Full wwPDB X-ray Structure Validation Report (i)

Nov 6, 2023 – 12:34 am GMT

PDB ID	:	6HNV
Title	:	Crystal structure of aminotransferase Aro9 from C. Albicans with ligands
Authors	:	Kiliszek, A.; Rzad, K.; Rypniewski, W.; Milewski, S.; Gabriel, I.
Deposited on	:	2018-09-17
Resolution	:	2.60 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 (i) 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 (i)) were used in the production of this report:

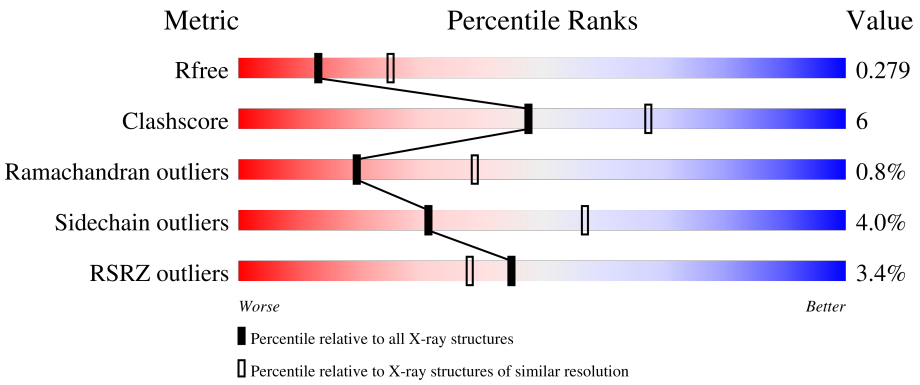
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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 <=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	524	3% 72% 15% • 11%
1	B	524	3% 76% 12% • 11%

2 Entry composition (i)

There are 9 unique types of molecules in this entry. The entry contains 7849 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.

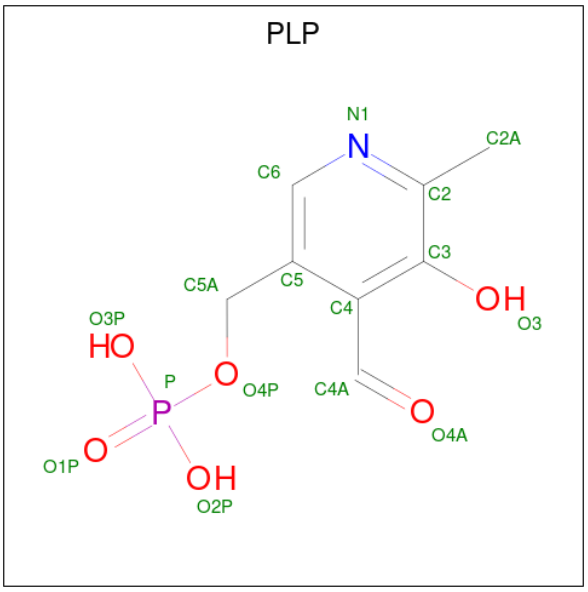
- Molecule 1 is a protein called Aromatic-amino-acid:2-oxoglutarate transaminase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	467	3765	2441	617	701	6	0	0	0
1	B	467	3772	2445	620	701	6	0	1	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	524	HIS	-	expression tag	UNP A0A1D8PMC5
B	524	HIS	-	expression tag	UNP A0A1D8PMC5

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: $C_8H_{10}NO_6P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	15	8	1	5	1	0	0

