



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:10 PM GMT

PDB ID : 1ECC
Title : ESCHERICHIA COLI GLUTAMINE PHOSPHORIBOSYLPYROPHOSPHATE (PRPP) AMIDOTRANSFERASE COMPLEXED WITH MN-CPRPP AND 5-OXO-NORLEUCINE
Authors : Krahn, J.M.; Smith, J.L.
Deposited on : 1997-07-09
Resolution : 2.40 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

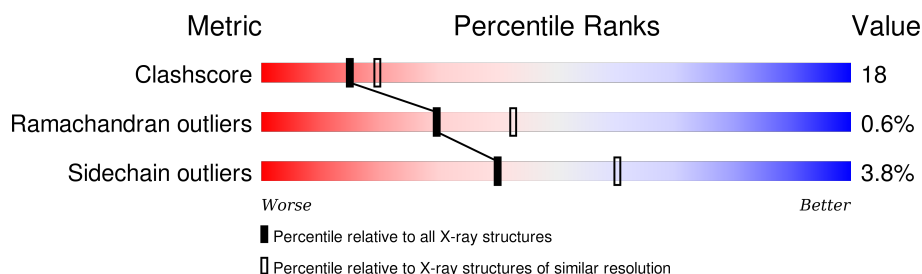
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)

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	504	
1	B	504	

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
3	ONL	A	2377	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 8092 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.

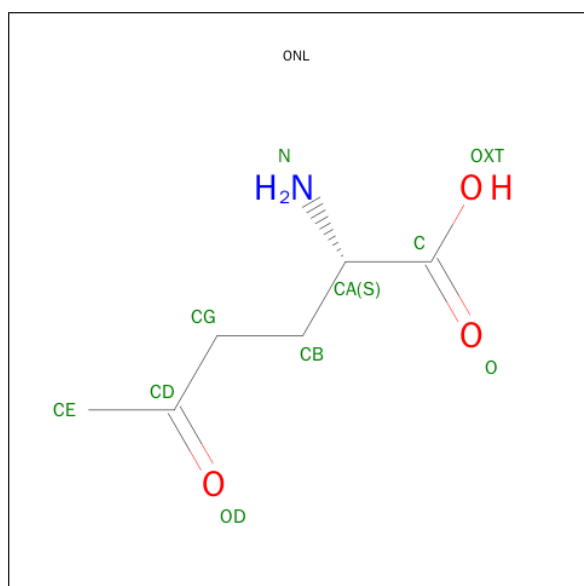
- Molecule 1 is a protein called GLUTAMINE PHOSPHORIBOSYLPYROPHOSPHATE AMIDOTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	492	Total	C	N	O	S	0	0	0
			3866	2426	692	731	17			
1	B	492	Total	C	N	O	S	0	0	0
			3866	2426	692	731	17			

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

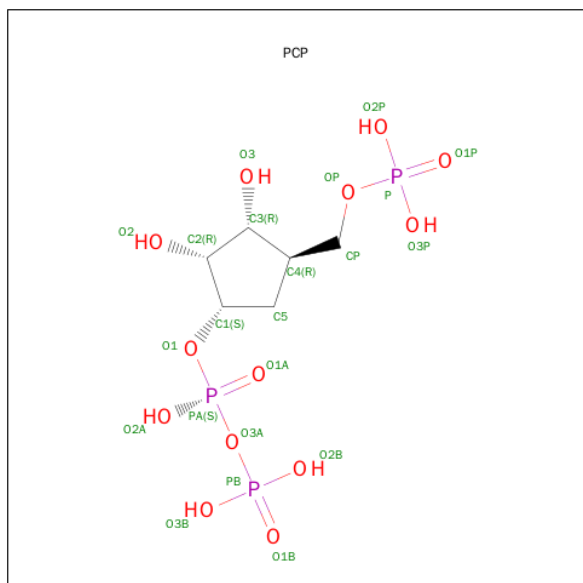
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	3	Total	Mn	0	0
			3	3		
2	A	2	Total	Mn	0	0
			2	2		

- Molecule 3 is 5-OXO-L-NORLEUCINE (three-letter code: ONL) (formula: C₆H₁₁NO₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			10	6	1	3		
3	B	1	Total	C	N	O	0	0
			10	6	1	3		

- Molecule 4 is 1-ALPHA-PYROPHOSPHORYL-2-ALPHA,3-ALPHA-DIHYDROXY-4-BETA-CYCLOPENTANE-METHANOL-5-PHOSPHATE (three-letter code: PCP) (formula: $C_6H_{15}O_{13}P_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	P	0	0
			22	6	13	3		
4	B	1	Total	C	O	P	0	0
			22	6	13	3		

- Molecule 5 is water.

