



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:37 PM GMT

PDB ID : 1BQD
Title : ASPARTATE AMINOTRANSFERASE P138A/P195A DOUBLE MUTANT
Authors : Malashkevich, V.N.; Jansonius, J.N.
Deposited on : 1998-08-14
Resolution : 2.10 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
<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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

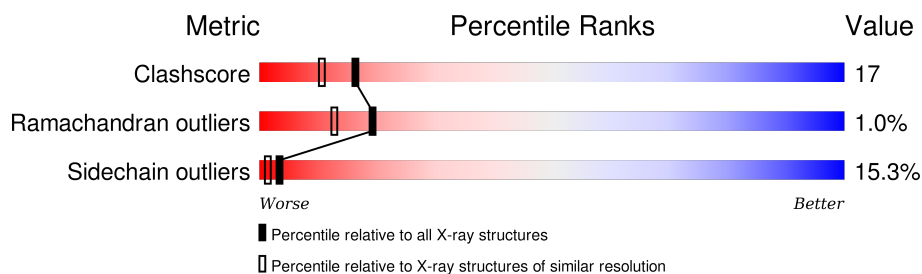
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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))
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	396	
1	B	396	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6848 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 ASPARTATE AMINOTRANSFERASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	396	Total	C	N	O	P	S	0	0	0
			3080	1940	537	589	1	13			
1	B	396	Total	C	N	O	P	S	0	0	0
			3080	1940	537	589	1	13			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	138	ALA	PRO	ENGINEERED	UNP P00509
A	195	ALA	PRO	ENGINEERED	UNP P00509
A	258	LLP	LYS	MODIFIED RESIDUE	UNP P00509
B	138	ALA	PRO	ENGINEERED	UNP P00509
B	195	ALA	PRO	ENGINEERED	UNP P00509
B	258	LLP	LYS	MODIFIED RESIDUE	UNP P00509

- Molecule 2 is water.

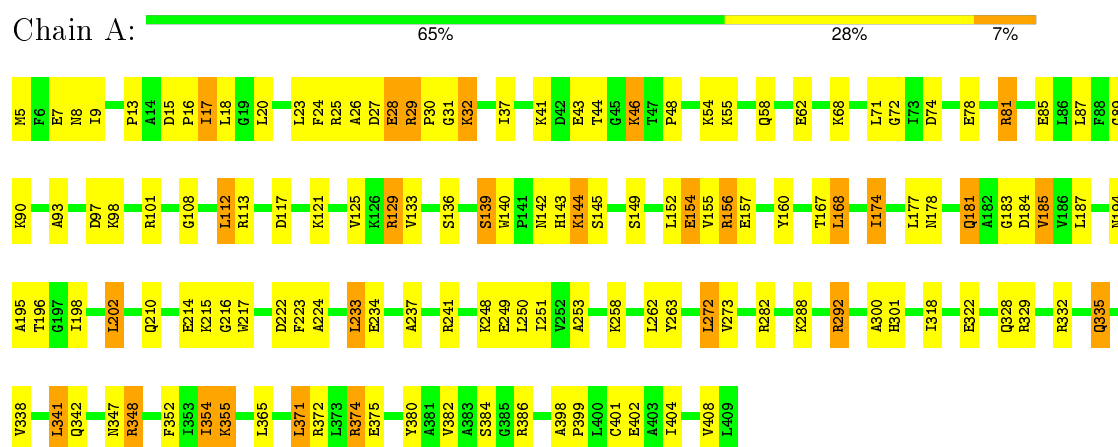
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	371	Total	O	0	0
			371	371		
2	B	317	Total	O	0	0
			317	317		