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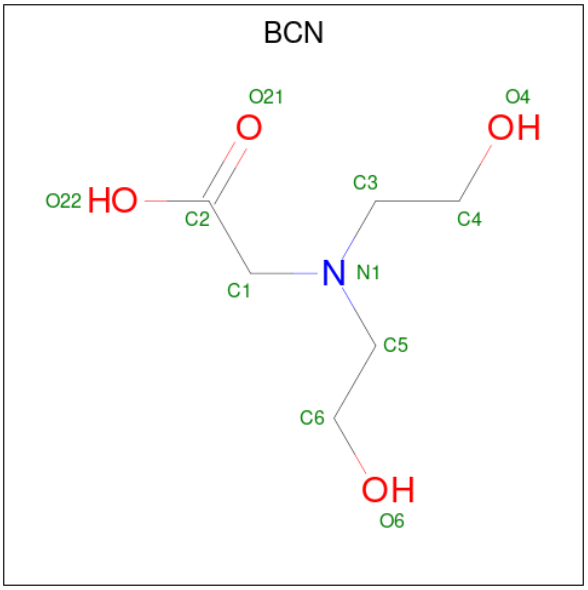
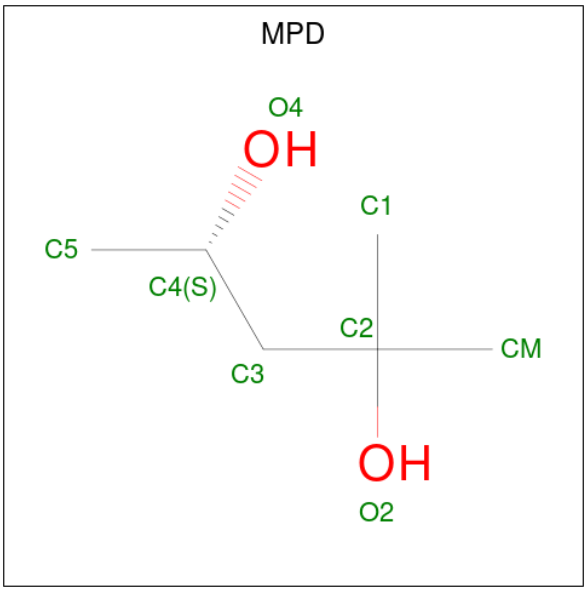
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	B	1	15	8	1	5	1	0	0

- Molecule 3 is BICINE (three-letter code: BCN) (formula: $C_6H_{13}NO_4$).



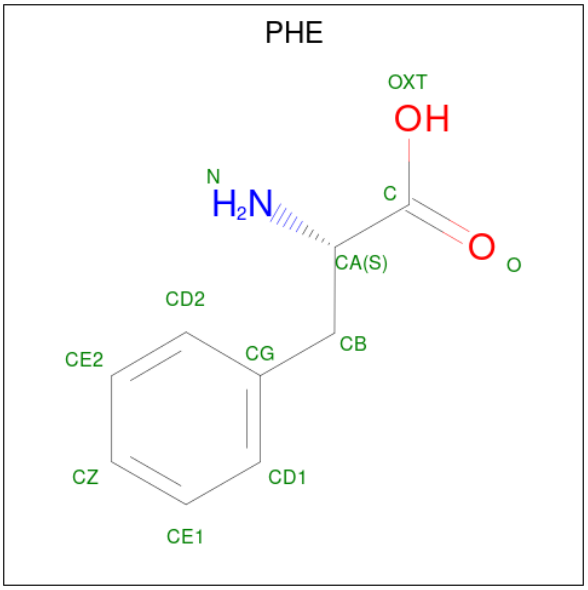
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	11	6	1	4	0	0
			Total	C	N	O		
3	B	1	11	6	1	4	0	0
			Total	C	N	O		
3	B	1	11	6	1	4	0	0

- Molecule 4 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



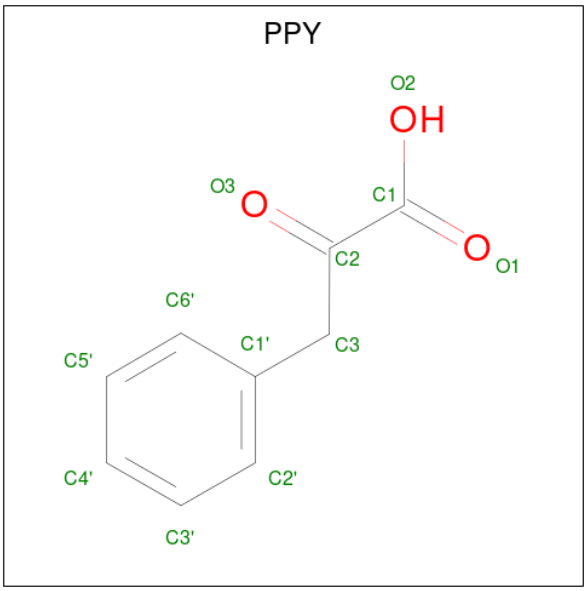
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
4	A	1	8	6	2	0	0

- Molecule 5 is PHENYLALANINE (three-letter code: PHE) (formula: $C_9H_{11}NO_2$).



