



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:27 AM GMT

PDB ID : 3P1U
Title : Crystal structure of a SusD homolog (BDI_0600) from Parabacteroides distasonis ATCC 8503 AT 2.05 Å resolution
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2010-09-30
Resolution : 2.05 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

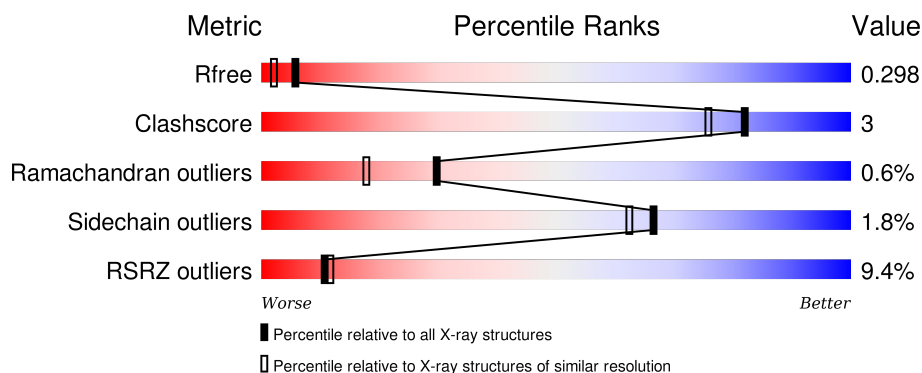
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	1192 (2.04-2.04)
Clashscore	102246	1269 (2.04-2.04)
Ramachandran outliers	100387	1258 (2.04-2.04)
Sidechain outliers	100360	1258 (2.04-2.04)
RSRZ outliers	91569	1194 (2.04-2.04)

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	529	 13% 87% 10% •
1	B	529	 5% 89% 8% •

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 8729 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 SusD homolog.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	515	Total	C	N	O	S	Se	0	7	0
			4067	2588	686	782	4	7			
1	B	515	Total	C	N	O	S	Se	0	5	0
			4065	2583	681	790	4	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	leader sequence	UNP A6L9L2
B	0	GLY	-	leader sequence	UNP A6L9L2

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

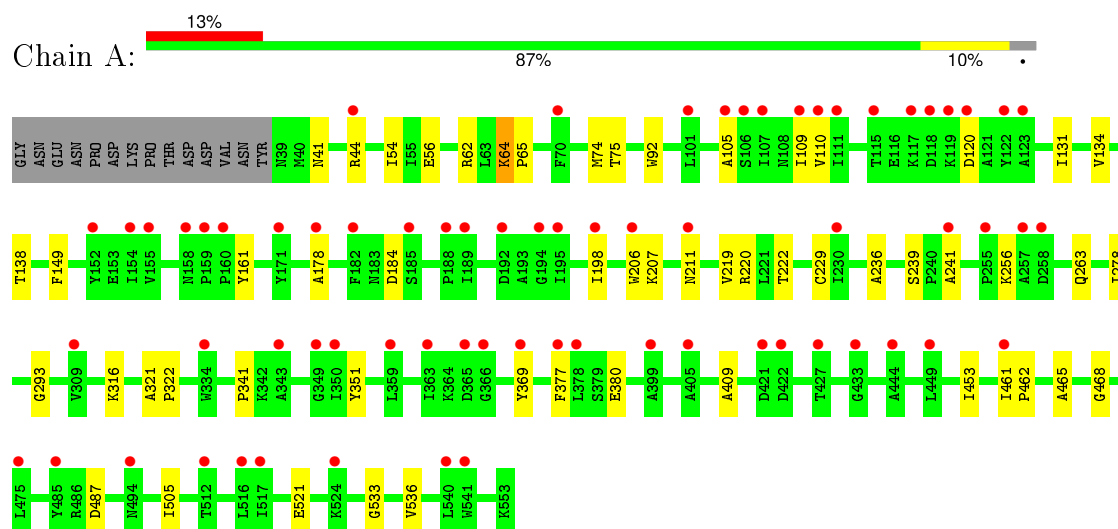
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	298	Total 298	O 298	0	0
3	B	294	Total 294	O 294	0	0

