wwPDB X-ray Structure Validation Summary Report (i)

May 13, 2020 – 08:09 am BST

PDB ID	:	4BWG
Title	:	Structural basis of subtilase cytotoxin SubAB assembly
Authors	:	Le Nours, J.; Paton, A.W.; Byres, E.; Troy, S.; Herdman, B.P.; Johnson, M.D.; Paton, J.C.; Rossjohn, J.; Beddoe, T.
Deposited on	:	2013-07-02
Resolution	:	2.60 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

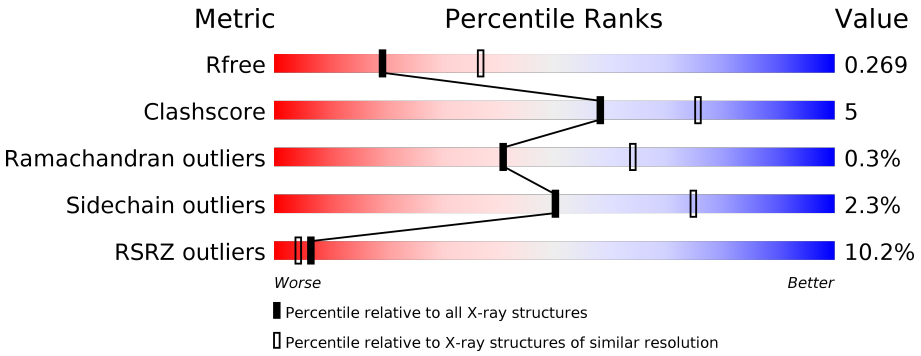
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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 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 <=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	347	83% 9% 7%
1	G	347	% 81% 11% 8%
2	B	120	13% 69% 11% 20%
2	C	120	5% 86% 8% • 5%
2	D	120	18% 68% 13% 18%
2	E	120	6% 79% 10% 11%

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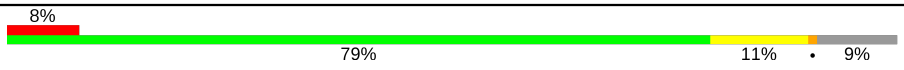
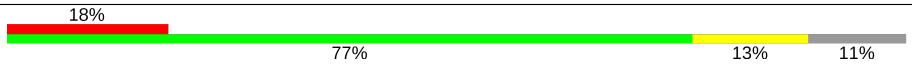
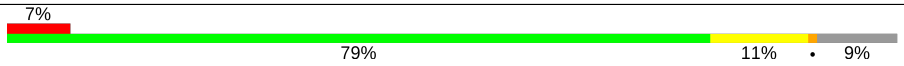
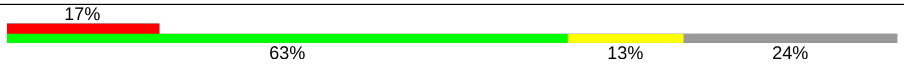
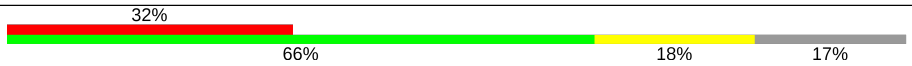
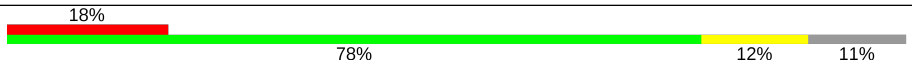
Mol	Chain	Length	Quality of chain
2	F	120	
2	H	120	
2	I	120	
2	J	120	
2	K	120	
2	L	120	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12621 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 SUBA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	321	Total	C	N	O	S	0	1	0
			2378	1503	416	450	9			
1	G	320	Total	C	N	O	S	0	0	0
			2366	1496	412	450	8			