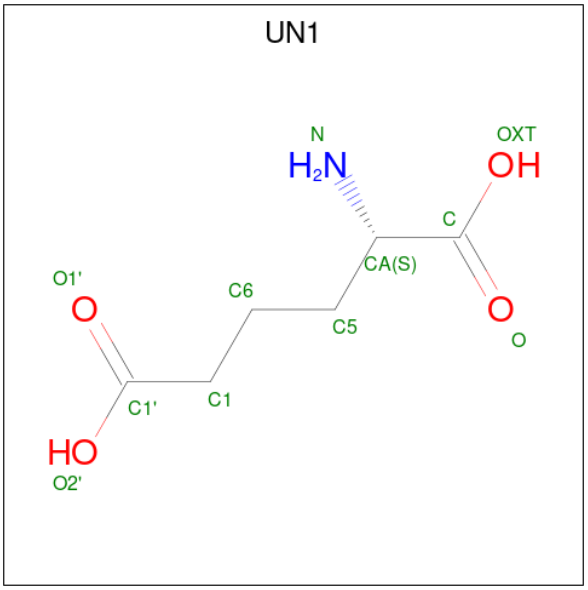
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
5	A	1	12	9	1	2	0	1

- Molecule 6 is 3-PHENYLPYRUVIC ACID (three-letter code: PPY) (formula: $C_9H_8O_3$).



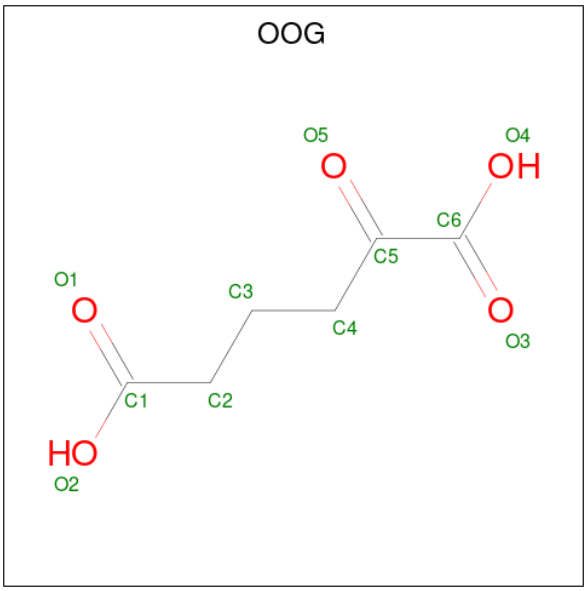
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
6	A	1	12	9	3	0	1

- Molecule 7 is 2-AMINOHEXANEDIOIC ACID (three-letter code: UN1) (formula: $C_6H_{11}NO_4$).



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

- Molecule 8 is 2-OXOADIPIC ACID (three-letter code: OOG) (formula: $C_6H_8O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
8	B	1	11	6	5	0	1

- Molecule 9 is water.

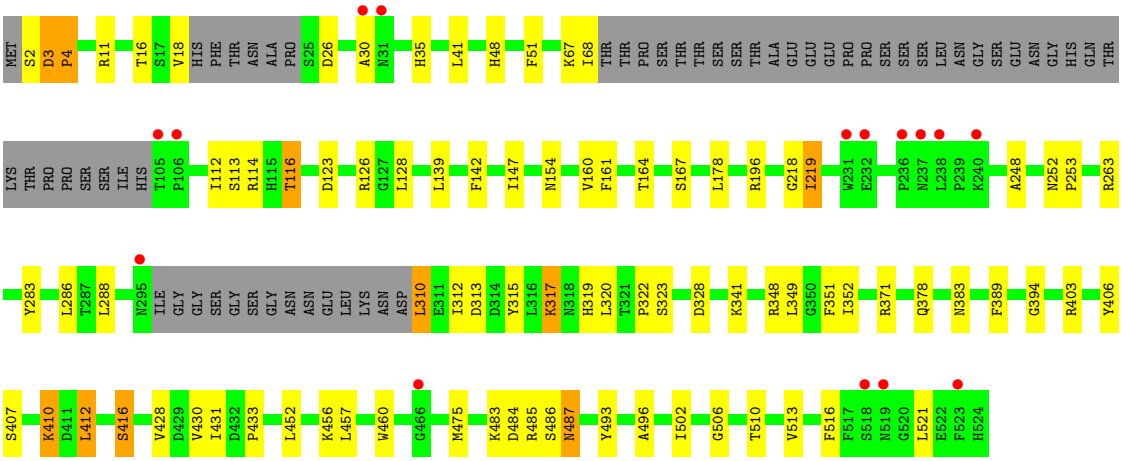
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
9	A	89	89	89	0	0
			Total	O		
9	B	106	106	106	0	0

