



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:30 AM GMT

PDB ID : 2R25
Title : Complex of YPD1 and SLN1-R1 with bound Mg²⁺ and BeF₃-
Authors : Copeland, D.M.; Zhao, X.; Soares, A.S.; West, A.H.
Deposited on : 2007-08-24
Resolution : 1.70 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
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) : trunk26865

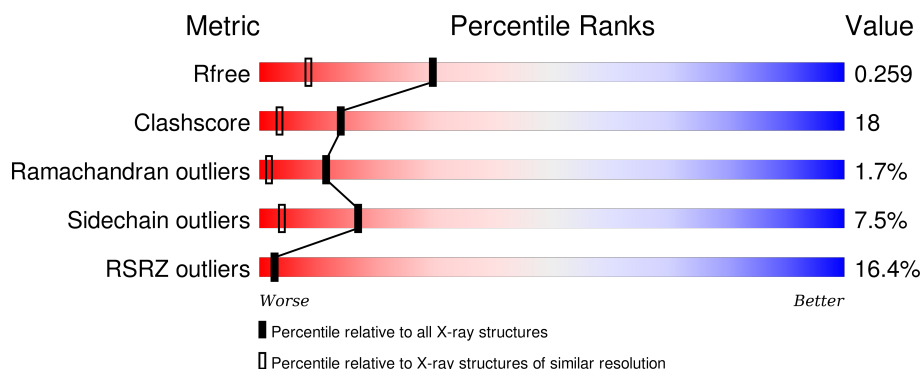
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	3190 (1.70-1.70)
Clashscore	102246	3585 (1.70-1.70)
Ramachandran outliers	100387	3527 (1.70-1.70)
Sidechain outliers	100360	3527 (1.70-1.70)
RSRZ outliers	91569	3200 (1.70-1.70)

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	167	<div> <div>14%</div> <div>82%</div> <div>12%</div> <div>5%</div> <div>..</div> </div>
2	B	133	<div> <div>20%</div> <div>75%</div> <div>20%</div> <div>5%</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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	NA	B	5	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 2723 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 Phosphorelay intermediate protein YPD1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	166	Total	C	N	O	S	0	6	0
			1383	871	232	275	5			

- Molecule 2 is a protein called Osmosensing histidine protein kinase SLN1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	133	Total	C	N	O	S	0	8	0
			1096	694	182	210	10			

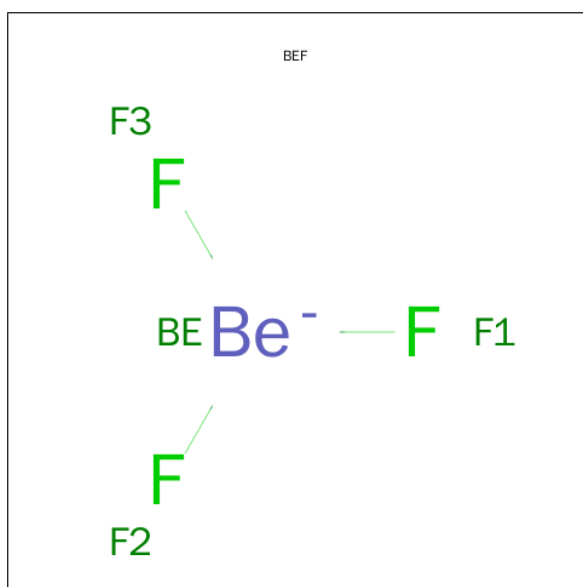
- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Na	0	0
			1	1		

- Molecule 5 is BERYLLIUM TRIFLUORIDE ION (three-letter code: BEF) (formula: BeF₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	Be	F	0	0
			4	1	3		