- Molecule 2 is a protein called SUBTILASE CYTOTOXIN, SUBUNIT B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	96	Total	C	N	O	S	0	2	1
			735	475	115	138	7			
2	C	114	Total	C	N	O	S	0	2	0
			868	556	136	168	8			
2	D	98	Total	C	N	O	S	0	2	1
			750	483	117	143	7			
2	E	107	Total	C	N	O	S	0	2	1
			818	520	129	161	8			
2	F	109	Total	C	N	O	S	0	2	1
			802	515	128	151	8			
2	H	107	Total	C	N	O	S	0	0	1
			780	504	124	146	6			
2	I	109	Total	C	N	O	S	0	2	1
			829	533	129	159	8			
2	J	91	Total	C	N	O	S	0	0	1
			656	422	105	124	5			
2	K	100	Total	C	N	O	S	0	2	1
			753	484	119	143	7			
2	L	107	Total	C	N	O	S	0	2	1
			791	511	127	145	8			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	119	LEU	-	expression tag	UNP Q3ZTX8

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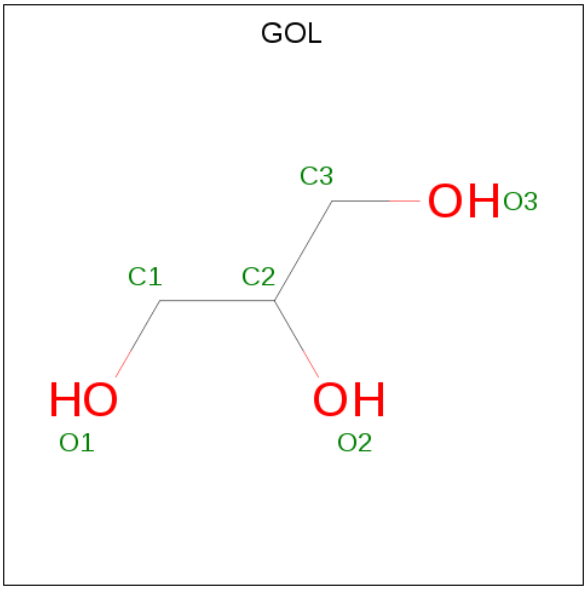
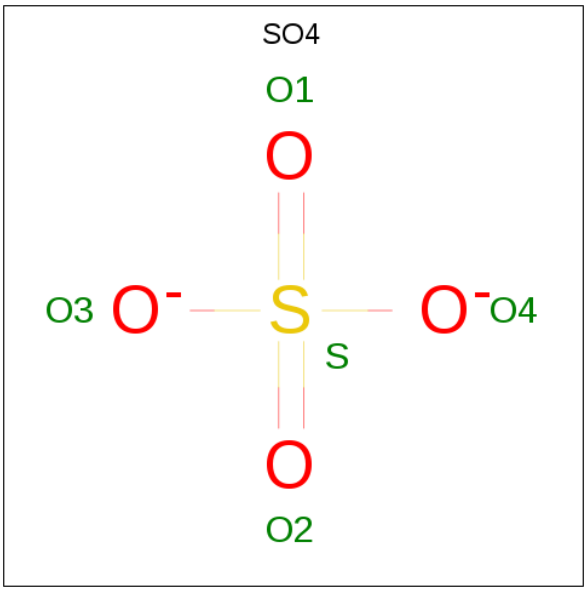
Chain	Residue	Modelled	Actual	Comment	Reference
B	120	GLU	-	expression tag	UNP Q3ZTX8
C	119	LEU	-	expression tag	UNP Q3ZTX8
C	120	GLU	-	expression tag	UNP Q3ZTX8
D	119	LEU	-	expression tag	UNP Q3ZTX8
D	120	GLU	-	expression tag	UNP Q3ZTX8
E	119	LEU	-	expression tag	UNP Q3ZTX8
E	120	GLU	-	expression tag	UNP Q3ZTX8
F	119	LEU	-	expression tag	UNP Q3ZTX8
F	120	GLU	-	expression tag	UNP Q3ZTX8
H	119	LEU	-	expression tag	UNP Q3ZTX8
H	120	GLU	-	expression tag	UNP Q3ZTX8
I	119	LEU	-	expression tag	UNP Q3ZTX8
I	120	GLU	-	expression tag	UNP Q3ZTX8
J	119	LEU	-	expression tag	UNP Q3ZTX8
J	120	GLU	-	expression tag	UNP Q3ZTX8
K	119	LEU	-	expression tag	UNP Q3ZTX8
K	120	GLU	-	expression tag	UNP Q3ZTX8
L	119	LEU	-	expression tag	UNP Q3ZTX8
L	120	GLU	-	expression tag	UNP Q3ZTX8