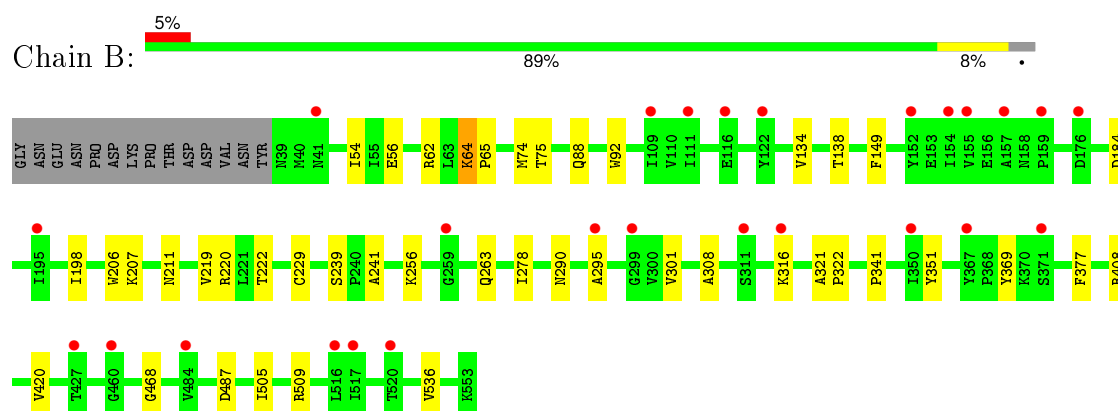
3 Residue-property plots [i](#)

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



• Molecule 1: SusD homolog



4 Data and refinement statistics

Property	Value	Source
Space group	P 6	Depositor
Cell constants a, b, c, α , β , γ	130.24Å 130.24Å 114.82Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	28.70 – 2.05 28.71 – 2.05	Depositor EDS
% Data completeness (in resolution range)	(Not available) (28.70-2.05) 100.0 (28.71-2.05)	Depositor EDS
R_{merge}	0.25	Depositor
R_{sym}	0.25	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.96 (at 2.04Å)	Xtriage
Refinement program	BUSTER 2.8.0	Depositor
R, R_{free}	0.243 , 0.265 0.271 , 0.298	Depositor DCC
R_{free} test set	3505 reflections (5.06%)	DCC
Wilson B-factor (Å ²)	16.4	Xtriage
Anisotropy	0.364	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 47.7	EDS
Estimated twinning fraction	0.042 for h,-h-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 69269 reflections	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	8729	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.06% 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: 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	0.50	0/4174	0.23	0/5651
1	B	0.50	0/4168	0.23	0/5646
All	All	0.50	0/8342	0.23	0/11297

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	4067	0	3943	29	0
1	B	4065	0	3925	22	0
2	B	5	0	0	0	0
3	A	298	0	0	2	0
3	B	294	0	0	0	0
All	All	8729	0	7868	51	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 3.

