



Full wwPDB X-ray Structure Validation Report

Oct 13, 2014 – 06:59 PM EDT

PDB ID : 4O13
Title : The crystal structure of NAMPT in complex with GNE-618
Authors : Oh, A.; Coons, M.; Brillantes, B.; Wang, W.
Deposited on : 2013-12-15
Resolution : 1.75 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

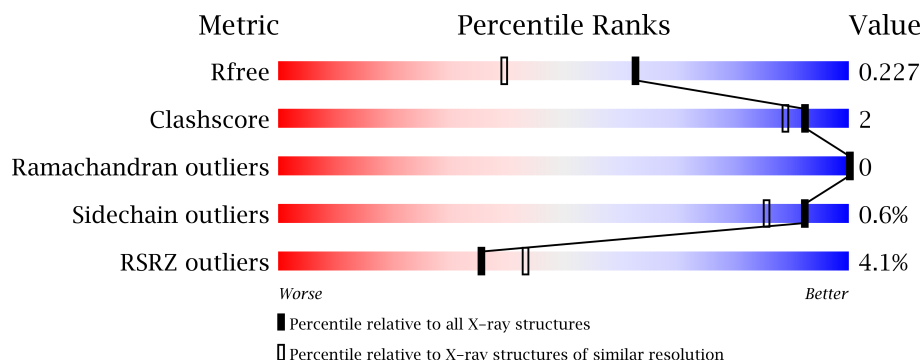
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.16 November 2013
Xtriage (Phenix) : dev-1439
EDS : stable23828
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable23828

1 Overall quality at a glance

The reported resolution of this entry is 1.75 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}	66092	1134 (1.76-1.76)
Clashscore	79885	1304 (1.76-1.76)
Ramachandran outliers	78287	1288 (1.76-1.76)
Sidechain outliers	78261	1288 (1.76-1.76)
RSRZ outliers	66119	1135 (1.76-1.76)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	501	
1	B	501	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	PO4	A	602	-	X
3	PO4	A	603	-	X
3	PO4	B	603	-	X
4	EDO	A	604	-	X
4	EDO	A	605	-	X
4	EDO	A	606	-	X

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Geometry	Electron density
4	EDO	B	604	-	X
4	EDO	B	605	-	X
4	EDO	B	606	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8573 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

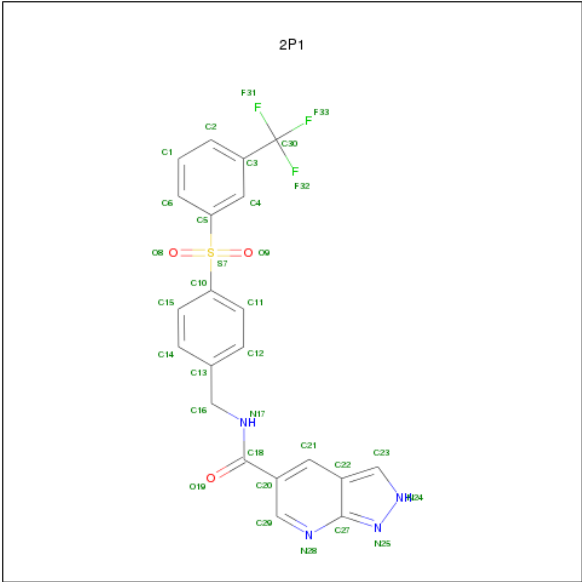
- Molecule 1 is a protein called Nicotinamide phosphoribosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	471	Total	C	N	O	S	0	1	0
			3773	2428	623	715	7			
1	B	469	Total	C	N	O	S	0	0	0
			3755	2416	620	712	7			

There are 20 discrepancies between the modelled and reference sequences:

