



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:32 PM GMT

PDB ID : 1PII
Title : THREE-DIMENSIONAL STRUCTURE OF THE BIFUNCTIONAL ENZYME PHOSPHORIBOSYLANTHRANILATE ISOMERASE: INDOLE GLYCEROLPHOSPHATE SYNTHASE FROM ESCHERICHIA COLI REFINED AT 2.0 ANGSTROMS RESOLUTION
Authors : Wilmanns, M.; Priestle, J.P.; Jansonius, J.N.
Deposited on : 1991-06-21
Resolution : 2.00 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

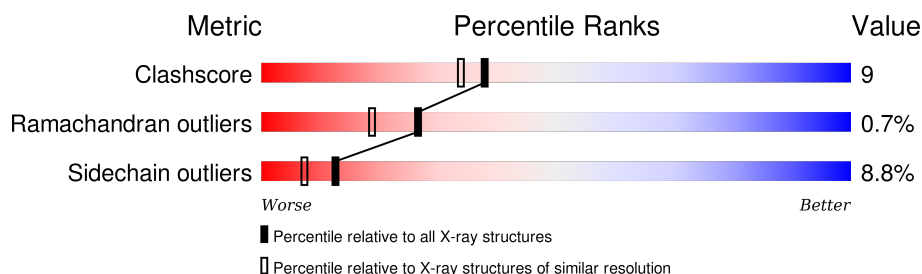
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)

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	452	

2 Entry composition [i](#)

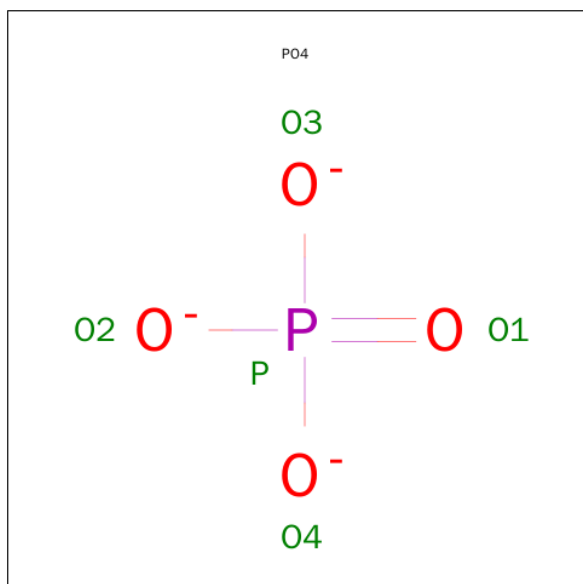
There are 3 unique types of molecules in this entry. The entry contains 4162 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 N-(5'PHOSPHORIBOSYL)ANTHRANILATE ISOMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	452	Total	C	N	O	S	0	5	0
			3524	2225	621	665	13			

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is water.

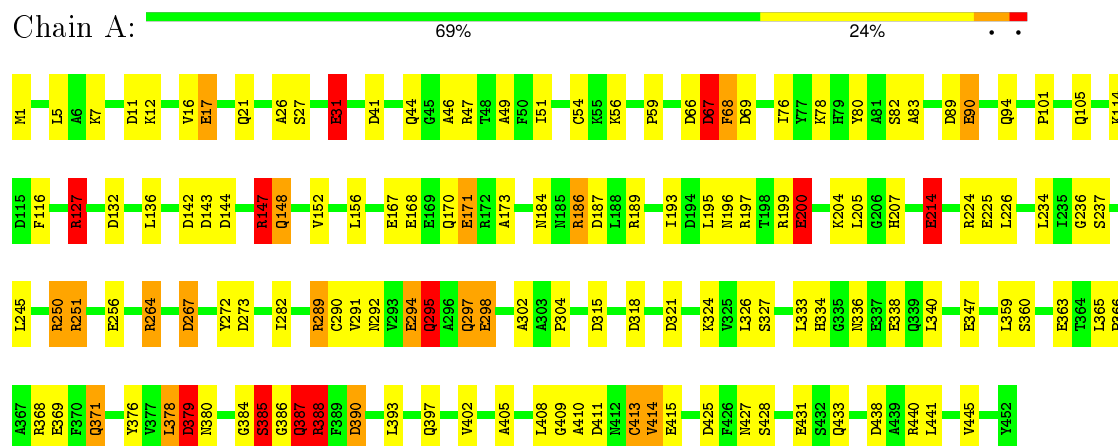
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	628	Total	O	8	0
			628	628		

