



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 15, 2017 – 05:25 am GMT

PDB ID : 5FQ7
Title : Crystal structure of the SusCD complex BT2261-2264 from *Bacteroides thetaiotaomicron*
Authors : Glenwright, A.J.; Pothula, K.R.; Chorev, D.S.; Basle, A.; Robinson, C.V.; Kleinekathoefer, U.; Bolam, D.N.; van den Berg, B.
Deposited on : 2015-12-07
Resolution : 3.40 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

<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
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

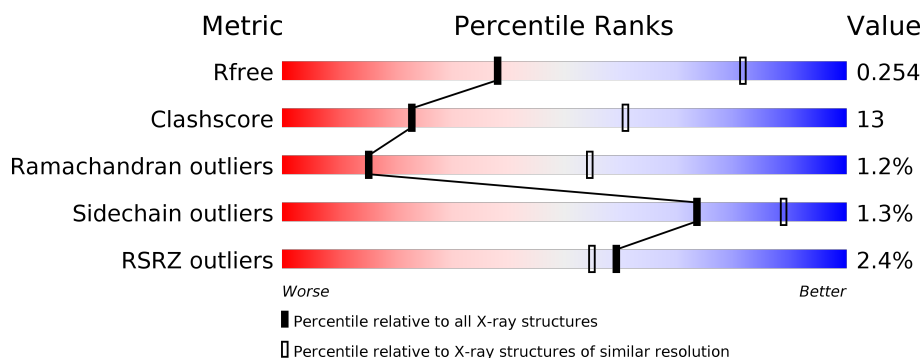
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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.



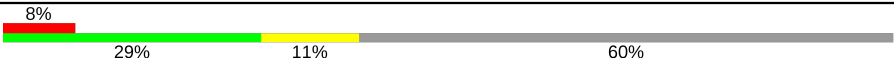

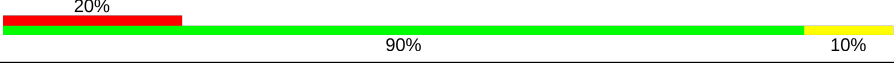

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	1679 (3.50-3.30)
Clashscore	112137	1832 (3.50-3.30)
Ramachandran outliers	110173	1789 (3.50-3.30)
Sidechain outliers	110143	1789 (3.50-3.30)
RSRZ outliers	101464	1709 (3.50-3.30)

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	480	<div> <div>70%</div> <div>29%</div> <div>.</div> </div>
1	C	480	<div> <div>71%</div> <div>28%</div> <div>.</div> </div>
2	B	984	<div> <div>2%</div> <div>66%</div> <div>29%</div> <div>..</div> </div>
2	D	984	<div> <div>%</div> <div>68%</div> <div>26%</div> <div>..</div> </div>
3	E	148	<div> <div>4%</div> <div>84%</div> <div>14%</div> <div>.</div> </div>
3	F	148	<div> <div>8%</div> <div>76%</div> <div>22%</div> <div>..</div> </div>

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Mol	Chain	Length	Quality of chain
4	G	212	
4	H	212	
5	I	10	
5	P	10	

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
7	NA	B	1988	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 25886 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 BT_2263.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	480	Total	C	N	O	S	0	0	0
			3744	2370	616	741	17			
1	C	480	Total	C	N	O	S	0	0	0
			3744	2370	616	741	17			

- Molecule 2 is a protein called BT_2264.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	943	Total	C	N	O	S	0	0	0
			7342	4643	1224	1445	30			
2	D	943	Total	C	N	O	S	0	0	0
			7337	4642	1224	1441	30			

- Molecule 3 is a protein called BT_2261.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	145	Total	C	N	O	S	0	0	0
			1134	717	178	234	5			
3	F	145	Total	C	N	O	S	0	0	0
			1134	717	178	234	5			