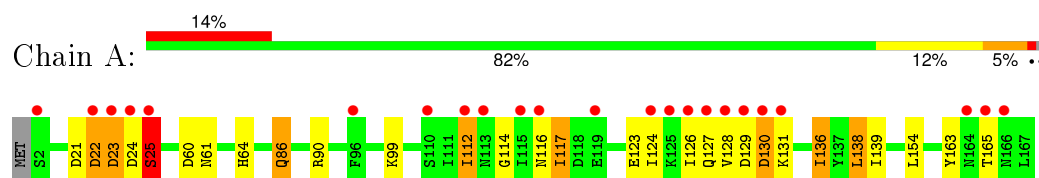
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	157	Total	O	0	0
			157	157		
6	B	81	Total	O	0	0
			81	81		

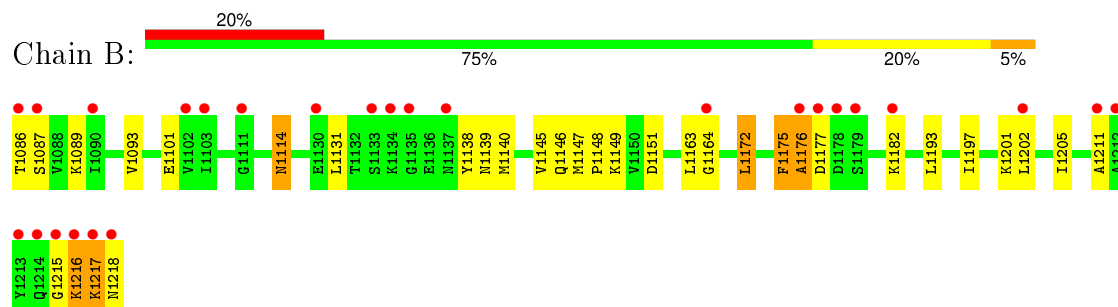
3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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: Phosphorelay intermediate protein YPD1



- Molecule 2: Osmosensing histidine protein kinase SLN1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	52.01Å 75.54Å 98.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	17.75 – 1.70 17.75 – 1.70	Depositor EDS
% Data completeness (in resolution range)	99.2 (17.75-1.70) 99.2 (17.75-1.70)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.61 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.196 , 0.240 0.231 , 0.259	Depositor DCC
R_{free} test set	2168 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	30.7	Xtriage
Anisotropy	0.101	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 47.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 42989 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2723	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.19% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: NA, MG, BEF

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	0.87	0/1406	0.87	2/1894 (0.1%)
2	B	0.76	1/1111 (0.1%)	0.80	2/1490 (0.1%)
All	All	0.82	1/2517 (0.0%)	0.84	4/3384 (0.1%)

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
2	B	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1138	TYR	C-O	5.41	1.33	1.23

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1140	MET	CB-CG-SD	-6.49	92.92	112.40
1	A	138	LEU	CB-CG-CD1	6.11	121.39	111.00
1	A	138	LEU	CA-CB-CG	5.64	128.26	115.30
2	B	1138	TYR	N-CA-C	-5.09	97.26	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	1164	GLY	Peptide
2	B	1175	PHE	Peptide

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	1383	0	1377	33	0
2	B	1096	0	1124	60	1
3	B	1	0	0	0	0
4	B	1	0	0	0	0
5	B	4	0	0	0	0
6	A	157	0	0	11	1
6	B	81	0	0	2	0
All	All	2723	0	2501	92	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 18.

All (92) 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)
2:B:1216:LYS:CB	2:B:1217:LYS:HB2	1.24	1.63
2:B:1216:LYS:HA	2:B:1217:LYS:CD	1.34	1.58
2:B:1216:LYS:CA	2:B:1217:LYS:HD3	1.51	1.37
2:B:1216:LYS:HB3	2:B:1217:LYS:CB	1.59	1.28
2:B:1216:LYS:CB	2:B:1217:LYS:CB	2.19	1.15
1:A:128:VAL:H	1:A:129:ASP:HB2	0.99	1.14
2:B:1172:LEU:CD2	2:B:1197[B]:ILE:HD11	1.81	1.10
2:B:1216:LYS:CA	2:B:1217:LYS:HB2	1.81	1.09
2:B:1172:LEU:HD21	2:B:1197[B]:ILE:HD11	1.33	1.08
1:A:128:VAL:N	1:A:129:ASP:HB2	1.80	0.96
1:A:24:ASP:O	1:A:25:SER:HB3	1.68	0.94
2:B:1172:LEU:HD21	2:B:1197[B]:ILE:CD1	1.96	0.94
2:B:1216:LYS:HA	2:B:1217:LYS:CG	2.00	0.92
2:B:1216:LYS:HB2	2:B:1217:LYS:HB2	1.50	0.91
2:B:1216:LYS:CA	2:B:1217:LYS:CB	2.44	0.91

