



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 13, 2026 – 11:33 PM UTC

PDB ID : 8IR4 / pdb_00008ir4
Title : Crystal structure of the SLF1 BRCT domain in complex with a Rad18 peptide containing pS442
Authors : Xiang, S.; Huang, W.; Qiu, F.
Deposited on : 2023-03-17
Resolution : 1.62 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 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

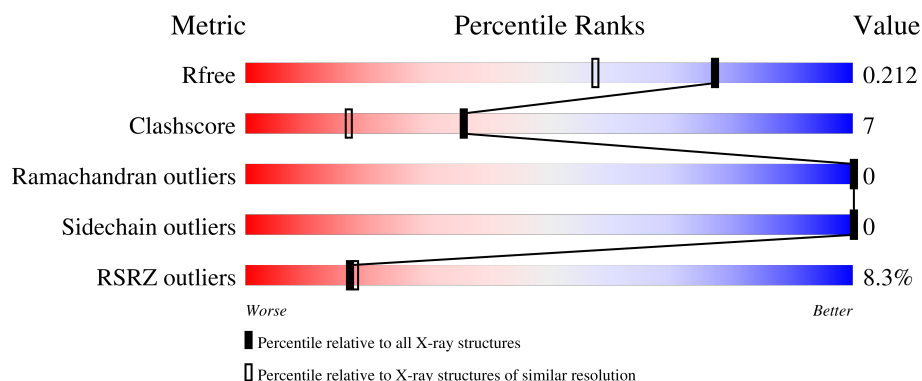
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.62 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	6728 (1.64-1.60)
Clashscore	190562	7023 (1.64-1.60)
Ramachandran outliers	187476	6898 (1.64-1.60)
Sidechain outliers	187428	6896 (1.64-1.60)
RSRZ outliers	180081	6727 (1.64-1.60)

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	207	3% 87% 7% 5%
1	B	207	10% 78% 17% 5%
2	C	17	29% 76% 18% 6%
2	D	17	24% 24% 59% 18%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4004 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 SMC5-SMC6 complex localization factor protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	196	Total	C	N	O	S	0	0	0
			1594	1028	278	281	7			
1	B	197	Total	C	N	O	S	0	2	0
			1615	1041	282	285	7			

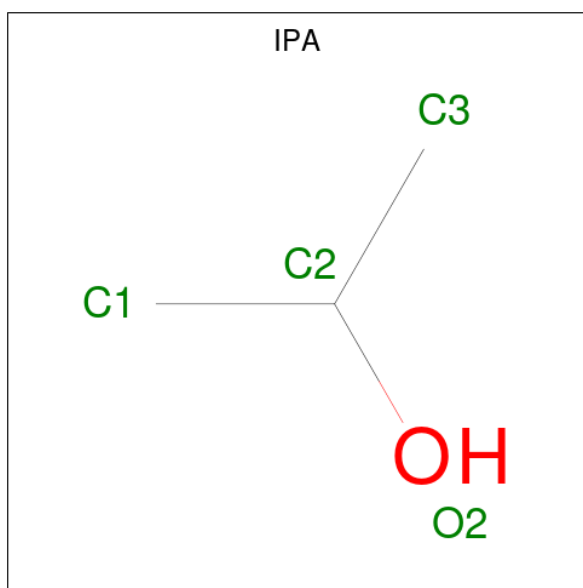
There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	200	LEU	-	expression tag	UNP Q9BQI6
A	201	GLU	-	expression tag	UNP Q9BQI6
A	202	HIS	-	expression tag	UNP Q9BQI6
A	203	HIS	-	expression tag	UNP Q9BQI6
A	204	HIS	-	expression tag	UNP Q9BQI6
A	205	HIS	-	expression tag	UNP Q9BQI6
A	206	HIS	-	expression tag	UNP Q9BQI6
A	207	HIS	-	expression tag	UNP Q9BQI6
B	200	LEU	-	expression tag	UNP Q9BQI6
B	201	GLU	-	expression tag	UNP Q9BQI6
B	202	HIS	-	expression tag	UNP Q9BQI6
B	203	HIS	-	expression tag	UNP Q9BQI6
B	204	HIS	-	expression tag	UNP Q9BQI6
B	205	HIS	-	expression tag	UNP Q9BQI6
B	206	HIS	-	expression tag	UNP Q9BQI6
B	207	HIS	-	expression tag	UNP Q9BQI6

- Molecule 2 is a protein called SER-ASP-SER-CYS-ASN-SER-SEP-SER-SER-ASP-ILE-ILE-ARG-ASP-LEU-LEU-GLU.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	C	16	Total 124	C 69	N 20	O 33	P 1	S 1	0	0	0
2	D	14	Total 107	C 60	N 18	O 27	P 1	S 1	0	0	0

- Molecule 3 is ISOPROPYL ALCOHOL (CCD ID: IPA) (formula: C_3H_8O).



