



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:19 PM UTC

PDB ID : 5JWG / pdb_00005jwg
Title : Crystal structure of Porphyromonas endodontalis DPP11 in complex with dipeptide Arg-Asp
Authors : Bezerra, G.A.; Fedosyuk, S.; Ohara-Nemoto, Y.; Nemoto, T.K.; Djinic-Carugo, K.
Deposited on : 2016-05-12
Resolution : 2.20 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

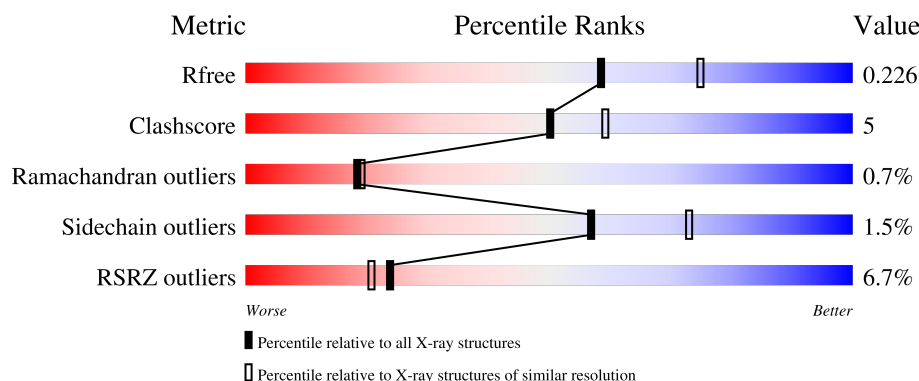
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	703	3% 87% 12% .
1	B	703	11% 86% 12% ..

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11235 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.

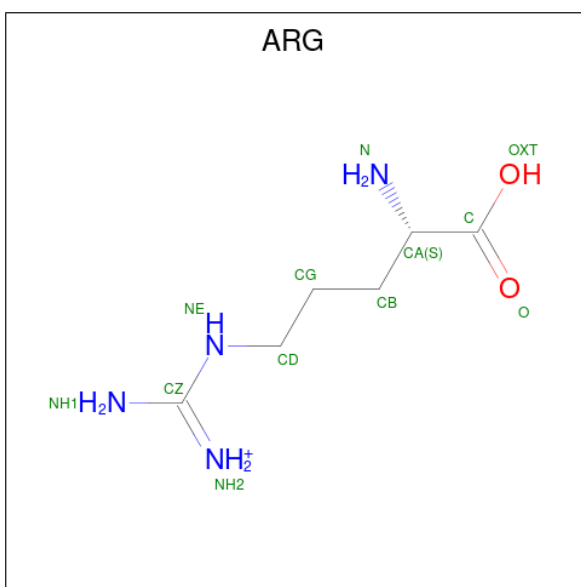
- Molecule 1 is a protein called Asp/Glu-specific dipeptidyl-peptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	703	Total	C	N	O	S	0	1	0
			5558	3517	967	1048	26			
1	B	695	Total	C	N	O	S	0	0	0
			5211	3313	896	979	23			

There are 16 discrepancies between the modelled and reference sequences:

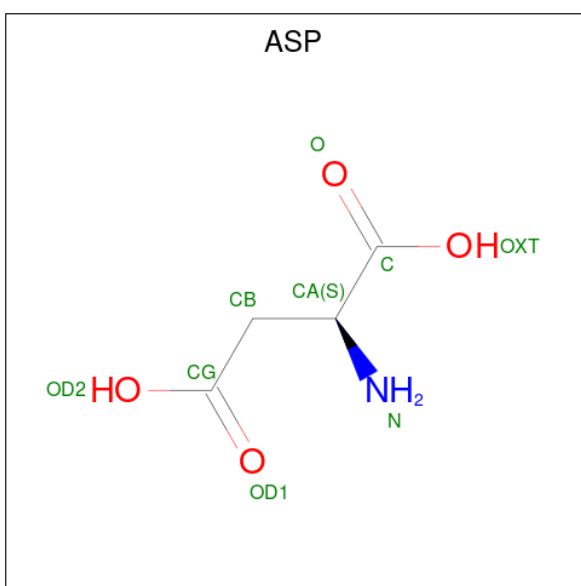
Chain	Residue	Modelled	Actual	Comment	Reference
A	21	MET	-	initiating methionine	UNP F8WQK8
A	652	ALA	SER	engineered mutation	UNP F8WQK8
A	718	HIS	-	expression tag	UNP F8WQK8
A	719	HIS	-	expression tag	UNP F8WQK8
A	720	HIS	-	expression tag	UNP F8WQK8
A	721	HIS	-	expression tag	UNP F8WQK8
A	722	HIS	-	expression tag	UNP F8WQK8
A	723	HIS	-	expression tag	UNP F8WQK8
B	21	MET	-	initiating methionine	UNP F8WQK8
B	652	ALA	SER	engineered mutation	UNP F8WQK8
B	718	HIS	-	expression tag	UNP F8WQK8
B	719	HIS	-	expression tag	UNP F8WQK8
B	720	HIS	-	expression tag	UNP F8WQK8
B	721	HIS	-	expression tag	UNP F8WQK8
B	722	HIS	-	expression tag	UNP F8WQK8
B	723	HIS	-	expression tag	UNP F8WQK8