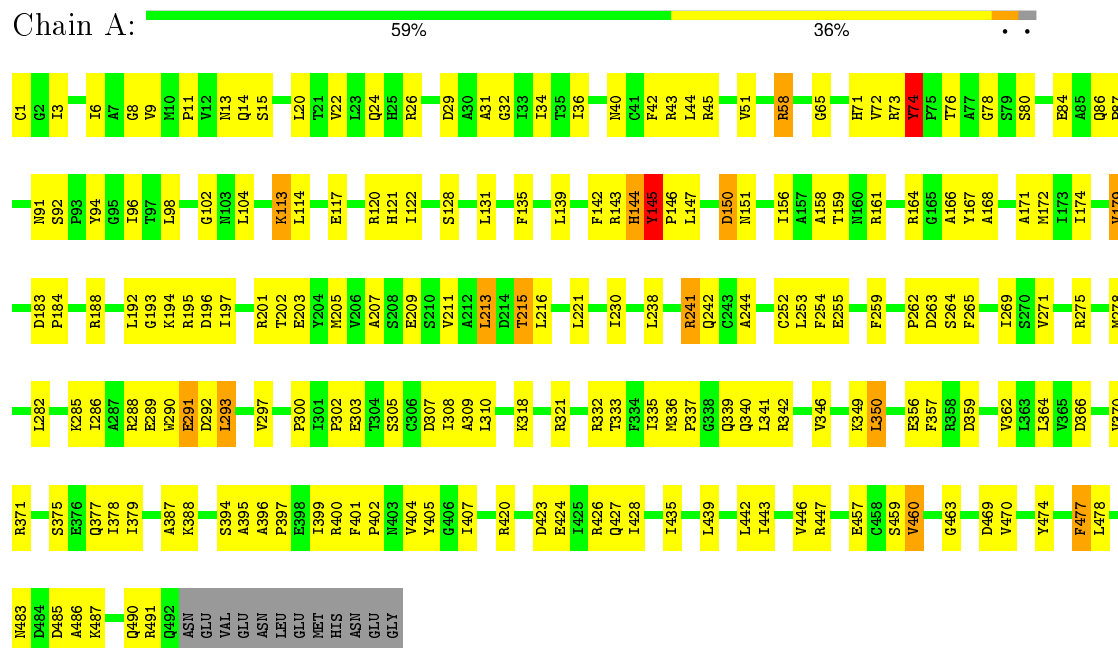
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	141	Total	O	0	0
			141	141		
5	B	150	Total	O	0	0
			150	150		

3 Residue-property plots

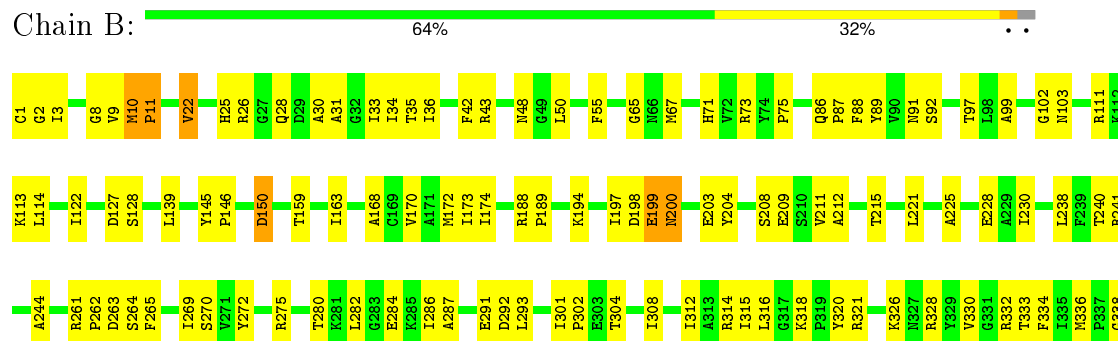
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 ($\text{RSRZ} > 2$). 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: GLUTAMINE PHOSPHORIBOSYLPYROPHOSPHATE AMIDOTRANSFERASE



- Molecule 1: GLUTAMINE PHOSPHORIBOSYLPYROPHOSPHATE AMIDOTRANSFERASE



R339	R340	R341	R342	R343	R344	S345	R349	L350	M353	R354	F357	Y365	D366	D367	Y370	Q377	I378	I379	E380	R381	A387	Y390	Y391	L392	A393	S394	A395	I399	R400	F401	P402	M403	M409	P410	E414	L415	T416	A417	D423	E424	I425	R426	Q427	I428	I429	L434
D438	L439	N440	D441	V446	R447	A448	E449	N450	P451	F456	N462	G463	V464	D469	L475	D476	F477	T480	L481	R482	N483	D484	D485	A486	V489	Q492	ASN	GLU	VAL	GLU	ASN	LEU	GLU	MET	HIS	ASN	GLU	GLY								

