



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:10 PM GMT

PDB ID : 4BQ0
Title : Pseudomonas aeruginosa beta-alanine:pyruvate aminotransferase holoenzyme
without divalent cations on dimer-dimer interface
Authors : Isupov, M.N.; Lebedev, A.A.; Westlake, A.; Sayer, C.; Littlechild, J.A.
Deposited on : 2013-05-29
Resolution : 1.77 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) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

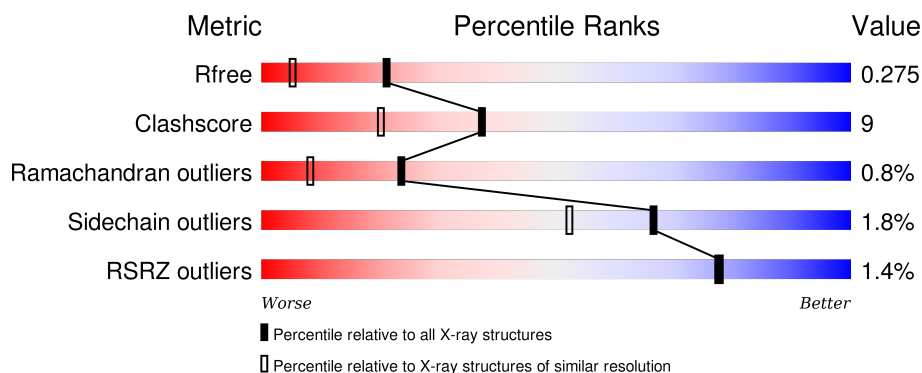
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.77 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}	91344	6655 (1.80-1.76)
Clashscore	102246	7658 (1.80-1.76)
Ramachandran outliers	100387	7570 (1.80-1.76)
Sidechain outliers	100360	7569 (1.80-1.76)
RSRZ outliers	91569	6671 (1.80-1.76)

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.

Mol	Chain	Length	Quality of chain
1	A	448	<div> <div>87%</div> <div>11% .</div> </div>
1	B	448	<div> <div>83%</div> <div>14% .</div> </div>
1	C	448	<div> <div>3%</div> <div>75%</div> <div>20% . .</div> </div>
1	D	448	<div> <div>2%</div> <div>86%</div> <div>10% . .</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15871 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 BETA-ALANINE--PYRUVATE TRANSAMINASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	439	Total	C	N	O	S	0	34	0
			3544	2260	619	646	19			
1	B	436	Total	C	N	O	S	0	37	0
			3536	2252	617	650	17			
1	C	435	Total	C	N	O	S	0	44	0
			3566	2279	620	647	20			
1	D	435	Total	C	N	O	S	0	30	0
			3483	2215	609	641	18			

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Cl	0	0
			1	1		
2	A	1	Total	Cl	0	0
			1	1		
2	D	1	Total	Cl	0	0
			1	1		
2	C	1	Total	Cl	0	0
			1	1		

- Molecule 3 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: C₈H₁₀NO₆P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			16	8	1	6	1		
3	B	1	Total	C	N	O	P	0	0
			16	8	1	6	1		
3	C	1	Total	C	N	O	P	0	0
			16	8	1	6	1		
3	D	1	Total	C	N	O	P	0	0
			16	8	1	6	1		

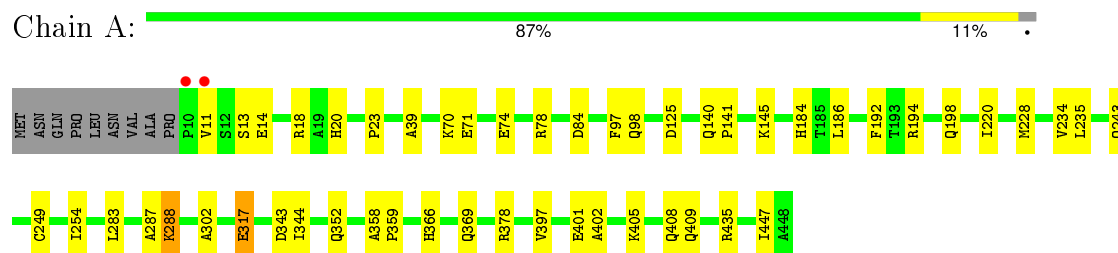
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	557	Total	O	0	0
			557	557		
4	B	439	Total	O	0	0
			439	439		
4	C	378	Total	O	0	0
			378	378		
4	D	300	Total	O	0	0
			300	300		

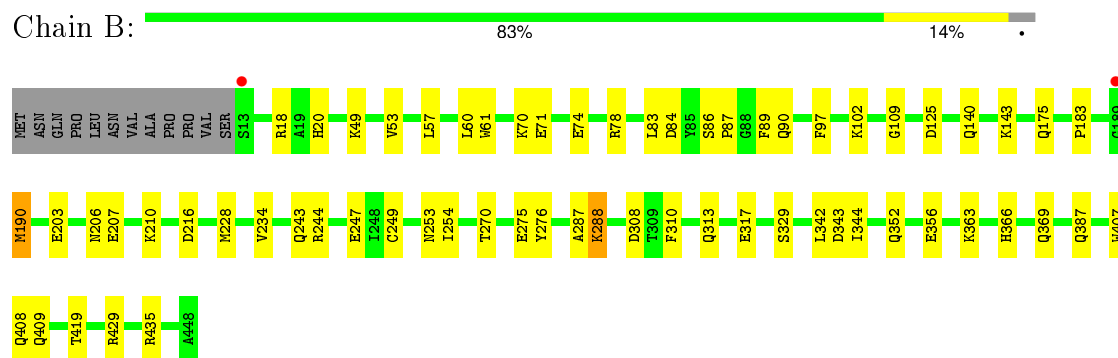
3 Residue-property plots [i](#)

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.