3 Residue-property plots

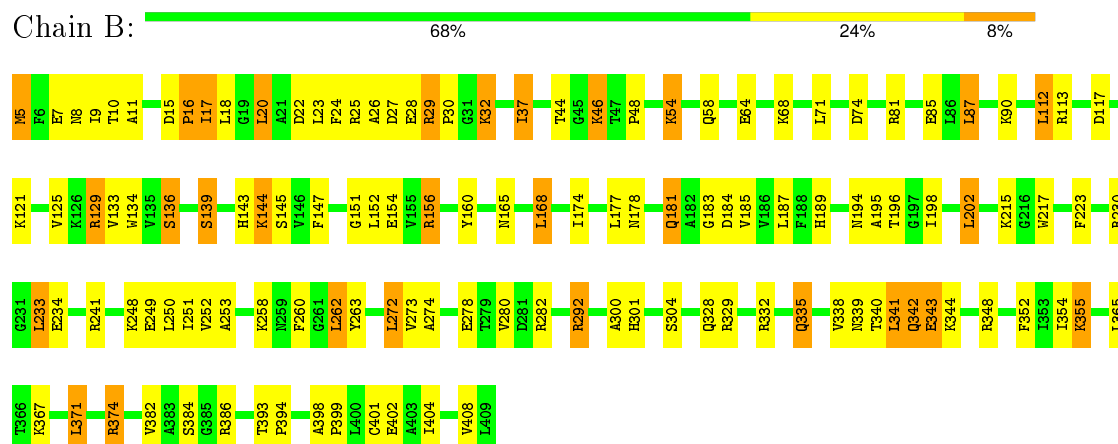
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

• Molecule 1: ASPARTATE AMINOTRANSFERASE



• Molecule 1: ASPARTATE AMINOTRANSFERASE



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	88.49 Å 79.88 Å 89.68 Å 90.00° 119.58° 90.00°	Depositor
Resolution (Å)	20.00 – 2.10	Depositor
% Data completeness (in resolution range)	92.1 (20.00-2.10)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT V. 5-D	Depositor
R, R_{free}	0.204 , 0.240	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6848	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	0.56	0/3114	0.59	0/4216
1	B	0.58	0/3114	0.60	0/4216
All	All	0.57	0/6228	0.60	0/8432

There are no bond length outliers.

There are no bond angle outliers.

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	3080	0	3019	115	0
1	B	3080	0	3019	109	0
2	A	371	0	0	20	0
2	B	317	0	0	14	0
All	All	6848	0	6038	204	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 17.