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.

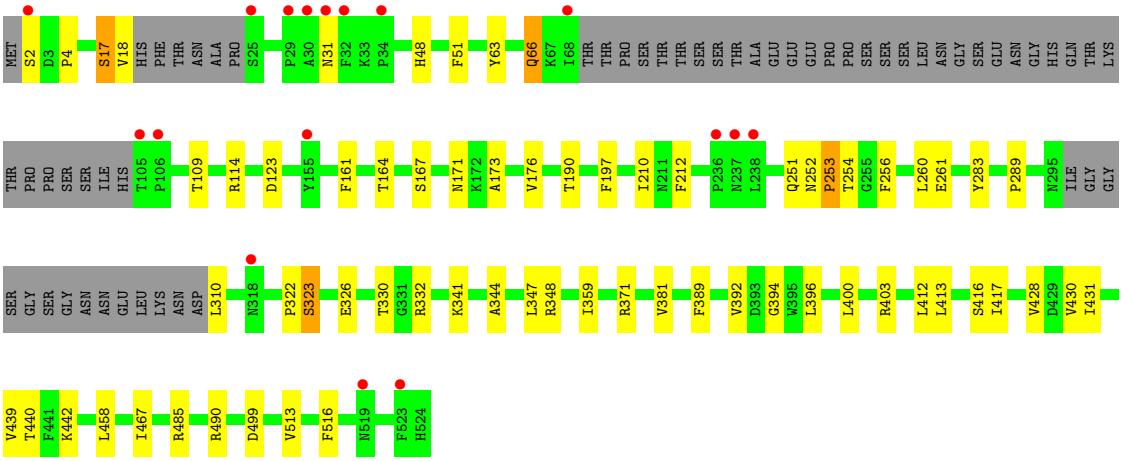
• Molecule 1: Aromatic-amino-acid:2-oxoglutarate transaminase

Chain A:



• Molecule 1: Aromatic-amino-acid:2-oxoglutarate transaminase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	74.57Å 88.54Å 160.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.53 – 2.60 46.53 – 2.60	Depositor EDS
% Data completeness (in resolution range)	94.0 (46.53-2.60) 94.0 (46.53-2.60)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.93 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.8.0232	Depositor
R, R_{free}	0.199 , 0.280 0.204 , 0.279	Depositor DCC
R_{free} test set	948 reflections (3.00%)	wwPDB-VP
Wilson B-factor (Å^2)	37.5	Xtriage
Anisotropy	0.397	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 47.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	7849	wwPDB-VP
Average B, all atoms (Å^2)	50.0	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 28.38 % 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 1.8628e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: OOG, PPY, UN1, PLP, BCN, MPD

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	0.71	0/3861	0.86	0/5240
1	B	0.71	0/3872	0.87	0/5254
All	All	0.71	0/7733	0.87	0/10494

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 (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3765	0	3715	55	0
1	B	3772	0	3724	43	0
2	A	15	0	6	0	0
2	B	15	0	6	0	0
3	A	11	0	12	1	0
3	B	22	0	24	1	0
4	A	8	0	14	0	0
5	A	12	0	8	1	0
6	A	12	0	7	2	0
7	B	11	0	4	1	0
8	B	11	0	6	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	A	89	0	0	3	0
9	B	106	0	0	1	0
All	All	7849	0	7526	93	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 6.