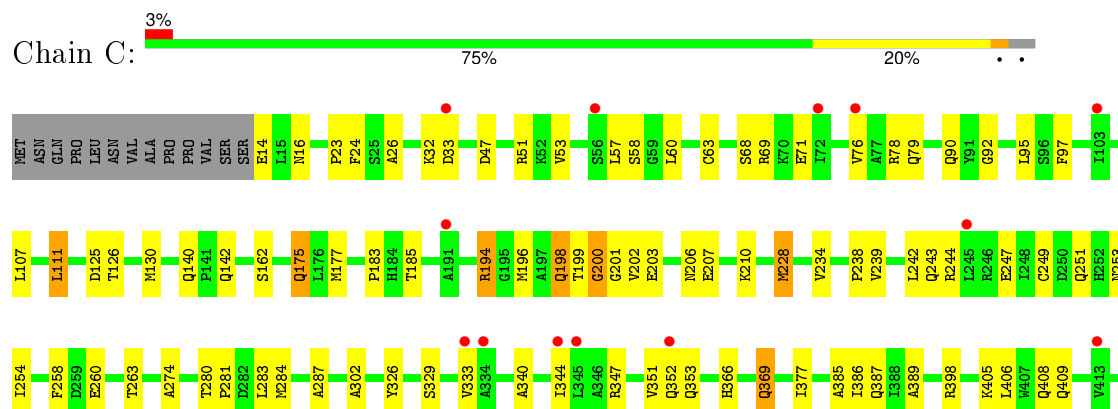
• Molecule 1: BETA-ALANINE--PYRUVATE TRANSAMINASE



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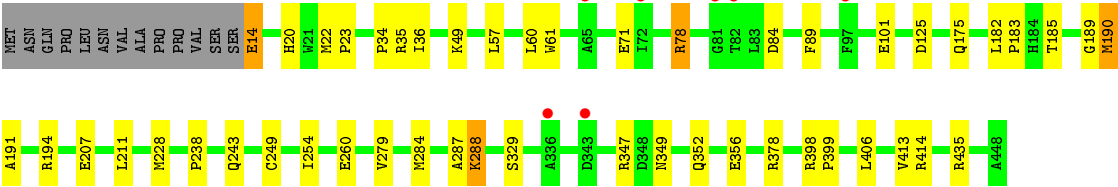
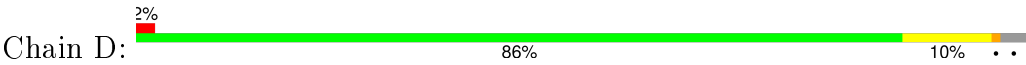


• Molecule 1: BETA-ALANINE--PYRUVATE TRANSAMINASE





● Molecule 1: BETA-ALANINE--PYRUVATE TRANSAMINASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	111.95Å 192.18Å 76.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	63.37 – 1.77 63.29 – 1.77	Depositor EDS
% Data completeness (in resolution range)	99.0 (63.37-1.77) 99.0 (63.29-1.77)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 1.77Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.216 , 0.276 0.215 , 0.275	Depositor DCC
R_{free} test set	7987 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	16.4	Xtriage
Anisotropy	1.281	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 54.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Outliers	4 of 159283 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	15871	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 70.35 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 3.1679e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: CL, PLP

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	0.57	0/3717	0.71	1/5026 (0.0%)
1	B	0.56	0/3717	0.71	1/5024 (0.0%)
1	C	0.50	0/3762	0.67	0/5086
1	D	0.49	0/3643	0.68	1/4927 (0.0%)
All	All	0.53	0/14839	0.69	3/20063 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	84	ASP	CB-CG-OD1	6.29	123.96	118.30
1	A	84	ASP	CB-CG-OD1	5.75	123.47	118.30
1	D	84	ASP	CB-CG-OD1	5.10	122.89	118.30

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	3544	0	3586	64	0
1	B	3536	0	3572	62	0
1	C	3566	0	3638	103	0
1	D	3483	0	3496	51	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	16	0	7	2	0
3	B	16	0	7	1	0
3	C	16	0	7	0	0
3	D	16	0	7	0	0
4	A	557	0	0	24	0
4	B	439	0	0	28	0
4	C	378	0	0	33	0
4	D	300	0	0	17	0
All	All	15871	0	14320	265	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 9.