- Molecule 4 is a protein called BT_2262.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	G	84	Total	C	N	O	S	0	0	0
			665	422	102	140	1			
4	H	89	Total	C	N	O	S	0	0	0
			698	441	108	148	1			

- Molecule 5 is a protein called PEPTIDE.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	I	10	Total	C	N	O	0	0	0
			40	20	10	10			
5	P	10	Total	C	N	O	0	0	0
			40	20	10	10			

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Mg	0	0
			1	1		
6	D	1	Total	Mg	0	0
			1	1		

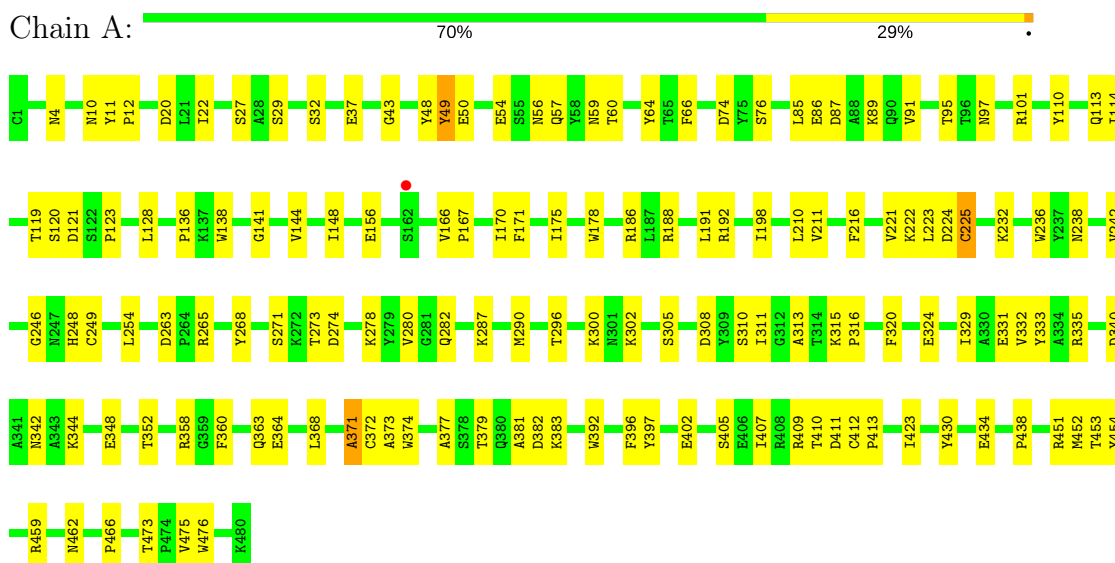
- Molecule 7 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	3	Total	Na	0	0
			3	3		
7	D	2	Total	Na	0	0
			2	2		
7	C	1	Total	Na	0	0
			1	1		

