



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:34 PM UTC

PDB ID : 7M11 / pdb_00007m11
Title : L-asparaginase from Escherichia coli with bound Aspartic acid
Authors : Souza, T.A.C.B.; Morais, S.B.; Morini, F.S.; Meissner, G.O.
Deposited on : 2021-03-11
Resolution : 1.83 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

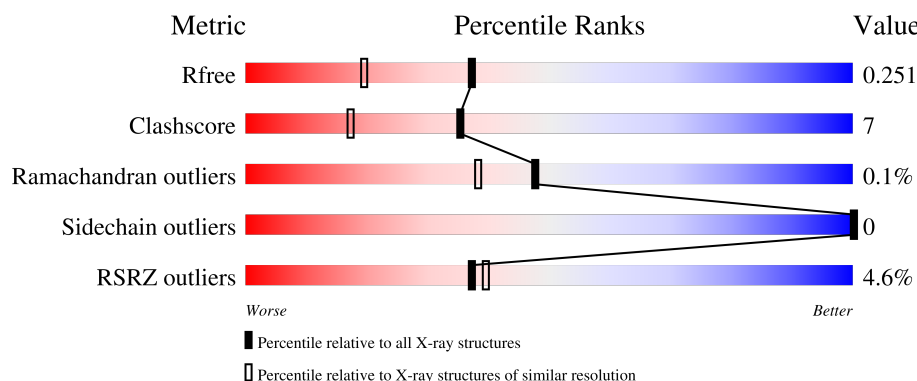
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

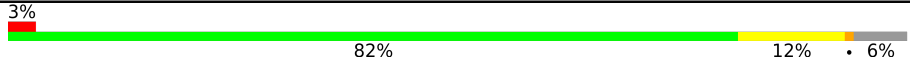
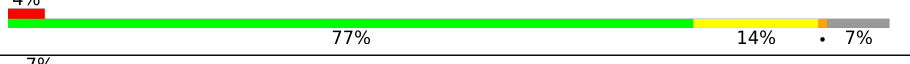
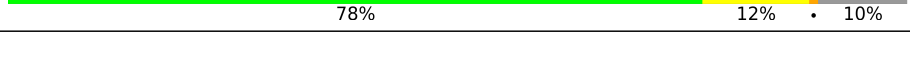
The reported resolution of this entry is 1.83 Å.

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	1296 (1.84-1.84)
Clashscore	190562	1329 (1.84-1.84)
Ramachandran outliers	187476	1318 (1.84-1.84)
Sidechain outliers	187428	1318 (1.84-1.84)
RSRZ outliers	180081	1296 (1.84-1.84)

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	326	
1	B	326	
1	C	326	
1	D	326	

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
2	ASP	C	401	-	-	X	-

2 Entry composition [i](#)

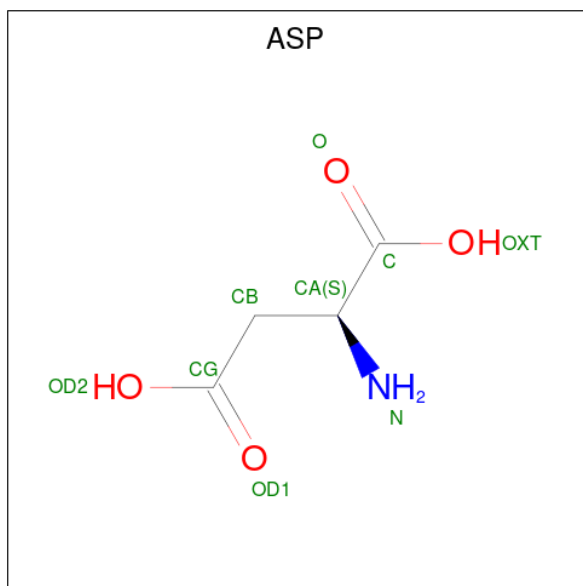
There are 3 unique types of molecules in this entry. The entry contains 9779 atoms, of which 0 are hydrogens and 0 are deuteriums.