All (265) 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:194[A]:ARG:CZ	4:C:2220:HOH:O	1.74	1.31
1:C:198[B]:GLN:NE2	4:C:2228:HOH:O	1.63	1.30
1:D:435[B]:ARG:HD3	4:D:2297:HOH:O	1.27	1.30
1:B:409[B]:GLN:NE2	4:B:2410:HOH:O	1.58	1.27
1:A:243[A]:GLN:OE1	4:A:2393:HOH:O	1.55	1.24
1:B:210[B]:LYS:NZ	4:B:2266:HOH:O	1.69	1.22
4:B:2269:HOH:O	1:C:210[B]:LYS:NZ	1.67	1.22
1:B:409[B]:GLN:OE1	4:B:2408:HOH:O	1.57	1.21
1:A:243[A]:GLN:HG2	4:A:2392:HOH:O	1.35	1.19
1:A:409[B]:GLN:OE1	4:A:2527:HOH:O	1.56	1.18
1:B:102[A]:LYS:HE3	4:B:2122:HOH:O	1.40	1.17
1:C:32[B]:LYS:HA	1:C:32[B]:LYS:HE2	1.29	1.15
1:C:185[B]:THR:HG21	1:C:238:PRO:HD3	1.29	1.11
1:D:435[A]:ARG:NH2	4:D:2295:HOH:O	1.82	1.11
1:D:435[A]:ARG:NE	4:D:2295:HOH:O	1.83	1.11
1:C:194[A]:ARG:CG	1:C:194[A]:ARG:HH11	1.62	1.10
1:A:369[B]:GLN:NE2	4:A:2338:HOH:O	1.65	1.10
1:B:409[A]:GLN:OE1	4:B:2407:HOH:O	1.68	1.09
1:C:194[A]:ARG:NH2	4:C:2220:HOH:O	1.78	1.09
4:B:2225:HOH:O	1:C:175[A]:GLN:NE2	1.85	1.08
1:B:71[A]:GLU:CD	4:B:2071:HOH:O	1.92	1.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:352[B]:GLN:CD	4:A:2476:HOH:O	1.90	1.07
1:C:243[B]:GLN:OE1	4:C:2254:HOH:O	1.71	1.07
1:C:194[A]:ARG:HH11	1:C:194[A]:ARG:HG3	1.00	1.07
1:A:435[B]:ARG:HH11	1:A:435[B]:ARG:HA	1.14	1.06
1:D:435[A]:ARG:CZ	4:D:2295:HOH:O	2.02	1.03
1:A:70[B]:LYS:HD2	1:A:74[B]:GLU:CD	1.78	1.02
1:D:435[B]:ARG:CD	4:D:2297:HOH:O	1.83	1.02
1:A:435[B]:ARG:NH1	1:A:435[B]:ARG:HA	1.75	1.02
1:B:409[B]:GLN:CD	4:B:2410:HOH:O	1.85	1.02
1:D:349[B]:ASN:OD1	4:D:2248:HOH:O	1.77	1.00
1:A:14[A]:GLU:OE1	4:A:2010:HOH:O	1.77	1.00
1:C:33[B]:ASP:OD1	4:C:2027:HOH:O	1.77	1.00
1:C:198[B]:GLN:CD	4:C:2228:HOH:O	1.90	0.99
1:A:369[B]:GLN:CG	4:A:2338:HOH:O	2.08	0.99
1:A:435[B]:ARG:NH1	4:A:2548:HOH:O	1.96	0.99
1:B:71[A]:GLU:OE1	4:B:2071:HOH:O	1.81	0.98
1:C:353[B]:GLN:CG	4:C:2312:HOH:O	2.09	0.98
1:C:347[A]:ARG:NH1	4:C:2077:HOH:O	1.63	0.97
1:B:102[A]:LYS:CE	4:B:2122:HOH:O	1.97	0.96
1:B:409[B]:GLN:OE1	4:B:2410:HOH:O	1.75	0.96
1:B:102[B]:LYS:NZ	1:B:343[B]:ASP:OD1	1.99	0.94
1:C:340:ALA:O	1:C:344[B]:ILE:HD13	1.67	0.94
1:B:203[B]:GLU:OE2	4:B:2262:HOH:O	1.88	0.92
1:C:194[A]:ARG:NH1	1:C:194[A]:ARG:HG3	1.80	0.92
1:A:186[B]:LEU:HG	1:A:235:LEU:HD22	1.53	0.91
1:C:352[B]:GLN:HG3	4:C:2313:HOH:O	1.70	0.90
1:C:242:LEU:HD21	4:C:2169:HOH:O	1.70	0.90
1:A:18[B]:ARG:NH1	4:A:2020:HOH:O	2.04	0.90
1:B:408[A]:GLN:OE1	4:B:2407:HOH:O	1.89	0.89
1:C:71[B]:GLU:OE1	1:C:344[B]:ILE:CD1	2.21	0.88
1:C:198[B]:GLN:OE1	4:C:2228:HOH:O	1.88	0.87
1:B:317[B]:GLU:CD	1:B:317[B]:GLU:H	1.78	0.87
1:D:260:GLU:HB2	1:D:284[A]:MET:HE2	1.57	0.84
1:A:70[B]:LYS:CD	1:A:74[B]:GLU:CD	2.47	0.83
1:C:194[A]:ARG:CG	1:C:194[A]:ARG:NH1	2.34	0.82
1:C:32[B]:LYS:HA	1:C:32[B]:LYS:CE	2.02	0.81
1:C:142[B]:GLN:OE1	4:C:2150:HOH:O	1.99	0.80
1:A:317[A]:GLU:CD	1:A:317[A]:GLU:H	1.85	0.80
1:C:71[B]:GLU:OE1	1:C:344[B]:ILE:HD11	1.82	0.79
1:C:428:ALA:O	4:C:2316:HOH:O	2.01	0.79
1:B:102[A]:LYS:NZ	4:B:2122:HOH:O	2.06	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:68:SER:HB2	4:C:2038:HOH:O	1.84	0.78