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	78.20 Å 78.20 Å 308.80 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.40	Depositor
% Data completeness (in resolution range)	88.3 (25.00-2.40)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.176 , 0.265	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	8092	wwPDB-VP
Average B, all atoms (Å ²)	34.0	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: PCP, MN, ONL

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	0.60	0/3934	0.81	4/5324 (0.1%)
1	B	0.62	0/3934	0.80	3/5324 (0.1%)
All	All	0.61	0/7868	0.80	7/10648 (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

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	263	ASP	CB-CG-OD1	5.87	123.58	118.30
1	B	353	ASN	N-CA-C	-5.84	95.24	111.00
1	B	73	ARG	NE-CZ-NH2	-5.41	117.59	120.30
1	A	168	ALA	N-CA-C	-5.32	96.64	111.00
1	A	241	ARG	NE-CZ-NH2	-5.27	117.66	120.30
1	B	168	ALA	N-CA-C	-5.17	97.05	111.00
1	A	74	TYR	N-CA-C	-5.09	97.27	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	145	TYR	Sidechain

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	3866	0	3831	150	0
1	B	3866	0	3831	125	0
2	A	2	0	0	0	0
2	B	3	0	0	0	0
3	A	10	0	9	6	0
3	B	10	0	9	3	0
4	A	22	0	8	1	0
4	B	22	0	8	2	0
5	A	141	0	0	14	0
5	B	150	0	0	7	0
All	All	8092	0	7696	273	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 18.