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	308	Total	C	N	O	S	13	0	0
			2299	1439	394	458	8			
1	B	302	Total	C	N	O	S	9	0	0
			2271	1420	386	457	8			
1	C	326	Total	C	N	O	S	23	0	0
			2431	1517	415	491	8			
1	D	295	Total	C	N	O	S	18	0	0
			2212	1382	377	445	8			

- 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		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
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		

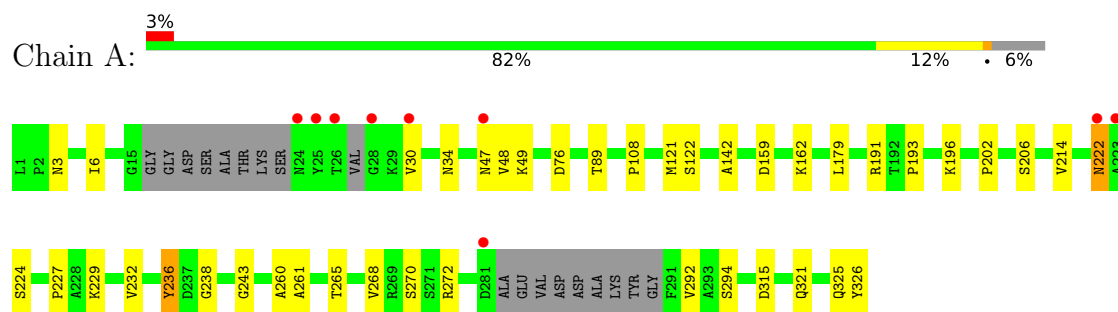
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	134	Total	O	0	0
			134	134		
3	B	131	Total	O	0	0
			131	131		
3	C	136	Total	O	0	0
			136	136		
3	D	129	Total	O	0	0
			129	129		

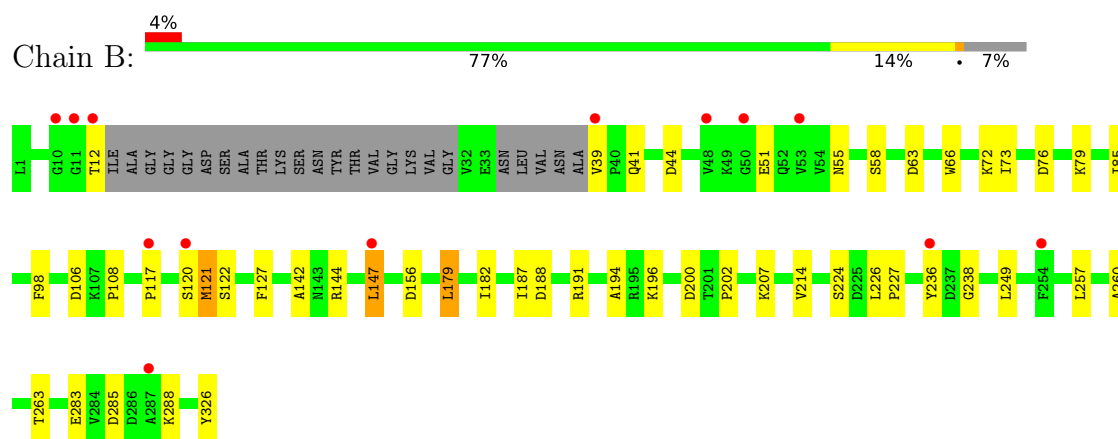
3 Residue-property plots [i](#)

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.