1:D:189[B]:GLY:O	4:D:2148:HOH:O	2.02	0.78
1:A:402:ALA:HB2	1:A:447[B]:ILE:HD11	1.65	0.78
1:A:352[A]:GLN:NE2	4:A:2476:HOH:O	2.17	0.77
4:C:2107:HOH:O	1:D:22:MET:HE2	1.84	0.77
1:C:71[B]:GLU:OE1	1:C:344[B]:ILE:HD12	1.85	0.76
1:B:140[B]:GLN:HG2	1:B:143:LYS:HG3	1.66	0.76
1:C:353[B]:GLN:HG2	4:C:2312:HOH:O	1.80	0.75
1:C:185[B]:THR:CG2	1:C:238:PRO:HD3	2.11	0.75
1:D:189[B]:GLY:O	1:D:191:ALA:N	2.16	0.75
1:C:183:PRO:O	4:C:2201:HOH:O	2.04	0.75
1:C:260:GLU:HG3	1:C:284[A]:MET:HG3	1.72	0.72
1:B:71[A]:GLU:OE2	4:B:2071:HOH:O	1.99	0.72
1:B:243:GLN:OE1	4:B:2301:HOH:O	2.08	0.72
1:B:270[B]:THR:HG21	1:B:276:TYR:HB2	1.71	0.72
1:B:210[A]:LYS:NZ	4:B:2263:HOH:O	2.22	0.71
1:B:49:LYS:HE3	4:B:2040:HOH:O	1.90	0.71
1:A:352[B]:GLN:NE2	4:A:2476:HOH:O	2.17	0.70
1:C:353[B]:GLN:HE21	1:C:427:ASN:HA	1.57	0.70
1:A:243[A]:GLN:CD	4:A:2393:HOH:O	2.14	0.69
1:D:260:GLU:CB	1:D:284[A]:MET:HE2	2.21	0.69
1:C:107:LEU:HD13	1:C:284[B]:MET:SD	2.32	0.69
1:C:353[B]:GLN:HB2	4:C:2312:HOH:O	1.93	0.69
1:A:70[B]:LYS:CD	1:A:74[B]:GLU:OE2	2.41	0.68
1:D:352[A]:GLN:OE1	1:D:356:GLU:HG3	1.93	0.68
1:B:243:GLN:HE21	1:B:247:GLU:HG3	1.59	0.67
1:C:366:HIS:O	1:C:369[B]:GLN:HG2	1.95	0.67
1:C:274:ALA:O	4:C:2273:HOH:O	2.12	0.67
1:A:343[B]:ASP:OD2	4:A:2182:HOH:O	2.12	0.66
1:C:366:HIS:O	1:C:369[B]:GLN:CG	2.43	0.66
1:A:70[B]:LYS:HD3	1:A:74[B]:GLU:OE2	1.96	0.66
1:A:243[A]:GLN:NE2	4:A:2393:HOH:O	2.26	0.66
1:D:435[B]:ARG:NH1	4:D:2297:HOH:O	1.62	0.66
1:A:243[A]:GLN:CG	4:A:2392:HOH:O	2.13	0.65
1:A:288:LYS:HE2	3:A:1450:PLP:H4A	1.78	0.65
1:A:366:HIS:O	1:A:369[B]:GLN:HG3	1.98	0.63
1:D:57:LEU:HD13	1:D:414:ARG:HB3	1.81	0.63
1:D:34:PRO:HB2	1:D:36[B]:ILE:HD11	1.81	0.63
1:B:366:HIS:O	1:B:369[B]:GLN:HG3	1.98	0.63
1:C:247:GLU:O	1:C:251[A]:GLN:HG2	1.99	0.62
1:C:24:PHE:N	4:C:2016:HOH:O	2.32	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:71[A]:GLU:HB2	1:C:344[A]:ILE:HD11	1.80	0.62
1:B:102[A]:LYS:NZ	4:B:2121:HOH:O	2.32	0.61
1:A:402:ALA:HB2	1:A:447[B]:ILE:CD1	2.30	0.61
1:C:249:CYS:HB3	1:C:254:ILE:O	2.00	0.61
1:C:405:LYS:HG2	4:C:2377:HOH:O	2.01	0.61
1:B:70[B]:LYS:NZ	1:B:74[B]:GLU:OE1	2.34	0.61
1:B:109:GLY:O	4:B:2130:HOH:O	2.16	0.61
1:C:51[B]:ARG:HH21	1:C:53:VAL:HG21	1.65	0.60
1:C:71[B]:GLU:CD	1:C:344[B]:ILE:HD11	2.21	0.60
1:C:206:ASN:ND2	4:C:2230:HOH:O	2.35	0.59
1:A:435[B]:ARG:CZ	4:A:2548:HOH:O	2.43	0.58
1:C:283:LEU:HD23	1:C:302:ALA:HA	1.83	0.58
1:A:369[B]:GLN:HG3	4:A:2338:HOH:O	1.89	0.57
1:A:352[B]:GLN:CG	4:A:2476:HOH:O	2.43	0.57
1:D:194[B]:ARG:NH1	4:D:2153:HOH:O	2.38	0.57
1:B:429:ARG:HD3	4:B:2420:HOH:O	2.04	0.56
1:A:98[B]:GLN:CD	4:A:2174:HOH:O	2.42	0.56
1:C:198[A]:GLN:HE21	1:C:239:VAL:HG21	1.71	0.56
1:C:377:ILE:HG12	1:C:386[B]:ILE:HG13	1.87	0.55
1:A:71[A]:GLU:HB2	1:A:344:ILE:HD11	1.88	0.55
1:A:70[B]:LYS:CD	1:A:74[B]:GLU:OE1	2.54	0.55
1:A:70[B]:LYS:HD2	1:A:74[B]:GLU:OE1	2.04	0.55
1:C:126:THR:O	1:C:130[B]:MET:HG3	2.07	0.55
1:D:435[B]:ARG:HG3	4:D:2297:HOH:O	2.04	0.54
1:B:140[B]:GLN:CG	1:B:143:LYS:HG3	2.36	0.54
1:B:270[B]:THR:CG2	1:B:276:TYR:HB2	2.36	0.54
1:D:182:LEU:HD22	4:D:2167:HOH:O	2.07	0.54