- Molecule 2 is ARGININE (CCD ID: ARG) (formula: C₆H₁₅N₄O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			11	6	4	1		
2	B	1	Total	C	N	O	0	0
			11	6	4	1		

- Molecule 3 is ASPARTIC ACID (CCD ID: ASP) (formula: C₄H₇NO₄).



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

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Na	0	0
			1	1		

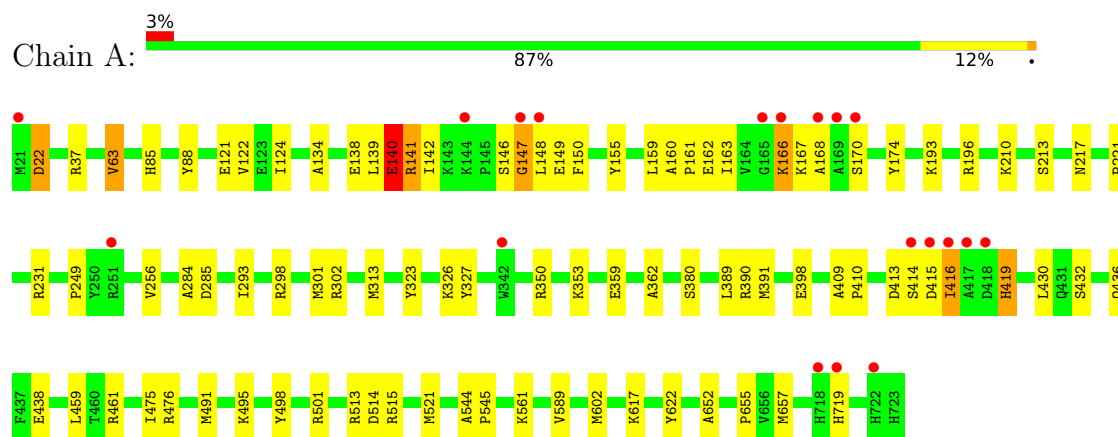
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	283	Total	O	0	0
			283	283		
6	B	138	Total	O	0	0
			138	138		

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: Asp/Glu-specific dipeptidyl-peptidase



• Molecule 1: Asp/Glu-specific dipeptidyl-peptidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	111.81Å 114.40Å 147.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.24 – 2.20 45.24 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.5 (45.24-2.20) 99.5 (45.24-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 2.20Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.181 , 0.227 0.186 , 0.226	Depositor DCC
R_{free} test set	4811 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	46.3	Xtriage
Anisotropy	0.264	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 43.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.021 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11235	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.82% 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

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

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	0.54	0/5687	0.84	4/7693 (0.1%)
1	B	0.51	0/5325	0.87	5/7232 (0.1%)
All	All	0.53	0/11012	0.85	9/14925 (0.1%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	21	MET	CB-CG-SD	-9.65	83.74	112.70
1	B	389	LEU	N-CA-C	-5.93	104.73	111.14
1	B	213	SER	N-CA-C	5.33	117.50	111.11
1	A	415	ASP	N-CA-C	-5.28	103.05	110.50
1	A	561	LYS	N-CA-C	-5.24	103.12	110.35
1	A	719	HIS	N-CA-C	-5.17	103.71	110.53
1	B	713	ILE	CA-C-N	-5.09	116.51	122.00
1	B	713	ILE	C-N-CA	-5.09	116.51	122.00
1	A	419	HIS	N-CA-C	5.03	121.52	110.80

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	5558	0	5374	63	0
1	B	5211	0	4841	49	0
2	A	11	0	12	0	0
2	B	11	0	12	0	0
3	A	9	0	4	1	0
3	B	9	0	4	1	0
4	A	4	0	6	2	0
5	A	1	0	0	0	0
6	A	283	0	0	5	0
6	B	138	0	0	3	0
All	All	11235	0	10253	112	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 5.