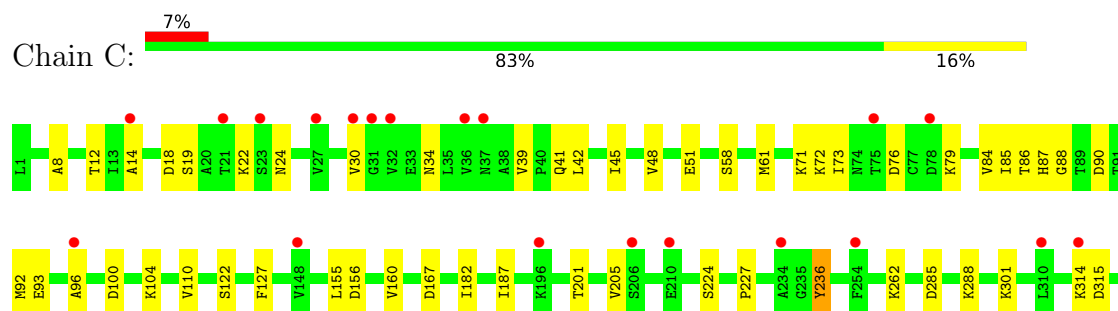
• Molecule 1: L-asparaginase 2



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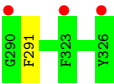
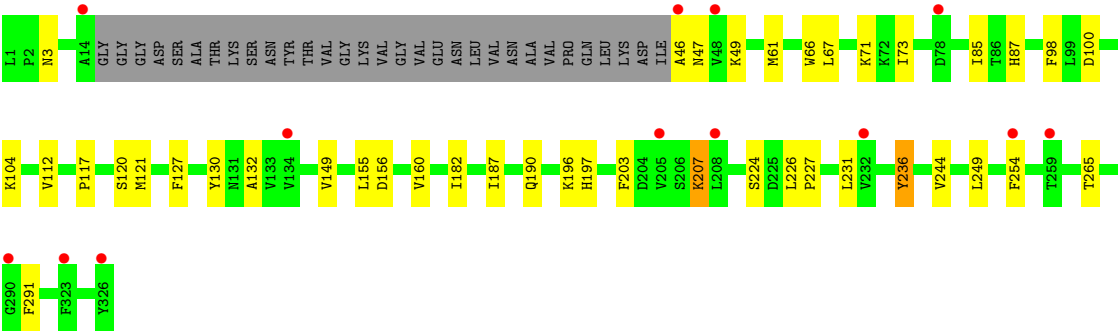
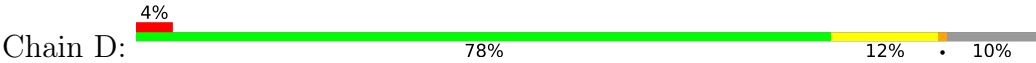


• Molecule 1: L-asparaginase 2





● Molecule 1: L-asparaginase 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.84Å 126.55Å 68.90Å 90.00° 107.36° 90.00°	Depositor
Resolution (Å)	47.18 – 1.83 47.18 – 1.83	Depositor EDS
% Data completeness (in resolution range)	87.8 (47.18-1.83) 87.7 (47.18-1.83)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	30.75 (at 1.83Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.206 , 0.252 0.206 , 0.251	Depositor DCC
R_{free} test set	2000 reflections (2.37%)	wwPDB-VP
Wilson B-factor (Å ²)	27.0	Xtriage
Anisotropy	0.300	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 39.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9779	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.84% 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

5.1 Standard geometry

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	0.41	0/2332	0.72	2/3172 (0.1%)
1	B	0.41	0/2305	0.74	5/3137 (0.2%)
1	C	0.42	1/2468 (0.0%)	0.72	0/3361
1	D	0.45	1/2246 (0.0%)	0.69	1/3058 (0.0%)
All	All	0.42	2/9351 (0.0%)	0.72	8/12728 (0.1%)

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
1	A	0	1
1	C	0	1
1	D	0	1
All	All	0	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	207	LYS	CG-CD	6.22	1.71	1.52
1	C	321	GLN	CD-NE2	5.51	1.44	1.33

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	147	LEU	CB-CG-CD2	-9.45	82.36	110.70
1	D	196	LYS	CA-CB-CG	-6.70	100.70	114.10
1	B	179	LEU	CB-CG-CD2	-5.80	93.28	110.70
1	B	147	LEU	CB-CG-CD1	5.74	127.92	110.70
1	B	147	LEU	CD1-CG-CD2	-5.72	98.21	110.80
1	B	121	MET	CB-CG-SD	-5.48	96.25	112.70
1	A	206	SER	CA-C-N	-5.40	114.06	122.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	206	SER	C-N-CA	-5.40	114.06	122.37

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	236	TYR	Sidechain
1	C	236	TYR	Sidechain
1	D	236	TYR	Sidechain

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	2299	0	2299	34	2
1	B	2271	0	2262	39	2
1	C	2431	0	2424	49	1
1	D	2212	0	2201	32	1
2	A	9	0	3	1	0
2	B	9	0	3	3	0
2	C	9	0	3	7	0
2	D	9	0	3	0	0
3	A	134	0	0	7	2
3	B	131	0	0	6	1
3	C	136	0	0	9	1
3	D	129	0	0	2	0
All	All	9779	0	9198	136	5

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 7.