1:C:71[B]:GLU:CD	1:C:344[B]:ILE:CD1	2.76	0.54
1:C:51[B]:ARG:HH21	1:C:53:VAL:CG2	2.21	0.53
1:C:203[A]:GLU:HG2	4:C:2231:HOH:O	2.07	0.53
1:B:71[B]:GLU:HB2	1:B:344:ILE:HD11	1.88	0.53
1:A:184:HIS:CE1	1:A:186[B]:LEU:HB2	2.44	0.53
1:C:258:PHE:CD2	1:C:281:PRO:HG3	2.44	0.53
1:C:443:ALA:HA	4:C:2377:HOH:O	2.09	0.52
1:B:206:ASN:OD1	1:B:244:ARG:NE	2.34	0.52
1:A:194[B]:ARG:NH2	4:A:2342:HOH:O	2.42	0.52
1:B:183:PRO:HG3	1:B:207[B]:GLU:OE2	2.10	0.52
1:C:183:PRO:HG3	1:C:207:GLU:HG3	1.92	0.52
1:B:18[B]:ARG:NH1	4:B:2007:HOH:O	2.40	0.52
1:C:71[A]:GLU:OE2	1:C:347[A]:ARG:NH2	2.43	0.52
4:C:2107:HOH:O	1:D:22:MET:CE	2.47	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74[A]:GLU:HG3	1:A:78[A]:ARG:HH11	1.75	0.51
1:A:18[B]:ARG:NH2	1:B:308:ASP:OD1	2.44	0.51
1:A:98[B]:GLN:NE2	4:A:2172:HOH:O	2.43	0.51
1:D:435[A]:ARG:HG2	4:D:2297:HOH:O	2.11	0.50
1:C:280:THR:O	4:C:2260:HOH:O	2.20	0.50
1:C:201:GLY:HA2	1:C:238:PRO:HG2	1.94	0.50
1:A:74[A]:GLU:HA	1:A:74[A]:GLU:OE1	2.12	0.50
1:C:202:VAL:HG13	1:C:244:ARG:HG2	1.94	0.50
1:B:408[A]:GLN:HG2	4:B:2407:HOH:O	2.11	0.50
1:C:353[B]:GLN:CB	4:C:2312:HOH:O	2.42	0.49
1:A:397:VAL:O	1:A:401:GLU:HG3	2.12	0.49
1:D:57:LEU:HD23	1:D:60:LEU:HA	1.93	0.49
1:D:243[B]:GLN:HG3	4:D:2054:HOH:O	2.12	0.49
1:C:408:GLN:OE1	1:C:409:GLN:NE2	2.43	0.49
1:C:23:PRO:HG3	1:D:329:SER:HB2	1.93	0.49
1:D:260:GLU:CD	1:D:284[A]:MET:HE3	2.33	0.49
1:B:57:LEU:HD23	1:B:60:LEU:HA	1.95	0.49
1:C:329:SER:HB2	1:D:23:PRO:HG3	1.95	0.49
1:A:140:GLN:N	1:A:141:PRO:HD3	2.28	0.48
1:C:385:ALA:C	1:C:386[B]:ILE:HD12	2.34	0.48
1:B:190[B]:MET:HG3	1:B:190[B]:MET:H	1.37	0.48
1:B:310:PHE:O	1:B:313[B]:GLN:HG2	2.13	0.48
1:C:111:LEU:HD13	1:C:284[B]:MET:SD	2.54	0.48
1:B:352[B]:GLN:HG2	4:B:2362:HOH:O	2.14	0.47
1:B:207[B]:GLU:CD	4:B:2267:HOH:O	2.52	0.47
1:B:102[B]:LYS:NZ	1:B:343[B]:ASP:CG	2.66	0.47
1:C:199[B]:THR:O	1:C:201:GLY:N	2.48	0.47
1:D:185:THR:HB	1:D:238:PRO:HD3	1.96	0.47
1:A:283:LEU:HD23	1:A:302:ALA:HA	1.96	0.47
1:C:202:VAL:HG12	4:C:2230:HOH:O	2.14	0.46
1:B:356[B]:GLU:HA	1:B:356[B]:GLU:OE1	2.15	0.46
1:C:389:ALA:O	1:C:398:ARG:NH1	2.48	0.46
1:C:436:LEU:O	1:C:440:VAL:HG23	2.15	0.46
1:C:47:ASP:OD2	1:C:51[B]:ARG:NE	2.48	0.46
1:D:406:LEU:HD13	1:D:413:VAL:HG21	1.97	0.46
1:D:71[B]:GLU:OE2	1:D:347:ARG:NH1	2.48	0.46
1:D:260:GLU:CB	1:D:284[A]:MET:CE	2.94	0.46
1:C:33[B]:ASP:CG	4:C:2026:HOH:O	2.54	0.46
1:D:49:LYS:HE2	4:D:2019:HOH:O	2.16	0.46
1:C:71[B]:GLU:HB3	1:C:344[B]:ILE:HD11	1.98	0.45
1:D:14:GLU:O	1:D:14:GLU:HG3	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:435:ARG:HG3	4:B:2422:HOH:O	2.16	0.45
1:A:288:LYS:CE	3:A:1450:PLP:H4A	2.47	0.45
1:A:186[B]:LEU:CG	1:A:235:LEU:HD22	2.38	0.45
1:D:435[B]:ARG:HG3	4:D:2294:HOH:O	2.17	0.45
1:B:53:VAL:HG11	1:B:407:TRP:CZ3	2.52	0.45
1:C:274:ALA:C	4:C:2273:HOH:O	2.53	0.44
1:B:89:PHE:O	1:B:90:GLN:HB2	2.17	0.44
1:A:70[B]:LYS:HE3	1:A:70[B]:LYS:HB3	1.56	0.44
1:C:351:VAL:HB	4:C:2311:HOH:O	2.16	0.44
1:B:288:LYS:HE2	3:B:1449:PLP:C4A	2.48	0.44
1:A:97:PHE:CD1	1:B:20:HIS:HB3	2.53	0.44
1:A:194[B]:ARG:NH1	4:A:2340:HOH:O	2.49	0.44