All (112) 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:313:MET:SD	6:A:1133:HOH:O	2.29	0.91
1:A:657:MET:SD	6:A:1158:HOH:O	2.30	0.89
1:A:416:ILE:HA	1:A:513:ARG:HE	1.39	0.86
1:B:388:ASP:HA	1:B:391:MET:HB3	1.56	0.85
1:A:313:MET:HE1	1:A:327:TYR:HB2	1.61	0.80
1:A:147:GLY:O	1:A:148:LEU:HG	1.85	0.76
1:A:298:ARG:HG3	1:A:302:ARG:HD3	1.68	0.74
1:A:22:ASP:HB2	1:A:37:ARG:NH2	2.05	0.72
1:B:310:LEU:HA	1:B:313:MET:HE2	1.77	0.67
1:A:514:ASP:O	1:A:515:ARG:HB3	1.97	0.65
1:B:22:ASP:O	6:B:901:HOH:O	2.15	0.65
1:B:298:ARG:HG3	1:B:302:ARG:HD3	1.79	0.63
1:A:146:SER:O	1:A:147:GLY:C	2.43	0.62
1:A:313:MET:HE2	1:A:323:TYR:HB3	1.83	0.61
1:A:617:LYS:NZ	6:A:904:HOH:O	2.34	0.59
1:B:213:SER:O	1:B:217:ASN:HB2	2.01	0.59
1:A:166:LYS:O	1:A:170:SER:HB3	2.02	0.59
1:A:655:PRO:HB2	1:A:657:MET:HE3	1.85	0.59
1:B:96:SER:OG	1:B:98:GLN:O	2.14	0.59
1:A:301:MET:HE1	1:A:461:ARG:HG2	1.85	0.58
1:B:389:LEU:HD11	1:B:538:GLU:HB3	1.84	0.58
1:A:138:GLU:HG2	1:A:159:LEU:HD22	1.84	0.58
1:A:231:ARG:HD3	1:A:249:PRO:HB3	1.86	0.58
1:B:340:ALA:O	1:B:344:ILE:HG12	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:148:LEU:HD12	1:A:544:ALA:O	2.04	0.58
1:B:22:ASP:OD2	1:B:37:ARG:NE	2.33	0.56
1:B:114:GLU:OE2	1:B:708:ARG:NH2	2.37	0.56
1:A:148:LEU:HD22	1:A:150:PHE:CE1	2.41	0.56
1:A:139:LEU:C	1:A:140:GLU:O	2.45	0.55
1:B:344:ILE:HD13	1:B:349:LEU:HD12	1.87	0.55
1:A:166:LYS:HG2	1:A:167:LYS:N	2.19	0.55
1:B:301:MET:HE1	1:B:461:ARG:HG2	1.88	0.55
1:A:134:ALA:HB1	1:A:163:ILE:HD12	1.89	0.54
1:B:123:GLU:HG2	1:B:194:VAL:HG22	1.90	0.54
1:A:146:SER:HB3	1:A:149:GLU:OE1	2.07	0.53
1:A:652:ALA:HB2	3:A:802:ASP:C	2.33	0.53
1:A:544:ALA:HB3	1:A:545:PRO:HD3	1.89	0.53
1:B:655:PRO:HB2	1:B:657:MET:HE3	1.89	0.53
1:A:301:MET:HE1	1:A:461:ARG:CG	2.39	0.53
1:A:213:SER:O	1:A:217:ASN:HB2	2.07	0.53
1:A:256:VAL:HG11	1:A:657:MET:HE2	1.91	0.52
1:A:491:MET:HE2	1:A:495:LYS:HD2	1.91	0.52
1:A:210:LYS:HE2	1:A:213:SER:HB2	1.92	0.52
1:A:432:SER:O	1:A:436:GLN:HG2	2.09	0.52
1:A:147:GLY:C	1:A:148:LEU:HG	2.35	0.51
1:A:142:ILE:HD11	1:A:155:TYR:OH	2.10	0.51
1:A:438:GLU:OE2	1:A:501:ARG:NH2	2.44	0.51
1:B:355:ALA:O	1:B:359:GLU:HG2	2.11	0.50
1:B:301:MET:HE1	1:B:461:ARG:CG	2.41	0.50
1:A:359:GLU:HA	4:A:803:EDO:H11	1.93	0.50
1:A:459:LEU:HD11	1:A:475:ILE:HD13	1.94	0.49
1:B:285:ASP:OD2	1:B:380:SER:OG	2.24	0.49
1:B:21:MET:HA	1:B:681:TYR:OH	2.12	0.49
1:A:313:MET:HE2	1:A:323:TYR:CB	2.43	0.49
1:B:288:THR:O	1:B:292:GLU:HG3	2.13	0.49
1:A:148:LEU:HD22	1:A:150:PHE:CD1	2.49	0.48