All (93) 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:323:SER:N	1:B:326:GLU:OE1	2.26	0.68
1:A:167:SER:HB3	1:B:371:ARG:HD3	1.75	0.67
1:B:212:PHE:O	1:B:442:LYS:NZ	2.28	0.67
1:A:428:VAL:CG1	1:A:513:VAL:HG11	2.25	0.66
1:B:413:LEU:O	1:B:416:SER:HB3	1.96	0.65
1:B:389:PHE:O	1:B:394:GLY:HA3	1.96	0.64
1:B:347:LEU:HD13	1:B:381:VAL:HG21	1.79	0.64
1:B:256:PHE:CE1	1:B:431:ILE:HG21	2.34	0.63
1:A:506:GLY:O	1:A:510:THR:HG23	1.99	0.62
1:A:160:VAL:HA	1:A:352:ILE:O	2.00	0.61
1:A:116:THR:HG22	1:A:123:ASP:HB3	1.83	0.60
1:A:487:ASN:HD22	1:A:487:ASN:H	1.49	0.60
1:A:112:ILE:HG12	1:A:383:ASN:HD22	1.67	0.59
1:B:210:ILE:HG21	1:B:254:THR:HG21	1.85	0.59
1:A:371:ARG:HH21	6:A:605[B]:PPY:C3	2.15	0.59
1:B:371:ARG:HH21	7:B:602[A]:UN1:CA	2.16	0.58
1:A:167:SER:HB3	1:B:371:ARG:CD	2.35	0.57
1:A:428:VAL:HG12	1:A:513:VAL:HG11	1.86	0.57
1:A:41:LEU:HG	1:B:467:ILE:HD11	1.87	0.57
5:A:604[A]:PHE:N	9:A:701:HOH:O	0.72	0.57
1:B:430:VAL:HG13	1:B:439:VAL:CG1	2.34	0.57
1:A:313:ASP:O	1:A:317:LYS:HB2	2.05	0.56
1:A:167:SER:CB	1:B:371:ARG:HD3	2.35	0.56
1:B:63:TYR:CZ	1:B:66:GLN:HB2	2.41	0.55
1:A:286:LEU:HD13	1:A:406:TYR:CD2	2.42	0.55
1:B:254:THR:O	1:B:440:THR:OG1	2.20	0.55
1:A:389:PHE:O	1:A:394:GLY:HA3	2.08	0.53
1:A:483:LYS:HG2	1:A:484:ASP:N	2.24	0.53
1:B:289:PRO:HB3	9:B:797:HOH:O	2.08	0.53
1:A:48:HIS:HA	1:A:51:PHE:CE2	2.45	0.52
1:A:496:ALA:HB3	1:A:502:ILE:HD11	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:412:LEU:HD21	1:B:499:ASP:HB3	1.92	0.51
1:A:18:VAL:HG12	1:A:18:VAL:O	2.10	0.51
1:A:35:HIS:NE2	9:A:702:HOH:O	2.26	0.51
1:B:171:ASN:HA	1:B:197:PHE:HB3	1.92	0.51
1:A:431:ILE:O	1:A:433:PRO:HD3	2.11	0.50
1:B:428:VAL:CG1	1:B:513:VAL:HG11	2.42	0.49
1:A:315:TYR:CZ	1:A:320:LEU:HD13	2.48	0.49
1:B:114:ARG:O	1:B:123:ASP:HB2	2.11	0.49
1:B:260:LEU:HD13	1:B:322:PRO:HG2	1.95	0.48
1:B:253:PRO:CB	1:B:490:ARG:HD3	2.44	0.48
1:B:253:PRO:HB3	1:B:490:ARG:HD3	1.95	0.48
1:A:371:ARG:HH21	6:A:605[B]:PPY:H31	1.78	0.48
1:A:263:ARG:NH1	1:A:322:PRO:O	2.47	0.47
3:A:602:BCN:H62	3:A:602:BCN:C2	2.43	0.47
1:A:310:LEU:HD12	1:A:403:ARG:NH1	2.30	0.47
1:B:371:ARG:HH21	8:B:605[B]:OOG:C2	2.28	0.47
1:A:248:ALA:HB3	9:A:740:HOH:O	2.15	0.46
