



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 02:21 PM GMT

PDB ID : 3WWQ
Title : Crystal structure of FAAP20 UBZ domain in complex with Lys63-linked diubiquitin
Authors : Sato, Y.; Fukai, S.
Deposited on : 2014-06-23
Resolution : 1.90 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
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

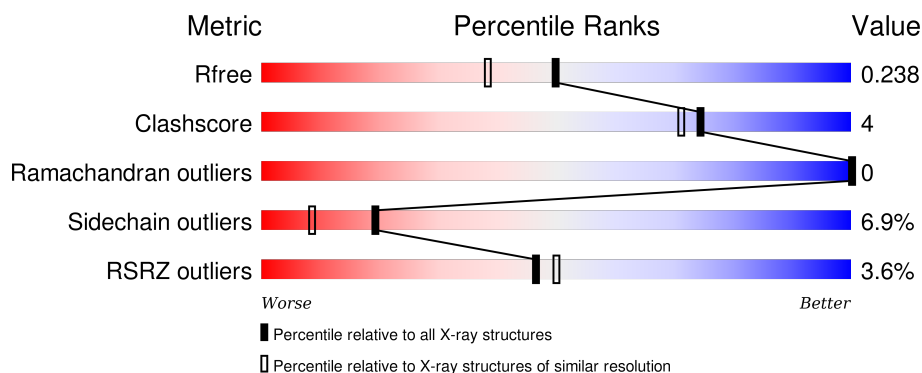
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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	76	<div> <div>84%</div> <div>16%</div> </div>
1	D	76	<div> <div>86%</div> <div>13%</div> </div>
1	G	76	<div> <div>83%</div> <div>16%</div> </div>
1	J	76	<div> <div>8%</div> <div>91%</div> <div>7%</div> </div>
2	B	77	<div> <div>87%</div> <div>9%</div> </div>

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Mol	Chain	Length	Quality of chain
2	E	77	<div><div></div><div>6%</div><div>82%</div><div>13%</div><div></div><div></div></div>
2	H	77	<div><div></div><div>%</div><div>84%</div><div>10%</div><div></div><div></div></div>
2	K	77	<div><div></div><div>6%</div><div>86%</div><div>10%</div><div></div><div></div></div>
3	C	44	<div><div></div><div>2%</div><div>75%</div><div>7%</div><div>5%</div><div>14%</div></div>
3	F	44	<div><div></div><div>7%</div><div>73%</div><div>16%</div><div></div><div>11%</div></div>
3	I	44	<div><div></div><div>5%</div><div>84%</div><div></div><div></div><div>14%</div></div>
3	L	44	<div><div></div><div>5%</div><div>70%</div><div>14%</div><div></div><div>16%</div></div>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6359 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 Ubiquitin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	76	Total	C	N	O	S	0	1	0
			608	381	108	118	1			
1	D	76	Total	C	N	O	S	0	0	0
			603	378	107	117	1			
1	G	76	Total	C	N	O	S	0	0	0
			603	378	107	117	1			
1	J	76	Total	C	N	O	S	0	0	0
			603	378	107	117	1			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	63	ARG	LYS	ENGINEERED MUTATION	UNP P0CG50
D	63	ARG	LYS	ENGINEERED MUTATION	UNP P0CG50
G	63	ARG	LYS	ENGINEERED MUTATION	UNP P0CG50
J	63	ARG	LYS	ENGINEERED MUTATION	UNP P0CG50

- Molecule 2 is a protein called Ubiquitin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	76	Total	C	N	O	S	0	0	0
			601	378	105	117	1			
2	E	76	Total	C	N	O	S	0	0	0
			601	378	105	117	1			
2	H	75	Total	C	N	O	S	0	0	0
			597	376	104	116	1			
2	K	76	Total	C	N	O	S	0	0	0
			601	378	105	117	1			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	77	ASP	-	SEE REMARK 999	UNP P0CG50
E	77	ASP	-	SEE REMARK 999	UNP P0CG50
H	77	ASP	-	SEE REMARK 999	UNP P0CG50
K	77	ASP	-	SEE REMARK 999	UNP P0CG50