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:VAL:H	1:A:129:ASP:CB	1.86	0.88
2:B:1216:LYS:HB3	2:B:1217:LYS:HB2	0.88	0.86
1:A:22:ASP:HA	6:A:314:HOH:O	1.73	0.86
2:B:1175:PHE:HB3	2:B:1176:ALA:HB3	1.63	0.81
2:B:1175:PHE:HA	2:B:1176:ALA:HB2	1.63	0.80
2:B:1086[A]:THR:HG22	2:B:1087[A]:SER:H	1.47	0.79
1:A:124:ILE:HD11	6:A:270:HOH:O	1.82	0.79
2:B:1216:LYS:CA	2:B:1217:LYS:CD	2.31	0.78
1:A:86:GLN:HE22	1:A:90:ARG:HD2	1.49	0.76
2:B:1172:LEU:HD23	2:B:1197[B]:ILE:HD11	1.69	0.75
2:B:1217:LYS:O	2:B:1218:ASN:OD1	2.05	0.75
2:B:1172:LEU:HD21	2:B:1197[B]:ILE:CG1	2.18	0.73
2:B:1086[B]:THR:HG23	2:B:1087[B]:SER:N	2.04	0.72
6:A:292:HOH:O	2:B:1176:ALA:HA	1.88	0.72
2:B:1086[A]:THR:O	2:B:1087[A]:SER:CB	2.41	0.69
2:B:1215:GLY:O	2:B:1217:LYS:CE	2.41	0.69
2:B:1216:LYS:HA	2:B:1217:LYS:CB	2.17	0.68
1:A:23:ASP:HB3	6:A:296:HOH:O	1.94	0.68
1:A:117:ILE:HG22	6:A:282:HOH:O	1.93	0.67
2:B:1101[A]:GLU:OE1	6:B:238:HOH:O	2.13	0.67
1:A:112:ILE:CD1	1:A:154:LEU:HD21	2.25	0.66
2:B:1215:GLY:O	2:B:1217:LYS:HE2	1.97	0.65
2:B:1216:LYS:HB3	2:B:1217:LYS:CG	2.26	0.65
2:B:1215:GLY:O	2:B:1217:LYS:HD3	1.97	0.64
2:B:1216:LYS:HA	2:B:1217:LYS:HD3	0.67	0.64
1:A:99:LYS:HD3	1:A:117:ILE:CG1	2.28	0.62
2:B:1193:LEU:HD21	2:B:1205:ILE:HG13	1.80	0.62
2:B:1217:LYS:C	2:B:1218:ASN:OD1	2.39	0.61
1:A:116:ASN:HA	6:A:282:HOH:O	1.99	0.61
1:A:86:GLN:NE2	1:A:90:ARG:HD2	2.15	0.60
2:B:1216:LYS:N	2:B:1217:LYS:HD3	2.16	0.60
2:B:1175:PHE:CA	2:B:1176:ALA:CB	2.80	0.59
2:B:1175:PHE:CA	2:B:1176:ALA:HB2	2.31	0.59
1:A:22:ASP:OD1	1:A:23:ASP:N	2.38	0.56
2:B:1215:GLY:O	2:B:1217:LYS:CD	2.53	0.55
2:B:1086[B]:THR:HG23	2:B:1087[B]:SER:H	1.69	0.55
2:B:1175:PHE:CB	2:B:1176:ALA:HB3	2.37	0.54
1:A:21:ASP:O	1:A:22:ASP:CG	2.46	0.54