All (204) 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:B:352:PHE:HA	1:B:355:LYS:HD3	1.35	1.05
1:A:352:PHE:HA	1:A:355:LYS:HD3	1.36	1.04
1:A:17:ILE:HG12	1:A:18:LEU:HD12	1.54	0.88
1:B:17:ILE:HG13	1:B:18:LEU:N	1.96	0.81
1:A:292:ARG:HG3	1:A:292:ARG:HH11	1.47	0.80
1:B:17:ILE:HG13	1:B:18:LEU:HD12	1.65	0.77
1:A:117:ASP:O	1:A:121:LYS:HG2	1.85	0.76
1:A:78:GLU:HB3	2:A:625:HOH:O	1.85	0.76
1:A:183:GLY:O	1:B:5:MET:HE1	1.86	0.75
1:A:282:ARG:HD2	1:B:9:ILE:O	1.87	0.75
1:B:352:PHE:CA	1:B:355:LYS:HD3	2.17	0.74
1:A:90:LYS:HG3	2:A:752:HOH:O	1.87	0.74
1:B:117:ASP:O	1:B:121:LYS:HG2	1.88	0.73
1:A:5:MET:HE1	1:B:183:GLY:O	1.88	0.73
1:B:328:GLN:O	1:B:332:ARG:HD3	1.89	0.73
1:B:24:PHE:HE1	1:B:32:LYS:HB2	1.53	0.72
1:A:258:LLP:H4'1	2:A:453:HOH:O	1.91	0.71
1:B:352:PHE:HA	1:B:355:LYS:CD	2.17	0.70
1:A:43:GLU:HB2	2:A:531:HOH:O	1.93	0.68
1:B:129:ARG:HH21	1:B:181:GLN:HG3	1.60	0.67
1:A:157:GLU:HB3	2:A:646:HOH:O	1.95	0.67
1:B:398:ALA:HB3	1:B:399:PRO:HD3	1.76	0.66
1:A:142:ASN:HB2	2:A:544:HOH:O	1.95	0.66
1:A:44:THR:OG1	1:A:46:LYS:HG2	1.95	0.66
1:B:189:HIS:CD2	1:B:194:ASN:H	2.14	0.66
1:A:258:LLP:H4'2	1:A:258:LLP:OP4	1.96	0.66
1:B:24:PHE:CE1	1:B:32:LYS:HB2	2.30	0.66
1:B:382:VAL:HG12	1:B:384:SER:H	1.61	0.66
1:A:5:MET:N	1:A:7:GLU:OE1	2.30	0.65
1:A:5:MET:CE	1:B:183:GLY:HA2	2.26	0.64
1:A:216:GLY:HA2	2:A:706:HOH:O	1.96	0.64
1:A:129:ARG:HG3	1:A:154:GLU:HB3	1.79	0.64
1:B:27:ASP:OD1	1:B:29:ARG:HB2	1.96	0.64
1:A:81:ARG:NH1	2:A:567:HOH:O	2.30	0.64
1:B:129:ARG:HG3	1:B:154:GLU:HB3	1.80	0.64
1:A:355:LYS:NZ	1:A:355:LYS:HB2	2.11	0.64
1:A:17:ILE:CG1	1:A:18:LEU:HD12	2.28	0.63
1:B:258:LLP:H4'2	1:B:258:LLP:OP4	1.99	0.63
1:A:292:ARG:HG3	1:A:292:ARG:NH1	2.13	0.62
1:B:136:SER:O	1:B:139:SER:HB2	2.00	0.62
1:A:139:SER:OG	1:A:140:TRP:N	2.31	0.62
1:B:260:PHE:HB3	1:B:262:LEU:HD22	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:189:HIS:HD2	1:B:194:ASN:H	1.46	0.62
1:A:160:TYR:O	1:A:168:LEU:HD23	2.00	0.61
1:B:44:THR:OG1	1:B:46:LYS:HG2	2.00	0.61
1:A:202:LEU:O	1:A:202:LEU:HD22	2.01	0.61
1:B:17:ILE:CG1	1:B:18:LEU:HD12	2.30	0.60
1:A:374:ARG:HD3	1:A:380:TYR:CE2	2.37	0.60
1:A:183:GLY:HA2	1:B:5:MET:HE2	1.82	0.60
1:A:382:VAL:HG12	1:A:384:SER:H	1.66	0.60
1:B:374:ARG:NH1	2:B:675:HOH:O	2.34	0.60
1:B:292:ARG:HH11	1:B:292:ARG:HG3	1.65	0.60
1:B:46:LYS:O	1:B:48:PRO:HD3	2.03	0.59
1:B:160:TYR:O	1:B:168:LEU:HD23	2.02	0.59
1:B:16:PRO:O	1:B:20:LEU:N	2.29	0.59
1:A:129:ARG:HH21	1:A:181:GLN:HG3	1.67	0.59
1:B:54:LYS:HE3	2:B:581:HOH:O	2.02	0.59
1:A:318:ILE:O	1:A:322:GLU:HG3	2.03	0.59
1:A:288:LYS:NZ	2:A:639:HOH:O	2.35	0.58
1:A:398:ALA:HB3	1:A:399:PRO:HD3	1.86	0.58
1:A:5:MET:HE1	1:B:183:GLY:C	2.24	0.57
1:B:18:LEU:HD22	2:B:654:HOH:O	2.04	0.57
1:A:338:VAL:HG13	1:A:342:GLN:HE21	1.69	0.57
1:A:183:GLY:HA2	1:B:5:MET:CE	2.35	0.57
1:A:9:ILE:O	1:B:282:ARG:HD2	2.05	0.57
1:B:18:LEU:HD13	2:B:654:HOH:O	2.04	0.57
1:B:71:LEU:HD11	1:B:300:ALA:HB2	1.87	0.57
1:B:341:LEU:HD13	1:B:401:CYS:SG	2.44	0.57
1:A:68:LYS:HD2	1:B:262:LEU:HD12	1.86	0.56
