Full wwPDB X-ray Structure Validation Report (i)

May 21, 2020 – 02:57 am BST

PDB ID	:	5MQ5
Title	:	A protease-resistant N24S Escherichia coli Asparaginase mutant with outstanding stability and enhanced anti-leukaemic activity
Authors	:	Maggi, M.; Mittelman, S.D.; Parmentier, J.H.; Whitmire, J.M.; Merrell, D.S.; Scotti, C.
Deposited on	:	2016-12-20
Resolution	:	1.60 Å(reported)

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

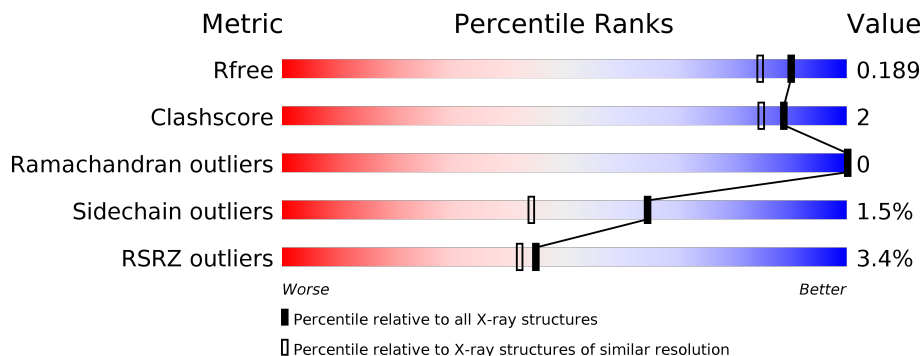|		
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation 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 1.60 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	Percentile Ranks	Value
Rfree		0.189
Clashscore		2
Ramachandran outliers		0
Sidechain outliers		1.5%
RSRZ outliers		3.4%

■ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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	332	4% 91% · 5%
1	B	332	3% 92% 6% ·
1	C	332	2% 92% 6% ··
1	D	332	4% 89% 5% 5%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 10835 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 L-asparaginase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	316	Total 2368	C 1480	N 407	O 473	S 8	0	0	0
1	C	326	Total 2429	C 1516	N 414	O 491	S 8	0	0	0
1	B	326	Total 2429	C 1516	N 414	O 491	S 8	0	0	0
1	D	314	Total 2348	C 1468	N 401	O 471	S 8	0	0	0

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	expression tag	UNP P00805
A	-4	HIS	-	expression tag	UNP P00805
A	-3	HIS	-	expression tag	UNP P00805
A	-2	HIS	-	expression tag	UNP P00805
A	-1	HIS	-	expression tag	UNP P00805
A	0	HIS	-	expression tag	UNP P00805
A	24	SER	ASN	engineered mutation	UNP P00805
C	-5	HIS	-	expression tag	UNP P00805
C	-4	HIS	-	expression tag	UNP P00805
C	-3	HIS	-	expression tag	UNP P00805
C	-2	HIS	-	expression tag	UNP P00805
C	-1	HIS	-	expression tag	UNP P00805
C	0	HIS	-	expression tag	UNP P00805
C	24	SER	ASN	engineered mutation	UNP P00805
B	-5	HIS	-	expression tag	UNP P00805
B	-4	HIS	-	expression tag	UNP P00805
B	-3	HIS	-	expression tag	UNP P00805
B	-2	HIS	-	expression tag	UNP P00805
B	-1	HIS	-	expression tag	UNP P00805
B	0	HIS	-	expression tag	UNP P00805
B	24	SER	ASN	engineered mutation	UNP P00805

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-5	HIS	-	expression tag	UNP P00805
D	-4	HIS	-	expression tag	UNP P00805
D	-3	HIS	-	expression tag	UNP P00805
D	-2	HIS	-	expression tag	UNP P00805
D	-1	HIS	-	expression tag	UNP P00805
D	0	HIS	-	expression tag	UNP P00805
D	24	SER	ASN	engineered mutation	UNP P00805