1:B:417:ILE:HG21	1:B:430:VAL:HG21	1.97	0.46
1:B:252:ASN:HA	1:B:253:PRO:HA	1.66	0.46
1:A:252:ASN:HA	1:A:253:PRO:HA	1.71	0.46
1:A:139:LEU:HD21	1:A:349:LEU:HD21	1.99	0.45
1:A:457:LEU:HD12	1:A:457:LEU:HA	1.88	0.45
1:B:458:LEU:HD13	1:B:516:PHE:CG	2.51	0.45
1:A:416:SER:OG	1:A:506:GLY:HA3	2.17	0.45
1:A:196:ARG:HH11	1:B:371:ARG:HH12	1.65	0.44
1:A:288:LEU:HD22	1:A:407:SER:HA	1.99	0.44
1:B:256:PHE:CE1	1:B:431:ILE:CG2	3.00	0.44
1:B:289:PRO:HD3	1:B:403:ARG:NH2	2.32	0.44
1:A:11:ARG:NH2	1:A:178:LEU:O	2.50	0.44
1:A:218:GLY:O	1:A:219:ILE:C	2.55	0.44
1:A:315:TYR:HA	1:A:319:HIS:HB2	2.00	0.44
1:A:487:ASN:H	1:A:487:ASN:ND2	2.14	0.43
1:B:396:LEU:O	1:B:400:LEU:HG	2.18	0.43
1:A:154:ASN:HB3	1:A:328:ASP:O	2.18	0.43
3:B:603:BCN:O21	3:B:603:BCN:H51	2.19	0.43
1:B:254:THR:HA	1:B:440:THR:OG1	2.19	0.43
1:B:164:THR:HA	1:B:348:ARG:O	2.19	0.43
1:A:164:THR:HA	1:A:348:ARG:O	2.18	0.43
1:B:161:PHE:HB2	1:B:359:ILE:HG23	2.01	0.43
1:B:17:SER:O	1:B:18:VAL:C	2.57	0.42
1:B:417:ILE:HG21	1:B:430:VAL:CG2	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:ARG:O	1:A:123:ASP:HB2	2.20	0.42
1:A:516:PHE:CE2	1:A:521:LEU:HA	2.55	0.42
1:B:63:TYR:CE1	1:B:66:GLN:CB	3.02	0.42
1:A:483:LYS:O	1:A:486:SER:HB2	2.20	0.42
1:A:452:LEU:O	1:A:456:LYS:HG3	2.19	0.42
1:A:123:ASP:OD2	1:A:126:ARG:NH2	2.45	0.41
1:A:253:PRO:HG3	1:A:475:MET:HE3	2.02	0.41
1:A:128:LEU:O	1:B:344:ALA:HB1	2.20	0.41
1:A:67:LYS:O	1:A:68:ILE:C	2.58	0.41
1:A:457:LEU:O	1:A:460:TRP:HB2	2.21	0.41
1:A:371:ARG:HD3	1:B:167:SER:CB	2.50	0.41
1:A:410:LYS:HB2	1:A:493:TYR:CZ	2.55	0.41
1:B:2:SER:HB3	1:B:330:THR:HG22	2.02	0.41
1:A:161:PHE:O	1:A:351:PHE:HA	2.21	0.41
1:A:371:ARG:CD	1:B:167:SER:HB3	2.51	0.41
1:B:48:HIS:HA	1:B:51:PHE:CE2	2.56	0.41
1:B:173:ALA:O	1:B:176:VAL:HG12	2.22	0.40
1:A:2:SER:OG	1:A:3:ASP:N	2.54	0.40
1:A:139:LEU:HD23	1:A:378:GLN:HG2	2.02	0.40
1:A:412:LEU:HD12	1:A:412:LEU:HA	1.92	0.40
1:A:3:ASP:OD1	1:A:3:ASP:C	2.60	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	459/524 (88%)	422 (92%)	33 (7%)	4 (1%)	17 35
1	B	460/524 (88%)	433 (94%)	24 (5%)	3 (1%)	22 43
All	All	919/1048 (88%)	855 (93%)	57 (6%)	7 (1%)	19 39