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:A:41:ASN:HD22	1:A:44[B]:ARG:HH12	1.30	0.76
1:A:533:GLY:HA2	3:A:676:HOH:O	1.87	0.73
1:B:64:LYS:HB3	1:B:65:PRO:CD	2.22	0.70
1:A:64:LYS:HB3	1:A:65:PRO:CD	2.23	0.68
1:A:105:ALA:O	1:A:109:ILE:HG12	2.03	0.58
1:A:207:LYS:HB3	1:A:241:ALA:HB1	1.88	0.55
1:B:256:LYS:HB2	1:B:263:GLN:HB3	1.89	0.55
1:B:207:LYS:HB3	1:B:241:ALA:HB1	1.87	0.55
1:A:256:LYS:HB2	1:A:263:GLN:HB3	1.88	0.54
1:B:295:ALA:HA	1:B:316[A]:LYS:HD3	1.90	0.53
1:B:64:LYS:HB3	1:B:65:PRO:HD3	1.90	0.53
1:A:64:LYS:HB3	1:A:65:PRO:HD3	1.92	0.52
1:A:41:ASN:HB3	1:A:110:VAL:HG13	1.92	0.51
1:B:211:ASN:HD21	1:B:239:SER:HB3	1.76	0.50
1:B:134:VAL:O	1:B:138:THR:HG23	2.13	0.49
1:A:131[B]:ILE:HD11	1:A:178:ALA:HB3	1.94	0.49
1:A:134:VAL:O	1:A:138:THR:HG23	2.13	0.48
1:A:211:ASN:HD21	1:A:239:SER:HB3	1.78	0.48
1:A:293:GLY:HA2	1:A:316:LYS:HD3	1.96	0.47
1:A:198:ILE:HD12	1:A:206:TRP:HH2	1.79	0.47
1:B:219:VAL:O	1:B:222:THR:HG23	2.14	0.47
1:B:505:ILE:HG13	1:B:536:VAL:HB	1.97	0.47
1:A:505:ILE:HG13	1:A:536:VAL:HB	1.97	0.47
1:A:219:VAL:O	1:A:222:THR:HG23	2.14	0.47
1:B:198:ILE:HD12	1:B:206:TRP:HH2	1.80	0.47
1:B:290:ASN:HA	1:B:316[A]:LYS:HB2	1.98	0.46
1:B:301:VAL:HG12	1:B:308:ALA:HA	1.97	0.46
1:A:64:LYS:HB3	1:A:65:PRO:HD2	1.99	0.44
1:B:64:LYS:HB3	1:B:65:PRO:HD2	1.98	0.44
1:B:408:ARG:HA	1:B:420:VAL:HG21	1.99	0.44
1:A:453:ILE:HG23	3:A:605:HOH:O	2.17	0.43
1:A:74:MSE:SE	1:A:465:ALA:HA	2.69	0.43
1:B:88:GLN:HG3	1:B:509:ARG:HG3	1.99	0.43
1:B:321:ALA:HB3	1:B:322:PRO:HD3	2.00	0.43
1:A:321:ALA:HB3	1:A:322:PRO:HD3	2.00	0.42
1:A:74:MSE:HE3	1:A:468:GLY:HA3	2.02	0.42
1:A:62:ARG:HG2	1:A:92:TRP:CE2	2.55	0.42
1:B:64:LYS:CB	1:B:65:PRO:CD	2.94	0.42
1:A:341:PRO:HG3	1:A:351:TYR:CD1	2.55	0.42
1:A:211:ASN:ND2	1:A:236:ALA:O	2.53	0.42
1:B:62:ARG:HG2	1:B:92:TRP:CE2	2.55	0.42
1:A:64:LYS:CB	1:A:65:PRO:CD	2.95	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:341:PRO:HG3	1:B:351:TYR:CD1	2.55	0.41
1:B:75:THR:HB	1:B:278:ILE:HG22	2.02	0.41
1:A:198:ILE:HD12	1:A:206:TRP:CH2	2.54	0.41
1:A:380:GLU:HG3	1:A:409:ALA:HB3	2.03	0.41
1:A:75:THR:HB	1:A:278:ILE:HG22	2.03	0.41
1:A:461:ILE:HA	1:A:462:PRO:HA	1.89	0.41
1:B:74:MSE:HE3	1:B:468:GLY:HA3	2.02	0.40
1:A:161:TYR:CD1	1:A:521:GLU:HB3	2.56	0.40
1:B:198:ILE:HD12	1:B:206:TRP:CH2	2.56	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	520/529 (98%)	496 (95%)	21 (4%)	3 (1%)	30	18
1	B	518/529 (98%)	494 (95%)	21 (4%)	3 (1%)	30	18
All	All	1038/1058 (98%)	990 (95%)	42 (4%)	6 (1%)	30	18

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	64	LYS
1	A	487	ASP
1	B	64	LYS
1	B	487	ASP
1	A	54	ILE
1	B	54	ILE

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	422/435 (97%)	413 (98%)	9 (2%)	61	56
1	B	425/435 (98%)	418 (98%)	7 (2%)	70	67
All	All	847/870 (97%)	831 (98%)	16 (2%)	66	60