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	6	3	3	0	0

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



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

- Molecule 5 is water.

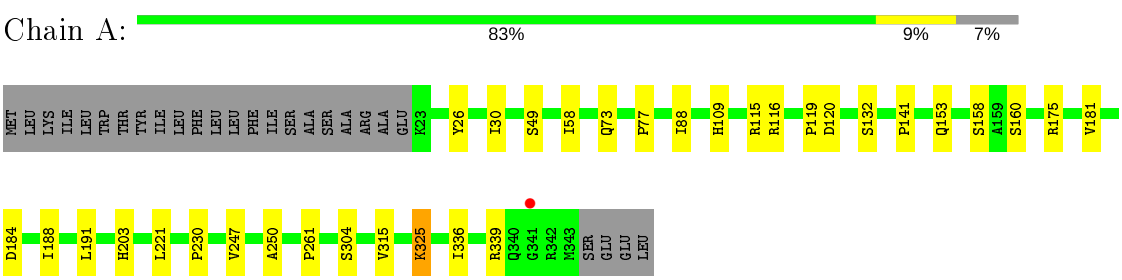
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	28	Total O 28 28	0	0
5	B	3	Total O 3 3	0	0
5	C	9	Total O 9 9	0	0
5	E	3	Total O 3 3	0	0
5	G	23	Total O 23 23	0	0
5	H	6	Total O 6 6	0	0
5	I	3	Total O 3 3	0	0
5	J	2	Total O 2 2	0	0
5	K	3	Total O 3 3	0	0
5	L	4	Total O 4 4	0	0

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.

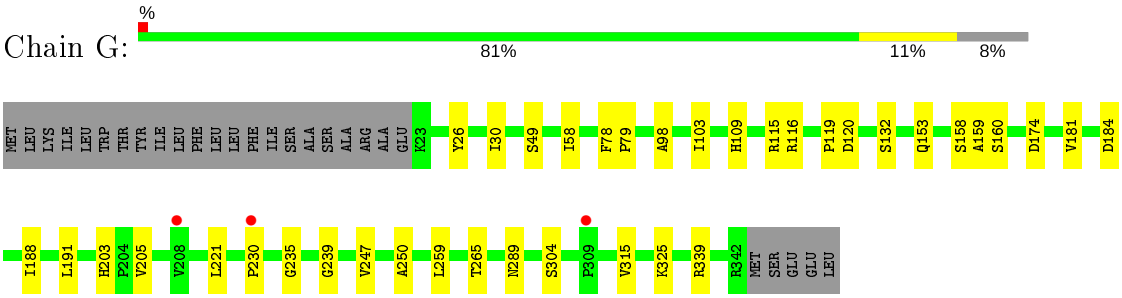
- Molecule 1: SUBA

Chain A:



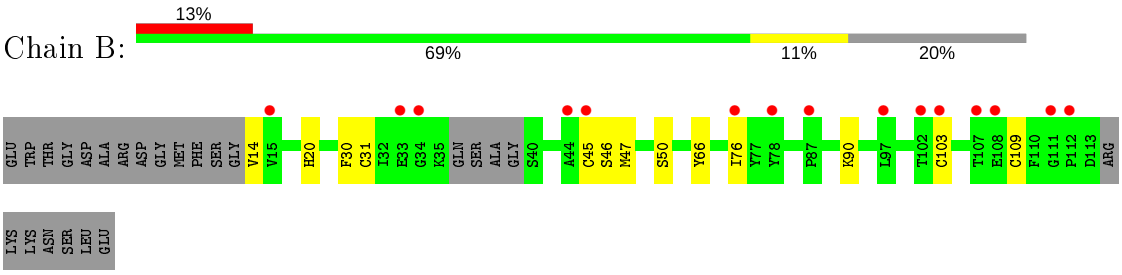
- Molecule 1: SUBA

Chain G:



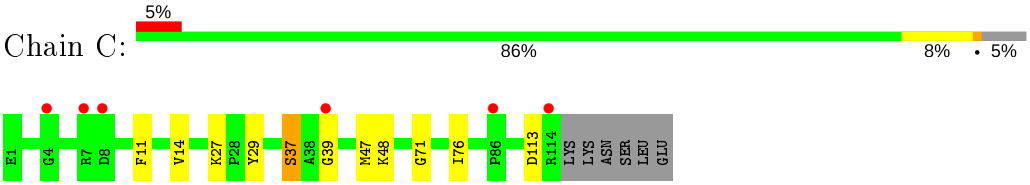
- Molecule 2: SUBTILASE CYTOTOXIN, SUBUNIT B

Chain B:



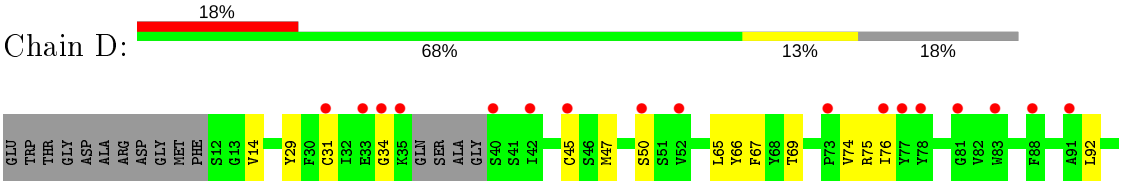
- Molecule 2: SUBTILASE CYTOTOXIN, SUBUNIT B

Chain C:



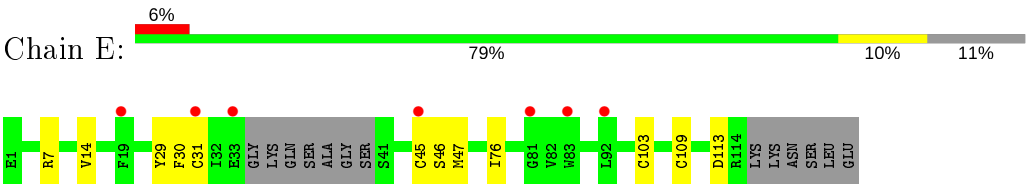
• Molecule 2: SUBTILASE CYTOTOXIN, SUBUNIT B

Chain D:



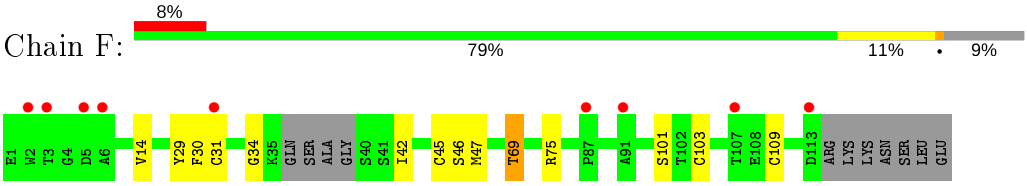
• Molecule 2: SUBTILASE CYTOTOXIN, SUBUNIT B

Chain E:



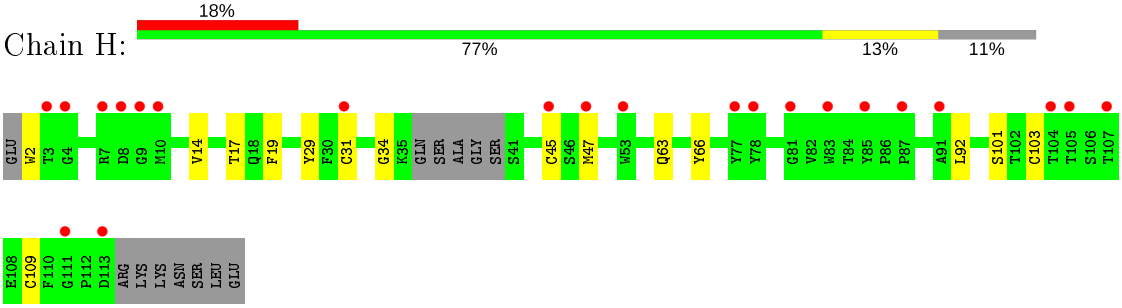
• Molecule 2: SUBTILASE CYTOTOXIN, SUBUNIT B

Chain F:



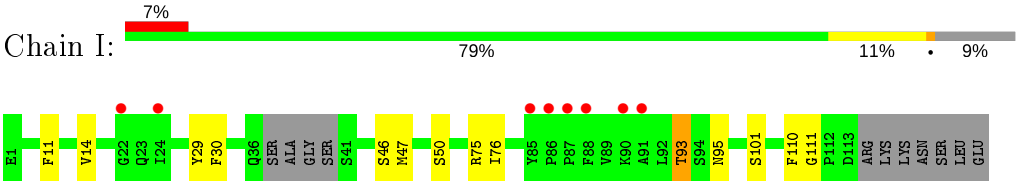
• Molecule 2: SUBTILASE CYTOTOXIN, SUBUNIT B

Chain H:



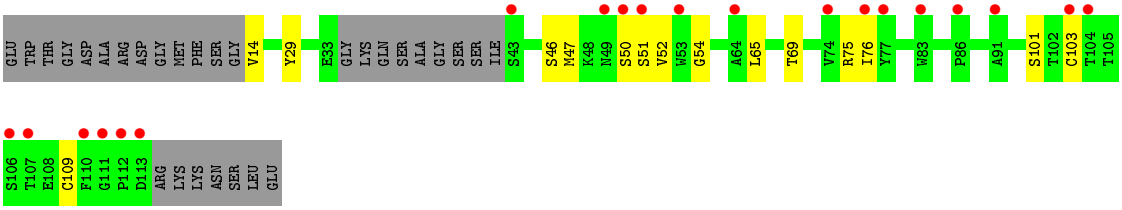
• Molecule 2: SUBTILASE CYTOTOXIN, SUBUNIT B

Chain I:



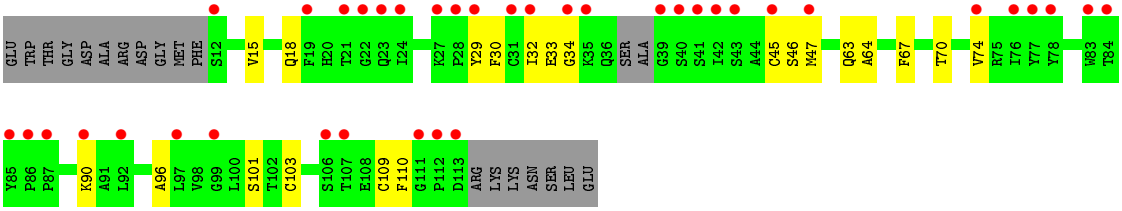
• Molecule 2: SUBTILASE CYTOTOXIN, SUBUNIT B

Chain J:



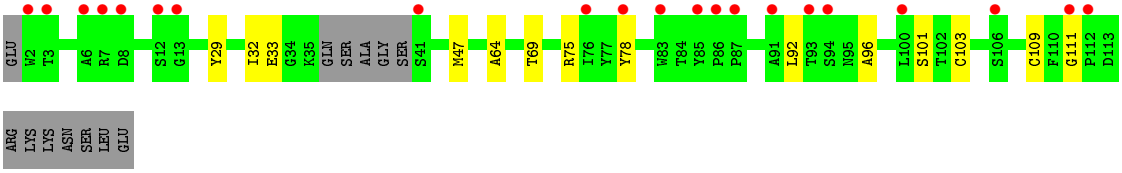
• Molecule 2: SUBTILASE CYTOTOXIN, SUBUNIT B