- Molecule 3 is a protein called Fanconi anemia-associated protein of 20 kDa.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	38	Total	C	N	O	S	0	0	0
			298	183	51	60	4			
3	F	39	Total	C	N	O	S	0	0	0
			303	186	52	61	4			
3	I	38	Total	C	N	O	S	0	0	0
			298	183	51	60	4			
3	L	37	Total	C	N	O	S	0	0	0
			293	180	50	59	4			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	137	GLY	-	EXPRESSION TAG	UNP Q6NZ36
C	138	PRO	-	EXPRESSION TAG	UNP Q6NZ36
C	139	GLY	-	EXPRESSION TAG	UNP Q6NZ36
C	140	HIS	-	EXPRESSION TAG	UNP Q6NZ36
C	141	MET	-	EXPRESSION TAG	UNP Q6NZ36
F	137	GLY	-	EXPRESSION TAG	UNP Q6NZ36
F	138	PRO	-	EXPRESSION TAG	UNP Q6NZ36
F	139	GLY	-	EXPRESSION TAG	UNP Q6NZ36
F	140	HIS	-	EXPRESSION TAG	UNP Q6NZ36
F	141	MET	-	EXPRESSION TAG	UNP Q6NZ36
I	137	GLY	-	EXPRESSION TAG	UNP Q6NZ36
I	138	PRO	-	EXPRESSION TAG	UNP Q6NZ36
I	139	GLY	-	EXPRESSION TAG	UNP Q6NZ36
I	140	HIS	-	EXPRESSION TAG	UNP Q6NZ36
I	141	MET	-	EXPRESSION TAG	UNP Q6NZ36
L	137	GLY	-	EXPRESSION TAG	UNP Q6NZ36
L	138	PRO	-	EXPRESSION TAG	UNP Q6NZ36
L	139	GLY	-	EXPRESSION TAG	UNP Q6NZ36
L	140	HIS	-	EXPRESSION TAG	UNP Q6NZ36
L	141	MET	-	EXPRESSION TAG	UNP Q6NZ36

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	I	1	Total Zn 1 1	0	0
4	L	1	Total Zn 1 1	0	0
4	C	1	Total Zn 1 1	0	0
4	F	1	Total Zn 1 1	0	0

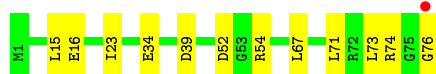
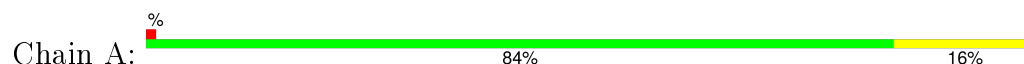
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	48	Total O 48 48	0	0
5	B	45	Total O 45 45	0	0
5	C	27	Total O 27 27	0	0
5	D	46	Total O 46 46	0	0
5	E	19	Total O 19 19	0	0
5	F	12	Total O 12 12	0	0
5	G	30	Total O 30 30	0	0
5	H	51	Total O 51 51	0	0
5	I	9	Total O 9 9	0	0
5	J	25	Total O 25 25	0	0
5	K	23	Total O 23 23	0	0
5	L	11	Total O 11 11	0	0

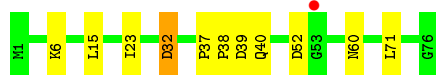
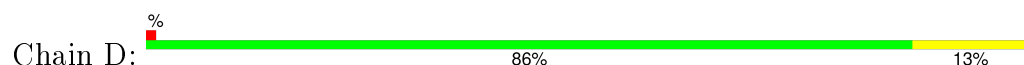
3 Residue-property plots [i](#)

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.

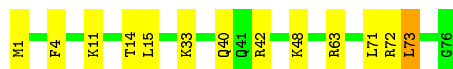
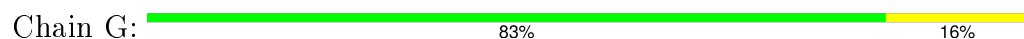
- Molecule 1: Ubiquitin



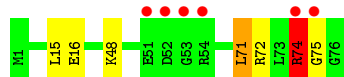
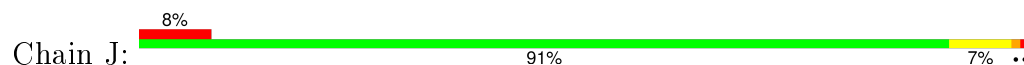
- Molecule 1: Ubiquitin