- Molecule 2 is ASPARTIC ACID (three-letter code: ASP) (formula: $C_4H_7NO_4$).
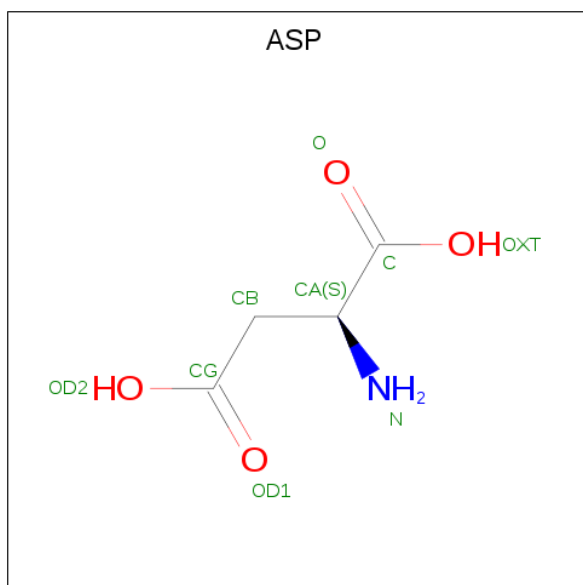
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 9	C 4	N 1	O 4	0	0
2	C	1	Total 9	C 4	N 1	O 4	0	0
2	B	1	Total 9	C 4	N 1	O 4	0	0
2	D	1	Total 9	C 4	N 1	O 4	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	333	Total 333	O 333	0	0
3	C	349	Total 349	O 349	0	0

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| Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	273	Total O 273 273	0	0
3	D	270	Total O 270 270	0	0

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.
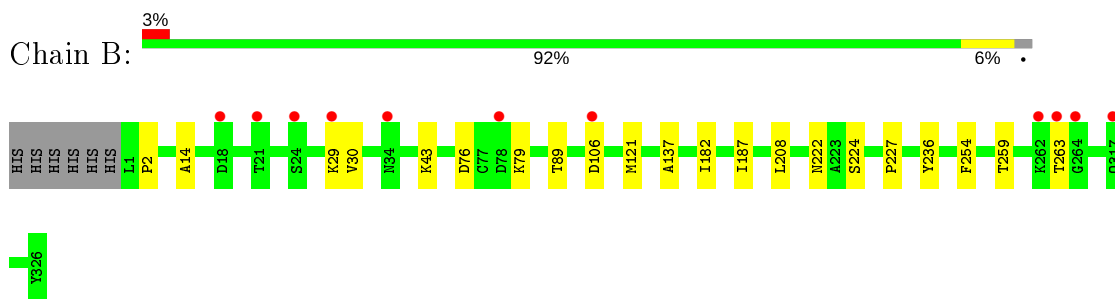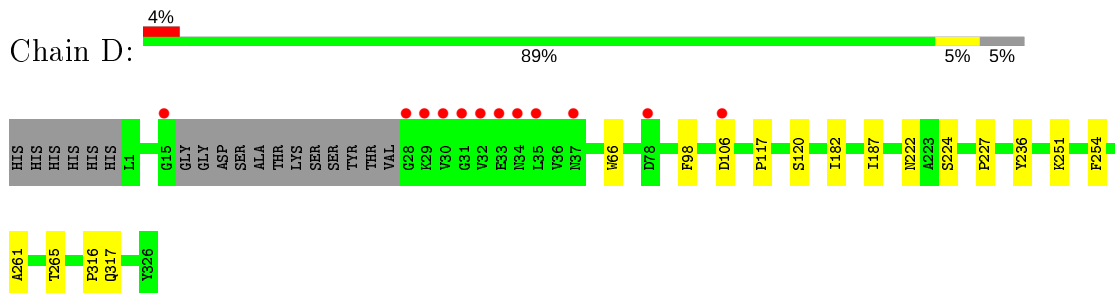
- Molecule 1: L-asparaginase 2
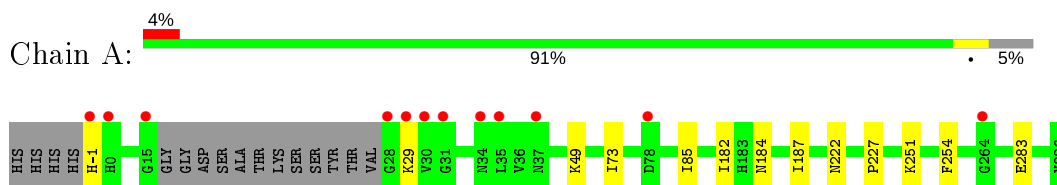
- Molecule 1: L-asparaginase 2
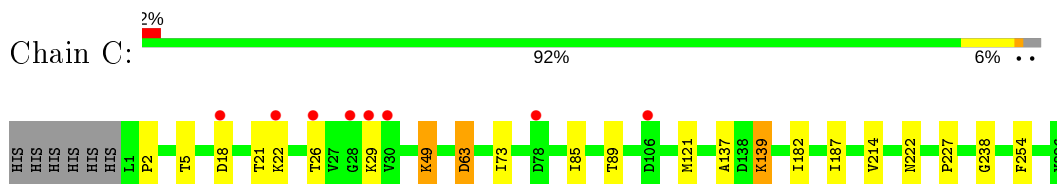
- Molecule 1: L-asparaginase 2