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	GLU
1	A	120[A]	ASP
1	A	120[B]	ASP
1	A	149	PHE
1	A	184	ASP
1	A	220	ARG
1	A	229	CYS
1	A	369	TYR
1	A	377	PHE
1	B	56	GLU
1	B	149	PHE
1	B	184	ASP
1	B	220	ARG
1	B	229	CYS
1	B	369	TYR
1	B	377	PHE

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	41	ASN
1	B	41	ASN
1	B	211	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 ⓘ

1 ligand is 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	SO4	B	554	-	4,4,4	0.19	0	6,6,6	0.09	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	SO4	B	554	-	-	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	507/529 (95%)	1.09	69 (13%) 4 4	10, 18, 33, 51	0
1	B	507/529 (95%)	0.62	26 (5%) 32 36	8, 16, 30, 48	0
All	All	1014/1058 (95%)	0.85	95 (9%) 11 12	8, 17, 32, 51	0

All (95) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	155	VAL	6.2
1	A	122	TYR	4.4
1	A	120[A]	ASP	4.2
1	A	516	LEU	4.1
1	A	185	SER	4.0
1	B	154	ILE	3.9
1	A	461	ILE	3.7
1	A	118	ASP	3.7
1	B	155	VAL	3.4
1	A	343	ALA	3.3
1	B	116	GLU	3.3
1	A	399	ALA	3.2
1	B	367	TYR	3.2
1	A	365	ASP	3.2
1	A	123	ALA	3.0
1	A	110	VAL	3.0
1	A	189	ILE	3.0
1	B	259	GLY	3.0
1	A	512	THR	3.0
1	B	157	ALA	3.0
1	A	117	LYS	2.9
1	A	158	ASN	2.9
1	A	230	ILE	2.9
1	A	309	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	363	ILE	2.8
1	A	109	ILE	2.8
1	A	198	ILE	2.8
1	A	433	GLY	2.8
1	A	111	ILE	2.8
1	B	484	VAL	2.8
1	B	371	SER	2.7
1	B	152	TYR	2.7
1	B	111	ILE	2.7
1	A	194	GLY	2.7
1	A	115	THR	2.6
1	A	405	ALA	2.6
1	B	295	ALA	2.6
1	A	359	LEU	2.6
1	A	427	THR	2.6
1	A	195	ILE	2.6
1	A	211	ASN	2.6
1	A	154	ILE	2.5
1	A	159	PRO	2.5
1	A	475	LEU	2.5
1	A	541	TRP	2.5
1	A	106	SER	2.4
1	A	192	ASP	2.4
1	A	119	LYS	2.4
1	B	122	TYR	2.4
1	A	160	PRO	2.4
1	A	449	LEU	2.4
1	B	516	LEU	2.4
1	A	178	ALA	2.4
1	A	182	PHE	2.4
1	B	41	ASN	2.4
1	B	176[A]	ASP	2.3
1	A	105	ALA	2.3
1	A	524[A]	LYS	2.3
1	B	460	GLY	2.3
1	A	101	LEU	2.2
1	B	195	ILE	2.2
1	B	311[A]	SER	2.2
1	A	241	ALA	2.2
1	A	257	ALA	2.2
1	A	444	ALA	2.2
1	B	299	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	107	ILE	2.2
1	A	517	ILE	2.2
1	A	70	PHE	2.2
1	A	188	PRO	2.2
1	A	378	LEU	2.1
1	A	206	TRP	2.1
1	B	109	ILE	2.1
1	A	421	ASP	2.1
1	B	316[A]	LYS	2.1
1	A	349	GLY	2.1
1	B	517	ILE	2.1
1	A	540	LEU	2.1
1	A	369	TYR	2.1
1	A	366	GLY	2.1
1	B	520	THR	2.1
1	A	255	PRO	2.1
1	A	377	PHE	2.1
1	A	171	TYR	2.1
1	A	485	TYR	2.1
1	B	427	THR	2.1
1	B	159	PRO	2.1
1	A	334	TRP	2.0
1	A	494	ASN	2.0
1	A	152	TYR	2.0
1	A	350	ILE	2.0
1	A	422	ASP	2.0
1	A	44[A]	ARG	2.0
1	A	258	ASP	2.0
1	B	350	ILE	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 [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. 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	SO4	B	554	5/5	0.95	0.15	-	35,39,40,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.