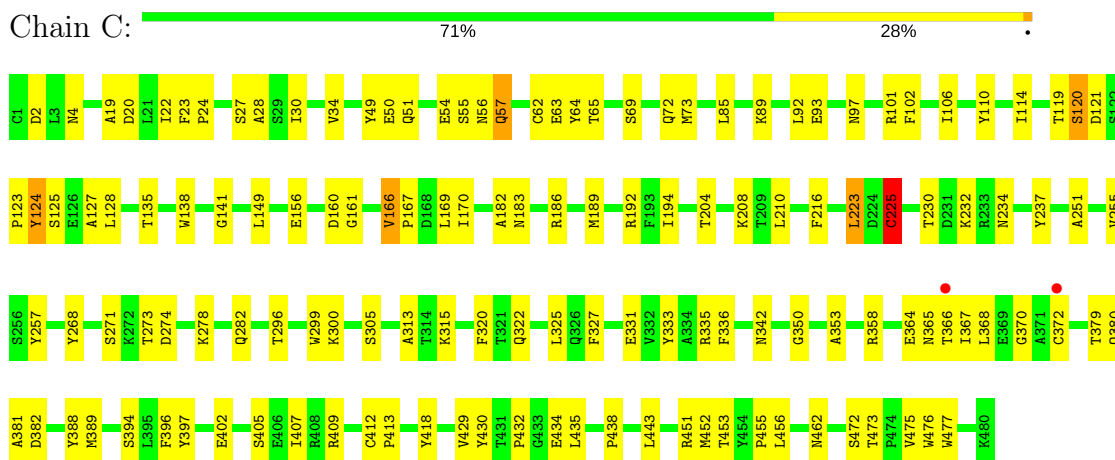
3 Residue-property plots

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: BT_2263

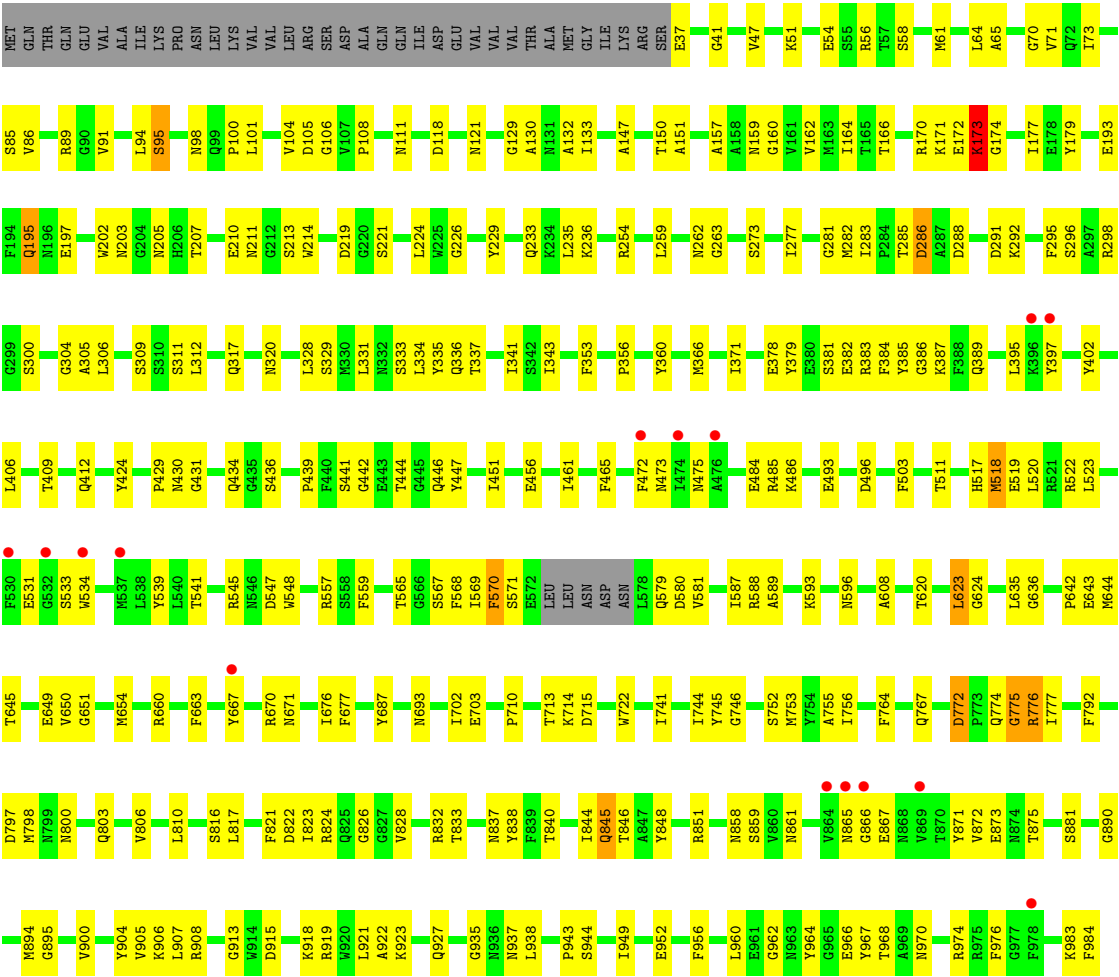


• Molecule 1: BT_2263

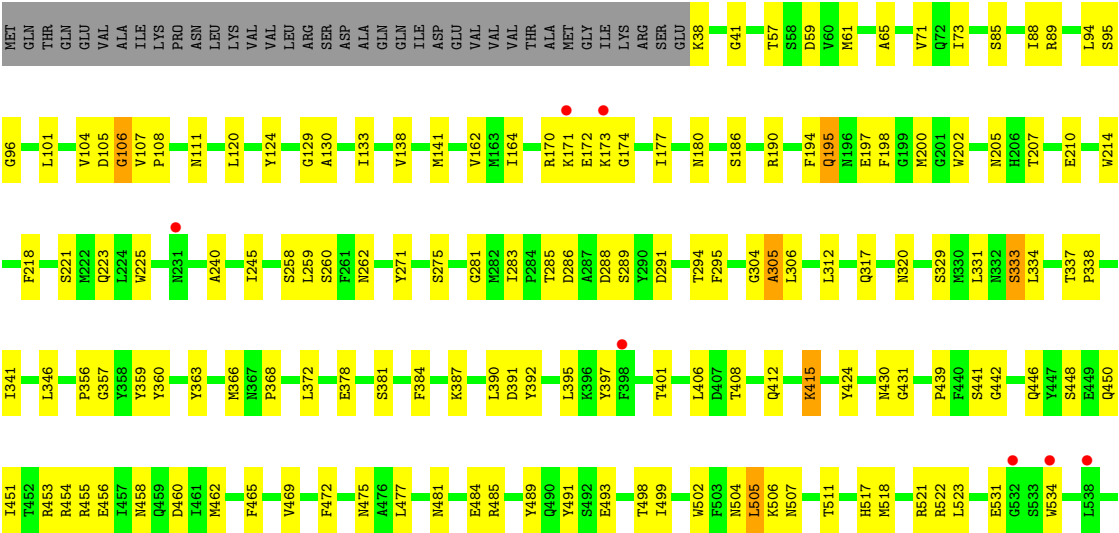


• Molecule 2: BT_2264





• Molecule 2: BT_2264

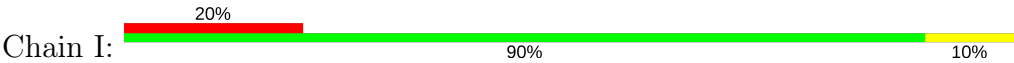


THR	ALA	SER	PRO	LEU	LYS	SER	GLY	ILE	TYR	THR	VAL	ALA	GLU	GLY	SER	LYS	ARG	THR	ALA	PRO	SER	VAL	VAL	ALA	PHE	SER	GLY	TYR	GLU	ILE	VAL	ILE	PHE	GLN	MET	LEU	GLU	PRO	GLY	ILE	PHE	THR	TYR	ILE	SER	ASP	PHE	LEU	GLY	GLY	TRP	TYR	ASP	GLN	ARG	ALA	GLY	TYR	GLY	PRO	ASP
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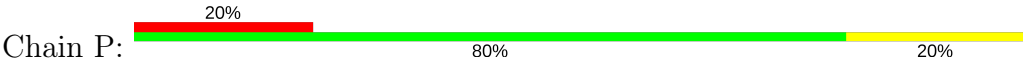
TYR	ALA	MET	VAL	GLY	LYS	PHE	GLU	LEU	ASN	ASP	ASP	ASN	THR	ILE	THR	PRO	LEU	GLU	SER	TYR	VAL	ALA	GLY	TRP	GLY	ASP	SER	MET	ASP	GLN	MET	THR	THR	ASN	LEU	LEU	ASP	PRO	ALA	ALA	THR	GLY	THR	LEU	LYS	TRP	THR	VAL	ALA	TYR	ALA	GLY	GLN	LEU	SER	PHE	ASP	ILE	VAL
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LYS	GLN
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● Molecule 5: PEPTIDE