- Molecule 1: L-asparaginase 2

4   Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	152.00Å 62.39Å 142.27Å 90.00° 117.92° 90.00°	Depositor
Resolution (Å)	12.97 – 1.60 12.97 – 1.60	Depositor EDS
% Data completeness (in resolution range)	96.9 (12.97-1.60) 96.9 (12.97-1.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$<I/\sigma(I)>$ [1]	1.75 (at 1.60Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.147 , 0.198 0.156 , 0.189	Depositor DCC
R_{free} test set	1545 reflections (1.03%)	wwPDB-VP
Wilson B-factor (Å^2)	19.3	Xtriage
Anisotropy	0.294	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.41 , 54.3	EDS
L-test for twinning[2]	$<|L|> = 0.49$, $<L^2> = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	10835	wwPDB-VP
Average B, all atoms (Å^2)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.52% of the height of the origin peak. No significant pseudotranslation is detected.*

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[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 (i)

5.1 Standard geometry (i)

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	0.47	1/2405 (0.0%)	0.58	0/3274
1	B	0.31	0/2466	0.53	0/3358
1	C	0.38	1/2466 (0.0%)	0.58	0/3358
1	D	0.42	2/2383 (0.1%)	0.54	0/3244
All	All	0.40	4/9720 (0.0%)	0.56	0/13234

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	-1	HIS	C-N	-14.52	1.00	1.34
1	D	316	PRO	C-N	-11.98	1.06	1.34
1	D	317	GLN	C-N	-6.46	1.19	1.34
1	C	63	ASP	C-N	-5.32	1.21	1.34

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	2368	0	2360	6	2
1	B	2429	0	2423	14	0
1	C	2429	0	2423	15	2
1	D	2348	0	2345	8	0
2	A	9	0	3	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	9	0	3	1	0
2	C	9	0	3	2	0
2	D	9	0	3	0	0
3	A	333	0	0	0	0
3	B	273	0	0	0	0
3	C	349	0	0	1	0
3	D	270	0	0	0	0
All	All	10835	0	9563	37	2

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 2.

All (37) 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:222:ASN:HD22	1:C:63:ASP:H	1.30	0.78
1:C:139:LYS:H	1:C:139:LYS:HE2	1.62	0.63
1:C:139:LYS:H	1:C:139:LYS:CE	2.15	0.59
1:B:29:LYS:O	1:B:29:LYS:HD2	2.03	0.58
1:A:222:ASN:ND2	1:C:63:ASP:H	1.99	0.58
1:B:76:ASP:HA	1:B:79:LYS:HE3	1.87	0.57
1:D:117:PRO:HG2	1:D:120:SER:HB3	1.86	0.56
1:D:182:ILE:HG12	1:D:187:ILE:HG12	1.88	0.55
1:B:182:ILE:HG12	1:B:187:ILE:HG12	1.89	0.55
1:B:227:PRO:HB3	1:D:227:PRO:HB3	1.90	0.54
1:C:139:LYS:HD2	3:C:1285:HOH:O	2.07	0.54
1:B:14:ALA:HA	1:B:30:VAL:HG22	1.92	0.52
1:A:182:ILE:HG12	1:A:187:ILE:HG12	1.92	0.51
1:C:182:ILE:HG12	1:C:187:ILE:HG12	1.93	0.51
1:C:26:THR:HB	1:C:29:LYS:NZ	2.27	0.49
1:D:106:ASP:OD1	1:D:106:ASP:N	2.44	0.48
1:C:89:THR:OG1	2:C:401:ASP:OD2	2.26	0.47
1:B:29:LYS:C	1:B:29:LYS:HD2	2.36	0.46
1:A:227:PRO:HB3	1:C:227:PRO:HB3	1.96	0.46
1:C:73:ILE:HD11	1:C:85:ILE:HD11	1.98	0.44
1:B:89:THR:OG1	2:B:401:ASP:OD2	2.30	0.44
1:C:21:THR:HG22	1:C:121:MET:CE	2.47	0.44
1:B:106:ASP:N	1:B:106:ASP:OD2	2.50	0.43
1:C:2:PRO:HG3	1:C:137:ALA:HB1	2.00	0.43
1:B:208:LEU:HA	1:B:208:LEU:HD23	1.90	0.43
1:B:224:SER:HB2	1:D:236:TYR:OH	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2:PRO:HG3	1:B:137:ALA:HB1	2.01	0.42
1:B:14:ALA:CA	1:B:30:VAL:HG22	2.50	0.42
1:C:214:VAL:HA	1:C:238:GLY:O	2.20	0.41
1:B:259:THR:O	1:B:263:THR:HG23	2.20	0.41
1:C:5:THR:HA	1:C:49:LYS:HG2	2.03	0.41
1:B:236:TYR:OH	1:D:224:SER:HB2	2.20	0.41
1:D:261:ALA:HA	1:D:265:THR:O	2.20	0.41
1:C:139:LYS:H	1:C:139:LYS:CD	2.33	0.41
1:D:66:TRP:HB3	1:D:98:PHE:CE2	2.56	0.40
1:A:73:ILE:HD11	1:A:85:ILE:HD11	2.03	0.40
1:A:283:GLU:OE2	2:C:401:ASP:N	2.55	0.40