6:A:292:HOH:O	2:B:1176:ALA:CB	2.55	0.54
1:A:112:ILE:HD11	1:A:154:LEU:HD21	1.89	0.54
1:A:21:ASP:O	1:A:22:ASP:OD1	2.26	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:LYS:HD3	1:A:117:ILE:HG12	1.90	0.52
1:A:114:GLY:CA	6:A:294:HOH:O	2.57	0.52
2:B:1193:LEU:CD2	2:B:1201:LYS:HE2	2.40	0.52
2:B:1197[B]:ILE:N	2:B:1197[B]:ILE:HD13	2.24	0.51
2:B:1172:LEU:HD21	2:B:1197[B]:ILE:HG13	1.93	0.51
2:B:1175:PHE:CB	2:B:1176:ALA:CB	2.90	0.50
2:B:1175:PHE:HD1	2:B:1176:ALA:HB2	1.77	0.50
1:A:64:HIS:HB2	1:A:86:GLN:HG3	1.94	0.50
2:B:1175:PHE:HB3	2:B:1176:ALA:CB	2.38	0.49
2:B:1217:LYS:O	2:B:1218:ASN:CG	2.51	0.49
2:B:1215:GLY:C	2:B:1217:LYS:HD3	2.32	0.49
1:A:21:ASP:OD2	1:A:25:SER:HA	2.12	0.49
1:A:86:GLN:HE22	1:A:90:ARG:CD	2.22	0.49
1:A:112:ILE:O	1:A:112:ILE:HG23	2.14	0.47
2:B:1145:VAL:O	2:B:1151:ASP:HB2	2.15	0.47
2:B:1086[A]:THR:HG22	2:B:1087[A]:SER:N	2.22	0.47
6:A:292:HOH:O	2:B:1176:ALA:HB2	2.14	0.46
2:B:1089:LYS:H	2:B:1139:ASN:ND2	2.13	0.45
1:A:163:TYR:HB3	1:A:165:THR:HG22	1.99	0.45
1:A:127:GLN:HB2	1:A:130:ASP:HB2	1.99	0.45
1:A:99:LYS:HD3	1:A:117:ILE:HG13	1.98	0.45
1:A:24:ASP:O	1:A:25:SER:CB	2.50	0.45
2:B:1086[B]:THR:CG2	2:B:1087[B]:SER:N	2.71	0.44
2:B:1172:LEU:CD2	2:B:1197[B]:ILE:CD1	2.66	0.44
1:A:114:GLY:N	6:A:294:HOH:O	2.46	0.44
2:B:1148:PRO:O	2:B:1149:LYS:HB2	2.18	0.43
1:A:136[A]:ILE:H	1:A:136[A]:ILE:HG13	1.62	0.43
2:B:1175:PHE:CD1	2:B:1176:ALA:HB2	2.53	0.43
1:A:86:GLN:C	1:A:86:GLN:HE21	2.22	0.43
2:B:1177:ASP:HB2	6:B:228:HOH:O	2.18	0.43
2:B:1093:VAL:HG12	2:B:1147:MET:SD	2.59	0.42
2:B:1139:ASN:HD22	2:B:1139:ASN:H	1.66	0.42
1:A:114:GLY:HA3	6:A:294:HOH:O	2.17	0.42
2:B:1089:LYS:HE2	2:B:1114:ASN:OD1	2.20	0.41
1:A:61:ASN:HA	2:B:1146:GLN:HE22	1.84	0.41
1:A:112:ILE:HD12	1:A:112:ILE:HA	1.56	0.41

