



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 03:29 pm BST

PDB ID : 2PWP
Title : Crystal structure of spermidine synthase from Plasmodium falciparum in complex with spermidine
Authors : Qiu, W.; Dong, A.; Ren, H.; Wu, H.; Zhao, Y.; Schapira, M.; Wasney, G.; Vedadi, M.; Lew, J.; Kozieradzki, I.; Edwards, A.M.; Arrowsmith, C.H.; Weigelt, J.; Sundstrom, M.; Plotnikov, A.N.; Bochkarev, A.; Hui, R.; Structural Genomics Consortium (SGC)
Deposited on : 2007-05-11
Resolution : 2.10 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	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
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)	:	2.11

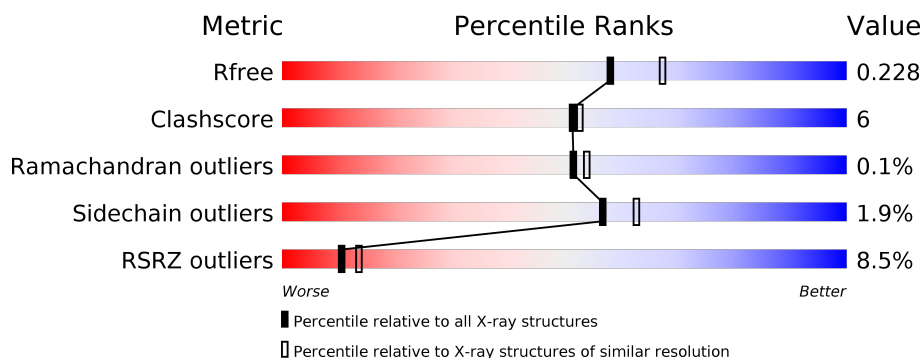
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	282	<div> <div>11%</div> <div> <div></div> <div>82%</div> <div>16%</div> <div></div> </div> </div>
1	B	282	<div> <div>7%</div> <div> <div></div> <div>83%</div> <div>13%</div> <div></div> </div> </div>
1	C	282	<div> <div>7%</div> <div> <div></div> <div>94%</div> <div>6%</div> <div></div> </div> </div>

2 Entry composition [i](#)

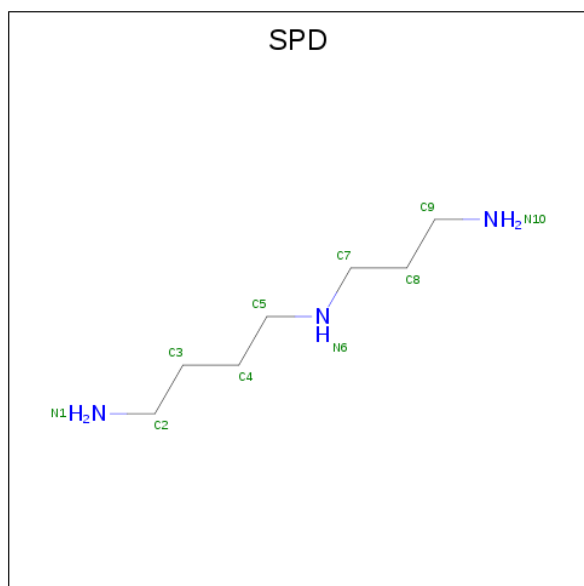
There are 5 unique types of molecules in this entry. The entry contains 7110 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Spermidine synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	281	Total	C	N	O	S	1	3	0
			2276	1476	356	430	14			
1	B	273	Total	C	N	O	S	6	5	0
			2217	1438	344	420	15			
1	C	281	Total	C	N	O	S	0	1	0
			2260	1468	353	425	14			

- Molecule 2 is SPERMIDINE (three-letter code: SPD) (formula: C₇H₁₉N₃).