All (2) 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)
1:A:184:ASN:OD1	1:C:22:LYS:NZ[2_757]	1.81	0.39
1:A:184:ASN:CG	1:C:22:LYS:NZ[2_757]	2.08	0.12

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	312/332 (94%)	307 (98%)	5 (2%)	0	100	100
1	B	324/332 (98%)	319 (98%)	5 (2%)	0	100	100
1	C	324/332 (98%)	321 (99%)	3 (1%)	0	100	100
1	D	310/332 (93%)	305 (98%)	5 (2%)	0	100	100
All	All	1270/1328 (96%)	1252 (99%)	18 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	259/272 (95%)	255 (98%)	4 (2%)	65	44
1	B	266/272 (98%)	262 (98%)	4 (2%)	65	44
1	C	266/272 (98%)	261 (98%)	5 (2%)	57	34
1	D	257/272 (94%)	254 (99%)	3 (1%)	71	54
All	All	1048/1088 (96%)	1032 (98%)	16 (2%)	65	44

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

Mol	Chain	Res	Type
1	A	29	LYS
1	A	49	LYS
1	A	251	LYS
1	A	254	PHE
1	C	18	ASP
1	C	49	LYS
1	C	139	LYS
1	C	222	ASN
1	C	254	PHE
1	B	43	LYS
1	B	121	MET
1	B	222	ASN
1	B	254	PHE
1	D	222	ASN
1	D	251	LYS
1	D	254	PHE

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

Mol	Chain	Res	Type
1	A	222	ASN

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)

4 ligands are modelled in this entry.

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)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	D	2
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	D	317:GLN	C	318:GLN	N	1.19

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Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	D	316:PRO	C	317:GLN	N	1.06
1	A	-1:HIS	C	0:HIS	N	1.00

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	316/332 (95%)	-0.40	12 (3%)	40	37	13, 20, 42, 73	0
1	B	326/332 (98%)	-0.29	11 (3%)	45	42	18, 27, 48, 74	0
1	C	326/332 (98%)	-0.45	8 (2%)	57	55	13, 20, 45, 62	0
1	D	314/332 (94%)	-0.29	12 (3%)	40	37	18, 26, 49, 98	0
All	All	1282/1328 (96%)	-0.36	43 (3%)	45	42	13, 23, 47, 98	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	-1	HIS	9.1
1	D	15	GLY	7.7
1	D	30	VAL	6.1
1	A	28	GLY	5.4
1	A	30	VAL	5.2
1	D	28	GLY	5.1
1	A	37	ASN	4.3
1	D	37	ASN	4.3
1	A	0	HIS	4.1
1	D	29	LYS	4.1
1	C	18	ASP	4.1
1	A	29	LYS	4.1
1	C	28	GLY	4.1
1	A	15	GLY	3.8
1	B	106	ASP	3.6
1	B	21	THR	3.5
1	D	34	ASN	3.3
1	B	29	LYS	3.2
1	D	31	GLY	3.1
1	D	35	LEU	3.1
1	A	34	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	317	GLN	2.8
1	B	18	ASP	2.7
1	D	33	GLU	2.6
1	B	263	THR	2.5
1	D	32	VAL	2.4
1	B	24	SER	2.3
1	A	31	GLY	2.3
1	C	30	VAL	2.3
1	B	78	ASP	2.3
1	D	106	ASP	2.3
1	D	78	ASP	2.2
1	C	26	THR	2.2
1	B	34	ASN	2.2
1	C	106	ASP	2.1
1	B	264	GLY	2.1
1	C	22	LYS	2.1
1	B	262	LYS	2.1
1	A	264	GLY	2.0
1	A	35	LEU	2.0
1	C	29	LYS	2.0
1	C	78	ASP	2.0
1	A	78	ASP	2.0

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. 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	B-factors($Å^2$)	Q<0.9
2	ASP	D	401	9/9	0.79	0.24	22,22,23,23	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	ASP	A	401	9/9	0.80	0.19	17,17,17,17	0
2	ASP	B	401	9/9	0.84	0.16	20,20,21,21	0
2	ASP	C	401	9/9	0.85	0.13	18,18,18,19	0

6.5 Other polymers (i)

There are no such residues in this entry.