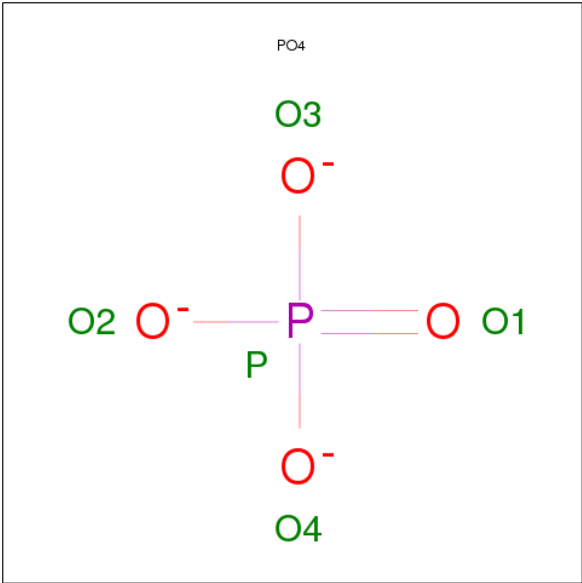
Chain	Residue	Modelled	Actual	Comment	Reference
A	492	LEU	-	EXPRESSION TAG	UNP P43490
A	493	GLU	-	EXPRESSION TAG	UNP P43490
A	494	HIS	-	EXPRESSION TAG	UNP P43490
A	495	HIS	-	EXPRESSION TAG	UNP P43490
A	496	HIS	-	EXPRESSION TAG	UNP P43490
A	497	HIS	-	EXPRESSION TAG	UNP P43490
A	498	HIS	-	EXPRESSION TAG	UNP P43490
A	499	HIS	-	EXPRESSION TAG	UNP P43490
A	500	HIS	-	EXPRESSION TAG	UNP P43490
A	501	HIS	-	EXPRESSION TAG	UNP P43490
B	492	LEU	-	EXPRESSION TAG	UNP P43490
B	493	GLU	-	EXPRESSION TAG	UNP P43490
B	494	HIS	-	EXPRESSION TAG	UNP P43490
B	495	HIS	-	EXPRESSION TAG	UNP P43490
B	496	HIS	-	EXPRESSION TAG	UNP P43490
B	497	HIS	-	EXPRESSION TAG	UNP P43490
B	498	HIS	-	EXPRESSION TAG	UNP P43490
B	499	HIS	-	EXPRESSION TAG	UNP P43490
B	500	HIS	-	EXPRESSION TAG	UNP P43490
B	501	HIS	-	EXPRESSION TAG	UNP P43490

- Molecule 2 is N-(4-{[3-(TRIFLUOROMETHYL)PHENYL]SULFONYL}BENZYL)-2H-PYRAZOLO[3,4-B]PYRIDINE-5-CARBOXAMIDE (three-letter code: 2P1) (formula: C₂₁H₁₅F₃N₄O₃S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	S	0	0
			32	21	3	4	3	1		
2	B	1	Total	C	F	N	O	S	0	0
			32	21	3	4	3	1		

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



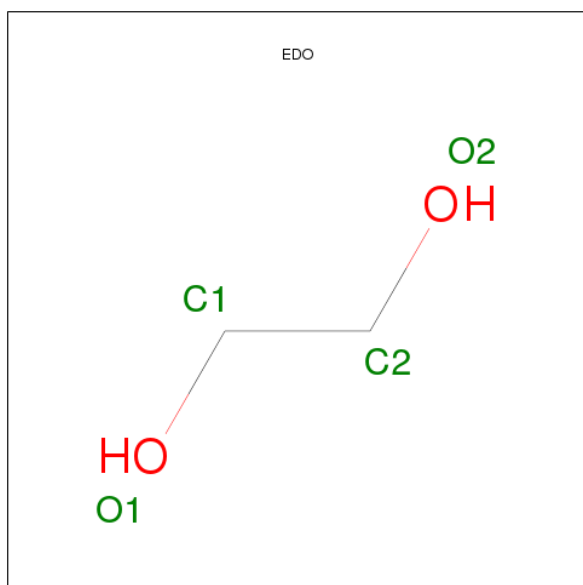
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	O P	0	0
			5	4 1		
3	A	1	Total	O P	0	0
			5	4 1		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	476	Total 476	O 476	0	0
5	B	453	Total 453	O 453	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	60.56Å 106.21Å 82.95Å 90.00° 96.39° 90.00°	Depositor
Resolution (Å)	34.86 – 1.75 46.22 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.1 (34.86-1.75) 99.1 (46.22-1.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 1.75Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.197 , 0.227 0.196 , 0.227	Depositor DCC
R_{free} test set	5226 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	16.5	Xtriage
Anisotropy	0.405	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 37.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 104764 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8573	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.01% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 2P1, PO4, EDO

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	0.37	0/3861	0.54	0/5231
1	B	0.36	0/3843	0.52	0/5206
All	All	0.37	0/7704	0.53	0/10437

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3773	0	3749	9	0
1	B	3755	0	3735	16	0
2	A	32	0	0	0	0
2	B	32	0	0	0	0
3	A	10	0	0	1	0
3	B	10	0	0	1	0
4	A	16	0	24	1	0
4	B	16	0	24	4	0
5	A	476	0	0	6	1
5	B	453	0	0	2	1

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	8573	0	7532	27	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 2.