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)
2:B:1101[B]:GLU:OE2	6:A:203:HOH:O[4_455]	2.07	0.13

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	170/167 (102%)	160 (94%)	8 (5%)	2 (1%)	16	3
2	B	138/133 (104%)	131 (95%)	4 (3%)	3 (2%)	8	1
All	All	308/300 (103%)	291 (94%)	12 (4%)	5 (2%)	11	2

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	1176	ALA
2	B	1211	ALA
2	B	1217	LYS
1	A	25	SER
1	A	22	ASP

5.3.2 Protein sidechains [i](#)

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	156/152 (103%)	142 (91%)	14 (9%)	12	2
2	B	124/118 (105%)	117 (94%)	7 (6%)	26	9
All	All	280/270 (104%)	259 (92%)	21 (8%)	17	4

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

Mol	Chain	Res	Type
1	A	23	ASP
1	A	25	SER
1	A	60	ASP
1	A	86	GLN
1	A	112	ILE
1	A	117	ILE
1	A	123	GLU
1	A	126	ILE
1	A	130	ASP
1	A	131	LYS
1	A	136[A]	ILE
1	A	136[B]	ILE
1	A	138	LEU
1	A	139	ILE
2	B	1114	ASN
2	B	1131	LEU
2	B	1163	LEU
2	B	1172	LEU
2	B	1182	LYS
2	B	1202	LEU
2	B	1216	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	GLN
1	A	86	GLN
1	A	94	HIS
1	A	113	ASN
2	B	1139	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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	$\# Z > 2$
5	BEF	B	2	2	0,3,3	0.00	-	0,3,3	0.00	-

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	BEF	B	2	2	-	0/0/0/0	0/0/0/0

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

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	166/167 (99%)	0.89	23 (13%) 4 4	22, 30, 57, 78	5 (3%)
2	B	133/133 (100%)	1.20	26 (19%) 1 1	21, 37, 68, 78	5 (3%)
All	All	299/300 (99%)	1.03	49 (16%) 2 2	21, 33, 67, 78	10 (3%)

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	126	ILE	13.7
2	B	1218	ASN	10.9
1	A	128	VAL	10.5
2	B	1217	LYS	8.2
2	B	1135	GLY	7.2
1	A	130	ASP	7.2
1	A	129	ASP	6.9
2	B	1086[A]	THR	6.5
2	B	1212	ALA	5.7
2	B	1215	GLY	5.5
2	B	1179	SER	4.8
1	A	125	LYS	4.8
1	A	22	ASP	4.8
1	A	127	GLN	4.7
2	B	1211	ALA	4.6
1	A	23	ASP	4.6
1	A	113	ASN	4.4
1	A	24	ASP	4.3
1	A	115	ILE	4.2
2	B	1214	GLN	4.2
1	A	124	ILE	4.0
2	B	1216	LYS	3.9
1	A	164	ASN	3.7
2	B	1177	ASP	3.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	B	1178	ASP	3.3
2	B	1134	LYS	3.3
2	B	1130	GLU	3.2
1	A	25	SER	3.2
2	B	1164	GLY	3.0
2	B	1103	ILE	3.0
1	A	131	LYS	2.9
2	B	1090	ILE	2.8
2	B	1213	TYR	2.8
1	A	110	SER	2.6
2	B	1137	ASN	2.5
2	B	1087[A]	SER	2.5
1	A	2	SER	2.5
1	A	116	ASN	2.4
2	B	1202	LEU	2.4
1	A	166[A]	ASN	2.4
1	A	165	THR	2.4
2	B	1176	ALA	2.2
1	A	96	PHE	2.2
2	B	1111	GLY	2.2
1	A	112	ILE	2.2
2	B	1102	VAL	2.2
2	B	1133	SER	2.1
1	A	119	GLU	2.1
2	B	1182	LYS	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
4	NA	B	5	1/1	0.64	0.22	3.63	39,39,39,39	1
5	BEF	B	2	4/4	0.94	0.09	-0.66	25,25,26,28	0
3	MG	B	1	1/1	1.00	0.02	-4.21	25,25,25,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.