All (273) 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:333:THR:HA	1:A:336:MET:SD	2.22	0.79
1:A:275:ARG:HA	1:A:278:MET:HE3	1.66	0.78
1:A:285:LYS:HD3	1:A:435:ILE:HD11	1.64	0.78
1:A:404:VAL:HG12	1:A:478:LEU:HD11	1.68	0.74
1:A:1:CYS:H1	3:A:2377:ONL:HE1	1.50	0.74
1:A:253:LEU:HD21	1:A:442:LEU:HD23	1.68	0.74
1:B:336:MET:HB2	1:B:342:ARG:HG2	1.70	0.72
1:A:282:LEU:HD22	1:A:395:ALA:HA	1.72	0.72
1:A:9:VAL:HG11	1:A:201:ARG:HD2	1.74	0.69
1:B:9:VAL:CG2	1:B:197:ILE:HG13	2.22	0.69
1:B:334:PHE:HE1	1:B:370:VAL:HG11	1.57	0.69
5:A:2383:HOH:O	1:B:326:LYS:HD3	1.93	0.69
1:B:113:LYS:HB3	1:B:113:LYS:NZ	2.08	0.68
1:B:350:LEU:CD1	1:B:378:ILE:HD11	2.23	0.68
1:A:487:LYS:O	1:A:491:ARG:HG3	1.94	0.68
1:B:139:LEU:HD21	1:B:173:ILE:HD13	1.76	0.67
1:B:379:ILE:HG23	1:B:390:VAL:HG21	1.74	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:9:VAL:HG21	1:B:197:ILE:HG13	1.75	0.67
1:A:333:THR:HG23	1:A:349:LYS:HD2	1.77	0.67
1:B:476:ASP:O	1:B:480:THR:HG23	1.94	0.67
1:A:36:ILE:O	1:A:65:GLY:HA3	1.96	0.66
1:A:282:LEU:O	1:A:286:ILE:HG13	1.97	0.65
1:B:447:ARG:HG2	1:B:456:PHE:HE2	1.61	0.65
1:B:199:GLU:O	1:B:200:ASN:HB2	1.97	0.65
1:A:135:PHE:O	1:A:139:LEU:HB2	1.98	0.64
1:B:30:ALA:HB2	1:B:48:ASN:HD22	1.63	0.64
1:A:58:ARG:HG2	5:A:2389:HOH:O	1.97	0.62
1:B:211:VAL:HG13	1:B:264:SER:OG	1.98	0.62
1:B:333:THR:HG22	1:B:345:SER:HB3	1.81	0.62
1:A:423:ASP:O	1:A:427:GLN:HG3	1.99	0.62
1:B:1:CYS:N	3:B:2450:ONL:HE1	2.14	0.62
1:B:334:PHE:CE1	1:B:370:VAL:HG11	2.35	0.62
1:B:334:PHE:HB2	4:B:505:PCP:H52	1.81	0.62
1:A:98:LEU:HD12	1:A:171:ALA:HB2	1.80	0.62
1:A:269:ILE:HD13	1:A:446:VAL:HG12	1.82	0.61
1:A:285:LYS:O	1:A:289:GLU:HG2	2.00	0.61
1:A:192:LEU:HG	1:A:221:LEU:HD12	1.83	0.61
1:A:1:CYS:N	3:A:2377:ONL:HE1	2.16	0.60
1:A:265:PHE:CZ	1:B:354:ARG:HD2	2.36	0.60
1:B:410:PRO:CB	1:B:482:ARG:HB2	2.31	0.60
1:B:350:LEU:HD12	1:B:378:ILE:HD11	1.83	0.60
1:A:86:GLN:HB3	1:A:87:PRO:HA	1.82	0.60
1:A:114:LEU:HD22	1:A:120:ARG:HG3	1.83	0.60
1:B:114:LEU:HD13	1:B:122:ILE:HD11	1.83	0.59
1:B:1:CYS:H3	3:B:2450:ONL:HE1	1.65	0.59
1:A:150:ASP:CG	1:A:241:ARG:HH22	2.04	0.59
1:A:443:ILE:O	1:A:447:ARG:HG3	2.03	0.59
1:B:36:ILE:HD13	1:B:174:ILE:HD13	1.83	0.59
1:A:24:GLN:OE1	1:A:51:VAL:HG23	2.03	0.59
1:A:253:LEU:CD2	1:A:442:LEU:HD23	2.33	0.59
1:A:339:GLN:OE1	1:A:371:ARG:HD2	2.03	0.59
1:B:462:ASN:CG	1:B:464:VAL:HG13	2.22	0.59
1:B:28:GLN:HE22	1:B:332:ARG:H	1.50	0.58
1:B:282:LEU:HD22	1:B:395:ALA:HA	1.84	0.58
1:A:474:TYR:O	1:A:477:PHE:HB3	2.04	0.57
1:B:410:PRO:HB2	1:B:482:ARG:HB2	1.86	0.57
1:B:36:ILE:O	1:B:65:GLY:HA3	2.04	0.57
1:A:76:THR:OG1	3:A:2377:ONL:HG2	2.03	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:285:LYS:CD	1:A:435:ILE:HD11	2.33	0.57
1:B:3:ILE:HG22	1:B:71:HIS:HD2	1.70	0.56
1:A:401:PHE:HB2	1:A:459:SER:O	2.04	0.56
1:A:346:VAL:HB	1:A:377:GLN:NE2	2.20	0.56
1:A:193:GLY:HA2	1:A:221:LEU:HG	1.86	0.56
1:B:423:ASP:O	1:B:427:GLN:HG3	2.05	0.56
1:A:164:ARG:HG3	5:A:2444:HOH:O	2.05	0.56
1:B:262:PRO:O	1:B:272:TYR:HB2	2.05	0.55
1:B:340:GLN:HG3	1:B:344:LYS:HE3	1.87	0.55
