH2	-5.53	117.53	120.30
1	B	179	ARG	NE-CZ-NH1	5.53	123.06	120.30
1	D	230	ARG	NE-CZ-NH2	-5.22	117.69	120.30
1	B	230	ARG	NE-CZ-NH2	5.21	122.90	120.30
1	C	230	ARG	NE-CZ-NH1	5.20	122.90	120.30
1	C	217	ARG	NE-CZ-NH1	5.04	122.82	120.30
1	C	118	ASP	CB-CG-OD1	5.01	122.81	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	188	ASN	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	1110	0	1072	21	0
1	B	1006	0	963	24	0
1	C	965	0	925	9	0
1	D	1083	0	1036	10	0
2	A	1	0	0	2	0
2	B	1	0	0	1	0
3	A	6	0	8	1	0
4	A	14	0	19	21	0
4	B	7	0	10	5	0
4	C	14	0	20	2	0
4	D	7	0	10	4	0
5	B	15	0	18	6	0
5	C	30	0	36	3	0
6	A	84	0	0	3	0
6	B	86	0	0	4	0
6	C	67	0	0	2	0
6	D	84	0	0	3	0
All	All	4580	0	4117	68	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.

All (68) 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:217:ARG:HB2	6:A:405:HOH:O	1.71	0.90
1:C:97:SER:N	6:C:401:HOH:O	2.13	0.81
1:A:109:GLN:OE1	6:A:401:HOH:O	2.02	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:TRP:CH2	4:A:303[B]:PEG:H42	2.22	0.73
1:D:106:TRP:CH2	4:D:301:PEG:H42	2.25	0.70
1:C:118:ASP:HB3	1:C:121:THR:HG22	1.76	0.67
1:D:146:GLN:OE1	6:D:401:HOH:O	2.12	0.66
1:A:212:MET:HB2	1:A:216:ALA:HB3	1.82	0.61
1:A:233:PRO:HG3	3:A:302:GOL:H32	1.81	0.61
4:A:303[A]:PEG:H12	1:B:219:THR:O	1.99	0.61
4:A:303[B]:PEG:C2	1:B:106:TRP:CH2	2.84	0.60
1:B:97:SER:N	6:B:403:HOH:O	2.34	0.60
1:B:137:GLU:HB2	4:B:303:PEG:H21	1.83	0.59
5:B:301:EPE:H92	6:B:401:HOH:O	2.01	0.59
4:A:303[B]:PEG:H22	1:B:106:TRP:CH2	2.38	0.58
4:A:303[B]:PEG:H32	1:B:221:PRO:HD2	1.86	0.57
1:A:219:THR:O	4:A:303[B]:PEG:O2	2.22	0.57
4:A:303[A]:PEG:O4	1:B:221:PRO:HG2	2.05	0.57
1:A:101:VAL:HG13	6:B:415:HOH:O	2.05	0.56
5:B:301:EPE:C9	6:B:401:HOH:O	2.54	0.55
1:A:106:TRP:CZ2	4:A:303[B]:PEG:H42	2.42	0.55
4:A:303[B]:PEG:H22	1:B:106:TRP:CZ2	2.42	0.55
1:B:230:ARG:NH1	1:C:232:ASP:OD2	2.40	0.54
4:C:302:PEG:H21	1:D:233:PRO:HG3	1.89	0.54
4:A:303[B]:PEG:H21	1:B:106:TRP:CH2	2.44	0.53
4:A:303[B]:PEG:H21	1:B:106:TRP:HH2	1.72	0.53
1:A:109:GLN:HG2	6:A:401:HOH:O	2.09	0.52
1:A:106:TRP:HZ2	4:A:303[B]:PEG:C1	2.22	0.52
1:A:173:LYS:O	1:A:173:LYS:HG2	2.10	0.52
1:B:105:TYR:O	1:B:109:GLN:HG2	2.10	0.51
1:B:116:THR:HG21	5:B:301:EPE:H52	1.93	0.51
4:A:303[B]:PEG:C2	1:B:106:TRP:HH2	2.25	0.49