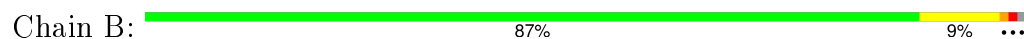
- Molecule 1: Ubiquitin



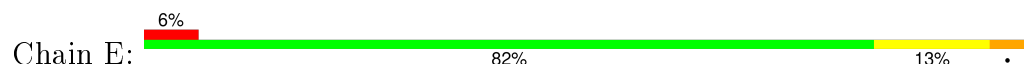
- Molecule 1: Ubiquitin

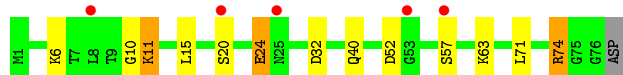


- Molecule 2: Ubiquitin

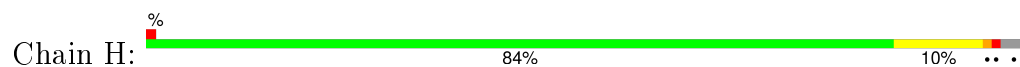


- Molecule 2: Ubiquitin

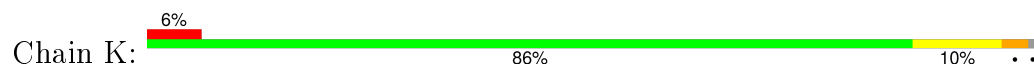




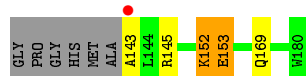
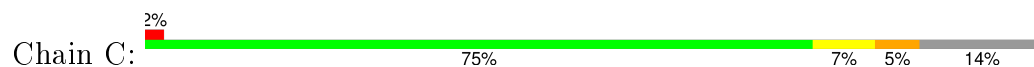
- Molecule 2: Ubiquitin



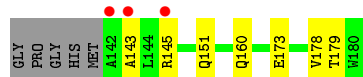
- Molecule 2: Ubiquitin



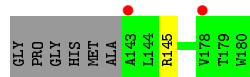
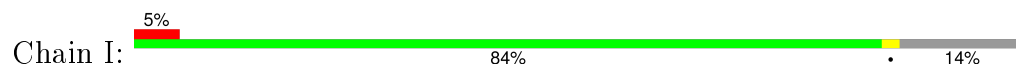
- Molecule 3: Fanconi anemia-associated protein of 20 kDa



- Molecule 3: Fanconi anemia-associated protein of 20 kDa



- Molecule 3: Fanconi anemia-associated protein of 20 kDa



- Molecule 3: Fanconi anemia-associated protein of 20 kDa



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.77Å 45.85Å 172.90Å 90.00° 98.06° 90.00°	Depositor
Resolution (Å)	50.00 – 1.90 38.47 – 1.90	Depositor EDS
% Data completeness (in resolution range)	95.0 (50.00-1.90) 94.8 (38.47-1.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.61 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.8.0069	Depositor
R, R_{free}	0.208 , 0.235 0.215 , 0.238	Depositor DCC
R_{free} test set	3533 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	24.1	Xtriage
Anisotropy	0.348	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 44.8	EDS
Estimated twinning fraction	0.022 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 70403 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6359	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.77% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	0.77	0/617	1.19	6/830 (0.7%)
1	D	0.71	0/609	1.03	3/819 (0.4%)
1	G	0.67	0/609	0.96	1/819 (0.1%)
1	J	0.65	0/609	0.97	1/819 (0.1%)
2	B	0.73	0/607	1.04	2/816 (0.2%)
2	E	0.63	0/607	1.06	3/816 (0.4%)
2	H	0.69	0/603	1.04	5/811 (0.6%)
2	K	0.66	0/607	1.06	2/816 (0.2%)
3	C	0.70	0/303	1.09	2/410 (0.5%)
3	F	0.75	0/308	0.90	0/417
3	I	0.57	0/303	0.96	2/410 (0.5%)
3	L	0.68	0/298	0.88	0/403
All	All	0.69	0/6080	1.03	27/8186 (0.3%)

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
3	F	0	1

There are no bond length outliers.

The worst 5 of 27 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	39	ASP	CB-CG-OD2	-9.91	109.38	118.30
1	A	39	ASP	CB-CG-OD1	9.63	126.96	118.30
3	C	145	ARG	NE-CZ-NH2	-9.55	115.52	120.30
2	E	32	ASP	CB-CG-OD1	-8.38	110.76	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	32	ASP	CB-CG-OD2	8.31	125.78	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	F	143	ALA	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	608	0	635	5	0
1	D	603	0	629	3	0
1	G	603	0	629	7	0
1	J	603	0	629	6	0
2	B	601	0	627	7	0
2	E	601	0	627	9	0
2	H	597	0	624	5	0
2	K	601	0	627	9	0
3	C	298	0	282	3	0
3	F	303	0	287	0	0
3	I	298	0	282	0	0
3	L	293	0	277	4	0
4	C	1	0	0	0	0
4	F	1	0	0	0	0
4	I	1	0	0	0	0
4	L	1	0	0	0	0
5	A	48	0	0	2	0
5	B	45	0	0	5	0
5	C	27	0	0	1	0
5	D	46	0	0	0	0
5	E	19	0	0	1	0
5	F	12	0	0	0	0
5	G	30	0	0	1	0
5	H	51	0	0	0	0
5	I	9	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	J	25	0	0	4	0
5	K	23	0	0	2	0
5	L	11	0	0	2	0
All	All	6359	0	6155	49	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 4.