● Molecule 5: PEPTIDE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	123.59Å 180.16Å 245.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.43 – 3.40 49.45 – 3.40	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.43-3.40) 94.7 (49.45-3.40)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.50 (at 3.40Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.201 , 0.255 0.198 , 0.254	Depositor DCC
R_{free} test set	7448 reflections (9.79%)	DCC
Wilson B-factor (Å ²)	80.6	Xtriage
Anisotropy	0.392	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 45.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	25886	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.79% 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: NA, MG

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	0.65	1/3835 (0.0%)	0.66	0/5220
1	C	0.68	2/3835 (0.1%)	0.65	0/5220
2	B	0.60	0/7518	0.70	3/10196 (0.0%)
2	D	0.62	1/7513 (0.0%)	0.71	4/10190 (0.0%)
3	E	0.49	0/1162	0.57	0/1578
3	F	0.43	0/1162	0.57	0/1578
4	G	0.36	0/677	0.55	0/918
4	H	0.55	0/711	0.65	0/966
5	I	0.59	0/39	1.00	0/47
5	P	0.55	0/39	1.03	0/47
All	All	0.61	4/26491 (0.0%)	0.67	7/35960 (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
2	B	0	2
2	D	0	1
All	All	0	3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	225	CYS	CB-SG	-6.56	1.71	1.82
1	A	225	CYS	CB-SG	-6.40	1.71	1.82
2	D	942	THR	C-N	-6.02	1.22	1.34
1	C	166	VAL	CB-CG1	-5.11	1.42	1.52

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	921	LEU	CB-CG-CD1	-7.91	97.56	111.00
2	D	624	GLY	N-CA-C	-7.29	94.87	113.10
2	B	624	GLY	N-CA-C	-6.89	95.88	113.10
2	D	623	LEU	CA-CB-CG	6.38	129.97	115.30
2	B	623	LEU	CA-CB-CG	5.51	127.98	115.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	623	LEU	Peptide
2	B	949	ILE	Peptide
2	D	623	LEU	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	3744	0	3562	96	0
1	C	3744	0	3562	95	0
2	B	7342	0	7013	208	0
2	D	7337	0	7014	213	0
3	E	1134	0	1050	12	0
3	F	1134	0	1050	21	0
4	G	665	0	626	16	0
4	H	698	0	656	21	0
5	I	40	0	32	1	0
5	P	40	0	32	1	0
6	B	1	0	0	0	0
6	D	1	0	0	0	0
7	B	3	0	0	0	0
7	C	1	0	0	0	0
7	D	2	0	0	0	0
All	All	25886	0	24597	653	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 13.

The worst 5 of 653 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:D:198:PHE:CD1	2:D:851:ARG:NH2	2.23	1.06
2:D:124:TYR:OH	2:D:415:LYS:NZ	1.93	1.02
2:B:172:GLU:HG2	2:B:173:LYS:HG2	1.43	0.98
2:D:198:PHE:HD1	2:D:851:ARG:NH2	1.59	0.97
2:D:579:GLN:HB2	2:D:581:VAL:HG12	1.46	0.93