All (136) 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:C:12:THR:OG1	2:C:401:ASP:O	1.88	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:112:VAL:HG11	1:D:132:ALA:HB2	1.51	0.90
1:C:58:SER:OG	2:C:401:ASP:OD2	1.91	0.88
1:A:196:LYS:NZ	1:A:325:GLN:OE1	2.07	0.86
1:C:87:HIS:HB3	1:C:92:MET:HE2	1.58	0.86
1:C:72:LYS:NZ	1:C:76:ASP:OD2	2.10	0.85
1:A:227:PRO:HB3	1:B:227:PRO:HB3	1.64	0.78
1:B:196:LYS:HD2	1:B:200:ASP:HB3	1.66	0.76
1:D:249:LEU:HB2	1:D:254:PHE:HB2	1.67	0.74
1:A:30:VAL:HG13	1:A:34:ASN:HB2	1.69	0.72
1:D:61:MET:HE3	1:D:87:HIS:NE2	2.04	0.72
1:B:12:THR:N	2:B:401:ASP:O	2.24	0.70
1:B:191:ARG:NH2	1:D:156:ASP:OD2	2.27	0.68
1:A:261:ALA:O	3:A:501:HOH:O	2.12	0.67
1:D:265:THR:OG1	3:D:501:HOH:O	2.06	0.67
1:B:58:SER:OG	2:B:401:ASP:OD2	2.08	0.66
1:A:3:ASN:OD1	3:A:502:HOH:O	2.14	0.65
1:B:51:GLU:OE2	1:B:72:LYS:NZ	2.27	0.63
1:D:46:ALA:N	3:D:506:HOH:O	2.31	0.63
1:C:72:LYS:NZ	3:C:508:HOH:O	2.31	0.63
1:B:196:LYS:HB2	1:B:326:TYR:CD2	2.34	0.63
1:B:121:MET:HE2	1:C:41:GLN:CD	2.24	0.63
1:C:227:PRO:HB3	1:D:227:PRO:HB3	1.81	0.63
1:A:179:LEU:HD21	1:A:193:PRO:HD3	1.80	0.62
1:C:262:LYS:NZ	3:C:510:HOH:O	2.32	0.61
1:C:34:ASN:ND2	3:C:511:HOH:O	2.33	0.61
1:D:112:VAL:HG12	1:D:149:VAL:O	2.02	0.58
1:A:76:ASP:OD2	3:A:504:HOH:O	2.17	0.58
1:C:236:TYR:OH	1:D:224:SER:HB2	2.03	0.57
1:B:144:ARG:HB2	1:B:147:LEU:HD11	1.87	0.57
1:D:67:LEU:O	1:D:71:LYS:HG3	2.05	0.57
1:B:182:ILE:HG12	1:B:187:ILE:HG12	1.88	0.56
1:B:260:ALA:O	1:B:263:THR:HG22	2.06	0.56
1:B:55:ASN:O	3:B:501:HOH:O	2.18	0.54
1:C:61:MET:HE3	1:C:87:HIS:NE2	2.22	0.54
1:D:66:TRP:HB3	1:D:98:PHE:CE2	2.43	0.54
1:A:159:ASP:OD1	3:A:505:HOH:O	2.18	0.54
1:C:24:ASN:O	3:C:501:HOH:O	2.19	0.53
1:B:66:TRP:HB3	1:B:98:PHE:CE1	2.44	0.53
1:A:47:ASN:OD1	3:A:502:HOH:O	2.19	0.52
1:B:156:ASP:HB3	1:B:179:LEU:CD1	2.40	0.52
1:B:214:VAL:HA	1:B:238:GLY:O	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:232:VAL:HG22	1:A:260:ALA:HB2	1.92	0.51
1:B:106:ASP:OD2	3:B:502:HOH:O	2.19	0.51
1:D:112:VAL:CG1	1:D:132:ALA:HB2	2.34	0.51
1:C:76:ASP:OD1	1:C:79:LYS:HE2	2.09	0.51
1:A:224:SER:HB2	1:B:236:TYR:OH	2.11	0.50
1:C:88:GLY:HA3	2:C:401:ASP:OD2	2.12	0.50
1:C:51:GLU:HB3	3:C:506:HOH:O	2.12	0.50
1:D:155:LEU:HD13	1:D:160:VAL:HG23	1.93	0.50
1:C:156:ASP:OD2	3:C:502:HOH:O	2.20	0.49
1:C:90:ASP:CG	2:C:401:ASP:CB	2.85	0.49
1:B:156:ASP:HB3	1:B:179:LEU:HD13	1.93	0.49
1:D:3:ASN:HA	1:D:49:LYS:HZ1	1.78	0.49
1:C:14:ALA:HA	1:C:30:VAL:CG1	2.43	0.48
1:B:76:ASP:OD1	1:B:79:LYS:HE2	2.13	0.48
1:C:12:THR:OG1	2:C:401:ASP:C	2.56	0.48
1:A:30:VAL:CG1	1:A:34:ASN:HB2	2.39	0.48
1:B:283:GLU:HB2	3:B:526:HOH:O	2.13	0.48
1:D:100:ASP:OD1	1:D:197:HIS:HB2	2.13	0.48
1:A:89:THR:OG1	1:A:162:LYS:NZ	2.45	0.48
1:D:73:ILE:HD11	1:D:85:ILE:HD11	1.96	0.47
1:B:285:ASP:OD2	1:B:288:LYS:HB2	2.13	0.47
1:C:39:VAL:O	1:C:39:VAL:HG13	2.13	0.47
1:B:188:ASP:OD2	3:B:503:HOH:O	2.21	0.47
1:C:61:MET:HB2	1:C:61:MET:HE2	1.70	0.47
1:C:224:SER:HB2	1:D:236:TYR:OH	2.14	0.47