1:A:358:ALA:HB3	1:A:359:PRO:HD3	2.00	0.44
1:C:79:GLN:NE2	1:C:95[B]:LEU:HD11	2.33	0.44
1:A:405:LYS:NZ	1:A:408:GLN:HE22	2.16	0.43
1:D:211:LEU:HD23	1:D:211:LEU:HA	1.83	0.43
1:C:57:LEU:HD23	1:C:60[A]:LEU:HA	2.00	0.43
1:C:353[B]:GLN:HG3	4:C:2312:HOH:O	1.94	0.43
1:C:387:GLN:HA	1:C:419:THR:HG22	1.99	0.43
1:C:435[B]:ARG:HA	1:C:435[B]:ARG:CZ	2.48	0.43
1:A:39:ALA:HA	1:B:83:LEU:HB2	1.99	0.43
1:B:78[A]:ARG:HG2	1:B:78[A]:ARG:HH11	1.83	0.43
1:C:69:ARG:NH2	1:C:344[A]:ILE:HG23	2.34	0.43
1:D:57:LEU:HD13	1:D:414:ARG:CB	2.46	0.43
1:D:183:PRO:HG3	1:D:207:GLU:HG3	2.00	0.43
1:C:194[A]:ARG:NH1	1:C:194[A]:ARG:HG2	2.30	0.43
1:A:145:LYS:HB3	1:A:220:ILE:HA	2.01	0.43
1:C:97:PHE:CD1	1:D:20:HIS:HB3	2.54	0.43
1:C:196:MET:HE1	1:C:239:VAL:HA	2.01	0.43
1:A:186[B]:LEU:HD21	1:A:378:ARG:HD3	2.01	0.43
1:B:140[B]:GLN:HG2	1:B:143:LYS:CG	2.42	0.43
1:C:326:TYR:HB2	1:C:329:SER:OG	2.19	0.42
1:C:58[B]:SER:OG	1:C:63:CYS:N	2.53	0.42
1:A:23:PRO:HG3	1:B:329:SER:HB2	2.01	0.42
1:B:249:CYS:HB3	1:B:254:ILE:O	2.18	0.42
1:A:435[B]:ARG:NH1	1:A:435[B]:ARG:CA	2.66	0.42
1:C:57:LEU:HD23	1:C:60[B]:LEU:HA	2.01	0.42
1:C:140[B]:GLN:HA	1:C:142[B]:GLN:NE2	2.35	0.42
1:B:387:GLN:HA	1:B:419:THR:HG22	2.02	0.42
1:B:102[B]:LYS:HE2	1:B:342:LEU:HB2	2.01	0.42
1:C:228[B]:MET:SD	1:C:263:THR:HG22	2.59	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:51[B]:ARG:HB2	1:C:51[B]:ARG:HE	1.41	0.42
1:D:61:TRP:HB2	1:D:288:LYS:HD3	2.01	0.42
1:B:86:SER:O	1:B:87:PRO:C	2.57	0.42
1:A:192:PHE:CE1	1:A:369[A]:GLN:HG3	2.55	0.42
1:A:186[B]:LEU:HA	1:A:186[B]:LEU:HD23	1.78	0.41
1:A:14[A]:GLU:HA	4:A:2006:HOH:O	2.19	0.41
1:B:270[B]:THR:OG1	1:B:275:GLU:CD	2.57	0.41
1:C:26:ALA:HB1	4:C:2020:HOH:O	2.20	0.41
1:D:243[A]:GLN:NE2	4:D:2194:HOH:O	1.77	0.41
1:D:398:ARG:HB2	1:D:399:PRO:HD3	2.02	0.41
1:B:344:ILE:HG13	4:B:2071:HOH:O	2.19	0.41
1:A:14[B]:GLU:HA	4:A:2006:HOH:O	2.20	0.41
1:C:199[B]:THR:HB	1:C:200[B]:GLY:H	1.67	0.41
1:A:141:PRO:CB	1:C:142[B]:GLN:HG3	2.51	0.41
1:D:378:ARG:HA	4:D:2151:HOH:O	2.21	0.41
1:C:76:VAL:HG13	1:C:333:VAL:HG13	2.03	0.41
1:C:92:GLY:O	1:D:36[B]:ILE:HA	2.21	0.41
1:D:34:PRO:CB	1:D:36[B]:ILE:HD11	2.50	0.41
1:C:385:ALA:O	1:C:386[B]:ILE:HD12	2.21	0.41
1:C:16:ASN:ND2	1:D:101[B]:GLU:OE1	2.54	0.41
1:B:175:GLN:NE2	1:D:175[A]:GLN:OE1	2.54	0.41
1:C:92:GLY:O	1:D:36[A]:ILE:HA	2.21	0.41
1:D:243[A]:GLN:HE21	1:D:279:VAL:HG22	1.86	0.41
1:B:363[B]:LYS:HB2	1:B:363[B]:LYS:HE3	1.72	0.40
1:D:34:PRO:HB2	1:D:36[B]:ILE:CD1	2.50	0.40
1:B:61:TRP:HB2	1:B:288:LYS:HD3	2.03	0.40
1:C:162:SER:OG	1:C:177[B]:MET:HG3	2.21	0.40
1:D:249:CYS:HB3	1:D:254:ILE:O	2.20	0.40
1:C:60[A]:LEU:HD21	1:D:89:PHE:CE2	2.56	0.40
1:A:249:CYS:HB3	1:A:254:ILE:O	2.20	0.40
1:C:90:GLN:O	1:D:35:ARG:NH2	2.44	0.40
1:A:20:HIS:HB3	1:B:97:PHE:CD1	2.56	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	471/448 (105%)	450 (96%)	18 (4%)	3 (1%)	30	13
1	B	471/448 (105%)	447 (95%)	20 (4%)	4 (1%)	24	8
1	C	477/448 (106%)	447 (94%)	25 (5%)	5 (1%)	19	5
1	D	463/448 (103%)	441 (95%)	18 (4%)	4 (1%)	21	6
All	All	1882/1792 (105%)	1785 (95%)	81 (4%)	16 (1%)	24	6