1:B:128:SER:OG	3:B:2450:ONL:HA	2.07	0.55
1:B:287:ALA:O	1:B:291:GLU:HB3	2.07	0.55
1:B:269:ILE:HG23	1:B:449:GLU:HB2	1.88	0.55
1:A:357:PHE:HB3	1:A:387:ALA:HB2	1.88	0.55
1:B:275:ARG:O	1:B:308:ILE:HG12	2.07	0.55
1:A:336:MET:HB3	1:A:341:LEU:HD23	1.89	0.54
1:B:10:MET:HB3	1:B:11:PRO:HD2	1.88	0.54
1:B:350:LEU:HD13	1:B:378:ILE:HD11	1.88	0.54
1:B:485:ASP:O	1:B:489:VAL:HG23	2.08	0.54
1:B:269:ILE:HG23	1:B:449:GLU:CB	2.38	0.54
1:B:36:ILE:HG12	1:B:42:PHE:CE1	2.42	0.54
1:B:67:MET:HG2	1:B:172:MET:SD	2.48	0.54
1:B:139:LEU:HD21	1:B:173:ILE:CD1	2.38	0.53
1:B:308:ILE:HG21	1:B:395:ALA:HB2	1.91	0.53
1:A:370:VAL:HG22	1:A:399:ILE:HD12	1.89	0.53
1:A:26:ARG:HD2	1:A:209:GLU:OE2	2.09	0.53
1:B:379:ILE:HD11	1:B:429:ILE:HD12	1.90	0.53
1:A:271:VAL:O	1:A:275:ARG:HG3	2.08	0.53
1:A:74:TYR:O	3:A:2377:ONL:HE3	2.09	0.53
1:B:333:THR:HG22	1:B:345:SER:CB	2.39	0.53
1:A:332:ARG:HB2	1:A:335:ILE:HG12	1.91	0.53
1:A:275:ARG:CA	1:A:278:MET:HE3	2.36	0.52
1:A:167:TYR:CE1	1:A:183:ASP:HA	2.44	0.52
1:A:308:ILE:HD13	1:A:395:ALA:CB	2.40	0.52
1:A:34:ILE:HB	1:A:44:LEU:HD12	1.91	0.52
1:A:241:ARG:HG2	1:A:242:GLN:N	2.24	0.52
1:A:254:PHE:HD2	1:A:407:ILE:HD13	1.75	0.52
1:A:15:SER:HB3	1:A:216:LEU:HD13	1.92	0.51
1:A:184:PRO:HB3	1:A:244:ALA:HB1	1.92	0.51
1:A:102:GLY:HA2	1:A:188:ARG:CD	2.41	0.51
1:A:8:GLY:HA2	1:A:203:GLU:HG2	1.92	0.51
1:A:194:LYS:HA	1:A:203:GLU:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:164:ARG:HB3	1:A:469:ASP:OD2	2.10	0.50
1:B:189:PRO:HB2	1:B:208:SER:O	2.11	0.50
1:A:3:ILE:O	1:A:207:ALA:HA	2.10	0.50
1:A:29:ASP:HB2	1:A:73:ARG:O	2.12	0.50
1:A:364:LEU:HD13	1:A:378:ILE:HG22	1.93	0.50
1:A:156:ILE:O	1:A:159:THR:HB	2.11	0.50
1:B:31:ALA:HA	1:B:71:HIS:O	2.12	0.50
1:B:225:ALA:O	1:B:228:GLU:HB3	2.12	0.50
1:A:262:PRO:HA	1:A:271:VAL:HB	1.93	0.50
1:A:404:VAL:HG21	1:A:470:VAL:HG21	1.93	0.50
1:B:113:LYS:HB3	1:B:113:LYS:HZ2	1.76	0.50
1:A:31:ALA:HA	1:A:71:HIS:O	2.12	0.49
1:A:300:PRO:HB3	1:A:309:ALA:HB3	1.93	0.49
1:A:375:SER:O	1:A:379:ILE:HG13	2.12	0.49
1:A:290:TRP:O	1:A:293:LEU:HB2	2.11	0.49
1:B:377:GLN:O	1:B:381:MET:HG3	2.12	0.49
1:A:6:ILE:HG12	1:A:205:MET:HB2	1.95	0.49
1:B:145:TYR:CD1	1:B:146:PRO:HA	2.48	0.49
1:A:194:LYS:HB2	1:A:202:THR:CG2	2.42	0.49
1:A:424:GLU:O	1:A:428:ILE:HG13	2.13	0.49
1:A:121:HIS:CE1	5:A:2515:HOH:O	2.66	0.49
1:A:102:GLY:HA2	1:A:188:ARG:HD2	1.94	0.48
1:B:282:LEU:HD23	1:B:312:ILE:HD11	1.94	0.48
1:A:337:PRO:HB2	1:A:491:ARG:NH1	2.29	0.48
1:A:307:ASP:HB2	5:A:2506:HOH:O	2.13	0.48
1:B:338:GLY:O	1:B:342:ARG:HG3	2.12	0.48
1:A:401:PHE:O	1:A:460:VAL:HG12	2.13	0.48
1:A:166:ALA:HB1	1:A:188:ARG:HG2	1.96	0.48
1:A:11:PRO:HA	1:A:65:GLY:O	2.14	0.48
1:A:359:ASP:O	1:A:388:LYS:HD3	2.14	0.48
1:B:330:VAL:HG11	1:B:349:LYS:HB2	1.95	0.48
1:B:447:ARG:NH1	1:B:451:PRO:O	2.47	0.48
1:A:357:PHE:CD1	1:A:362:VAL:HG11	2.49	0.48
1:A:404:VAL:HG22	5:A:2398:HOH:O	2.13	0.48
1:A:188:ARG:HD2	1:A:405:TYR:O	2.13	0.48
1:B:86:GLN:HB3	1:B:87:PRO:HA	1.96	0.48
1:A:44:LEU:HG	1:A:45:ARG:N	2.29	0.47
1:B:293:LEU:HD11	1:B:391:TYR:CE1	2.49	0.47
1:A:102:GLY:O	3:A:2377:ONL:HB1	2.14	0.47