1:C:90:ASP:CG	2:C:401:ASP:HB3	2.39	0.47
1:A:122:SER:HB2	1:D:127:PHE:HB2	1.96	0.47
1:B:41:GLN:OE1	1:B:41:GLN:N	2.41	0.46
1:B:226:LEU:HB2	1:B:227:PRO:HD3	1.97	0.46
1:C:18:ASP:N	3:C:501:HOH:O	2.26	0.46
1:C:155:LEU:HD13	1:C:160:VAL:HG23	1.98	0.46
1:C:93:GLU:HB2	1:C:301:LYS:HE2	1.97	0.46
1:C:314:LYS:HG3	3:C:531:HOH:O	2.16	0.46
1:B:39:VAL:N	3:B:518:HOH:O	2.49	0.46
1:B:122:SER:HB2	1:C:127:PHE:HB2	1.97	0.46
1:C:14:ALA:C	1:C:30:VAL:HG12	2.41	0.45
1:C:71:LYS:HE3	1:C:205:VAL:O	2.15	0.45
1:A:6:ILE:HD12	1:A:48:VAL:CG1	2.47	0.45
1:B:156:ASP:OD2	3:B:504:HOH:O	2.21	0.45
1:C:14:ALA:O	1:C:30:VAL:HG12	2.17	0.45
1:C:76:ASP:CG	1:C:79:LYS:HE2	2.42	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:8:ALA:HA	1:C:86:THR:OG1	2.18	0.44
1:D:254:PHE:CE1	1:D:291:PHE:HE2	2.35	0.44
1:B:127:PHE:HB2	1:C:122:SER:HB2	1.99	0.44
1:D:182:ILE:HG12	1:D:187:ILE:HG12	2.00	0.44
1:A:108:PRO:HD3	1:A:142:ALA:HB2	2.00	0.44
1:A:268:VAL:HG22	1:A:292:VAL:HB	2.00	0.44
1:B:73:ILE:HD11	1:B:85:ILE:HD11	2.00	0.44
1:A:236:TYR:OH	1:B:224:SER:HB2	2.18	0.44
1:D:117:PRO:HG2	1:D:120:SER:HB3	1.99	0.44
1:B:117:PRO:O	1:B:120:SER:HB2	2.18	0.43
1:C:39:VAL:HG13	1:C:42:LEU:HG	2.01	0.43
1:A:179:LEU:HG	1:A:191:ARG:O	2.18	0.43
1:C:19:SER:HB2	1:C:22:LYS:H	1.83	0.43
1:A:6:ILE:HD12	1:A:48:VAL:HG13	2.01	0.43
1:A:270:SER:HA	1:A:294:SER:OG	2.19	0.43
1:D:104:LYS:HA	1:D:104:LYS:HD3	1.84	0.43
1:A:202:PRO:HD3	1:A:326:TYR:CZ	2.53	0.42
1:C:96:ALA:O	1:C:100:ASP:HB2	2.20	0.42
1:C:90:ASP:OD2	2:C:401:ASP:HB3	2.20	0.42
1:C:167:ASP:OD2	3:C:503:HOH:O	2.22	0.42
1:B:108:PRO:HD3	1:B:142:ALA:HB2	2.00	0.42
1:C:42:LEU:O	1:C:45:ILE:HG12	2.19	0.42
1:A:321:GLN:O	1:A:325:GLN:HG2	2.19	0.42
1:C:73:ILE:HD11	1:C:85:ILE:HD11	2.00	0.42
1:C:90:ASP:HB3	1:D:244:VAL:HB	2.02	0.42
1:B:194:ALA:HB3	1:D:190:GLN:HA	2.02	0.42
1:C:182:ILE:HG12	1:C:187:ILE:HG12	2.01	0.42
1:B:202:PRO:HD3	1:B:326:TYR:OH	2.19	0.41
1:C:84:VAL:HA	1:C:110:VAL:O	2.20	0.41
1:D:121:MET:HE3	1:D:121:MET:HB2	1.77	0.41
1:A:222:ASN:HD22	1:B:63:ASP:CG	2.28	0.41
1:A:121:MET:HE1	1:D:130:TYR:CD1	2.55	0.41
1:A:232:VAL:CG2	1:A:260:ALA:HB2	2.50	0.41
1:A:260:ALA:HB1	1:A:265:THR:HB	2.02	0.41
1:B:249:LEU:CD2	1:B:257:LEU:HD12	2.50	0.41
1:D:231:LEU:HD22	1:D:236:TYR:CE2	2.55	0.41
1:A:121:MET:HE3	1:D:127:PHE:HD1	1.86	0.41
1:A:214:VAL:HA	1:A:238:GLY:O	2.21	0.41
1:B:12:THR:HG1	2:B:401:ASP:C	2.28	0.41
1:C:41:GLN:CD	1:C:41:GLN:H	2.28	0.41
1:C:285:ASP:OD2	1:C:288:LYS:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:47:ASN:O	1:D:49:LYS:NZ	2.52	0.41
1:A:121:MET:HE1	1:D:130:TYR:CG	2.56	0.41
1:A:243:GLY:O	1:A:272:ARG:HG2	2.21	0.41
1:B:121:MET:HE2	1:C:41:GLN:OE1	2.20	0.41
1:D:104:LYS:NZ	1:D:203:PHE:O	2.43	0.41
1:D:224:SER:OG	1:D:226:LEU:HB2	2.21	0.41
1:C:42:LEU:HD13	1:C:48:VAL:HG11	2.02	0.41
1:A:227:PRO:HD2	3:A:542:HOH:O	2.21	0.40
1:A:30:VAL:HG13	1:A:34:ASN:CB	2.44	0.40
2:A:401:ASP:N	1:B:283:GLU:OE1	2.54	0.40
1:C:104:LYS:NZ	1:C:201:THR:OG1	2.54	0.40
1:A:229:LYS:NZ	3:A:503:HOH:O	2.17	0.40