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

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

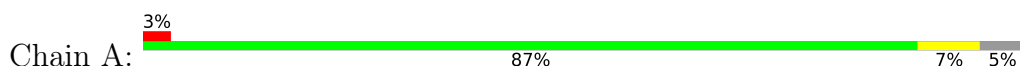
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	279	Total 279	O 279	0	0
6	B	234	Total 234	O 234	0	0
6	C	15	Total 15	O 15	0	0
6	D	21	Total 21	O 21	0	0

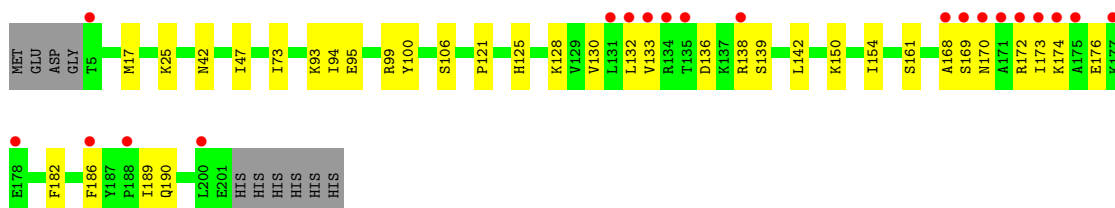
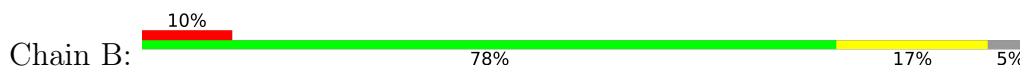
3 Residue-property plots [i](#)

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.

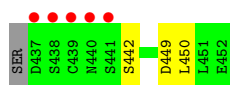
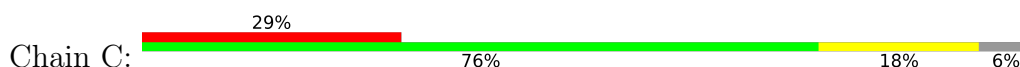
- Molecule 1: SMC5-SMC6 complex localization factor protein 1



- Molecule 1: SMC5-SMC6 complex localization factor protein 1



- Molecule 2: SER-ASP-SER-CYS-ASN-SER-SEP-SER-SER-ASP-ILE-ILE-ARG-ASP-LEU-LEU-GLU



- Molecule 2: SER-ASP-SER-CYS-ASN-SER-SEP-SER-SER-ASP-ILE-ILE-ARG-ASP-LEU-LEU-GLU



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	40.42Å 76.10Å 141.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.60 – 1.62 29.60 – 1.62	Depositor EDS
% Data completeness (in resolution range)	97.6 (29.60-1.62) 97.6 (29.60-1.62)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.32 (at 1.62Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.181 , 0.213 0.181 , 0.212	Depositor DCC
R_{free} test set	2828 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	18.5	Xtriage
Anisotropy	0.215	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 47.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4004	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.21% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: IPA, CL, MG, SEP

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	0.18	0/1630	0.41	0/2189
1	B	0.18	0/1654	0.45	0/2222
2	C	0.16	0/112	0.33	0/148
2	D	0.18	0/95	0.42	0/125
All	All	0.18	0/3491	0.43	0/4684

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	1594	0	1651	11	0
1	B	1615	0	1675	34	0
2	C	124	0	109	2	0
2	D	107	0	99	13	0
3	A	4	0	8	0	0
3	B	4	0	8	0	0
4	A	2	0	0	1	0
4	B	2	0	0	0	0
5	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	1	0	0	0	0
6	A	279	0	0	1	0
6	B	234	0	0	7	1
6	C	15	0	0	0	0
6	D	21	0	0	2	0
All	All	4004	0	3550	51	1

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 7.