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	478/480 (100%)	425 (89%)	45 (9%)	8 (2%)	11	47
1	C	478/480 (100%)	432 (90%)	40 (8%)	6 (1%)	14	53
2	B	939/984 (95%)	828 (88%)	99 (10%)	12 (1%)	14	53
2	D	939/984 (95%)	848 (90%)	80 (8%)	11 (1%)	15	54
3	E	143/148 (97%)	130 (91%)	11 (8%)	2 (1%)	13	51
3	F	143/148 (97%)	124 (87%)	19 (13%)	0	100	100
4	G	80/212 (38%)	77 (96%)	3 (4%)	0	100	100
4	H	87/212 (41%)	82 (94%)	5 (6%)	0	100	100
5	I	8/10 (80%)	7 (88%)	1 (12%)	0	100	100
5	P	8/10 (80%)	5 (62%)	2 (25%)	1 (12%)	0	4
All	All	3303/3668 (90%)	2958 (90%)	305 (9%)	40 (1%)	15	54

5 of 40 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	411	ASP

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Mol	Chain	Res	Type
2	B	714	LYS
2	B	943	PRO
1	C	57	GLN
1	C	120	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	392/392 (100%)	386 (98%)	6 (2%)	70	87
1	C	392/392 (100%)	387 (99%)	5 (1%)	73	89
2	B	796/836 (95%)	787 (99%)	9 (1%)	78	90
2	D	795/836 (95%)	784 (99%)	11 (1%)	71	88
3	E	121/124 (98%)	121 (100%)	0	100	100
3	F	121/124 (98%)	120 (99%)	1 (1%)	85	93
4	G	73/177 (41%)	71 (97%)	2 (3%)	50	80
4	H	76/177 (43%)	73 (96%)	3 (4%)	37	72
All	All	2766/3058 (90%)	2729 (99%)	37 (1%)	73	89

5 of 37 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	223	LEU
2	D	195	GLN
4	H	22	ASP
1	C	274	ASP
1	C	320	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
2	D	458	ASN

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Mol	Chain	Res	Type
3	E	123	GLN
2	D	800	ASN
2	B	490	GLN
2	D	803	GLN

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](#)

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	480/480 (100%)	-0.25	1 (0%) 94 94	52, 70, 91, 112	0
1	C	480/480 (100%)	-0.21	2 (0%) 92 90	56, 72, 93, 120	0
2	B	943/984 (95%)	-0.08	15 (1%) 72 67	55, 75, 109, 147	0
2	D	943/984 (95%)	0.00	13 (1%) 75 71	55, 74, 106, 146	0
3	E	145/148 (97%)	0.24	6 (4%) 38 34	76, 87, 103, 126	0
3	F	145/148 (97%)	0.34	12 (8%) 12 13	80, 99, 126, 151	0
4	G	84/212 (39%)	1.13	18 (21%) 1 1	84, 126, 152, 174	0
4	H	89/212 (41%)	0.71	8 (8%) 10 11	66, 93, 123, 179	0
5	I	10/10 (100%)	1.69	2 (20%) 1 1	53, 62, 75, 78	0
5	P	10/10 (100%)	1.41	2 (20%) 1 1	63, 68, 82, 93	0
All	All	3329/3668 (90%)	-0.00	79 (2%) 59 55	52, 75, 111, 179	0

The worst 5 of 79 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	F	25	ALA	4.4
4	G	82	THR	4.4
4	H	88	ALA	3.9
4	G	84	THR	3.7
3	F	133	TYR	3.6

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. 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
7	NA	B	1988	1/1	0.94	0.23	3.14	43,43,43,43	0
6	MG	B	1985	1/1	0.96	0.25	1.94	23,23,23,23	0
7	NA	D	1986	1/1	0.98	0.21	0.17	35,35,35,35	0
6	MG	D	1985	1/1	0.97	0.21	-0.26	38,38,38,38	0
7	NA	D	1987	1/1	0.97	0.15	-0.64	34,34,34,34	0
7	NA	C	1481	1/1	0.91	0.16	-0.64	44,44,44,44	0
7	NA	B	1987	1/1	0.97	0.13	-1.14	30,30,30,30	0
7	NA	B	1986	1/1	0.83	0.39	-	27,27,27,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.