



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:14 PM UTC

PDB ID : 5K3O / pdb_00005k3o
Title : Wolinella succinogenes L-asparaginase P121 and L-Aspartic acid
Authors : Nguyen, H.A.; Lave, A.
Deposited on : 2016-05-19
Resolution : 1.70 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

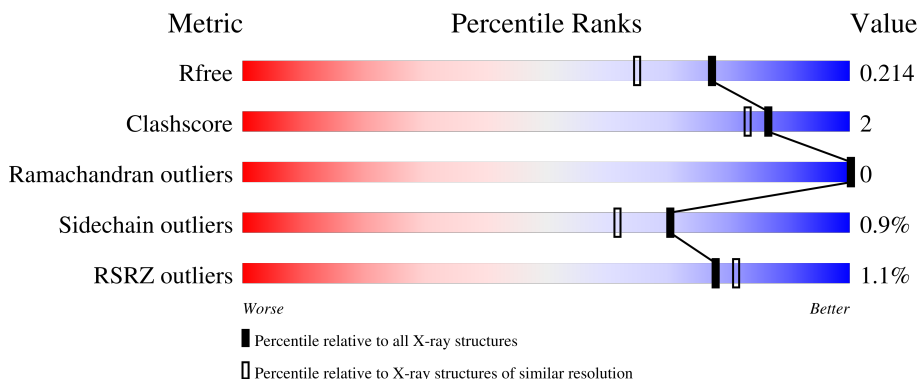
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

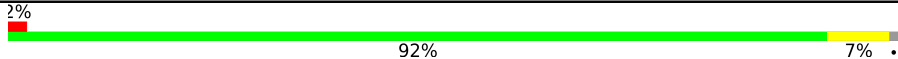
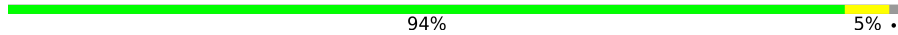
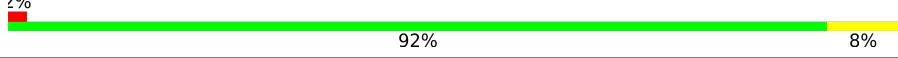
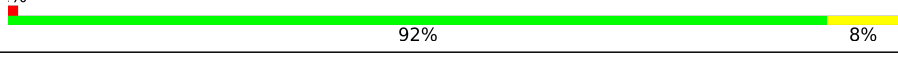
The reported resolution of this entry is 1.70 Å.

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}	180053	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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 of 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 $\leq 5\%$. 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	330	
1	B	330	
1	C	330	
1	D	330	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10710 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.

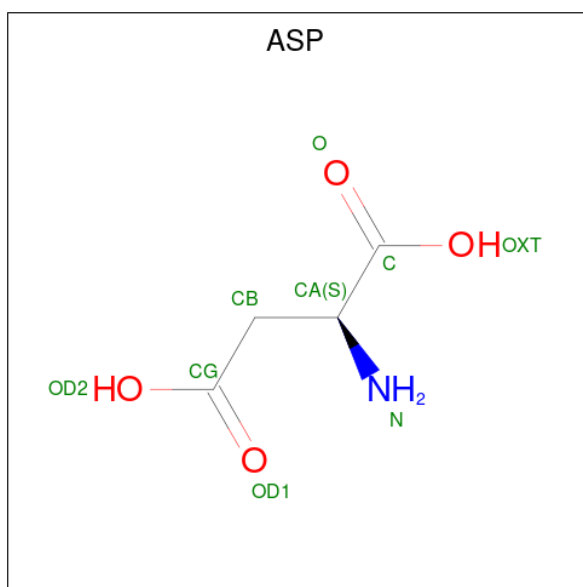
- Molecule 1 is a protein called L-asparaginase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	328	Total	C	N	O	S	0	1	0
			2443	1531	420	484	8			
1	B	328	Total	C	N	O	S	0	1	0
			2443	1531	420	484	8			
1	C	328	Total	C	N	O	S	0	1	0
			2443	1531	420	484	8			
1	D	328	Total	C	N	O	S	0	0	0
			2434	1526	418	482	8			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	121	PRO	SER	engineered mutation	UNP P50286
B	121	PRO	SER	engineered mutation	UNP P50286
C	121	PRO	SER	engineered mutation	UNP P50286
D	121	PRO	SER	engineered mutation	UNP P50286

- Molecule 2 is ASPARTIC ACID (CCD ID: ASP) (formula: C₄H₇NO₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			9	4	1	4		
2	B	1	Total	C	N	O	0	0
			9	4	1	4		
2	C	1	Total	C	N	O	0	0
			9	4	1	4		
2	D	1	Total	C	N	O	0	0
			9	4	1	4		