All (51) 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:B:133:VAL:HG21	1:B:189:ILE:HD11	1.41	1.02
1:B:190:GLN:HE21	2:D:451:LEU:HD11	1.42	0.82
1:B:138:ARG:NH1	2:D:449:ASP:OD2	2.13	0.81
1:A:18:GLU:OE1	6:A:1101:HOH:O	1.97	0.81
1:B:100:TYR:OH	6:B:401:HOH:O	1.99	0.80
1:B:133:VAL:HG13	6:B:403:HOH:O	1.84	0.76
1:B:42[B]:ASN:ND2	6:B:404:HOH:O	2.24	0.71
1:B:138:ARG:HB3	2:D:450:LEU:HD21	1.73	0.70
1:B:133:VAL:HG22	1:B:168:ALA:HA	1.72	0.70
1:B:173:ILE:HG12	1:B:186:PHE:HB3	1.74	0.69
1:A:128:LYS:HE3	1:A:154:ILE:HD11	1.75	0.68
2:D:445:ASP:HB3	2:D:448:ARG:NH1	2.08	0.67
1:B:190:GLN:NE2	2:D:451:LEU:HD11	2.12	0.64
2:D:445:ASP:HB3	2:D:448:ARG:HH11	1.61	0.64
1:B:25:LYS:NZ	6:B:402:HOH:O	2.17	0.63
1:B:128[B]:LYS:HE2	1:B:154:ILE:HD11	1.86	0.58
2:D:448:ARG:HH11	2:D:448:ARG:HG2	1.71	0.56
1:A:78:LYS:HD3	6:B:532:HOH:O	2.06	0.55
1:B:121:PRO:HB2	1:B:150:LYS:HG3	1.88	0.54
2:D:447:ILE:O	2:D:451:LEU:HD13	2.09	0.53
1:B:132:LEU:HD11	1:B:182:PHE:CE2	2.43	0.53
1:B:128[B]:LYS:NZ	1:B:161:SER:O	2.35	0.53
1:B:168:ALA:HB1	1:B:172:ARG:HG2	1.90	0.52
1:A:179:LYS:HE3	1:A:186:PHE:HB2	1.91	0.52
1:B:94:ILE:HD11	1:B:106:SER:HB3	1.92	0.51
1:B:169:SER:OG	6:B:403:HOH:O	2.19	0.51
1:A:174:LYS:HB2	1:A:179:LYS:NZ	2.28	0.49
1:B:136:ASP:H	1:B:139:SER:HB3	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:130:VAL:HG22	1:B:154:ILE:HG13	1.94	0.48
1:A:137:LYS:NZ	2:C:450:LEU:HA	2.31	0.46
1:B:142:LEU:HD11	2:D:450:LEU:HG	1.97	0.46
1:A:170:ASN:HB3	1:A:173:ILE:HD11	1.97	0.46
1:B:176:GLU:HB2	1:B:186:PHE:CE2	2.51	0.45
1:B:99:ARG:NH2	1:B:100:TYR:OH	2.50	0.44
1:B:170:ASN:O	1:B:174:LYS:HD3	2.18	0.44
1:A:131:LEU:HD23	1:A:167:ILE:HB	2.00	0.44
1:B:125:HIS:ND1	6:B:406:HOH:O	2.36	0.44
1:B:17:MET:CE	2:D:440:ASN:HB2	2.48	0.43
1:B:47:ILE:HG13	1:B:73:ILE:HD11	2.00	0.43
2:D:438:SER:N	6:D:502:HOH:O	2.51	0.43
1:B:138:ARG:HB3	2:D:450:LEU:CD2	2.45	0.43
1:A:137:LYS:NZ	2:C:449:ASP:O	2.37	0.43
1:B:93:LYS:HE3	1:B:95:GLU:OE2	2.19	0.42
1:B:169:SER:O	1:B:173:ILE:HG13	2.20	0.41
1:B:138:ARG:HD3	1:B:138:ARG:HA	1.77	0.41
1:B:170:ASN:O	1:B:173:ILE:HB	2.21	0.41
1:A:20:LYS:NZ	4:A:1002:CL:CL	2.77	0.41
2:D:446:ILE:O	2:D:450:LEU:HD23	2.21	0.41
1:B:17:MET:HE1	6:D:501:HOH:O	2.20	0.41
1:B:190:GLN:H	1:B:190:GLN:CD	2.28	0.40
1:A:146:LEU:HD21	1:A:192:LEU:HD21	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B:549:HOH:O	6:B:607:HOH:O[1_655]	2.12	0.08