Chain K:



• Molecule 2: SUBTILASE CYTOTOXIN, SUBUNIT B

Chain L:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	59.21Å 81.71Å 227.79Å 90.00° 100.24° 90.00°	Depositor
Resolution (Å)	38.07 – 2.60 66.03 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.9 (38.07-2.60) 99.9 (66.03-2.60)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.14 (at 2.62Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.219 , 0.250 0.237 , 0.269	Depositor DCC
R_{free} test set	3351 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å^2)	35.4	Xtriage
Anisotropy	0.311	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 60.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.025 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	12621	wwPDB-VP
Average B, all atoms (Å^2)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.22% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: GOL, 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	0.54	0/2427	0.72	0/3305
1	G	0.54	0/2414	0.71	0/3287
2	B	0.49	0/757	0.60	0/1037
2	C	0.53	0/894	0.66	0/1223
2	D	0.48	0/772	0.61	0/1056
2	E	0.46	0/842	0.61	0/1153
2	F	0.52	0/826	0.64	0/1133
2	H	0.54	0/804	0.65	0/1104
2	I	0.54	0/854	0.66	0/1168
2	J	0.58	0/677	0.65	0/931
2	K	0.54	0/775	0.59	0/1059
2	L	0.51	0/815	0.65	0/1117
All	All	0.53	0/12857	0.67	0/17573

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

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	2378	0	2330	19	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	2366	0	2332	17	0
2	B	735	0	685	8	0
2	C	868	0	786	5	0
2	D	750	0	699	13	0
2	E	818	0	732	6	0
2	F	802	0	709	9	0
2	H	780	0	699	10	0
2	I	829	0	751	9	0
2	J	656	0	557	11	0
2	K	753	0	684	12	0
2	L	791	0	702	12	0
3	A	6	0	8	3	0
4	G	5	0	0	0	0
5	A	28	0	0	0	0
5	B	3	0	0	0	0
5	C	9	0	0	0	0
5	E	3	0	0	0	0
5	G	23	0	0	1	0
5	H	6	0	0	0	0
5	I	3	0	0	0	0
5	J	2	0	0	0	0
5	K	3	0	0	0	0
5	L	4	0	0	0	0
All	All	12621	0	11674	114	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 5.

The worst 5 of 114 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:175:ARG:HE	3:A:1344:GOL:H12	1.35	0.90
2:D:31[B]:CYS:HG	2:D:45[B]:CYS:HG	1.23	0.82
2:J:51:SER:HA	2:J:54:GLY:H	1.47	0.79
1:G:339:ARG:HE	2:L:69:THR:HG21	1.50	0.75
2:I:93:THR:HG23	2:I:95:ASN:H	1.52	0.73

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	320/347 (92%)	308 (96%)	11 (3%)	1 (0%)	41 64
1	G	318/347 (92%)	306 (96%)	10 (3%)	2 (1%)	25 47
2	B	94/120 (78%)	93 (99%)	1 (1%)	0	100 100
2	C	114/120 (95%)	110 (96%)	3 (3%)	1 (1%)	17 35
2	D	96/120 (80%)	94 (98%)	2 (2%)	0	100 100
2	E	105/120 (88%)	102 (97%)	2 (2%)	1 (1%)	15 32
2	F	107/120 (89%)	104 (97%)	3 (3%)	0	100 100
2	H	103/120 (86%)	100 (97%)	3 (3%)	0	100 100
2	I	107/120 (89%)	105 (98%)	2 (2%)	0	100 100
2	J	87/120 (72%)	84 (97%)	3 (3%)	0	100 100
2	K	98/120 (82%)	96 (98%)	2 (2%)	0	100 100
2	L	105/120 (88%)	103 (98%)	2 (2%)	0	100 100
All	All	1654/1894 (87%)	1605 (97%)	44 (3%)	5 (0%)	41 64