1:B:34:ILE:HA	1:B:43:ARG:O	2.14	0.47
1:A:486:ALA:O	1:A:490:GLN:HG3	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:425:ILE:O	1:B:429:ILE:HG12	2.14	0.47
1:A:32:GLY:HA3	1:A:86:GLN:O	2.15	0.47
1:A:128:SER:OG	3:A:2377:ONL:HA	2.14	0.47
1:A:195:ARG:HB3	1:A:203:GLU:HB2	1.97	0.47
1:A:483:ASN:O	1:A:486:ALA:N	2.47	0.47
1:A:14:GLN:HB3	5:A:2395:HOH:O	2.15	0.47
1:A:22:VAL:HG11	1:A:215:THR:HG21	1.97	0.47
1:B:318:LYS:HA	1:B:318:LYS:HD3	1.69	0.47
1:A:195:ARG:HA	5:A:2453:HOH:O	2.14	0.47
1:A:446:VAL:HG23	5:A:2514:HOH:O	2.14	0.46
1:A:491:ARG:HD2	5:A:2491:HOH:O	2.14	0.46
1:B:321:ARG:HD3	1:B:321:ARG:HA	1.71	0.46
1:B:403:ASN:H	1:B:409:MET:HB2	1.79	0.46
1:A:131:LEU:HD23	1:A:131:LEU:O	2.15	0.46
1:B:75:PRO:HD2	5:B:2561:HOH:O	2.15	0.46
1:A:34:ILE:HD11	1:A:42:PHE:HB3	1.97	0.46
1:B:440:ASN:H	1:B:440:ASN:ND2	2.13	0.46
1:B:26:ARG:HD2	1:B:209:GLU:OE2	2.16	0.46
1:A:401:PHE:CD2	1:A:463:GLY:N	2.84	0.46
1:A:145:TYR:CD1	1:A:146:PRO:HA	2.51	0.46
1:B:261:ARG:HB3	1:B:263:ASP:OD1	2.16	0.46
1:A:252:CYS:HA	1:A:457:GLU:HB3	1.97	0.46
1:A:442:LEU:HG	5:A:2514:HOH:O	2.14	0.46
1:A:34:ILE:HA	1:A:43:ARG:O	2.16	0.46
1:B:302:PRO:HG3	5:B:2576:HOH:O	2.15	0.46
1:A:333:THR:O	1:A:342:ARG:HG2	2.16	0.46
1:A:13:ASN:OD1	1:A:14:GLN:N	2.48	0.46
1:B:212:ALA:O	1:B:215:THR:HG22	2.15	0.46
1:B:99:ALA:HB3	1:B:170:VAL:HG12	1.98	0.46
1:A:302:PRO:HA	1:A:303:GLU:HA	1.58	0.45
1:B:477:PHE:O	1:B:481:LEU:HD13	2.16	0.45
1:A:265:PHE:HA	1:A:269:ILE:O	2.17	0.45
1:A:113:LYS:O	1:A:117:GLU:HB2	2.16	0.45
1:B:392:LEU:HD12	1:B:393:ALA:N	2.32	0.45
1:B:301:ILE:HD12	1:B:366:ASP:HB2	1.98	0.45
5:A:2497:HOH:O	1:B:22:VAL:HG13	2.16	0.45
1:B:9:VAL:HG22	1:B:197:ILE:HG13	1.95	0.45
1:A:400:ARG:HD3	5:A:2394:HOH:O	2.16	0.45
1:B:475:LEU:HD23	1:B:475:LEU:HA	1.82	0.45
1:B:150:ASP:HB2	5:B:2572:HOH:O	2.16	0.45
1:B:198:ASP:O	1:B:200:ASN:N	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:288:ARG:HB2	1:A:288:ARG:HE	1.52	0.45
1:B:42:PHE:HD2	1:B:89:TYR:CE2	2.35	0.44
1:A:350:LEU:HD13	1:A:378:ILE:HD11	1.98	0.44
1:A:426:ARG:HG2	1:A:426:ARG:HH11	1.82	0.44
4:A:505:PCP:O3B	1:B:328:ARG:NH1	2.51	0.44
1:A:211:VAL:HG13	1:A:264:SER:OG	2.18	0.44
1:B:447:ARG:HG2	1:B:456:PHE:CE2	2.45	0.44
1:A:401:PHE:HA	1:A:402:PRO:HD3	1.77	0.44
1:B:392:LEU:O	1:B:434:LEU:HA	2.18	0.44
1:B:357:PHE:O	1:B:387:ALA:HA	2.17	0.44
1:B:486:ALA:HB1	5:B:2579:HOH:O	2.18	0.44
1:A:310:LEU:HD21	1:B:320:TYR:O	2.17	0.44
1:A:94:TYR:CE2	1:A:147:LEU:HG	2.52	0.44
1:A:321:ARG:HG3	1:A:356:GLU:OE1	2.18	0.44
1:A:96:ILE:HA	1:A:172:MET:O	2.17	0.44
1:B:462:ASN:OD1	1:B:464:VAL:HG13	2.17	0.43
1:B:282:LEU:O	1:B:286:ILE:HG13	2.18	0.43
1:A:357:PHE:HD1	1:A:362:VAL:HG11	1.83	0.43
1:B:304:THR:HB	1:B:367:ASP:OD1	2.18	0.43
1:B:2:GLY:O	1:B:71:HIS:HA	2.18	0.43
1:B:25:HIS:CE1	1:B:261:ARG:HG3	2.54	0.43
1:A:94:TYR:O	1:A:174:ILE:HB	2.18	0.43
1:B:240:THR:O	1:B:241:ARG:HB2	2.18	0.43
1:B:462:ASN:ND2	1:B:464:VAL:HG13	2.33	0.43
1:A:196:ASP:HA	1:A:202:THR:HA	2.00	0.43
1:A:255:GLU:HA	1:A:259:PHE:HB3	1.98	0.43
1:A:378:ILE:HA	1:A:378:ILE:HD13	1.76	0.43
1:A:213:LEU:HD23	1:A:213:LEU:N	2.32	0.43