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	30	ALA
1	B	251	GLN
1	A	4	PRO
1	A	219	ILE
1	A	312	ILE
1	B	392	VAL
1	B	4	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	413/462 (89%)	394 (95%)	19 (5%)	27	51
1	B	414/462 (90%)	401 (97%)	13 (3%)	40	66
All	All	827/924 (90%)	795 (96%)	32 (4%)	31	58

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	ASP
1	A	4	PRO
1	A	16	THR
1	A	26	ASP
1	A	113	SER
1	A	116	THR
1	A	142	PHE
1	A	147	ILE
1	A	283	TYR
1	A	310	LEU
1	A	317	LYS
1	A	323	SER
1	A	341	LYS
1	A	410	LYS
1	A	412	LEU
1	A	416	SER

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Mol	Chain	Res	Type
1	A	430	VAL
1	A	485	ARG
1	A	487	ASN
1	B	17	SER
1	B	31	ASN
1	B	66	GLN
1	B	109	THR
1	B	190	THR
1	B	253	PRO
1	B	261	GLU
1	B	283	TYR
1	B	310	LEU
1	B	323	SER
1	B	332	ARG
1	B	341	LYS
1	B	485	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	487	ASN
1	B	524	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

10 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
5	PHE	A	604[A]	-	11,12,12	0.66	0	14,15,15	1.17	2 (14%)
2	PLP	A	601	1	15,15,16	0.59	0	20,22,23	1.02	2 (10%)
7	UN1	B	602[A]	-	9,10,10	0.94	0	11,12,12	1.16	0
8	OOG	B	605[B]	-	10,10,10	5.44	2 (20%)	12,12,12	1.54	2 (16%)
2	PLP	B	601	1	15,15,16	0.77	0	20,22,23	0.91	0
6	PPY	A	605[B]	-	12,12,12	4.77	3 (25%)	15,15,15	1.88	4 (26%)
3	BCN	B	603	-	10,10,10	0.66	0	11,11,11	1.00	1 (9%)
3	BCN	A	602	-	10,10,10	0.83	0	11,11,11	1.00	0
4	MPD	A	603	-	7,7,7	0.16	0	9,10,10	0.50	0
3	BCN	B	604	-	10,10,10	0.77	0	11,11,11	0.69	0

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
5	PHE	A	604[A]	-	-	2/8/8/8	0/1/1/1
2	PLP	A	601	1	-	0/6/6/8	0/1/1/1
7	UN1	B	602[A]	-	-	5/10/10/10	-
8	OOG	B	605[B]	-	-	7/10/10/10	-
2	PLP	B	601	1	-	1/6/6/8	0/1/1/1
6	PPY	A	605[B]	-	-	3/8/8/8	0/1/1/1
3	BCN	B	603	-	-	4/10/10/10	-
3	BCN	A	602	-	-	4/10/10/10	-
4	MPD	A	603	-	-	1/5/5/5	-
3	BCN	B	604	-	-	4/10/10/10	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	B	605[B]	OOG	C5-C6	-16.92	1.30	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	605[B]	PPY	C2-C1	-16.23	1.31	1.53
8	B	605[B]	OOG	O4-C6	-2.22	1.24	1.30
6	A	605[B]	PPY	O2-C1	-2.15	1.24	1.30
6	A	605[B]	PPY	C3-C2	2.10	1.53	1.51

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	A	605[B]	PPY	C1'-C3-C2	4.67	119.70	113.64
6	A	605[B]	PPY	C3-C2-C1	3.65	124.01	117.85
8	B	605[B]	OOG	C4-C5-C6	3.58	122.62	115.97
5	A	604[A]	PHE	OXT-C-O	-3.18	116.87	124.09
6	A	605[B]	PPY	O1-C1-C2	-2.98	117.74	121.72
5	A	604[A]	PHE	OXT-C-CA	2.71	122.62	113.38
8	B	605[B]	OOG	O3-C6-C5	-2.34	118.59	121.72
3	B	603	BCN	C5-N1-C3	2.31	116.94	111.44
2	A	601	PLP	C4A-C4-C5	2.20	123.20	120.94
2	A	601	PLP	O3P-P-O2P	2.17	115.92	107.64
6	A	605[B]	PPY	O2-C1-C2	2.12	119.77	113.97

There are no chirality outliers.