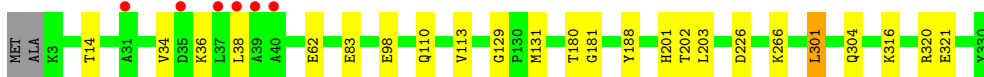
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	215	Total	O	0	0
			215	215		
3	B	226	Total	O	0	0
			226	226		
3	C	216	Total	O	0	0
			216	216		
3	D	254	Total	O	0	0
			254	254		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

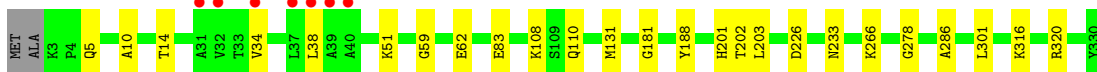
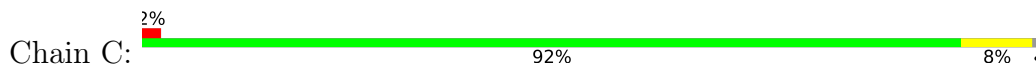
- Molecule 1: L-asparaginase



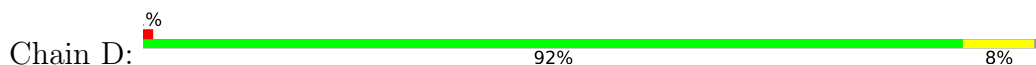
- Molecule 1: L-asparaginase



- Molecule 1: L-asparaginase



- Molecule 1: L-asparaginase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	144.75Å 71.46Å 141.50Å 90.00° 118.07° 90.00°	Depositor
Resolution (Å)	30.00 – 1.70 30.00 – 1.70	Depositor EDS
% Data completeness (in resolution range)	98.8 (30.00-1.70) 97.1 (30.00-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.26 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.8.0107	Depositor
R, R_{free}	0.175 , 0.207 0.185 , 0.214	Depositor DCC
R_{free} test set	6531 reflections (4.63%)	wwPDB-VP
Wilson B-factor (Å ²)	21.1	Xtriage
Anisotropy	0.301	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 36.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.032 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	10710	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% 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](#)

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	1.18	5/2476 (0.2%)	1.01	2/3356 (0.1%)
1	B	1.20	3/2476 (0.1%)	0.99	2/3356 (0.1%)
1	C	1.30	5/2476 (0.2%)	1.05	3/3356 (0.1%)
1	D	1.34	5/2467 (0.2%)	1.05	4/3344 (0.1%)
All	All	1.26	18/9895 (0.2%)	1.03	11/13412 (0.1%)

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	286	ALA	CA-C	7.90	1.56	1.52
1	C	108	LYS	CA-C	-6.70	1.45	1.53
1	D	180	THR	C-N	-6.32	1.25	1.33
1	B	301	LEU	CA-C	6.29	1.59	1.53
1	C	266	LYS	N-CA	6.04	1.53	1.46
1	C	233	ASN	CA-C	-6.04	1.45	1.52
1	A	266	LYS	N-CA	5.90	1.53	1.46
1	C	10	ALA	CA-C	-5.63	1.45	1.52
1	D	69	LEU	CA-C	-5.58	1.45	1.52
1	B	255	PRO	N-CA	5.58	1.54	1.47
1	D	16	ALA	CA-CB	5.30	1.60	1.53
1	A	301	LEU	CA-C	5.29	1.58	1.53
1	D	298	THR	N-CA	-5.29	1.39	1.46
1	A	113	VAL	N-CA	-5.13	1.39	1.46
1	A	301	LEU	C-O	5.12	1.29	1.24
1	B	245	HIS	N-CA	-5.11	1.40	1.46
1	A	180	THR	C-N	-5.04	1.27	1.33
1	D	261	LEU	CA-C	5.01	1.59	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	181	GLY	N-CA-C	-7.59	99.67	110.75
1	B	181	GLY	N-CA-C	-7.38	99.98	110.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	181	GLY	N-CA-C	-7.24	100.17	110.75
1	B	277	VAL	N-CA-C	6.79	117.57	110.72
1	C	181	GLY	N-CA-C	-6.55	101.19	110.75
1	A	129	GLY	O-C-N	6.04	123.25	121.07
1	D	170	THR	N-CA-C	5.36	121.37	114.56
1	D	278	GLY	N-CA-C	5.18	119.12	113.58
1	C	59	GLY	N-CA-C	-5.18	105.30	111.36
1	D	180	THR	N-CA-C	-5.17	99.80	110.80
1	C	278	GLY	N-CA-C	5.11	119.05	113.58