1:B:209:PRO:O	1:B:212:MET:HG2	2.12	0.49
1:B:134:SER:OG	4:B:303:PEG:H32	2.13	0.49
4:C:302:PEG:H11	6:C:413:HOH:O	2.13	0.48
1:D:219:THR:O	4:D:301:PEG:O2	2.31	0.48
1:D:179:ARG:NE	1:D:234:GLU:OE2	2.46	0.48
1:C:121:THR:HG21	5:C:304:EPE:C3	2.45	0.47
1:A:221:PRO:HD2	4:A:303[A]:PEG:H22	1.96	0.47
1:A:173:LYS:CG	1:A:173:LYS:O	2.63	0.46
1:B:125:ASP:OD2	5:B:301:EPE:H51	2.16	0.46
1:B:134:SER:OG	4:B:303:PEG:C3	2.64	0.46
1:D:129:VAL:HG11	1:D:141:TYR:CG	2.50	0.46
4:B:303:PEG:C4	1:C:173:LYS:HB2	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:98:TYR:OH	6:D:402:HOH:O	2.20	0.45
1:A:220:ASN:HA	2:A:301:CL:CL	2.54	0.45
4:A:303[B]:PEG:H41	1:B:106:TRP:HZ2	1.80	0.44
1:D:106:TRP:CZ2	4:D:301:PEG:H42	2.52	0.44
1:B:220:ASN:OD1	2:B:302:CL:CL	2.72	0.44
1:A:220:ASN:OD1	2:A:301:CL:CL	2.72	0.44
1:C:105:TYR:O	1:C:109:GLN:HG2	2.18	0.44
1:D:221:PRO:HD2	4:D:301:PEG:H12	2.00	0.43
1:B:137:GLU:N	4:B:303:PEG:H22	2.34	0.43
1:A:106:TRP:HZ2	4:A:303[B]:PEG:H11	1.84	0.43
5:B:301:EPE:N1	5:B:301:EPE:C7	2.80	0.42
5:B:301:EPE:H72	5:B:301:EPE:N1	2.34	0.42
4:A:303[B]:PEG:H41	1:B:106:TRP:CZ2	2.54	0.42
1:C:139:GLN:HG2	5:C:301:EPE:O3S	2.20	0.42
1:A:154:GLN:HE21	1:A:154:GLN:HB3	1.71	0.42
1:A:219:THR:O	4:A:303[A]:PEG:O2	2.38	0.41
1:A:106:TRP:HZ2	4:A:303[B]:PEG:H12	1.84	0.41
1:D:151:LEU:N	6:D:414:HOH:O	2.54	0.41
1:A:208:ASN:OD1	1:A:208:ASN:C	2.59	0.41
4:A:303[B]:PEG:H32	1:B:221:PRO:CD	2.49	0.41
1:A:117:TYR:OH	1:A:152:ASP:HA	2.21	0.40
4:A:303[A]:PEG:C1	1:B:219:THR:O	2.67	0.40
1:C:125:ASP:OD1	5:C:304:EPE:H32	2.21	0.40
1:C:182:LYS:O	1:C:195:THR:HA	2.22	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	136/142 (96%)	134 (98%)	2 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	122/142 (86%)	118 (97%)	3 (2%)	1 (1%)	24	11
1	C	115/142 (81%)	115 (100%)	0	0	100	100
1	D	132/142 (93%)	128 (97%)	4 (3%)	0	100	100
All	All	505/568 (89%)	495 (98%)	9 (2%)	1 (0%)	52	43

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	188	ASN

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	122/124 (98%)	116 (95%)	6 (5%)	31	15
1	B	112/124 (90%)	108 (96%)	4 (4%)	42	28
1	C	107/124 (86%)	102 (95%)	5 (5%)	32	16
1	D	120/124 (97%)	114 (95%)	6 (5%)	30	14
All	All	461/496 (93%)	440 (95%)	21 (5%)	33	17