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	194/207 (94%)	192 (99%)	2 (1%)	0	100	100
1	B	197/207 (95%)	194 (98%)	3 (2%)	0	100	100
2	C	13/17 (76%)	11 (85%)	2 (15%)	0	100	100
2	D	11/17 (65%)	9 (82%)	2 (18%)	0	100	100
All	All	415/448 (93%)	406 (98%)	9 (2%)	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	171/181 (94%)	171 (100%)	0	100	100
1	B	174/181 (96%)	174 (100%)	0	100	100
2	C	15/16 (94%)	15 (100%)	0	100	100
2	D	13/16 (81%)	13 (100%)	0	100	100
All	All	373/394 (95%)	373 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	190	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	SEP	D	442	2	8,9,10	1.48	1 (12%)	7,12,14	0.71	0
2	SEP	C	442	2	8,9,10	1.53	1 (12%)	7,12,14	0.84	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	SEP	D	442	2	-	2/6/8/10	-
2	SEP	C	442	2	-	1/6/8/10	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	442	SEP	P-O1P	3.51	1.61	1.50
2	D	442	SEP	P-O1P	3.24	1.60	1.50

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	442	SEP	C-CA-CB-OG
2	D	442	SEP	C-CA-CB-OG
2	D	442	SEP	CB-OG-P-O1P

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 7 are monoatomic - leaving 2 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	IPA	B	301	-	3,3,3	0.53	0	3,3,3	0.32	0
3	IPA	A	1001	-	3,3,3	0.53	0	3,3,3	0.34	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.

No monomer is involved in short contacts.

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	196/207 (94%)	0.00	6 (3%) 51 55	12, 20, 40, 59	0
1	B	197/207 (95%)	0.36	20 (10%) 12 12	12, 21, 81, 133	2 (1%)
2	C	15/17 (88%)	1.53	5 (33%) 1 1	21, 34, 83, 95	0
2	D	13/17 (76%)	1.70	4 (30%) 1 1	24, 40, 64, 84	0
All	All	421/448 (93%)	0.28	35 (8%) 17 18	12, 21, 69, 133	2 (0%)

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	133	VAL	7.3
1	B	173	ILE	5.7
1	B	132	LEU	5.6
1	B	171	ALA	5.5
2	C	439	CYS	4.6
1	B	175	ALA	4.4
1	B	168	ALA	4.4
2	D	439	CYS	4.3
1	B	169	SER	4.3
2	C	440	ASN	4.2
2	D	450	LEU	3.9
1	B	170	ASN	3.9
2	D	451	LEU	3.6
1	B	186	PHE	3.5
1	B	174	LYS	3.4
1	B	138	ARG	3.3
1	B	134	ARG	3.3
1	A	135	THR	3.1
1	B	5	THR	3.1
1	A	6	PRO	3.1
2	C	437	ASP	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	177	LYS	2.8
1	B	131	LEU	2.8
1	B	135	THR	2.7
1	B	178	GLU	2.6
1	B	200	LEU	2.5
1	A	138	ARG	2.4
2	C	441	SER	2.4
1	B	172	ARG	2.3
1	A	175	ALA	2.3
1	B	188	PRO	2.2
1	A	173	ILE	2.2
1	A	200	LEU	2.1
2	D	438	SER	2.1
2	C	438	SER	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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($\text{\AA}^2$)	Q<0.9
2	SEP	D	442	10/11	0.98	0.07	18,20,25,32	0
2	SEP	C	442	10/11	0.99	0.06	15,18,22,28	0

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($\text{\AA}^2$)	Q<0.9
3	IPA	B	301	4/4	0.90	0.13	20,26,27,32	0
5	MG	A	1005	1/1	0.95	0.14	38,38,38,38	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	MG	A	1004	1/1	0.96	0.20	36,36,36,36	0
3	IPA	A	1001	4/4	0.96	0.08	19,20,24,29	0
4	CL	A	1003	1/1	0.97	0.08	34,34,34,34	0
4	CL	B	303	1/1	0.98	0.05	27,27,27,27	0
5	MG	B	304	1/1	0.98	0.10	24,24,24,24	0
4	CL	A	1002	1/1	1.00	0.03	19,19,19,19	0
4	CL	B	302	1/1	1.00	0.02	17,17,17,17	0

6.5 Other polymers [i](#)

There are no such residues in this entry.