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: N-(5'PHOSPHORIBOSYL)ANTHRANILATE ISOMERASE



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	104.70 Å 104.70 Å 68.00 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) (15.00-2.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT, X-PLOR	Depositor
R, R_{free}	0.173 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4162	wwPDB-VP
Average B, all atoms (Å ²)	24.0	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: PO4

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	0.86	15/3587 (0.4%)	1.74	84/4863 (1.7%)

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

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	17	GLU	CD-OE2	7.33	1.33	1.25
1	A	347	GLU	CD-OE1	6.48	1.32	1.25
1	A	363	GLU	CD-OE2	5.97	1.32	1.25
1	A	294	GLU	CD-OE1	5.83	1.32	1.25
1	A	171	GLU	CD-OE2	5.77	1.31	1.25
1	A	168	GLU	CD-OE2	5.67	1.31	1.25
1	A	200	GLU	CD-OE2	5.60	1.31	1.25
1	A	298	GLU	CD-OE1	5.59	1.31	1.25
1	A	369	GLU	CD-OE1	5.53	1.31	1.25
1	A	214	GLU	CD-OE1	5.36	1.31	1.25
1	A	415	GLU	CD-OE2	5.30	1.31	1.25
1	A	90	GLU	CD-OE2	5.26	1.31	1.25
1	A	31	GLU	CD-OE1	5.21	1.31	1.25
1	A	167	GLU	CD-OE1	5.18	1.31	1.25
1	A	338	GLU	CD-OE2	5.02	1.31	1.25