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	200[A]	GLY
1	C	200[B]	GLY
1	D	190[A]	MET
1	D	190[B]	MET
1	A	287	ALA
1	A	288	LYS
1	B	287	ALA
1	D	288	LYS
1	B	288	LYS
1	D	287	ALA
1	B	216	ASP
1	C	111	LEU
1	C	287	ALA
1	A	234	VAL
1	B	234	VAL
1	C	234	VAL

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	376/350 (107%)	368 (98%)	8 (2%)	61	44

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	375/350 (107%)	369 (98%)	6 (2%)	70	57
1	C	381/350 (109%)	366 (96%)	15 (4%)	39	19
1	D	367/350 (105%)	360 (98%)	7 (2%)	65	49
All	All	1499/1400 (107%)	1463 (98%)	36 (2%)	66	38

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	VAL
1	A	13	SER
1	A	125	ASP
1	A	198[A]	GLN
1	A	198[B]	GLN
1	A	228	MET
1	A	317[A]	GLU
1	A	317[B]	GLU
1	B	125	ASP
1	B	190[A]	MET
1	B	190[B]	MET
1	B	228	MET
1	B	253[A]	ASN
1	B	253[B]	ASN
1	C	14	GLU
1	C	125	ASP
1	C	175[A]	GLN
1	C	175[B]	GLN
1	C	194[A]	ARG
1	C	194[B]	ARG
1	C	198[A]	GLN
1	C	198[B]	GLN
1	C	228[A]	MET
1	C	228[B]	MET
1	C	253[A]	ASN
1	C	253[B]	ASN
1	C	369[A]	GLN
1	C	369[B]	GLN
1	C	406	LEU
1	D	14	GLU
1	D	78[A]	ARG
1	D	78[B]	ARG
1	D	125	ASP