1:B:88:PHE:O	1:B:97:THR:HA	2.19	0.43
1:A:34:ILE:HG22	1:A:87:PRO:HG3	2.00	0.43
1:B:269:ILE:HD13	1:B:446:VAL:HG12	2.00	0.43
1:B:244:ALA:HB1	5:B:2473:HOH:O	2.19	0.43
1:B:438:ASP:O	1:B:441:ASP:HB2	2.18	0.43
1:A:349:LYS:NZ	5:A:2512:HOH:O	2.52	0.43
1:B:103:ASN:HA	1:B:127:ASP:OD2	2.19	0.43
1:B:401:PHE:HA	1:B:402:PRO:HD3	1.84	0.43
1:A:104:LEU:HD21	1:A:131:LEU:HD12	2.01	0.43
1:B:8:GLY:HA2	1:B:203:GLU:HG2	2.00	0.43
1:A:139:LEU:HD23	1:A:139:LEU:HA	1.59	0.42
1:B:308:ILE:HG13	5:B:2508:HOH:O	2.19	0.42
1:A:144:HIS:CE1	1:A:151:ASN:OD1	2.72	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:404:VAL:HG12	1:A:478:LEU:CD1	2.42	0.42
1:B:199:GLU:O	1:B:200:ASN:CB	2.65	0.42
1:A:142:PHE:C	1:A:143:ARG:HG3	2.39	0.42
1:B:354:ARG:HG2	5:B:2582:HOH:O	2.18	0.42
1:B:102:GLY:HA2	1:B:188:ARG:HD2	2.01	0.42
1:A:300:PRO:O	1:A:302:PRO:HD3	2.19	0.42
1:B:33:ILE:HG13	1:B:55:PHE:CZ	2.54	0.42
1:B:312:ILE:HD12	1:B:365:VAL:HG11	2.00	0.42
1:A:305:SER:HB3	1:A:366:ASP:HA	2.01	0.42
1:B:367:ASP:OD1	4:B:505:PCP:O2	2.37	0.42
1:B:343:ARG:O	1:B:377:GLN:NE2	2.52	0.42
1:A:179:VAL:HA	1:A:230:ILE:O	2.20	0.42
1:B:265:PHE:CE2	1:B:270:SER:HB2	2.54	0.42
1:A:366:ASP:O	1:A:394:SER:HA	2.19	0.42
1:A:114:LEU:HD13	1:A:122:ILE:HD11	2.01	0.42
1:A:340:GLN:HA	1:A:340:GLN:NE2	2.34	0.42
1:B:333:THR:HA	1:B:336:MET:SD	2.60	0.42
1:B:230:ILE:HG23	1:B:238:LEU:HD11	2.02	0.42
1:A:333:THR:CG2	1:A:349:LYS:HD2	2.49	0.41
1:A:3:ILE:HG22	1:A:71:HIS:HD2	1.84	0.41
1:A:364:LEU:HD13	1:A:378:ILE:CG2	2.50	0.41
1:A:297:VAL:HG12	1:A:362:VAL:HG12	2.02	0.41
1:B:314:ARG:NH2	1:B:315:ILE:HD11	2.35	0.41
1:A:230:ILE:HG23	1:A:238:LEU:HD11	2.03	0.41
1:A:318:LYS:HD3	1:A:318:LYS:HA	1.88	0.41
1:B:379:ILE:CD1	1:B:429:ILE:HD12	2.49	0.41
1:B:34:ILE:HD11	1:B:42:PHE:HB3	2.02	0.41
1:B:28:GLN:HB3	1:B:50:LEU:HD21	2.03	0.41
1:B:280:THR:O	1:B:284:GLU:HG3	2.21	0.41
1:A:20:LEU:HD22	1:A:72:VAL:HG13	2.02	0.41
1:B:42:PHE:CD2	1:B:89:TYR:CE2	3.09	0.41
1:B:312:ILE:O	1:B:316:LEU:HG	2.21	0.41
1:A:300:PRO:HB3	1:A:309:ALA:CB	2.51	0.41
1:A:143:ARG:C	1:A:145:TYR:H	2.24	0.41
1:A:158:ALA:O	1:A:161:ARG:HB2	2.21	0.41
1:B:410:PRO:HG3	1:B:482:ARG:HB3	2.03	0.41
1:A:439:LEU:O	1:A:443:ILE:HD12	2.21	0.41
1:A:420:ARG:NE	1:A:428:ILE:HD11	2.36	0.41
1:A:396:ALA:HA	1:A:397:PRO:HD3	1.85	0.41
1:B:194:LYS:HB3	1:B:204:TYR:CD1	2.56	0.41
1:B:221:LEU:HD23	1:B:221:LEU:HA	1.83	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:35:THR:OG1	1:B:43:ARG:HB2	2.22	0.40
1:A:142:PHE:CD2	1:A:151:ASN:HB3	2.56	0.40
1:B:399:ILE:HG22	1:B:415:LEU:HD13	2.02	0.40
1:B:159:THR:O	1:B:163:ILE:HG12	2.21	0.40
1:A:166:ALA:CB	1:A:405:TYR:HB3	2.51	0.40
1:A:197:ILE:HB	1:A:201:ARG:O	2.21	0.40
1:A:9:VAL:CG2	1:A:197:ILE:HG13	2.51	0.40
1:A:78:GLY:HA3	1:A:84:GLU:OE1	2.20	0.40
1:B:417:ALA:HA	1:B:425:ILE:HD11	2.03	0.40
1:A:370:VAL:HG22	1:A:399:ILE:CD1	2.52	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	490/504 (97%)	460 (94%)	27 (6%)	3 (1%)	30	43
1	B	490/504 (97%)	457 (93%)	30 (6%)	3 (1%)	30	43
All	All	980/1008 (97%)	917 (94%)	57 (6%)	6 (1%)	30	43