All (84) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	289	ARG	NE-CZ-NH1	12.64	126.62	120.30
1	A	186	ARG	NE-CZ-NH1	11.97	126.29	120.30
1	A	440	ARG	NE-CZ-NH1	11.76	126.18	120.30
1	A	315	ASP	CB-CG-OD2	-11.57	107.88	118.30
1	A	315	ASP	CB-CG-OD1	10.52	127.76	118.30
1	A	365	LEU	CB-CA-C	-9.82	91.55	110.20
1	A	199	ARG	NE-CZ-NH1	-9.23	115.69	120.30
1	A	66	ASP	C-N-CA	9.16	144.60	121.70
1	A	388	ARG	N-CA-C	9.16	135.72	111.00
1	A	318	ASP	CB-CG-OD2	8.82	126.24	118.30
1	A	387	GLN	N-CA-C	8.64	134.32	111.00
1	A	386	GLY	C-N-CA	8.47	142.88	121.70
1	A	189	ARG	NE-CZ-NH1	8.41	124.51	120.30
1	A	132	ASP	CB-CG-OD2	8.33	125.79	118.30
1	A	411	ASP	CB-CG-OD1	8.21	125.69	118.30
1	A	273	ASP	CB-CG-OD1	8.12	125.61	118.30
1	A	380	ASN	N-CA-C	-8.05	89.26	111.00
1	A	187	ASP	CB-CG-OD1	7.84	125.35	118.30
1	A	385	SER	N-CA-C	-7.64	90.37	111.00
1	A	11	ASP	CB-CG-OD2	-7.63	111.44	118.30
1	A	187	ASP	CB-CG-OD2	-7.42	111.62	118.30
1	A	289	ARG	CD-NE-CZ	7.27	133.78	123.60
1	A	318	ASP	CB-CG-OD1	-7.22	111.80	118.30
1	A	273	ASP	CB-CG-OD2	-7.18	111.84	118.30
1	A	440	ARG	NE-CZ-NH2	-7.09	116.75	120.30
1	A	387	GLN	C-N-CA	7.05	139.32	121.70
1	A	142	ASP	CB-CG-OD1	6.97	124.58	118.30
1	A	384	GLY	N-CA-C	-6.77	96.17	113.10
1	A	390	ASP	CB-CG-OD1	6.71	124.33	118.30
1	A	390	ASP	CB-CG-OD2	-6.69	112.28	118.30
1	A	289	ARG	NE-CZ-NH2	-6.65	116.97	120.30
1	A	379	ASP	CB-CG-OD2	-6.65	112.32	118.30
1	A	67	ASP	CB-CA-C	6.62	123.64	110.40
1	A	250	ARG	NE-CZ-NH1	6.53	123.56	120.30
1	A	363	GLU	N-CA-CB	6.43	122.18	110.60
1	A	411	ASP	CB-CG-OD2	-6.26	112.67	118.30
1	A	173	ALA	N-CA-CB	6.24	118.83	110.10
1	A	267	ASP	CB-CG-OD2	-6.22	112.70	118.30
1	A	371	GLN	CA-CB-CG	6.22	127.08	113.40
1	A	388	ARG	NE-CZ-NH2	-6.18	117.21	120.30
1	A	127	ARG	CD-NE-CZ	-6.09	115.07	123.60
1	A	295	GLN	CB-CA-C	6.04	122.47	110.40
1	A	11	ASP	CB-CG-OD1	6.03	123.73	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	68	PHE	N-CA-C	6.03	127.27	111.00
1	A	143	ASP	CB-CG-OD1	-6.02	112.88	118.30
1	A	142	ASP	CB-CG-OD2	-5.99	112.91	118.30
1	A	425	ASP	CB-CG-OD1	-5.96	112.94	118.30
1	A	205	LEU	C-N-CA	-5.95	109.80	122.30
1	A	80	TYR	CB-CA-C	-5.95	98.50	110.40
1	A	59	PRO	C-N-CA	-5.92	106.90	121.70
1	A	360	SER	CB-CA-C	5.88	121.28	110.10
1	A	297	GLN	CB-CG-CD	5.81	126.70	111.60
1	A	47	ARG	NE-CZ-NH1	5.78	123.19	120.30
1	A	69	ASP	CB-CG-OD2	5.78	123.50	118.30
1	A	132	ASP	CB-CG-OD1	-5.71	113.16	118.30
1	A	47	ARG	CA-CB-CG	5.63	125.79	113.40
1	A	144	ASP	CB-CG-OD1	-5.63	113.23	118.30
1	A	387	GLN	N-CA-CB	-5.62	100.49	110.60
1	A	413	CYS	N-CA-CB	-5.62	100.48	110.60
1	A	388	ARG	NE-CZ-NH1	5.57	123.08	120.30
1	A	41	ASP	CB-CG-OD1	-5.55	113.30	118.30
1	A	267	ASP	CB-CG-OD1	5.54	123.28	118.30
1	A	89	ASP	CB-CG-OD1	5.54	123.28	118.30
1	A	292	ASN	N-CA-CB	5.44	120.39	110.60
1	A	49	ALA	CB-CA-C	-5.40	102.00	110.10
1	A	186	ARG	NE-CZ-NH2	-5.40	117.60	120.30
1	A	207	HIS	N-CA-CB	5.39	120.31	110.60
1	A	440	ARG	CD-NE-CZ	5.32	131.04	123.60
1	A	47	ARG	CB-CA-C	-5.30	99.80	110.40
1	A	66	ASP	CB-CG-OD2	-5.29	113.53	118.30
1	A	264	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	A	147	ARG	CD-NE-CZ	5.24	130.94	123.60
1	A	46	ALA	C-N-CA	5.23	134.78	121.70
1	A	16	VAL	CA-CB-CG2	-5.20	103.09	110.90
1	A	69	ASP	CB-CG-OD1	-5.18	113.64	118.30
1	A	438	ASP	CB-CG-OD2	-5.14	113.67	118.30
1	A	297	GLN	CB-CA-C	5.12	120.64	110.40
1	A	251	ARG	NE-CZ-NH2	5.09	122.84	120.30
1	A	438	ASP	CB-CG-OD1	5.08	122.88	118.30