1:A:63:VAL:HG22	1:A:124:ILE:HG12	1.96	0.48
1:A:409:ALA:HB2	1:A:521:MET:HE3	1.96	0.48
1:B:652:ALA:HB2	3:B:802:ASP:C	2.39	0.48
1:A:63:VAL:HG13	1:A:122:VAL:HG13	1.96	0.47
1:A:140:GLU:O	1:A:141:ARG:HB2	2.13	0.47
1:B:93:LYS:O	1:B:443:LYS:NZ	2.47	0.47
1:A:498:TYR:HB2	1:A:521:MET:CE	2.45	0.47
1:B:422:ASP:C	1:B:424:ALA:H	2.23	0.47
1:B:468:ALA:HA	1:B:471:GLN:CD	2.40	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:MET:HA	1:A:301:MET:HE2	1.95	0.47
1:A:416:ILE:CA	1:A:513:ARG:HE	2.20	0.47
1:B:301:MET:HA	1:B:301:MET:HE2	1.96	0.46
1:A:302:ARG:HD2	1:A:398:GLU:OE1	2.16	0.46
1:B:231:ARG:HD3	1:B:249:PRO:HB3	1.96	0.46
1:B:544:ALA:HB3	1:B:545:PRO:HD3	1.97	0.45
1:A:121:GLU:HG2	1:A:196:ARG:HG2	1.98	0.45
1:A:413:ASP:CG	1:A:414:SER:H	2.25	0.45
1:A:221:PRO:HG2	6:A:1174:HOH:O	2.16	0.45
1:A:284:ALA:O	1:A:350:ARG:NH1	2.50	0.45
1:B:177:GLU:OE2	1:B:179:LYS:HE3	2.17	0.44
1:A:409:ALA:CB	1:A:521:MET:HE3	2.47	0.44
1:A:602:MET:HG3	1:A:622:TYR:CD2	2.52	0.44
1:A:140:GLU:O	1:A:141:ARG:CB	2.63	0.44
1:A:389:LEU:H	1:A:389:LEU:HD12	1.83	0.44
1:A:168:ALA:C	1:A:170:SER:H	2.26	0.44
1:A:174:TYR:OH	1:A:193:LYS:NZ	2.51	0.44
1:B:63:VAL:HG13	1:B:122:VAL:HG13	1.99	0.44
1:A:362:ALA:HB3	4:A:803:EDO:H22	1.99	0.44
1:A:410:PRO:HB3	1:A:436:GLN:HG3	1.99	0.44
1:B:465:ARG:HH11	1:B:465:ARG:HG3	1.83	0.44
1:B:586:PRO:HG3	1:B:592:TYR:CE2	2.53	0.44
1:B:388:ASP:OD1	1:B:388:ASP:C	2.60	0.43
1:A:160:ALA:N	1:A:161:PRO:HD2	2.32	0.43
1:B:309:MET:O	1:B:313:MET:HG3	2.18	0.43
1:A:293:ILE:HG21	1:A:390:ARG:HG2	2.00	0.43
1:B:404:ILE:HG22	1:B:406:MET:HG2	2.00	0.43
1:B:117:ASN:HA	1:B:118:PRO:HD2	1.93	0.43
1:B:409:ALA:HB2	1:B:521:MET:HE3	2.01	0.42
1:B:44:GLN:HG3	1:B:660:ARG:HG2	2.01	0.42
1:B:699:ILE:O	1:B:703:PHE:HB3	2.20	0.42
1:B:660:ARG:HD3	6:B:903:HOH:O	2.20	0.42
1:B:302:ARG:NH1	1:B:402:MET:HE1	2.34	0.42
1:B:125:VAL:HG22	1:B:192:LYS:HE2	2.01	0.42
1:A:85:HIS:HA	1:A:88:TYR:HB2	2.01	0.41
1:B:536:ALA:O	1:B:539:VAL:HG12	2.21	0.41
1:A:285:ASP:OD2	1:A:380:SER:OG	2.34	0.41
1:B:459:LEU:HD23	1:B:459:LEU:HA	1.80	0.41
1:B:28:MET:HE3	1:B:51:TYR:HB2	2.03	0.41
1:A:353:LYS:HA	1:A:353:LYS:HD2	1.89	0.41
1:B:32:ASN:HB2	6:B:959:HOH:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:389:LEU:HD22	1:B:539:VAL:HB	2.03	0.41
1:A:326:LYS:NZ	6:A:918:HOH:O	2.54	0.41
1:B:83:ASN:HB2	1:B:86:CYS:SG	2.60	0.41
1:B:461:ARG:HD2	1:B:465:ARG:NH1	2.36	0.41
1:B:369:ILE:HG22	1:B:371:THR:HB	2.03	0.41
1:B:76:ASN:HB3	1:B:255:TRP:CH2	2.57	0.40