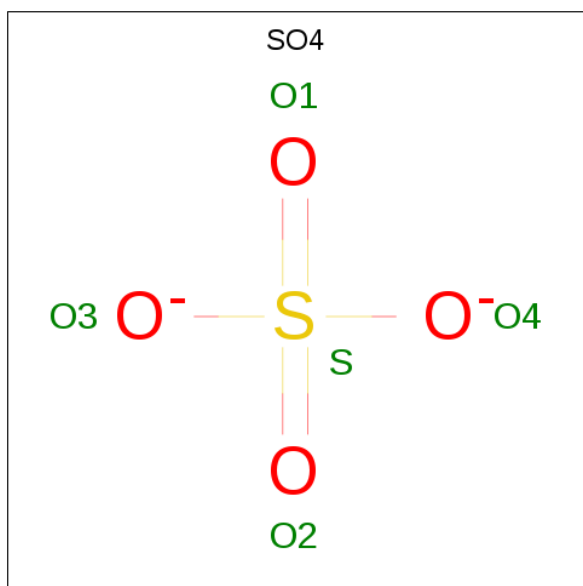
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			10	7	3		
2	B	1	Total	C	N	0	0
			10	7	3		
2	C	1	Total	C	N	0	0
			10	7	3		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

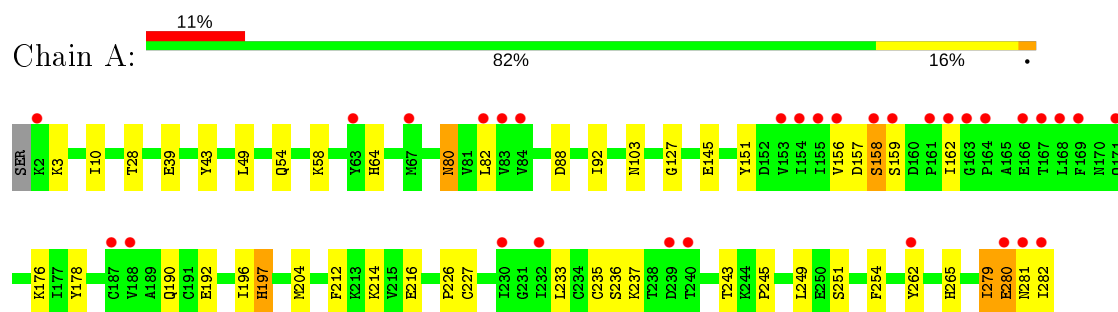
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	63	Total	O	0	0
			63	63		
5	B	92	Total	O	0	0
			92	92		
5	C	149	Total	O	0	0
			149	149		

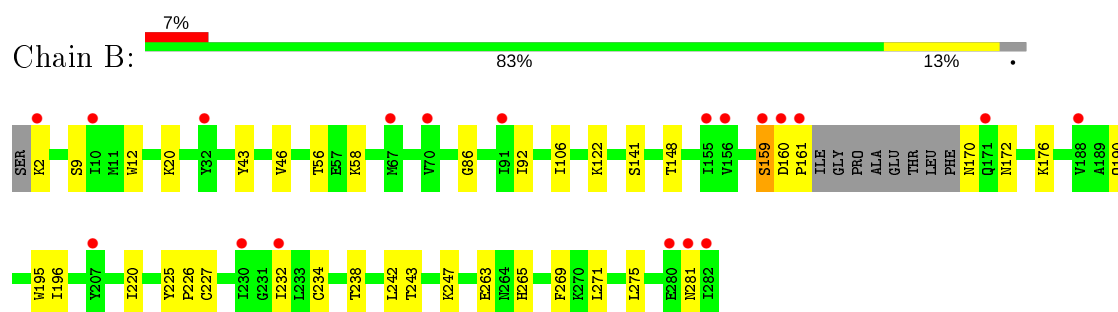
3 Residue-property plots [i](#)

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.