1	A	386	GLY	CA-C-N	-5.07	106.05	117.20
1	A	224	ARG	NE-CZ-NH2	5.05	122.82	120.30
1	A	321	ASP	CB-CG-OD1	5.01	122.81	118.30
1	A	148[A]	GLN	N-CA-CB	-5.00	101.60	110.60
1	A	148[B]	GLN	N-CA-CB	-5.00	101.60	110.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	197[A]	ARG	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	3524	0	3503	65	0
2	A	10	0	0	2	0
3	A	628	0	0	21	0
All	All	4162	0	3503	65	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 (65) 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:186:ARG:NH2	1:A:193:ILE:HD11	1.76	1.01
1:A:147:ARG:HG2	1:A:147:ARG:HH11	1.44	0.82
1:A:101:PRO:O	1:A:105:GLN:HG2	1.84	0.78
1:A:414:VAL:HG23	3:A:953:HOH:O	1.88	0.74
1:A:186:ARG:CZ	1:A:193:ILE:HD11	2.19	0.71
1:A:295:GLN:HG2	3:A:663:HOH:O	1.93	0.67
1:A:31:GLU:HG3	3:A:908:HOH:O	1.92	0.67
1:A:289:ARG:NH1	1:A:431:GLU:OE1	2.31	0.64
1:A:184:ASN:HA	1:A:214:GLU:HG3	1.81	0.62
1:A:27:SER:HB3	3:A:646:HOH:O	1.99	0.62
1:A:76:ILE:HD13	1:A:245:LEU:HB3	1.81	0.61
1:A:67:ASP:HA	3:A:834:HOH:O	2.00	0.61
1:A:237:SER:OG	2:A:453:PO4:O2	2.11	0.60
1:A:304:PRO:HB2	3:A:562:HOH:O	1.99	0.60
1:A:147:ARG:HG2	1:A:147:ARG:NH1	2.15	0.58
1:A:371:GLN:OE1	3:A:467:HOH:O	2.18	0.56
1:A:359:LEU:HB3	1:A:366:PRO:HG3	1.86	0.56
1:A:378:LEU:HD22	1:A:402:VAL:HG11	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:ARG:HD3	1:A:267:ASP:OD2	2.05	0.56
1:A:379:ASP:HB3	1:A:405:ALA:HB3	1.89	0.55
1:A:371:GLN:HG3	3:A:1042:HOH:O	2.06	0.55
1:A:368:ARG:NH2	1:A:397:GLN:HB3	2.22	0.55
1:A:433:GLN:NE2	3:A:764:HOH:O	2.40	0.54
1:A:379:ASP:CB	1:A:405:ALA:HB3	2.37	0.54
1:A:171:GLU:HG2	3:A:784:HOH:O	2.08	0.54
1:A:148[B]:GLN:O	1:A:152:VAL:HG23	2.08	0.54
1:A:186:ARG:HH22	1:A:193:ILE:HD11	1.70	0.52
1:A:433:GLN:HG2	3:A:967:HOH:O	2.10	0.52
1:A:148[A]:GLN:O	1:A:152:VAL:HG23	2.10	0.52
1:A:385:SER:OG	3:A:605:HOH:O	2.19	0.51
1:A:225:GLU:OE1	3:A:716:HOH:O	2.19	0.51
1:A:410:ALA:HA	1:A:445:VAL:HG12	1.93	0.51
1:A:427:ASN:HB3	2:A:454:PO4:O2	2.11	0.50
1:A:105:GLN:HG3	3:A:555:HOH:O	2.12	0.50
1:A:26:ALA:HA	3:A:922:HOH:O	2.13	0.48
1:A:409:GLY:O	1:A:413:CYS:HB2	2.13	0.48
1:A:272:TYR:HE1	3:A:693:HOH:O	1.95	0.48
1:A:56:LYS:HE3	1:A:94:GLN:OE1	2.14	0.48
1:A:56:LYS:HE2	3:A:715:HOH:O	2.13	0.47
1:A:368:ARG:HB3	1:A:376:TYR:CZ	2.51	0.46
1:A:390:ASP:O	1:A:393:LEU:HB2	2.16	0.46
1:A:250:ARG:NH1	1:A:272:TYR:O	2.48	0.46
1:A:196:ASN:O	1:A:200:GLU:HG2	2.16	0.46
1:A:368:ARG:HH22	1:A:397:GLN:HB3	1.80	0.45
1:A:127:ARG:HG3	1:A:156:LEU:HD13	1.98	0.45
1:A:251:ARG:NH1	1:A:256:GLU:OE1	2.50	0.45
1:A:291:VAL:HG12	1:A:295:GLN:CG	2.46	0.45
1:A:428:SER:O	1:A:431:GLU:HB2	2.17	0.45
1:A:326:LEU:HD23	1:A:326:LEU:HA	1.74	0.44
1:A:12:LYS:HD2	1:A:116:PHE:HB2	2.00	0.43
1:A:136:LEU:HD23	1:A:136:LEU:HA	1.69	0.43
1:A:282:ILE:O	1:A:290:CYS:HA	2.18	0.43
1:A:245:LEU:HD12	1:A:245:LEU:HA	1.77	0.43
1:A:90:GLU:O	1:A:94:GLN:HA	2.18	0.43
1:A:195:LEU:HD13	1:A:226:LEU:HD11	2.01	0.43
1:A:234:LEU:HD23	1:A:234:LEU:C	2.39	0.42
1:A:387:GLN:HG2	1:A:387:GLN:H	1.52	0.42
1:A:82:SER:CB	3:A:825:HOH:O	2.68	0.42
1:A:359:LEU:HD13	1:A:366:PRO:HG2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:ILE:HG12	1:A:83:ALA:HB3	2.02	0.42
1:A:94:GLN:OE1	3:A:715:HOH:O	2.22	0.42
1:A:302:ALA:HB1	3:A:1039:HOH:O	2.19	0.42
1:A:54[B]:CYS:SG	1:A:68:PHE:CZ	3.14	0.41
1:A:324[A]:LYS:NZ	3:A:609:HOH:O	2.54	0.40
1:A:408:LEU:HA	1:A:408:LEU:HD23	1.77	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	455/452 (101%)	440 (97%)	12 (3%)	3 (1%)	26 19