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	153	LYS
1	A	154	GLN
1	A	194	GLU
1	A	212	MET
1	A	214	GLU
1	A	215	SER
1	B	159	VAL
1	B	212	MET
1	B	215	SER
1	B	230	ARG

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Mol	Chain	Res	Type
1	C	99	ASP
1	C	136	SER
1	C	160	ARG
1	C	212	MET
1	C	215	SER
1	D	123	GLN
1	D	173	LYS
1	D	185	LYS
1	D	186	ARG
1	D	213	SER
1	D	214	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	109	GLN
1	A	154	GLN
1	B	188	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 2 are monoatomic - leaving 10 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	A	302	-	5,5,5	0.87	0	5,5,5	1.09	0
4	PEG	A	303[A]	-	6,6,6	0.56	0	5,5,5	1.30	1 (20%)
4	PEG	A	303[B]	-	6,6,6	0.66	0	5,5,5	0.55	0
5	EPE	B	301	-	15,15,15	2.22	1 (6%)	19,20,20	2.49	5 (26%)
4	PEG	B	303	-	6,6,6	0.49	0	5,5,5	0.87	0
5	EPE	C	301	-	15,15,15	2.05	1 (6%)	19,20,20	3.00	6 (31%)
4	PEG	C	302	-	6,6,6	0.70	0	5,5,5	1.00	0
4	PEG	C	303	-	6,6,6	1.08	0	5,5,5	1.09	0
5	EPE	C	304	-	15,15,15	2.13	1 (6%)	19,20,20	1.52	5 (26%)
4	PEG	D	301	-	6,6,6	1.13	0	5,5,5	1.38	1 (20%)

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	A	302	-	-	0/4/4/4	0/0/0/0
4	PEG	A	303[A]	-	-	0/4/4/4	0/0/0/0
4	PEG	A	303[B]	-	-	0/4/4/4	0/0/0/0
5	EPE	B	301	-	-	0/9/19/19	0/1/1/1
4	PEG	B	303	-	-	0/4/4/4	0/0/0/0
5	EPE	C	301	-	-	0/9/19/19	0/1/1/1
4	PEG	C	302	-	-	0/4/4/4	0/0/0/0
4	PEG	C	303	-	-	0/4/4/4	0/0/0/0
5	EPE	C	304	-	-	0/9/19/19	0/1/1/1
4	PEG	D	301	-	-	0/4/4/4	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	301	EPE	C10-S	-8.17	1.65	1.77
5	C	304	EPE	C10-S	-7.79	1.65	1.77
5	C	301	EPE	C10-S	-7.34	1.66	1.77

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	301	EPE	C5-C6-N1	-5.68	99.57	110.65
5	C	301	EPE	O2S-S-C10	-4.35	103.79	106.87
5	C	301	EPE	O3S-S-O2S	-3.46	103.61	111.26
5	B	301	EPE	C3-C2-N1	-2.85	105.10	110.65
4	A	303[A]	PEG	O2-C3-C4	-2.56	99.56	110.25
5	C	301	EPE	C8-C7-N4	-2.53	103.56	113.44
5	C	304	EPE	O2S-S-O1S	-2.28	107.52	113.96
5	C	304	EPE	C2-C3-N4	2.21	114.96	110.65
5	C	304	EPE	O3S-S-C10	2.25	109.67	104.99
5	B	301	EPE	O3S-S-C10	2.43	110.04	104.99
4	D	301	PEG	O2-C2-C1	2.56	120.96	110.25
5	C	304	EPE	C7-N4-C3	2.78	117.30	111.25
5	C	301	EPE	C7-N4-C3	3.03	117.83	111.25
5	C	301	EPE	O3S-S-C10	3.30	111.84	104.99
5	C	304	EPE	O1S-S-C10	3.56	109.39	106.87
5	B	301	EPE	C9-N1-C2	4.15	120.28	111.25
5	B	301	EPE	O1S-S-C10	6.23	111.27	106.87
5	C	301	EPE	O1S-S-C10	9.87	113.84	106.87