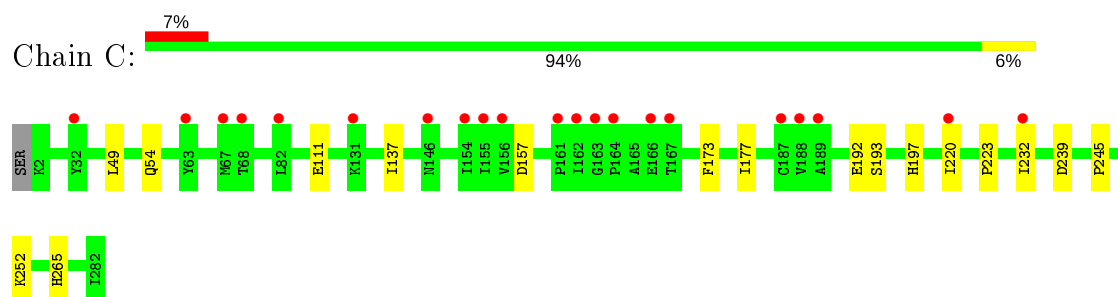
• Molecule 1: Spermidine synthase



• Molecule 1: Spermidine synthase



• Molecule 1: Spermidine synthase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	198.97 Å 134.50 Å 48.71 Å 90.00° 95.88° 90.00°	Depositor
Resolution (Å)	30.00 – 2.10 29.62 – 2.09	Depositor EDS
% Data completeness (in resolution range)	100.0 (30.00-2.10) 99.3 (29.62-2.09)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 2.10 Å)	Xtriage
Refinement program	REFMAC 5.1.19	Depositor
R, R_{free}	0.195 , 0.232 0.192 , 0.228	Depositor DCC
R_{free} test set	3765 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	39.0	Xtriage
Anisotropy	0.059	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 52.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7110	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.88% 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: GOL, SPD, SO4

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	0.33	0/2330	0.52	0/3147
1	B	0.39	0/2276	0.55	1/3072 (0.0%)
1	C	0.41	0/2313	0.53	0/3122
All	All	0.37	0/6919	0.53	1/9341 (0.0%)

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
1	B	0	1
All	All	0	3

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	161	PRO	N-CA-C	-5.72	97.24	112.10

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	157	ASP	Peptide
1	A	279	ILE	Peptide
1	B	160	ASP	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	2276	0	2275	37	0
1	B	2217	0	2220	27	0
1	C	2260	0	2278	15	0
2	A	10	0	19	5	0
2	B	10	0	19	2	0
2	C	10	0	19	1	0
3	A	6	0	8	0	0
3	B	6	0	8	0	0
3	C	6	0	8	0	0
4	B	5	0	0	0	0
5	A	63	0	0	0	0
5	B	92	0	0	1	0
5	C	149	0	0	1	0
All	All	7110	0	6854	77	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 6.