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Mol	Chain	Res	Type
1	D	190[A]	MET
1	D	190[B]	MET
1	D	228	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	175	GLN
1	A	408	GLN
1	B	73	GLN
1	B	175	GLN
1	B	243	GLN
1	B	251	GLN
1	B	373	ASN
1	C	79	GLN
1	C	252	HIS
1	D	73	GLN
1	D	79	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	$\# Z > 2$
3	PLP	A	1450	-	16,16,16	3.75	4 (25%)	21,23,23	1.60	3 (14%)
3	PLP	B	1449	-	16,16,16	3.78	4 (25%)	21,23,23	1.30	3 (14%)
3	PLP	C	1449	-	16,16,16	3.84	3 (18%)	21,23,23	1.23	2 (9%)
3	PLP	D	1449	-	16,16,16	3.37	3 (18%)	21,23,23	1.38	4 (19%)

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
3	PLP	A	1450	-	-	0/8/8/8	0/1/1/1
3	PLP	B	1449	-	-	0/8/8/8	0/1/1/1
3	PLP	C	1449	-	-	0/8/8/8	0/1/1/1
3	PLP	D	1449	-	-	0/8/8/8	0/1/1/1

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1449	PLP	P-O3P	-2.64	1.45	1.54
3	A	1450	PLP	P-O3P	-2.51	1.45	1.54
3	A	1450	PLP	C4-C5	4.37	1.47	1.42
3	C	1449	PLP	C4-C5	4.87	1.48	1.42
3	B	1449	PLP	C4-C3	5.03	1.47	1.40
3	D	1449	PLP	C4-C5	5.24	1.49	1.42
3	B	1449	PLP	C4-C5	5.47	1.49	1.42
3	A	1450	PLP	C4-C3	5.73	1.48	1.40
3	D	1449	PLP	C4-C3	5.90	1.48	1.40
3	C	1449	PLP	C4-C3	6.07	1.48	1.40
3	D	1449	PLP	C3-C2	10.56	1.48	1.40
3	A	1450	PLP	C3-C2	12.53	1.49	1.40
3	B	1449	PLP	C3-C2	12.60	1.49	1.40
3	C	1449	PLP	C3-C2	12.92	1.49	1.40

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1450	PLP	C3-C4-C5	-4.24	114.93	118.11
3	C	1449	PLP	C3-C4-C5	-3.22	115.69	118.11
3	D	1449	PLP	O4A-C4A-C4	-2.79	119.47	125.11
3	D	1449	PLP	O2P-P-O4P	-2.17	100.32	106.56
3	B	1449	PLP	O2P-P-O4P	-2.09	100.55	106.56
3	A	1450	PLP	O3P-P-O4P	2.06	112.50	106.56
3	A	1450	PLP	O3P-P-O1P	2.25	117.82	110.58
3	D	1449	PLP	C6-N1-C2	2.26	123.89	119.28
3	B	1449	PLP	O3P-P-O1P	2.34	118.12	110.58
3	D	1449	PLP	O3-C3-C2	2.45	121.92	117.66
3	B	1449	PLP	O3-C3-C2	2.53	122.06	117.66
3	C	1449	PLP	O3-C3-C2	2.72	122.39	117.66

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1450	PLP	2	0
3	B	1449	PLP	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	439/448 (97%)	-0.09	2 (0%) 91 91	8, 19, 33, 86	0
1	B	436/448 (97%)	-0.08	2 (0%) 91 91	7, 19, 33, 69	0
1	C	435/448 (97%)	0.50	14 (3%) 51 49	17, 27, 40, 70	0
1	D	435/448 (97%)	0.21	7 (1%) 74 74	15, 25, 41, 68	0
All	All	1745/1792 (97%)	0.13	25 (1%) 78 77	7, 23, 38, 86	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	11	VAL	6.6
1	A	10	PRO	4.9
1	C	191	ALA	4.2
1	C	72	ILE	4.0
1	C	103	ILE	3.2
1	B	13	SER	3.2
1	C	333	VAL	3.1
1	D	72	ILE	3.0
1	C	413	VAL	2.8
1	C	33[A]	ASP	2.7
1	D	336	ALA	2.7
1	D	97	PHE	2.7
1	D	65	ALA	2.7
1	C	344[A]	ILE	2.6
1	B	189[A]	GLY	2.5
1	C	345	LEU	2.3
1	D	81	GLY	2.3
1	C	352[A]	GLN	2.2
1	C	334	ALA	2.2
1	C	245	LEU	2.1
1	C	76	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	424	PRO	2.1
1	C	56	SER	2.1
1	D	82	THR	2.0
1	D	343[A]	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	PLP	A	1450	16/16	0.97	0.09	0.70	8,15,26,37	0
3	PLP	B	1449	16/16	0.98	0.08	0.33	10,17,31,43	0
2	CL	A	1449	1/1	0.96	0.08	-0.06	26,26,26,26	1
3	PLP	C	1449	16/16	0.97	0.10	-0.08	13,26,41,75	0
3	PLP	D	1449	16/16	0.97	0.08	-0.43	12,24,30,49	0
2	CL	D	1450	1/1	0.98	0.05	-2.34	30,30,30,30	1
2	CL	B	1450	1/1	0.98	0.07	-4.49	22,22,22,22	1
2	CL	C	1450	1/1	0.95	0.06	-5.52	28,28,28,28	1

6.5 Other polymers [i](#)

There are no such residues in this entry.