1:B:202:LEU:HD22	1:B:202:LEU:O	2.06	0.56
1:B:37:ILE:HB	2:B:521:HOH:O	2.06	0.56
1:A:18:LEU:N	1:A:18:LEU:HD12	2.21	0.56
1:B:113:ARG:NH2	2:B:595:HOH:O	2.39	0.55
1:B:22:ASP:HA	1:B:25:ARG:NH2	2.20	0.55
1:A:347:ASN:HB3	2:A:740:HOH:O	2.07	0.55
1:A:44:THR:OG1	1:A:46:LYS:HE2	2.05	0.55
1:B:398:ALA:HB3	1:B:399:PRO:CD	2.37	0.54
1:A:5:MET:HE2	1:B:183:GLY:HA2	1.88	0.54
1:B:194:ASN:OD1	1:B:195:ALA:HB2	2.08	0.53
1:B:25:ARG:NH1	1:B:25:ARG:HB2	2.23	0.53
1:A:341:LEU:HD13	1:A:401:CYS:SG	2.48	0.53
1:B:85:GLU:CD	1:B:90:LYS:HD2	2.28	0.53
1:A:404:ILE:O	1:A:408:VAL:HG22	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:GLY:HA3	2:A:602:HOH:O	2.08	0.53
1:A:112:LEU:HD13	1:A:253:ALA:CB	2.38	0.53
1:A:24:PHE:HE1	1:A:32:LYS:HB2	1.74	0.53
1:B:339:ASN:O	1:B:343:GLU:HG3	2.09	0.53
1:A:27:ASP:OD1	1:A:29:ARG:HB2	2.09	0.53
1:B:292:ARG:NH1	1:B:292:ARG:HG3	2.24	0.52
1:B:74:ASP:OD1	1:B:74:ASP:N	2.35	0.52
1:A:46:LYS:O	1:A:48:PRO:HD3	2.10	0.52
1:B:29:ARG:HG2	1:B:29:ARG:HH11	1.74	0.52
1:A:335:GLN:HA	1:A:354:ILE:CD1	2.40	0.52
1:A:78:GLU:CD	1:A:78:GLU:H	2.13	0.52
1:B:249:GLU:HA	1:B:273:VAL:O	2.10	0.52
1:A:248:LYS:NZ	2:A:470:HOH:O	2.41	0.51
1:B:134:TRP:CE3	1:B:156:ARG:HB2	2.46	0.51
1:B:25:ARG:HH11	1:B:25:ARG:HB2	1.74	0.51
1:B:251:ILE:HG12	1:B:272:LEU:HD23	1.92	0.51
1:A:15:ASP:O	1:A:17:ILE:N	2.44	0.51
1:A:328:GLN:O	1:A:332:ARG:HD3	2.11	0.51
1:A:194:ASN:OD1	1:A:195:ALA:HB2	2.10	0.51
1:A:282:ARG:HG3	1:B:11:ALA:HB2	1.92	0.50
1:B:113:ARG:O	1:B:113:ARG:HD2	2.10	0.50
1:B:348:ARG:HB3	2:B:604:HOH:O	2.11	0.50
1:A:68:LYS:O	1:B:263:TYR:HB2	2.11	0.50
1:B:29:ARG:HG2	1:B:29:ARG:NH1	2.27	0.50
1:A:93:ALA:O	1:A:97:ASP:HB2	2.10	0.50
1:A:17:ILE:HG13	1:A:18:LEU:N	2.25	0.50
1:A:5:MET:HE1	1:B:183:GLY:HA2	1.92	0.49
1:A:348:ARG:HG2	2:A:740:HOH:O	2.12	0.49
1:A:24:PHE:CE1	1:A:32:LYS:HB2	2.47	0.49
1:A:5:MET:HE1	1:B:183:GLY:CA	2.42	0.49
1:A:74:ASP:OD1	1:A:74:ASP:N	2.37	0.49
1:B:248:LYS:HG3	2:B:614:HOH:O	2.13	0.49
1:A:113:ARG:NH2	2:B:595:HOH:O	2.45	0.49
1:A:18:LEU:HD12	1:A:18:LEU:H	1.77	0.49
1:A:129:ARG:NH1	2:A:649:HOH:O	2.46	0.48
1:B:15:ASP:HB3	2:B:654:HOH:O	2.12	0.48
1:A:196:THR:HB	1:A:198:ILE:HD12	1.93	0.48
1:A:5:MET:HB2	1:B:249:GLU:OE1	2.14	0.48
1:A:195:ALA:HB1	1:A:386:ARG:HD2	1.95	0.48
1:A:250:LEU:HD23	1:A:251:ILE:N	2.28	0.48
1:B:335:GLN:HA	1:B:354:ILE:CD1	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:194:ASN:HA	1:A:195:ALA:HA	1.56	0.47
1:A:237:ALA:O	1:A:241:ARG:HD3	2.14	0.47
1:A:348:ARG:HD3	2:A:740:HOH:O	2.14	0.47
1:B:194:ASN:HA	1:B:195:ALA:HA	1.56	0.47
1:B:151:GLY:HA2	2:B:578:HOH:O	2.15	0.47
1:B:250:LEU:HD21	1:B:252:VAL:CG2	2.44	0.47
1:B:29:ARG:NH2	1:B:374:ARG:O	2.48	0.47
1:A:143:HIS:NE2	1:A:222:ASP:OD2	2.48	0.47
1:A:233:LEU:HD12	1:A:233:LEU:HA	1.76	0.47
1:A:144:LYS:HB2	1:A:144:LYS:HE3	1.40	0.46
1:B:338:VAL:HG13	1:B:342:GLN:HE21	1.79	0.46
1:B:87:LEU:O	1:B:241:ARG:HD2	2.15	0.46
1:A:183:GLY:C	1:B:5:MET:HE1	2.36	0.46
1:B:195:ALA:HB1	1:B:386:ARG:HD2	1.98	0.46