The worst 5 of 49 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:72:ARG:NH1	5:B:121:HOH:O	2.07	0.86
1:A:34:GLU:HG3	5:A:102:HOH:O	1.76	0.85
2:B:42:ARG:NH1	5:B:121:HOH:O	2.14	0.81
2:B:49:GLN:OE1	5:B:109:HOH:O	2.05	0.74
1:J:16:GLU:OE1	5:J:119:HOH:O	2.08	0.72

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	75/76 (99%)	75 (100%)	0	0	100	100
1	D	74/76 (97%)	74 (100%)	0	0	100	100
1	G	74/76 (97%)	74 (100%)	0	0	100	100
1	J	74/76 (97%)	74 (100%)	0	0	100	100
2	B	74/77 (96%)	74 (100%)	0	0	100	100
2	E	74/77 (96%)	74 (100%)	0	0	100	100
2	H	73/77 (95%)	73 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	K	74/77 (96%)	74 (100%)	0	0	100	100
3	C	36/44 (82%)	36 (100%)	0	0	100	100
3	F	37/44 (84%)	37 (100%)	0	0	100	100
3	I	36/44 (82%)	36 (100%)	0	0	100	100
3	L	35/44 (80%)	35 (100%)	0	0	100	100
All	All	736/788 (93%)	736 (100%)	0	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	69/68 (102%)	66 (96%)	3 (4%)	35	23
1	D	68/68 (100%)	63 (93%)	5 (7%)	17	7
1	G	68/68 (100%)	64 (94%)	4 (6%)	24	12
1	J	68/68 (100%)	64 (94%)	4 (6%)	24	12
2	B	68/69 (99%)	64 (94%)	4 (6%)	24	12
2	E	68/69 (99%)	61 (90%)	7 (10%)	9	3
2	H	68/69 (99%)	64 (94%)	4 (6%)	24	12
2	K	68/69 (99%)	64 (94%)	4 (6%)	24	12
3	C	34/37 (92%)	32 (94%)	2 (6%)	24	12
3	F	34/37 (92%)	28 (82%)	6 (18%)	2	0
3	I	34/37 (92%)	34 (100%)	0	100	100
3	L	34/37 (92%)	30 (88%)	4 (12%)	6	2
All	All	681/696 (98%)	634 (93%)	47 (7%)	19	8

5 of 47 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	F	145	ARG
3	F	179	THR
3	L	144	LEU
3	F	160	GLN
1	G	11	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 16 such sidechains are listed below:

Mol	Chain	Res	Type
1	G	60	ASN
2	H	25	ASN
2	K	40	GLN
2	E	40	GLN
2	K	49	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 4 ligands modelled in this entry, 4 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	76/76 (100%)	-0.06	1 (1%) 79 82	21, 31, 49, 71	1 (1%)
1	D	76/76 (100%)	-0.05	1 (1%) 79 82	22, 32, 43, 53	1 (1%)
1	G	76/76 (100%)	0.03	0 100 100	25, 39, 55, 58	2 (2%)
1	J	76/76 (100%)	0.50	6 (7%) 15 17	30, 48, 75, 81	1 (1%)
2	B	76/77 (98%)	0.18	0 100 100	23, 35, 48, 56	0
2	E	76/77 (98%)	0.63	5 (6%) 22 24	34, 50, 70, 74	0
2	H	75/77 (97%)	0.22	1 (1%) 79 82	22, 34, 45, 51	3 (4%)
2	K	76/77 (98%)	0.69	5 (6%) 22 24	24, 42, 66, 71	1 (1%)
3	C	38/44 (86%)	0.08	1 (2%) 59 63	23, 37, 55, 65	0
3	F	39/44 (88%)	0.53	3 (7%) 16 18	23, 36, 72, 82	1 (2%)
3	I	38/44 (86%)	0.24	2 (5%) 30 33	26, 45, 62, 77	0
3	L	37/44 (84%)	0.45	2 (5%) 29 33	29, 43, 68, 77	1 (2%)
All	All	759/788 (96%)	0.28	27 (3%) 46 50	21, 38, 64, 82	11 (1%)

The worst 5 of 27 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	F	142	ALA	5.8
3	I	143	ALA	4.4
3	F	143	ALA	3.9
2	E	20	SER	3.7
1	A	76	GLY	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
4	ZN	I	201	1/1	0.98	0.09	-0.80	28,28,28,28	0
4	ZN	C	201	1/1	0.99	0.07	-1.12	23,23,23,23	0
4	ZN	L	201	1/1	0.98	0.07	-1.25	28,28,28,28	0
4	ZN	F	201	1/1	0.99	0.05	-1.59	25,25,25,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.