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

9 monomers are involved in 42 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	302	GOL	1	0
4	A	303[A]	PEG	5	0
4	A	303[B]	PEG	16	0
5	B	301	EPE	6	0
4	B	303	PEG	5	0
5	C	301	EPE	1	0
4	C	302	PEG	2	0
5	C	304	EPE	2	0
4	D	301	PEG	4	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(Å ²)	Q<0.9
1	A	139/142 (97%)	0.22	15 (10%) 8 12	28, 40, 86, 108	0
1	B	126/142 (88%)	0.15	6 (4%) 34 45	28, 40, 80, 116	0
1	C	121/142 (85%)	0.24	10 (8%) 14 22	28, 39, 79, 96	0
1	D	136/142 (95%)	0.24	15 (11%) 7 12	29, 40, 86, 107	0
All	All	522/568 (91%)	0.21	46 (8%) 12 20	28, 40, 85, 116	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	120	TYR	7.3
1	C	159	VAL	6.5
1	B	159	VAL	5.3
1	B	186	ARG	4.8
1	A	119	TRP	4.1
1	D	120	TYR	4.0
1	D	190	THR	4.0
1	D	188	ASN	4.0
1	B	188	ASN	3.9
1	D	155	TYR	3.9
1	C	184	THR	3.8
1	D	119	TRP	3.5
1	B	187	THR	3.4
1	A	187	THR	3.4
1	A	147	GLY	3.3
1	D	189	GLU	3.1
1	C	119	TRP	3.1
1	D	186	ARG	3.1
1	B	190	THR	3.1
1	C	145	PHE	3.0
1	A	154	GLN	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	191	GLY	2.9
1	D	157	SER	2.9
1	D	153	LYS	2.9
1	C	120	TYR	2.8
1	A	148	ASP	2.7
1	A	155	TYR	2.7
1	A	153	LYS	2.7
1	D	191	GLY	2.6
1	D	154	GLN	2.6
1	C	194	GLU	2.6
1	C	117	TYR	2.6
1	A	186	ARG	2.4
1	A	149	LYS	2.3
1	D	187	THR	2.3
1	C	143	SER	2.3
1	A	157	SER	2.3
1	C	146	GLN	2.2
1	B	120	TYR	2.2
1	D	158	ASN	2.2
1	A	152	ASP	2.2
1	A	158	ASN	2.2
1	D	159	VAL	2.2
1	D	171	ASN	2.1
1	A	238	VAL	2.1
1	A	150	ALA	2.0

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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	CL	A	301	1/1	0.90	0.20	12.38	72,72,72,72	0
4	PEG	B	303	7/7	0.65	0.48	11.68	37,39,42,44	7
3	GOL	A	302	6/6	0.90	0.23	7.86	23,26,28,31	6
4	PEG	C	302	7/7	0.89	0.19	7.70	40,57,65,66	0
4	PEG	A	303[B]	7/7	0.92	0.21	7.08	21,23,25,26	7
4	PEG	A	303[A]	7/7	0.92	0.21	7.06	21,25,28,30	7
5	EPE	C	304	15/15	0.73	0.37	5.02	84,102,158,177	0
4	PEG	D	301	7/7	0.81	0.20	3.99	36,42,53,57	2
4	PEG	C	303	7/7	0.60	0.22	3.66	51,65,72,75	0
5	EPE	C	301	15/15	0.76	0.26	3.23	43,71,80,90	10
5	EPE	B	301	15/15	0.85	0.22	2.07	65,76,82,84	9
2	CL	B	302	1/1	0.98	0.07	-1.73	52,52,52,52	0

6.5 Other polymers [i](#)

There are no such residues in this entry.