All (77) 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:192:GLU:HB3	1:C:197:HIS:CD2	1.82	1.14
1:C:192:GLU:HB3	1:C:197:HIS:HD2	1.29	0.97
1:B:2:LYS:HD2	1:B:20:LYS:HE2	1.46	0.94
1:A:192:GLU:HB3	1:A:197[A]:HIS:CD2	2.11	0.86
1:A:80:ASN:HD21	1:A:103:ASN:HD22	1.33	0.77
1:C:193:SER:H	1:C:197:HIS:HD2	1.33	0.77
1:A:192:GLU:HB3	1:A:197[A]:HIS:NE2	2.05	0.71
1:A:192:GLU:CB	1:A:197[A]:HIS:CD2	2.80	0.64
1:A:43:TYR:CD2	1:A:58:LYS:HD3	2.36	0.61
1:A:158:SER:O	1:A:159:SER:HB3	2.02	0.60
1:A:245:PRO:HG2	1:A:262:TYR:CE1	2.37	0.60
1:C:239:ASP:HB2	5:C:588:HOH:O	2.02	0.60
1:B:170:ASN:HD21	1:B:172:ASN:HB3	1.65	0.59
1:A:88:ASP:OD2	2:A:401:SPD:H92	2.02	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:173:PHE:CZ	1:C:177:ILE:HD11	2.37	0.58
1:B:243:THR:HG22	1:B:269:PHE:CG	2.37	0.58
1:C:193:SER:H	1:C:197:HIS:CD2	2.19	0.58
1:C:111:GLU:HB3	1:C:137:ILE:HD12	1.86	0.57
1:A:178:TYR:CE1	1:A:237:LYS:HE2	2.39	0.57
1:C:173:PHE:O	1:C:177:ILE:HD13	2.05	0.55
1:B:86:GLY:CA	1:B:106:ILE:HD11	2.36	0.55
1:B:170:ASN:HD22	1:B:172:ASN:H	1.56	0.54
1:A:280:GLU:O	1:A:281:ASN:HB2	2.06	0.54
1:A:80:ASN:ND2	1:A:103:ASN:HD22	2.01	0.54
1:B:226:PRO:O	1:B:227:CYS:HB2	2.08	0.54
1:B:243:THR:HG22	1:B:269:PHE:CD1	2.43	0.53
1:B:43:TYR:CD2	1:B:58:LYS:HD3	2.45	0.52
1:A:190:GLN:HE22	2:A:401:SPD:H52	1.75	0.52
1:C:245:PRO:HG3	1:C:265:HIS:CG	2.44	0.51
1:C:245:PRO:HG3	1:C:265:HIS:ND1	2.25	0.51
1:B:220:ILE:HD11	1:B:232:ILE:HD11	1.93	0.50
1:A:280:GLU:C	1:A:282:ILE:H	2.15	0.50
1:B:170:ASN:ND2	1:B:172:ASN:HB3	2.26	0.50
1:B:225:TYR:CE1	2:B:401:SPD:H82	2.47	0.50
1:B:92:ILE:HD11	1:B:106:ILE:HD13	1.94	0.49
1:B:86:GLY:C	1:B:106:ILE:HD11	2.33	0.49
1:C:192:GLU:HB3	1:C:197:HIS:CG	2.40	0.48
1:B:86:GLY:HA3	1:B:106:ILE:HD11	1.97	0.47
1:A:54:GLN:O	2:A:401:SPD:H82	2.15	0.47
1:A:3:LYS:HB2	1:A:3:LYS:HE3	1.52	0.47
1:B:86:GLY:HA3	1:B:106:ILE:CD1	2.45	0.47
1:B:170:ASN:ND2	1:B:172:ASN:H	2.13	0.46
1:A:28:THR:OG1	1:B:122:LYS:HD2	2.16	0.46
1:B:195:TRP:CE2	1:C:223:PRO:HD3	2.50	0.46
1:A:279:ILE:HA	1:A:282:ILE:HD11	1.97	0.45
1:B:46:VAL:HG22	1:B:56:THR:HG22	1.97	0.45
1:C:157:ASP:OD1	2:C:401:SPD:H72	2.16	0.45
1:A:214:LYS:HB3	1:A:236:SER:HB2	1.99	0.45
1:A:204:MET:HB3	1:A:233:LEU:HD22	2.00	0.44
1:C:220:ILE:HD11	1:C:232:ILE:HD11	1.99	0.44
1:B:263:GLU:HG3	5:B:619:HOH:O	2.17	0.44
1:B:9:SER:HB3	1:B:12:TRP:CE2	2.53	0.43
1:A:64:HIS:NE2	2:A:401:SPD:N10	2.67	0.43
1:A:226:PRO:O	1:A:227:CYS:HB2	2.18	0.43
1:A:156:VAL:HG12	1:A:158:SER:HB2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:196:ILE:HG12	1:B:227:CYS:SG	2.58	0.43
1:A:216:GLU:OE2	1:A:243:THR:HG23	2.19	0.43
1:A:82:LEU:HB2	1:A:151:TYR:CE1	2.54	0.42
1:B:271:LEU:HD13	1:B:275:LEU:HD23	2.01	0.42
1:B:172:ASN:OD1	1:B:176:LYS:HE3	2.20	0.42
1:A:249:LEU:HB3	1:A:254:PHE:CG	2.55	0.42
1:A:49:LEU:HG	1:A:54:GLN:HG3	2.02	0.42
1:A:92:ILE:HG22	1:A:127:GLY:HA3	2.02	0.41
1:B:234:CYS:HB3	1:B:242:LEU:CD1	2.50	0.41
1:A:265:HIS:C	1:A:265:HIS:CD2	2.93	0.41
1:A:279:ILE:HA	1:A:282:ILE:CD1	2.50	0.41
1:A:196:ILE:HG12	1:A:227:CYS:SG	2.61	0.41
1:A:212:PHE:CG	1:A:235:CYS:HB3	2.56	0.41
1:A:10:ILE:HD11	1:A:162:ILE:HD13	2.01	0.41
1:B:265:HIS:CD2	1:B:265:HIS:C	2.94	0.41
1:A:145[A]:GLU:OE1	1:A:176:LYS:HE2	2.21	0.41
1:A:280:GLU:C	1:A:282:ILE:N	2.72	0.41
1:A:158:SER:O	1:A:159:SER:CB	2.68	0.41
1:C:49:LEU:HG	1:C:54:GLN:HG3	2.02	0.40
1:A:190:GLN:NE2	2:A:401:SPD:H52	2.36	0.40
1:B:225:TYR:OH	2:B:401:SPD:H31	2.21	0.40
1:A:280:GLU:HB2	1:A:281:ASN:H	1.78	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	282/282 (100%)	276 (98%)	6 (2%)	0	100	100
1	B	274/282 (97%)	265 (97%)	8 (3%)	1 (0%)	34	32
1	C	280/282 (99%)	271 (97%)	9 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	836/846 (99%)	812 (97%)	23 (3%)	1 (0%)	51 54