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	145	TYR
1	B	199	GLU
1	B	200	ASN
1	B	469	ASP
1	A	144	HIS
1	A	291	GLU

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	411/422 (97%)	393 (96%)	18 (4%)	35	53
1	B	411/422 (97%)	398 (97%)	13 (3%)	46	68
All	All	822/844 (97%)	791 (96%)	31 (4%)	40	60

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	ASN
1	A	58	ARG
1	A	74	TYR
1	A	80	SER
1	A	91	ASN
1	A	92	SER
1	A	113	LYS
1	A	150	ASP
1	A	179	VAL
1	A	213	LEU
1	A	215	THR
1	A	291	GLU
1	A	292	ASP
1	A	293	LEU
1	A	350	LEU
1	A	460	VAL
1	A	477	PHE
1	A	485	ASP
1	B	10	MET
1	B	11	PRO
1	B	22	VAL
1	B	91	ASN
1	B	92	SER
1	B	111	ARG
1	B	150	ASP
1	B	292	ASP
1	B	410	PRO

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Mol	Chain	Res	Type
1	B	414	GLU
1	B	440	ASN
1	B	464	VAL
1	B	483	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	48	ASN
1	A	141	ASN
1	A	144	HIS
1	B	28	GLN
1	B	48	ASN
1	B	144	HIS
1	B	185	ASN
1	B	427	GLN
1	B	440	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 5 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	ONL	A	2377	1	6,9,9	0.37	0	5,11,11	1.10	1 (20%)
4	PCP	A	505	2	20,22,22	1.82	5 (25%)	31,35,35	1.42	2 (6%)
3	ONL	B	2450	1	6,9,9	0.22	0	5,11,11	0.58	0
4	PCP	B	505	2	20,22,22	1.43	4 (20%)	31,35,35	1.50	6 (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	ONL	A	2377	1	-	0/5/9/9	0/0/0/0
4	PCP	A	505	2	-	0/17/33/33	0/1/1/1
3	ONL	B	2450	1	-	0/5/9/9	0/0/0/0
4	PCP	B	505	2	-	0/17/33/33	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	505	PCP	C4-C3	-2.60	1.46	1.54
4	B	505	PCP	P-O2P	-2.57	1.45	1.54
4	B	505	PCP	P-O3P	-2.34	1.46	1.54
4	A	505	PCP	C5-C4	-2.32	1.47	1.53
4	A	505	PCP	C3-C2	-2.12	1.47	1.53
4	A	505	PCP	P-O2P	-2.10	1.47	1.54
4	A	505	PCP	CP-C4	2.97	1.58	1.51
4	B	505	PCP	PA-O1	3.09	1.69	1.60
4	A	505	PCP	PA-O1	4.01	1.71	1.60

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	505	PCP	C4-C3-C2	-5.75	98.20	103.50
4	B	505	PCP	O3-C3-C4	-3.28	104.46	112.62
4	B	505	PCP	O3P-P-OP	-2.53	99.27	106.56
4	B	505	PCP	C4-C3-C2	-2.46	101.23	103.50
3	A	2377	ONL	CB-CG-CD	-2.45	111.78	114.52
4	A	505	PCP	O2A-PA-O3A	2.16	114.89	105.09

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Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
4	B	505	PCP	O3A-PA-O1	2.19	109.93	103.63
4	B	505	PCP	C5-C4-C3	2.32	107.85	102.24
4	B	505	PCP	OP-P-O1P	2.71	114.04	107.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	2377	ONL	6	0
4	A	505	PCP	1	0
3	B	2450	ONL	3	0
4	B	505	PCP	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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.