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	702/703 (100%)	673 (96%)	24 (3%)	5 (1%)	18	19
1	B	689/703 (98%)	655 (95%)	29 (4%)	5 (1%)	18	19
All	All	1391/1406 (99%)	1328 (96%)	53 (4%)	10 (1%)	18	19

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	416	ILE
1	A	589	VAL
1	B	450	GLU
1	B	452	ASP
1	B	589	VAL
1	A	140	GLU
1	A	147	GLY
1	A	419	HIS
1	B	169	ALA
1	B	423	PRO

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	568/583 (97%)	559 (98%)	9 (2%)	55	71
1	B	489/583 (84%)	482 (99%)	7 (1%)	59	75
All	All	1057/1166 (91%)	1041 (98%)	16 (2%)	57	73

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

Mol	Chain	Res	Type
1	A	22	ASP
1	A	63	VAL
1	A	140	GLU
1	A	141	ARG
1	A	162	GLU
1	A	166	LYS
1	A	391	MET
1	A	430	LEU
1	A	476	ARG
1	B	21	MET
1	B	63	VAL
1	B	69	CYS
1	B	77	GLN
1	B	302	ARG
1	B	348	SER
1	B	360	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	217	ASN
1	A	261	GLN
1	A	332	ASN
1	B	44	GLN
1	B	54	ASN

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 ⓘ

Of 6 ligands modelled in this entry, 1 is monoatomic - leaving 5 for Mogul analysis.

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	ARG	A	801	-	9,10,11	0.42	0	5,11,13	0.28	0
3	ASP	B	802	-	7,8,8	1.09	0	6,10,10	1.42	1 (16%)
2	ARG	B	801	-	9,10,11	0.74	0	5,11,13	0.34	0
4	EDO	A	803	-	3,3,3	0.43	0	2,2,2	0.38	0
3	ASP	A	802	-	7,8,8	1.16	0	6,10,10	1.35	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	ARG	A	801	-	-	0/8/9/11	-
3	ASP	B	802	-	-	5/8/8/8	-
2	ARG	B	801	-	-	0/8/9/11	-
4	EDO	A	803	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ASP	A	802	-	-	4/8/8/8	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	802	ASP	OXT-C-O	-2.45	118.52	124.08
3	B	802	ASP	OXT-C-O	-2.15	119.21	124.08

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	802	ASP	C-CA-CB-CG
3	B	802	ASP	C-CA-CB-CG
3	A	802	ASP	N-CA-CB-CG
4	A	803	EDO	O1-C1-C2-O2
3	A	802	ASP	OXT-C-CA-CB
3	B	802	ASP	N-CA-CB-CG
3	A	802	ASP	O-C-CA-CB
3	B	802	ASP	O-C-CA-CB
3	B	802	ASP	OXT-C-CA-CB
3	B	802	ASP	OXT-C-CA-N