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	184	ASP
1	G	184	ASP
2	E	113	ASP
2	C	37	SER
1	G	159	ALA

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	244/282 (86%)	237 (97%)	7 (3%)	42	68
1	G	245/282 (87%)	237 (97%)	8 (3%)	38	64
2	B	81/102 (79%)	80 (99%)	1 (1%)	71	87
2	C	90/102 (88%)	88 (98%)	2 (2%)	52	76
2	D	83/102 (81%)	82 (99%)	1 (1%)	71	87
2	E	87/102 (85%)	86 (99%)	1 (1%)	73	88
2	F	80/102 (78%)	78 (98%)	2 (2%)	47	73
2	H	78/102 (76%)	75 (96%)	3 (4%)	33	59
2	I	87/102 (85%)	85 (98%)	2 (2%)	50	75
2	J	63/102 (62%)	62 (98%)	1 (2%)	62	82
2	K	79/102 (78%)	78 (99%)	1 (1%)	69	86
2	L	77/102 (76%)	77 (100%)	0	100	100
All	All	1294/1584 (82%)	1265 (98%)	29 (2%)	50	76

5 of 29 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	F	69	THR
1	G	109	HIS
2	I	93	THR
1	G	49	SER
1	G	120	ASP

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

Mol	Chain	Res	Type
2	D	62	ASN

5.3.3 RNA

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)

2 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	$\#\|Z\| > 2$
3	GOL	A	1344	-	5,5,5	0.23	0	5,5,5	0.26	0
4	SO4	G	1343	-	4,4,4	0.14	0	6,6,6	0.28	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
3	GOL	A	1344	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1344	GOL	O1-C1-C2-C3
3	A	1344	GOL	O2-C2-C3-O3
3	A	1344	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
3	A	1344	GOL	C1-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1344	GOL	3	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, 95^{th} 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	321/347 (92%)	-0.02	1 (0%) 94 93	11, 38, 85, 109	0
1	G	320/347 (92%)	0.08	3 (0%) 84 82	12, 41, 83, 105	0
2	B	96/120 (80%)	0.99	15 (15%) 2 1	43, 68, 111, 138	0
2	C	114/120 (95%)	0.29	6 (5%) 26 20	14, 45, 83, 116	0
2	D	98/120 (81%)	1.31	21 (21%) 0 0	43, 77, 116, 133	0
2	E	107/120 (89%)	0.63	7 (6%) 18 14	22, 62, 87, 100	0
2	F	109/120 (90%)	0.52	9 (8%) 11 8	25, 57, 88, 101	0
2	H	107/120 (89%)	1.26	22 (20%) 1 0	49, 73, 100, 116	0
2	I	109/120 (90%)	0.44	8 (7%) 15 11	16, 52, 85, 108	0
2	J	91/120 (75%)	1.29	20 (21%) 0 0	44, 67, 92, 110	3 (3%)
2	K	100/120 (83%)	1.71	38 (38%) 0 0	30, 78, 99, 136	0
2	L	107/120 (89%)	1.02	21 (19%) 1 0	33, 68, 95, 114	0
All	All	1679/1894 (88%)	0.58	171 (10%) 6 4	11, 56, 95, 138	3 (0%)

The worst 5 of 171 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	K	22	GLY	6.9
2	K	78	TYR	6.1
2	J	76	ILE	5.6
2	L	76	ILE	4.9
2	H	111	GLY	4.7

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 (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, 95^{th} 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
3	GOL	A	1344	6/6	0.89	0.27	38,40,47,49	0
4	SO4	G	1343	5/5	0.94	0.19	62,63,67,67	0

6.5 Other polymers (i)

There are no such residues in this entry.