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	159	SER

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	255/254 (100%)	248 (97%)	7 (3%)	44 48
1	B	250/254 (98%)	243 (97%)	7 (3%)	43 47
1	C	254/254 (100%)	253 (100%)	1 (0%)	91 94
All	All	759/762 (100%)	744 (98%)	15 (2%)	57 60

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	GLU
1	A	80	ASN
1	A	158	SER
1	A	197[A]	HIS
1	A	197[B]	HIS
1	A	251	SER
1	A	280	GLU
1	B	141	SER
1	B	148	THR
1	B	159	SER
1	B	190	GLN
1	B	238	THR
1	B	247	LYS
1	B	281	ASN
1	C	252	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	GLN
1	A	80	ASN
1	A	134	ASN
1	A	179	ASN
1	A	190	GLN
1	B	103	ASN
1	B	170	ASN
1	B	246	ASN
1	C	134	ASN
1	C	171	GLN
1	C	172	ASN
1	C	197	HIS
1	C	246	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](#)

7 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
2	SPD	A	401	-	9,9,9	0.53	0	8,8,8	2.67	3 (37%)
2	SPD	C	401	-	9,9,9	0.39	0	8,8,8	0.88	0
3	GOL	B	501	-	5,5,5	0.39	0	5,5,5	0.31	0
3	GOL	A	501	-	5,5,5	0.35	0	5,5,5	0.30	0
3	GOL	C	501	-	5,5,5	0.33	0	5,5,5	0.41	0
4	SO4	B	601	-	4,4,4	0.14	0	6,6,6	0.13	0
2	SPD	B	401	-	9,9,9	0.49	0	8,8,8	0.55	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
2	SPD	A	401	-	-	0/7/7/7	-
2	SPD	C	401	-	-	0/7/7/7	-
3	GOL	B	501	-	-	0/4/4/4	-
3	GOL	A	501	-	-	2/4/4/4	-
3	GOL	C	501	-	-	0/4/4/4	-
2	SPD	B	401	-	-	0/7/7/7	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	SPD	C4-C5-N6	-5.01	98.62	112.14
2	A	401	SPD	C8-C7-N6	-4.46	100.10	112.14
2	A	401	SPD	C7-N6-C5	2.67	126.03	113.45