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	67	ASP
1	A	388	ARG
1	A	236	GLY

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	367/362 (101%)	335 (91%)	32 (9%)	13 7

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

Mol	Chain	Res	Type
1	A	1	MET
1	A	5	LEU
1	A	7	LYS
1	A	17	GLU
1	A	21	GLN
1	A	31	GLU
1	A	44	GLN
1	A	67	ASP
1	A	78	LYS
1	A	114	LYS
1	A	127	ARG
1	A	147	ARG
1	A	170	GLN
1	A	200	GLU
1	A	204	LYS
1	A	214	GLU
1	A	294	GLU
1	A	295	GLN
1	A	297	GLN
1	A	298	GLU
1	A	327	SER
1	A	333	LEU
1	A	334	HIS
1	A	336	ASN
1	A	340	LEU
1	A	378	LEU
1	A	379	ASP
1	A	385	SER
1	A	387	GLN
1	A	388	ARG
1	A	414	VAL
1	A	441	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	GLN

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Mol	Chain	Res	Type
1	A	29	GLN
1	A	44	GLN
1	A	295	GLN
1	A	372	HIS
1	A	387	GLN
1	A	418	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 ⓘ

2 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	$\# Z > 2$
2	PO4	A	453	-	4,4,4	2.40	3 (75%)	6,6,6	0.28	0
2	PO4	A	454	-	4,4,4	0.93	0	6,6,6	0.28	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
2	PO4	A	453	-	-	0/0/0/0	0/0/0/0
2	PO4	A	454	-	-	0/0/0/0	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	453	PO4	P-O2	2.11	1.61	1.53
2	A	453	PO4	P-O1	2.11	1.61	1.52
2	A	453	PO4	P-O3	3.58	1.66	1.53

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	453	PO4	1	0
2	A	454	PO4	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 ⓘ

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

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

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

6.3 Carbohydrates ⓘ

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

6.4 Ligands ⓘ

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

6.5 Other polymers ⓘ

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