1:B:234:GLU:OE1	1:B:241:ARG:NH2	2.49	0.45
1:B:129:ARG:NH2	1:B:181:GLN:HG3	2.27	0.45
1:A:224:ALA:HB1	1:A:258:LLP:HZ1	1.81	0.45
1:B:398:ALA:O	1:B:402:GLU:HG3	2.15	0.45
1:A:262:LEU:HD23	1:B:68:LYS:HD3	1.99	0.45
1:A:17:ILE:CG1	1:A:18:LEU:N	2.80	0.45
1:B:196:THR:HB	1:B:198:ILE:HD12	1.98	0.45
1:A:181:GLN:O	1:A:184:ASP:HB2	2.17	0.45
1:A:234:GLU:OE1	1:A:241:ARG:NH2	2.50	0.45
1:A:125:VAL:HG11	1:A:185:VAL:CG1	2.47	0.45
1:A:55:LYS:HE2	2:A:681:HOH:O	2.18	0.45
1:A:177:LEU:O	1:A:217:TRP:HZ2	2.00	0.44
1:B:233:LEU:HD12	1:B:233:LEU:HA	1.76	0.44
1:B:165:ASN:ND2	2:B:726:HOH:O	2.30	0.44
1:A:54:LYS:NZ	1:B:64:GLU:O	2.51	0.44
1:A:85:GLU:CD	1:A:90:LYS:HD2	2.38	0.44
1:A:28:GLU:H	1:A:28:GLU:HG2	1.60	0.44
1:B:24:PHE:CE1	1:B:32:LYS:HD2	2.53	0.44
1:B:250:LEU:HD23	1:B:250:LEU:C	2.38	0.44
1:B:25:ARG:NH1	1:B:25:ARG:CB	2.80	0.44
1:A:263:TYR:HB2	1:B:68:LYS:O	2.18	0.44
1:A:144:LYS:HE2	1:A:155:VAL:HG11	2.00	0.44
1:A:112:LEU:HD12	1:A:112:LEU:HA	1.86	0.43
1:A:174:ILE:N	1:A:174:ILE:HD13	2.34	0.43
1:A:129:ARG:HD2	1:A:154:GLU:OE1	2.19	0.43
1:B:5:MET:N	1:B:7:GLU:OE1	2.52	0.43
1:A:233:LEU:HD23	2:A:474:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:GLY:HA3	1:A:258:LLP:H5'1	2.01	0.43
1:B:371:LEU:O	1:B:371:LEU:HD22	2.19	0.43
1:B:20:LEU:HA	1:B:20:LEU:HD23	1.74	0.43
1:A:249:GLU:HA	1:A:273:VAL:O	2.19	0.43
1:B:18:LEU:HB2	2:B:654:HOH:O	2.19	0.42
1:A:156:ARG:HD3	2:A:649:HOH:O	2.19	0.42
1:B:404:ILE:O	1:B:408:VAL:HG22	2.19	0.42
1:A:174:ILE:HD12	1:A:174:ILE:HA	1.81	0.42
1:B:143:HIS:O	1:B:147:PHE:HD2	2.02	0.42
1:B:393:THR:HB	1:B:394:PRO:HD2	2.01	0.42
1:A:85:GLU:O	1:A:89:GLY:N	2.52	0.42
1:B:181:GLN:O	1:B:184:ASP:HB2	2.19	0.42
1:B:112:LEU:HD13	1:B:253:ALA:CB	2.50	0.42
1:B:230:ARG:NH1	2:B:450:HOH:O	2.41	0.42
1:A:58:GLN:NE2	1:A:62:GLU:OE2	2.53	0.42
1:B:187:LEU:C	1:B:187:LEU:HD23	2.40	0.42
1:A:72:GLY:HA3	2:A:764:HOH:O	2.20	0.42
1:B:177:LEU:HB3	1:B:217:TRP:CH2	2.55	0.41
1:A:249:GLU:OE1	1:B:5:MET:N	2.53	0.41
1:A:187:LEU:C	1:A:187:LEU:HD23	2.41	0.41
1:A:29:ARG:O	1:A:32:LYS:NZ	2.39	0.41
1:A:71:LEU:HD11	1:A:300:ALA:HB2	2.02	0.41
1:B:340:THR:O	1:B:344:LYS:HB2	2.19	0.41
1:A:58:GLN:HE21	1:A:62:GLU:HG3	1.84	0.41
1:A:371:LEU:HA	1:A:371:LEU:HD23	1.87	0.41
1:B:341:LEU:HA	1:B:341:LEU:HD12	1.82	0.41
1:B:125:VAL:HG11	1:B:185:VAL:CG1	2.50	0.41
1:A:18:LEU:CD1	1:A:18:LEU:H	2.34	0.41
1:B:144:LYS:HE3	1:B:144:LYS:HB2	1.21	0.41
1:B:250:LEU:HD21	1:B:252:VAL:HG23	2.02	0.41
1:A:46:LYS:C	1:A:48:PRO:HD3	2.41	0.41
1:B:274:ALA:HB3	1:B:280:VAL:HB	2.01	0.41
1:A:292:ARG:NH1	1:A:292:ARG:CG	2.81	0.40
1:A:25:ARG:C	1:A:27:ASP:H	2.24	0.40
1:A:13:PRO:HA	2:A:778:HOH:O	2.21	0.40
1:A:167:THR:OG1	1:A:168:LEU:N	2.54	0.40
1:A:251:ILE:HG12	1:A:272:LEU:HD23	2.04	0.40
1:B:262:LEU:HA	1:B:262:LEU:HD12	1.90	0.40
1:B:25:ARG:CZ	1:B:25:ARG:CB	3.00	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	393/396 (99%)	374 (95%)	15 (4%)	4 (1%)	19	13
1	B	393/396 (99%)	374 (95%)	15 (4%)	4 (1%)	19	13
All	All	786/792 (99%)	748 (95%)	30 (4%)	8 (1%)	19	13