All (31) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	603	BCN	C4-C3-N1-C5
3	B	603	BCN	N1-C1-C2-O21
3	B	603	BCN	N1-C1-C2-O22
4	A	603	MPD	C2-C3-C4-O4
5	A	604[A]	PHE	N-CA-CB-CG
5	A	604[A]	PHE	C-CA-CB-CG
6	A	605[B]	PPY	O2-C1-C2-C3
7	B	602[A]	UN1	C6-C5-CA-N
8	B	605[B]	OOG	C4-C5-C6-O4
3	A	602	BCN	C6-C5-N1-C1
3	B	604	BCN	N1-C5-C6-O6
8	B	605[B]	OOG	C1-C2-C3-C4
3	B	603	BCN	N1-C5-C6-O6
3	A	602	BCN	C2-C1-N1-C5
7	B	602[A]	UN1	CA-C5-C6-C1
3	A	602	BCN	N1-C5-C6-O6
3	B	604	BCN	N1-C3-C4-O4
7	B	602[A]	UN1	C6-C5-CA-C

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Mol	Chain	Res	Type	Atoms
3	A	602	BCN	C4-C3-N1-C1
8	B	605[B]	OOG	C3-C4-C5-C6
8	B	605[B]	OOG	O5-C5-C6-O3
6	A	605[B]	PPY	O1-C1-C2-C3
8	B	605[B]	OOG	C4-C5-C6-O3
3	B	604	BCN	C4-C3-N1-C5
2	B	601	PLP	C5A-O4P-P-O1P
6	A	605[B]	PPY	O2-C1-C2-O3
8	B	605[B]	OOG	O5-C5-C6-O4
8	B	605[B]	OOG	C2-C3-C4-C5
3	B	604	BCN	C2-C1-N1-C3
7	B	602[A]	UN1	C6-C1-C1'-O2'
7	B	602[A]	UN1	C6-C1-C1'-O1'

There are no ring outliers.

6 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	604[A]	PHE	1	0
7	B	602[A]	UN1	1	0
8	B	605[B]	OOG	1	0
6	A	605[B]	PPY	2	0
3	B	603	BCN	1	0
3	A	602	BCN	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 (i)

6.1 Protein, DNA and RNA chains (i)

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	467/524 (89%)	0.01	15 (3%)	47	40	30, 47, 72, 100	0
1	B	467/524 (89%)	0.03	17 (3%)	42	35	30, 48, 73, 127	0
All	All	934/1048 (89%)	0.02	32 (3%)	45	38	30, 47, 73, 127	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	236	PRO	5.5
1	B	236	PRO	4.4
1	B	237	ASN	4.2
1	B	105	THR	4.1
1	A	30	ALA	3.9
1	A	237	ASN	3.9
1	B	155	TYR	3.8
1	B	32	PHE	3.8
1	A	31	ASN	3.7
1	A	519	ASN	3.6
1	B	68	ILE	3.5
1	B	29	PRO	3.4
1	A	105	THR	3.3
1	A	518	SER	3.3
1	B	2	SER	3.3
1	A	295	ASN	3.3
1	B	30	ALA	3.3
1	A	232	GLU	3.2
1	B	25	SER	3.2
1	B	523	PHE	3.0
1	B	31	ASN	3.0
1	A	106	PRO	2.7
1	B	318	ASN	2.5
1	B	238	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	523	PHE	2.4
1	B	519	ASN	2.3
1	A	466	GLY	2.2
1	B	106	PRO	2.1
1	A	231	TRP	2.1
1	A	240	LYS	2.1
1	A	238	LEU	2.0
1	B	34	PRO	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 monosaccharides 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
3	BCN	A	602	11/11	0.61	0.27	80,94,106,107	0
5	PHE	A	604[A]	12/12	0.81	0.39	74,78,79,79	12
6	PPY	A	605[B]	12/12	0.82	0.35	70,75,77,77	12
3	BCN	B	604	11/11	0.84	0.34	85,96,100,102	0
4	MPD	A	603	8/8	0.84	0.28	83,91,104,109	0
3	BCN	B	603	11/11	0.87	0.17	68,81,96,97	0
7	UN1	B	602[A]	11/11	0.89	0.40	57,60,64,65	11
8	OOG	B	605[B]	11/11	0.89	0.39	47,57,66,66	11
2	PLP	A	601	15/16	0.98	0.18	37,42,49,52	0
2	PLP	B	601	15/16	0.98	0.17	39,46,50,54	0

6.5 Other polymers (i)

There are no such residues in this entry.