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	2443	0	2497	10	1
1	B	2443	0	2497	8	0
1	C	2443	0	2497	11	1
1	D	2434	0	2490	13	0
2	A	9	0	3	1	0
2	B	9	0	3	1	0
2	C	9	0	3	1	0
2	D	9	0	3	2	0
3	A	215	0	0	3	0
3	B	226	0	0	3	0
3	C	216	0	0	4	0
3	D	254	0	0	6	1
All	All	10710	0	9993	42	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 (42) 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:62:GLU:OE2	3:A:501:HOH:O	1.73	1.05
1:C:320:ARG:NH1	3:C:502:HOH:O	2.18	0.77
1:C:62:GLU:OE1	3:C:501:HOH:O	2.02	0.76
1:C:110[A]:GLN:HG3	1:C:203:LEU:HD12	1.84	0.58
1:B:316:LYS:NZ	3:B:502:HOH:O	2.36	0.57
1:B:201:HIS:HD2	1:B:202:THR:OG1	1.88	0.56
1:C:201:HIS:HD2	1:C:202:THR:OG1	1.88	0.56
1:D:201:HIS:HD2	1:D:202:THR:OG1	1.88	0.55
1:A:201:HIS:HE1	3:A:639:HOH:O	1.90	0.55
1:A:201:HIS:HD2	1:A:202:THR:OG1	1.89	0.55
1:A:110[A]:GLN:HG3	1:A:203:LEU:HD12	1.88	0.54
1:B:110[B]:GLN:HG3	1:B:203:LEU:HD12	1.90	0.53
1:B:301:LEU:C	3:B:566:HOH:O	2.52	0.53
1:A:131:MET:HE1	1:A:188:TYR:CE2	2.46	0.51
1:B:14:THR:OG1	2:B:400:ASP:CG	2.54	0.50
1:D:156:ASP:OD2	3:D:501:HOH:O	2.18	0.49
1:C:131:MET:HE1	1:C:188:TYR:CE2	2.47	0.49
1:C:201:HIS:HE1	3:C:590:HOH:O	1.95	0.49
1:D:131:MET:HE1	1:D:188:TYR:CE2	2.48	0.47
1:A:14:THR:OG1	2:A:400:ASP:CG	2.57	0.47
1:A:301:LEU:C	3:A:609:HOH:O	2.57	0.47
1:D:320:ARG:NH1	3:D:510:HOH:O	2.47	0.46
1:B:131:MET:HE1	1:B:188:TYR:CE2	2.51	0.46
1:C:14:THR:OG1	2:C:400:ASP:CG	2.58	0.46
1:D:3:LYS:N	3:D:511:HOH:O	2.49	0.45
1:C:5:GLN:NE2	1:C:51:LYS:HE2	2.32	0.45
1:A:34:VAL:O	1:A:38:LEU:HG	2.16	0.45
1:D:14:THR:OG1	2:D:400:ASP:CG	2.59	0.45
1:D:34:VAL:O	1:D:38:LEU:HG	2.17	0.45
1:D:201:HIS:HE1	3:D:619:HOH:O	1.99	0.44
1:B:34:VAL:O	1:B:38:LEU:HG	2.17	0.44
1:C:34:VAL:O	1:C:38:LEU:HG	2.17	0.44
1:C:301:LEU:C	3:C:581:HOH:O	2.60	0.44
1:D:119:MET:HE3	1:D:119:MET:HB3	1.92	0.43
1:C:5:GLN:CD	1:C:51:LYS:HE2	2.44	0.42
1:D:301:LEU:C	3:D:575:HOH:O	2.63	0.42
1:A:98:GLU:OE2	1:A:304:GLN:HB2	2.21	0.40
1:B:201:HIS:HE1	3:B:665:HOH:O	2.04	0.40
1:A:320:ARG:NH2	1:A:321:GLU:HG3	2.35	0.40
1:D:88:ILE:HD11	1:D:137:VAL:HA	2.03	0.40
1:D:13:GLY:HA2	2:D:400:ASP:OXT	2.22	0.40
1:D:110:GLN:HG3	3:D:627:HOH:O	2.21	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)
3:D:541:HOH:O	3:D:541:HOH:O[2_454]	1.67	0.53
1:A:83:GLU:OE2	1:C:83:GLU:OE1[4_445]	2.16	0.04

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	327/330 (99%)	318 (97%)	9 (3%)	0	100	100
1	B	327/330 (99%)	318 (97%)	9 (3%)	0	100	100
1	C	327/330 (99%)	318 (97%)	9 (3%)	0	100	100
1	D	326/330 (99%)	317 (97%)	9 (3%)	0	100	100
All	All	1307/1320 (99%)	1271 (97%)	36 (3%)	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	265/265 (100%)	262 (99%)	3 (1%)	65	54
1	B	265/265 (100%)	262 (99%)	3 (1%)	65	54
1	C	265/265 (100%)	263 (99%)	2 (1%)	73	65