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	16	PRO
1	A	26	ALA
1	A	30	PRO
1	A	301	HIS
1	B	26	ALA
1	B	30	PRO
1	B	301	HIS
1	B	16	PRO

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	317/317 (100%)	267 (84%)	50 (16%)	3	1
1	B	317/317 (100%)	270 (85%)	47 (15%)	4	2
All	All	634/634 (100%)	537 (85%)	97 (15%)	3	1

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	17	ILE
1	A	20	LEU
1	A	23	LEU
1	A	28	GLU
1	A	29	ARG
1	A	32	LYS
1	A	37	ILE
1	A	41	LYS
1	A	46	LYS
1	A	81	ARG
1	A	87	LEU
1	A	98	LYS
1	A	101	ARG
1	A	112	LEU
1	A	129	ARG
1	A	133	VAL
1	A	136	SER
1	A	139	SER
1	A	144	LYS
1	A	145	SER
1	A	149	SER
1	A	152	LEU
1	A	154	GLU
1	A	156	ARG
1	A	168	LEU
1	A	174	ILE
1	A	178	ASN
1	A	181	GLN
1	A	185	VAL
1	A	202	LEU
1	A	210	GLN
1	A	214	GLU
1	A	215	LYS
1	A	223	PHE
1	A	233	LEU
1	A	272	LEU
1	A	292	ARG
1	A	329	ARG
1	A	335	GLN
1	A	341	LEU
1	A	348	ARG
1	A	354	ILE