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	501	GOL	O1-C1-C2-C3
3	A	501	GOL	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	SPD	5	0
2	C	401	SPD	1	0
2	B	401	SPD	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 ⓘ

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	281/282 (99%)	0.38	31 (11%) 5 7	30, 45, 80, 101	13 (4%)
1	B	273/282 (96%)	0.16	19 (6%) 16 20	21, 36, 66, 110	10 (3%)
1	C	281/282 (99%)	0.13	21 (7%) 14 18	21, 33, 62, 86	4 (1%)
All	All	835/846 (98%)	0.22	71 (8%) 10 13	21, 38, 72, 110	27 (3%)

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	162	ILE	6.2
1	B	161	PRO	6.1
1	C	162	ILE	5.4
1	B	282	ILE	5.2
1	B	281	ASN	4.4
1	A	155	ILE	4.4
1	A	167	THR	4.3
1	B	10	ILE	4.2
1	A	281	ASN	4.1
1	A	156	VAL	3.9
1	A	282	ILE	3.7
1	C	167	THR	3.7
1	A	169	PHE	3.7
1	C	166	GLU	3.7
1	A	166	GLU	3.6
1	B	171	GLN	3.6
1	A	164	PRO	3.6
1	A	188	VAL	3.4
1	B	232	ILE	3.3
1	A	163	GLY	3.3
1	A	240	THR	3.3
1	A	154	ILE	3.3
1	A	2	LYS	3.2

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Mol	Chain	Res	Type	RSRZ
1	C	161	PRO	3.1
1	A	82	LEU	3.1
1	B	32	TYR	3.0
1	C	163	GLY	3.0
1	C	188	VAL	2.9
1	B	188	VAL	2.9
1	C	68	THR	2.9
1	A	84	VAL	2.8
1	C	154	ILE	2.7
1	A	83	VAL	2.7
1	B	156	VAL	2.7
1	A	161	PRO	2.7
1	C	155	ILE	2.7
1	A	67	MET	2.7
1	A	262	TYR	2.7
1	B	280[A]	GLU	2.7
1	A	168	LEU	2.7
1	B	159	SER	2.6
1	B	67	MET	2.6
1	C	164	PRO	2.6
1	B	155	ILE	2.5
1	A	239	ASP	2.5
1	A	63	TYR	2.5
1	A	230	ILE	2.5
1	C	63	TYR	2.5
1	B	91	ILE	2.5
1	A	171	GLN	2.4
1	C	67	MET	2.4
1	C	156	VAL	2.4
1	A	153	VAL	2.4
1	C	131	LYS	2.4
1	B	160	ASP	2.4
1	B	230	ILE	2.4
1	A	187	CYS	2.3
1	A	232	ILE	2.3
1	C	220	ILE	2.3
1	A	159	SER	2.3
1	B	2	LYS	2.3
1	A	158	SER	2.2
1	C	187	CYS	2.2
1	C	32	TYR	2.2
1	C	82	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	189	ALA	2.2
1	B	207	TYR	2.1
1	B	70	VAL	2.1
1	C	232	ILE	2.1
1	C	146	ASN	2.1
1	A	280	GLU	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(\AA^2)	Q<0.9
2	SPD	B	401	10/10	0.91	0.40	27,46,55,59	0
2	SPD	A	401	10/10	0.93	0.31	32,43,63,68	0
3	GOL	A	501	6/6	0.94	0.18	51,57,62,74	0
3	GOL	C	501	6/6	0.95	0.11	42,50,61,63	0
2	SPD	C	401	10/10	0.95	0.35	25,39,54,54	0
4	SO4	B	601	5/5	0.96	0.13	54,56,75,92	0
3	GOL	B	501	6/6	0.96	0.15	29,43,47,52	0

6.5 Other polymers [i](#)

There are no such residues in this entry.