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	264/265 (100%)	262 (99%)	2 (1%)	73	65
All	All	1059/1060 (100%)	1049 (99%)	10 (1%)	70	62

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	LYS
1	A	226	ASP
1	A	316	LYS
1	B	26	SER
1	B	53	GLU
1	B	226	ASP
1	C	226	ASP
1	C	316	LYS
1	D	26	SER
1	D	226	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	GLN
1	A	201	HIS
1	A	252	ASN
1	A	259	ASN
1	B	5	GLN
1	B	81	GLN
1	B	201	HIS
1	B	252	ASN
1	B	259	ASN
1	C	201	HIS
1	C	258	GLN
1	D	201	HIS
1	D	252	ASN
1	D	259	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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	$\# Z > 2$
2	ASP	B	400	-	7,8,8	1.40	2 (28%)	6,10,10	1.17	1 (16%)
2	ASP	C	400	-	7,8,8	1.06	0	6,10,10	2.07	2 (33%)
2	ASP	A	400	-	7,8,8	1.35	1 (14%)	6,10,10	1.51	2 (33%)
2	ASP	D	400	-	7,8,8	1.22	1 (14%)	6,10,10	1.20	1 (16%)

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	ASP	B	400	-	-	2/8/8/8	-
2	ASP	C	400	-	-	4/8/8/8	-
2	ASP	A	400	-	-	4/8/8/8	-
2	ASP	D	400	-	-	4/8/8/8	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	400	ASP	OD1-CG	2.66	1.30	1.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	400	ASP	OD1-CG	2.49	1.30	1.22
2	B	400	ASP	OD1-CG	2.47	1.30	1.22
2	B	400	ASP	OXT-C	-2.07	1.24	1.30

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	400	ASP	OD2-CG-CB	3.22	124.03	114.00
2	C	400	ASP	OXT-C-O	-2.89	117.53	124.08
2	B	400	ASP	OXT-C-O	-2.70	117.94	124.08
2	A	400	ASP	OD2-CG-CB	2.39	121.43	114.00
2	D	400	ASP	OD1-CG-CB	-2.22	116.04	122.84
2	A	400	ASP	OXT-C-O	-2.06	119.41	124.08

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	400	ASP	O-C-CA-N
2	B	400	ASP	O-C-CA-N
2	C	400	ASP	O-C-CA-N
2	D	400	ASP	O-C-CA-N
2	C	400	ASP	OXT-C-CA-N
2	D	400	ASP	OXT-C-CA-N
2	A	400	ASP	OXT-C-CA-N
2	B	400	ASP	OXT-C-CA-N
2	D	400	ASP	O-C-CA-CB
2	C	400	ASP	O-C-CA-CB
2	C	400	ASP	OXT-C-CA-CB
2	A	400	ASP	O-C-CA-CB
2	A	400	ASP	OXT-C-CA-CB
2	D	400	ASP	OXT-C-CA-CB

There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	400	ASP	1	0
2	C	400	ASP	1	0
2	A	400	ASP	1	0
2	D	400	ASP	2	0

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	328/330 (99%)	-0.02	6 (1%) 67 72	15, 25, 48, 64	1 (0%)
1	B	328/330 (99%)	-0.06	0 100 100	15, 24, 45, 63	1 (0%)
1	C	328/330 (99%)	0.00	7 (2%) 63 68	12, 23, 47, 58	1 (0%)
1	D	328/330 (99%)	-0.17	2 (0%) 85 88	12, 21, 41, 60	0
All	All	1312/1320 (99%)	-0.06	15 (1%) 78 81	12, 23, 45, 64	3 (0%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	37	LEU	3.4
1	C	38	LEU	2.8
1	A	37	LEU	2.6
1	A	31	ALA	2.5
1	A	40	ALA	2.5
1	C	31	ALA	2.4
1	C	39	ALA	2.4
1	C	34	VAL	2.4
1	D	37	LEU	2.4
1	C	32	VAL	2.4
1	A	39	ALA	2.3
1	A	38	LEU	2.3
1	D	38	LEU	2.3
1	C	40	ALA	2.2
1	A	35	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 oligosaccharides 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($\text{\AA}^2$)	Q<0.9
2	ASP	C	400	9/9	0.94	0.07	25,26,30,32	0
2	ASP	B	400	9/9	0.95	0.06	20,23,26,27	0
2	ASP	A	400	9/9	0.95	0.06	23,25,26,27	0
2	ASP	D	400	9/9	0.97	0.06	19,21,25,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.