All (27) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:434:ARG:HG2	1:B:457:LEU:HD11	1.72	0.69
3:A:603:PO4:O3	5:A:814:HOH:O	2.13	0.64
1:B:396:ASN:HA	4:B:605:EDO:H22	1.79	0.63
1:A:177:LEU:HD22	1:A:484:ILE:HD11	1.82	0.62
5:A:956:HOH:O	1:B:206:ILE:HG21	2.08	0.52
1:B:451:GLU:OE2	1:B:455:GLN:NE2	2.44	0.51
1:B:396:ASN:HD22	4:B:605:EDO:H22	1.75	0.51
1:B:396:ASN:HD22	4:B:605:EDO:C2	2.23	0.50
5:A:956:HOH:O	1:B:157:TYR:HD2	1.93	0.50
1:B:187:GLU:H	1:B:187:GLU:CD	2.15	0.49
1:A:172:LEU:HD21	1:A:361:LEU:HD11	1.95	0.48
1:B:336:GLU:OE1	5:B:897:HOH:O	2.19	0.48
3:B:602:PO4:O2	5:B:839:HOH:O	2.21	0.46
1:B:114:ILE:HD11	1:B:144:LEU:HG	1.97	0.46
1:B:32:LYS:HB3	1:B:405:VAL:HB	1.97	0.46
1:B:70:LEU:HD11	1:B:151:ILE:HD11	1.99	0.45
1:B:37:PHE:CZ	1:B:397:CYS:HB3	2.53	0.44
1:A:259:LYS:NZ	5:A:1152:HOH:O	2.50	0.44
1:B:354:ASP:OD2	4:B:606:EDO:H11	2.18	0.43
1:A:414:PHE:CE2	1:A:427:LYS:HE2	2.54	0.43
1:B:175:TYR:HB3	1:B:375:ILE:HG13	2.01	0.42
1:A:354:ASP:OD1	5:A:732:HOH:O	2.22	0.41
1:A:396:ASN:HD22	4:A:604:EDO:C2	2.33	0.41
1:B:331:LYS:HD3	1:B:331:LYS:HA	1.90	0.41
1:A:175:TYR:HB3	1:A:375:ILE:HG13	2.02	0.41
1:A:199:SER:O	5:A:956:HOH:O	2.22	0.41
1:A:66:LEU:HD23	1:A:70:LEU:HD12	2.01	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
5:A:800:HOH:O	5:B:952:HOH:O[1_554]	2.10	0.10

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	468/501 (93%)	458 (98%)	10 (2%)	0	100	100
1	B	465/501 (93%)	454 (98%)	11 (2%)	0	100	100
All	All	933/1002 (93%)	912 (98%)	21 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	414/440 (94%)	412 (100%)	2 (0%)	94	88
1	B	413/440 (94%)	410 (99%)	3 (1%)	91	82
All	All	827/880 (94%)	822 (99%)	5 (1%)	92	85

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	TYR
1	A	195	TYR
1	B	18	TYR
1	B	94	ASP
1	B	195	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

14 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	2P1	A	601	-	35,35,35	1.19	3 (8%)	49,52,52	1.84	10 (20%)
3	PO4	A	602	-	4,4,4	0.20	0	6,6,6	0.29	0
3	PO4	A	603	-	4,4,4	0.31	0	6,6,6	0.29	0
4	EDO	A	604	-	3,3,3	0.46	0	2,2,2	0.47	0
4	EDO	A	605	-	3,3,3	0.58	0	2,2,2	0.24	0
4	EDO	A	606	-	3,3,3	0.49	0	2,2,2	0.56	0
4	EDO	A	607	-	3,3,3	0.48	0	2,2,2	0.49	0
2	2P1	B	601	-	35,35,35	1.18	4 (11%)	49,52,52	1.80	9 (18%)
3	PO4	B	602	-	4,4,4	0.27	0	6,6,6	0.29	0
3	PO4	B	603	-	4,4,4	0.32	0	6,6,6	0.30	0
4	EDO	B	604	-	3,3,3	0.51	0	2,2,2	0.42	0
4	EDO	B	605	-	3,3,3	0.47	0	2,2,2	0.49	0
4	EDO	B	606	-	3,3,3	0.50	0	2,2,2	0.39	0
4	EDO	B	607	-	3,3,3	0.51	0	2,2,2	0.21	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	2P1	A	601	-	-	0/27/27/27	0/4/4/4
3	PO4	A	602	-	-	0/0/0/0	0/0/0/0
3	PO4	A	603	-	-	0/0/0/0	0/0/0/0
4	EDO	A	604	-	-	0/1/1/1	0/0/0/0
4	EDO	A	605	-	-	0/1/1/1	0/0/0/0
4	EDO	A	606	-	-	0/1/1/1	0/0/0/0
4	EDO	A	607	-	-	0/1/1/1	0/0/0/0
2	2P1	B	601	-	-	0/27/27/27	0/4/4/4
3	PO4	B	602	-	-	0/0/0/0	0/0/0/0
3	PO4	B	603	-	-	0/0/0/0	0/0/0/0
4	EDO	B	604	-	-	0/1/1/1	0/0/0/0
4	EDO	B	605	-	-	0/1/1/1	0/0/0/0
4	EDO	B	606	-	-	0/1/1/1	0/0/0/0
4	EDO	B	607	-	-	0/1/1/1	0/0/0/0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	2P1	C23-N24	-3.29	1.30	1.33
2	A	601	2P1	C23-N24	-3.09	1.31	1.33
2	A	601	2P1	C27-N28	-2.76	1.32	1.37
2	A	601	2P1	C23-C22	2.61	1.46	1.40
2	B	601	2P1	C27-N28	-2.50	1.33	1.37
2	B	601	2P1	C23-C22	2.42	1.45	1.40
2	B	601	2P1	F31-C30	2.00	1.40	1.32