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Mol	Chain	Res	Type
1	A	355	LYS
1	A	365	LEU
1	A	371	LEU
1	A	372	ARG
1	A	374	ARG
1	A	375	GLU
1	A	402	GLU
1	B	5	MET
1	B	8	ASN
1	B	10	THR
1	B	17	ILE
1	B	20	LEU
1	B	23	LEU
1	B	28	GLU
1	B	29	ARG
1	B	32	LYS
1	B	37	ILE
1	B	46	LYS
1	B	54	LYS
1	B	58	GLN
1	B	81	ARG
1	B	87	LEU
1	B	112	LEU
1	B	129	ARG
1	B	133	VAL
1	B	136	SER
1	B	139	SER
1	B	144	LYS
1	B	145	SER
1	B	152	LEU
1	B	156	ARG
1	B	168	LEU
1	B	174	ILE
1	B	178	ASN
1	B	181	GLN
1	B	202	LEU
1	B	215	LYS
1	B	223	PHE
1	B	233	LEU
1	B	262	LEU
1	B	272	LEU
1	B	278	GLU

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Mol	Chain	Res	Type
1	B	292	ARG
1	B	304	SER
1	B	329	ARG
1	B	335	GLN
1	B	341	LEU
1	B	342	GLN
1	B	343	GLU
1	B	355	LYS
1	B	365	LEU
1	B	367	LYS
1	B	371	LEU
1	B	374	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	58	GLN
1	A	148	ASN
1	A	206	GLN
1	B	148	ASN
1	B	189	HIS
1	B	206	GLN
1	B	210	GLN
1	B	328	GLN
1	B	339	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	LLP	A	258	1	23,24,25	1.13	1 (4%)	28,32,34	1.45	5 (17%)
1	LLP	B	258	1	23,24,25	0.97	1 (4%)	28,32,34	1.52	8 (28%)

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
1	LLP	A	258	1	-	0/15/17/19	0/1/1/1
1	LLP	B	258	1	-	0/15/17/19	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	258	LLP	P-OP3	-2.18	1.46	1.54
1	A	258	LLP	CE-NZ	2.09	1.51	1.46

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	258	LLP	C5-C6-N1	-2.74	119.11	123.86
1	B	258	LLP	C3-C4-C5	-2.69	116.09	118.11
1	A	258	LLP	C3-C2-N1	-2.63	116.98	120.61
1	B	258	LLP	C5-C6-N1	-2.48	119.56	123.86
1	B	258	LLP	C3-C2-N1	-2.29	117.44	120.61
1	B	258	LLP	CD-CE-NZ	-2.21	107.36	110.98
1	B	258	LLP	C2'-C2-C3	2.06	123.52	121.04
1	B	258	LLP	C6-N1-C2	2.35	124.06	119.28
1	B	258	LLP	CE-NZ-C4'	2.42	125.96	118.97
1	B	258	LLP	OP3-P-OP1	2.44	118.45	110.58
1	A	258	LLP	OP3-P-OP1	2.48	118.58	110.58
1	A	258	LLP	C6-N1-C2	2.49	124.36	119.28
1	A	258	LLP	C2'-C2-C3	2.76	124.36	121.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	258	LLP	4	0
1	B	258	LLP	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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](#)

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.