All (5) 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:C:315:ASP:OD2	1:D:207:LYS:NZ[1_655]	1.87	0.33
1:A:315:ASP:OD2	1:B:207:LYS:NZ[1_655]	2.08	0.12
1:A:49:LYS:NZ	1:B:44:ASP:OD2[1_656]	2.10	0.10
3:A:634:HOH:O	3:C:636:HOH:O[2_555]	2.16	0.04
3:A:518:HOH:O	3:B:624:HOH:O[1_656]	2.17	0.03

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	300/326 (92%)	294 (98%)	5 (2%)	1 (0%)	36	25
1	B	296/326 (91%)	287 (97%)	9 (3%)	0	100	100
1	C	324/326 (99%)	315 (97%)	9 (3%)	0	100	100
1	D	291/326 (89%)	285 (98%)	6 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1211/1304 (93%)	1181 (98%)	29 (2%)	1 (0%)	48 38

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	222	ASN

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	252/266 (95%)	252 (100%)	0	100 100
1	B	250/266 (94%)	250 (100%)	0	100 100
1	C	266/266 (100%)	266 (100%)	0	100 100
1	D	242/266 (91%)	242 (100%)	0	100 100
All	All	1010/1064 (95%)	1010 (100%)	0	100 100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	ASN
1	A	183	HIS
1	C	3	ASN
1	C	47	ASN
1	C	222	ASN
1	C	317	GLN
1	D	248	ASN
1	D	325	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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	A	401	-	7,8,8	1.20	1 (14%)	6,10,10	1.24	0
2	ASP	B	401	-	7,8,8	1.30	2 (28%)	6,10,10	0.92	1 (16%)
2	ASP	C	401	-	7,8,8	1.78	2 (28%)	6,10,10	1.12	0
2	ASP	D	401	-	7,8,8	1.17	1 (14%)	6,10,10	1.36	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	A	401	-	-	5/8/8/8	-
2	ASP	B	401	-	-	0/8/8/8	-
2	ASP	C	401	-	-	3/8/8/8	-
2	ASP	D	401	-	-	2/8/8/8	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	401	ASP	CB-CG	-3.30	1.42	1.51
2	C	401	ASP	OD2-CG	-2.55	1.22	1.30
2	D	401	ASP	OXT-C	-2.28	1.23	1.30
2	A	401	ASP	OXT-C	-2.26	1.23	1.30
2	B	401	ASP	OXT-C	-2.21	1.23	1.30
2	B	401	ASP	OD2-CG	-2.11	1.23	1.30