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	2P1	C20-C29-N28	-5.63	120.80	125.25
2	B	601	2P1	C20-C29-N28	-5.46	120.93	125.25
2	A	601	2P1	C29-N28-C27	4.68	122.11	116.85
2	B	601	2P1	C21-C20-C29	4.14	120.53	116.59
2	B	601	2P1	C29-N28-C27	3.93	121.27	116.85
2	A	601	2P1	O9-S7-C5	-3.91	104.07	107.94
2	A	601	2P1	C21-C20-C29	3.80	120.20	116.59
2	B	601	2P1	C3-C4-C5	3.64	120.84	118.89
2	B	601	2P1	O8-S7-O9	3.32	123.88	119.11
2	B	601	2P1	C5-S7-C10	-2.98	99.36	104.58
2	A	601	2P1	C3-C4-C5	2.94	120.46	118.89

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	2P1	O8-S7-C10	2.92	110.84	107.94
2	A	601	2P1	O8-S7-O9	2.76	123.06	119.11
2	A	601	2P1	C5-S7-C10	-2.58	100.06	104.58
2	B	601	2P1	N25-C27-N28	2.44	132.49	125.81
2	A	601	2P1	N25-C27-N28	2.39	132.35	125.81
2	B	601	2P1	O9-S7-C5	-2.34	105.62	107.94
2	B	601	2P1	C20-C21-C22	-2.30	117.00	121.25
2	A	601	2P1	C20-C21-C22	-2.19	117.21	121.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	471/501 (94%)	0.58	21 (4%) 32 39	9, 16, 30, 57	0
1	B	469/501 (93%)	0.50	18 (3%) 38 47	9, 15, 28, 61	0
All	All	940/1002 (93%)	0.54	39 (4%) 35 43	9, 16, 29, 61	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	409	LEU	4.5
1	A	486	LEU	4.2
1	B	484	ILE	3.7
1	B	77	LYS	3.5
1	B	451	GLU	3.4
1	A	453	TYR	3.4
1	A	204	ALA	3.2
1	A	370	GLN	3.2
1	B	94	ASP	2.7
1	A	108	TYR	2.7
1	B	450	LEU	2.7
1	B	486	LEU	2.7
1	B	399	PHE	2.7
1	B	156	TRP	2.6
1	B	62	LEU	2.5
1	A	198	VAL	2.4
1	B	151	ILE	2.4
1	B	411	ILE	2.4
1	B	419	ALA	2.4
1	A	53	LYS	2.4
1	B	148	ILE	2.3
1	A	455	GLN	2.3
1	A	454	GLY	2.3
1	A	386	LEU	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	484	ILE	2.2
1	B	88	LYS	2.2
1	B	454	GLY	2.2
1	A	451	GLU	2.2
1	A	298	LEU	2.2
1	A	450	LEU	2.2
1	B	204	ALA	2.2
1	A	151[A]	ILE	2.2
1	A	156	TRP	2.1
1	A	43	LYS	2.1
1	A	159	ILE	2.1
1	A	62	LEU	2.1
1	A	255	LYS	2.1
1	A	161	VAL	2.1
1	B	108	TYR	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	PO4	B	603	5/5	0.42	19.21	22,42,46,55	0
3	PO4	A	602	5/5	0.41	11.89	24,36,47,48	0
4	EDO	B	605	4/4	0.37	7.80	34,38,38,41	0
4	EDO	B	606	4/4	0.26	7.01	31,33,34,37	0
4	EDO	A	604	4/4	0.25	5.61	35,36,36,37	0
4	EDO	A	606	4/4	0.20	3.72	34,35,35,41	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	EDO	B	604	4/4	0.15	3.01	15,17,17,21	0
3	PO4	A	603	5/5	0.18	2.98	33,36,41,46	0
4	EDO	A	605	4/4	0.16	2.82	21,22,28,28	0
3	PO4	B	602	5/5	0.19	1.55	35,36,39,40	0
2	2P1	B	601	32/32	0.11	0.75	12,17,20,20	0
4	EDO	B	607	4/4	0.16	0.70	27,27,31,36	0
2	2P1	A	601	32/32	0.12	-0.15	11,18,21,23	0
4	EDO	A	607	4/4	0.11	-0.44	15,19,21,22	0

6.5 Other polymers ⓘ

There are no such residues in this entry.