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	802	ASP	1	0
4	A	803	EDO	2	0
3	A	802	ASP	1	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

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	703/703 (100%)	-0.12	19 (2%) 56 53	32, 53, 81, 105	1 (0%)
1	B	695/703 (98%)	0.48	75 (10%) 11 8	31, 72, 144, 192	0
All	All	1398/1406 (99%)	0.18	94 (6%) 24 21	31, 58, 128, 192	1 (0%)

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	431	GLN	5.0
1	A	169	ALA	4.5
1	B	169	ALA	4.4
1	A	415	ASP	4.1
1	B	512	ASP	4.1
1	B	418	ASP	4.1
1	A	417	ALA	3.8
1	B	415	ASP	3.8
1	B	21	MET	3.7
1	B	507	PHE	3.7
1	A	148	LEU	3.6
1	B	718	HIS	3.6
1	B	449	VAL	3.6
1	A	416	ILE	3.6
1	A	147	GLY	3.6
1	B	430	LEU	3.6
1	A	719	HIS	3.6
1	B	530	TYR	3.5
1	B	510	ASN	3.5
1	B	473	ILE	3.4
1	B	414	SER	3.3
1	B	141	ARG	3.3
1	B	342	TRP	3.3
1	A	418	ASP	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	416	ILE	3.2
1	B	452	ASP	3.2
1	A	414	SER	3.1
1	B	516	LEU	3.1
1	B	159	LEU	3.1
1	B	479	ILE	3.0
1	B	484	SER	3.0
1	B	420	TRP	3.0
1	B	143	LYS	3.0
1	B	515	ARG	3.0
1	B	481	GLU	3.0
1	B	167	LYS	2.9
1	B	438	GLU	2.9
1	A	718	HIS	2.9
1	B	168	ALA	2.9
1	B	491	MET	2.9
1	B	417	ALA	2.8
1	B	489	VAL	2.8
1	B	403	GLY	2.8
1	B	562	GLY	2.8
1	B	714	GLN	2.8
1	B	399	GLY	2.8
1	B	441	PHE	2.8
1	B	483	GLY	2.7
1	B	423	PRO	2.7
1	B	497	ILE	2.7
1	B	400	ILE	2.7
1	B	459	LEU	2.7
1	B	505	GLU	2.7
1	A	21	MET	2.7
1	B	138	GLU	2.6
1	B	166	LYS	2.6
1	B	513	ARG	2.6
1	B	165	GLY	2.6
1	B	469	GLU	2.6
1	B	402	MET	2.6
1	A	166	LYS	2.5
1	A	170	SER	2.4
1	B	511	PRO	2.4
1	B	297	ILE	2.4
1	B	410	PRO	2.4
1	B	421	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	534	LYS	2.4
1	B	389	LEU	2.4
1	B	458	LEU	2.4
1	A	251	ARG	2.4
1	B	514	ASP	2.3
1	B	493	PHE	2.3
1	A	342	TRP	2.3
1	A	168	ALA	2.3
1	B	428	ALA	2.3
1	B	475	ILE	2.3
1	B	140	GLU	2.3
1	B	501	ARG	2.3
1	A	144	LYS	2.3
1	B	404	ILE	2.2
1	B	526	ALA	2.2
1	B	504	PHE	2.2
1	A	165	GLY	2.2
1	B	467	PRO	2.1
1	B	524	PHE	2.1
1	B	476	ARG	2.1
1	B	496	SER	2.1
1	B	523	ARG	2.1
1	B	172	PRO	2.1
1	A	722	HIS	2.0
1	B	436	GLN	2.0
1	B	429	GLY	2.0
1	B	508	MET	2.0
1	B	392	ARG	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(Å ²)	Q<0.9
4	EDO	A	803	4/4	0.87	0.21	61,62,64,64	0
3	ASP	B	802	9/9	0.94	0.09	40,45,62,68	0
5	NA	A	804	1/1	0.94	0.09	60,60,60,60	0
2	ARG	A	801	11/12	0.96	0.10	41,48,60,68	0
2	ARG	B	801	11/12	0.96	0.07	31,39,46,53	0
3	ASP	A	802	9/9	0.96	0.07	41,46,53,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.