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	401	ASP	OXT-C-O	-2.69	117.99	124.08
2	B	401	ASP	OXT-C-O	-2.10	119.33	124.08

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	ASP	O-C-CA-N
2	A	401	ASP	OXT-C-CA-N
2	C	401	ASP	O-C-CA-CB
2	C	401	ASP	OXT-C-CA-CB
2	A	401	ASP	OXT-C-CA-CB
2	A	401	ASP	O-C-CA-CB
2	D	401	ASP	CA-CB-CG-OD1
2	D	401	ASP	CA-CB-CG-OD2
2	A	401	ASP	CA-CB-CG-OD2
2	C	401	ASP	OXT-C-CA-N

There are no ring outliers.

3 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	ASP	1	0
2	B	401	ASP	3	0
2	C	401	ASP	7	0

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	308/326 (94%)	0.43	9 (2%) 53 58	18, 31, 44, 68	3 (0%)
1	B	302/326 (92%)	0.50	13 (4%) 40 42	18, 32, 46, 71	2 (0%)
1	C	326/326 (100%)	0.72	22 (6%) 24 25	24, 34, 51, 61	6 (1%)
1	D	295/326 (90%)	0.62	13 (4%) 39 42	21, 33, 45, 51	5 (1%)
All	All	1231/1304 (94%)	0.57	57 (4%) 37 39	18, 33, 47, 71	16 (1%)

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	25	TYR	5.2
1	C	23	SER	5.2
1	A	222	ASN	4.3
1	C	30	VAL	4.2
1	C	254	PHE	3.7
1	C	21	THR	3.4
1	A	28	GLY	3.1
1	B	11	GLY	3.1
1	C	321	GLN	3.0
1	B	287	ALA	2.9
1	C	234	ALA	2.9
1	B	236	TYR	2.8
1	C	31	GLY	2.8
1	C	14	ALA	2.7
1	D	254	PHE	2.7
1	C	206	SER	2.7
1	D	232	VAL	2.7
1	C	96	ALA	2.6
1	B	254	PHE	2.6
1	D	326	TYR	2.6
1	D	48	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	290	GLY	2.5
1	D	78	ASP	2.5
1	C	196	LYS	2.5
1	B	39	VAL	2.5
1	B	10	GLY	2.4
1	B	12	THR	2.4
1	B	117	PRO	2.4
1	A	223	ALA	2.4
1	D	14	ALA	2.4
1	C	148	VAL	2.4
1	C	210	GLU	2.3
1	C	310	LEU	2.3
1	A	26	THR	2.3
1	D	46	ALA	2.2
1	A	281	ASP	2.2
1	C	314	LYS	2.2
1	C	75	THR	2.2
1	D	259	THR	2.2
1	C	322	ILE	2.2
1	C	37	ASN	2.2
1	A	30	VAL	2.2
1	C	32	VAL	2.2
1	B	147	LEU	2.2
1	B	48	VAL	2.1
1	C	78	ASP	2.1
1	C	27	VAL	2.1
1	A	24	ASN	2.1
1	A	47	ASN	2.1
1	B	53	VAL	2.1
1	D	323	PHE	2.0
1	B	120	SER	2.0
1	D	134	VAL	2.0
1	D	208	LEU	2.0
1	B	50	GLY	2.0
1	C	36	VAL	2.0
1	D	205	VAL	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 [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	401	9/9	0.60	0.15	35,39,46,46	0
2	ASP	B	401	9/9	0.81	0.10	34,39,46,48	0
2	ASP	A	401	9/9	0.92	0.07	27,30,33,34	0
2	ASP	D	401	9/9	0.94	0.06	28,31,38,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.