



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 10, 2019 – 01:04 PM EST

PDB ID : 6CBK
Title : X-ray structure of NeoB from *Streptomyces fradiae* in complex with PMP
Authors : Thoden, J.B.; Dow, G.T.; Holden, H.M.
Deposited on : 2018-02-03
Resolution : 1.75 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 ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix) : 1.13
EDS : rb-20031633
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20031633

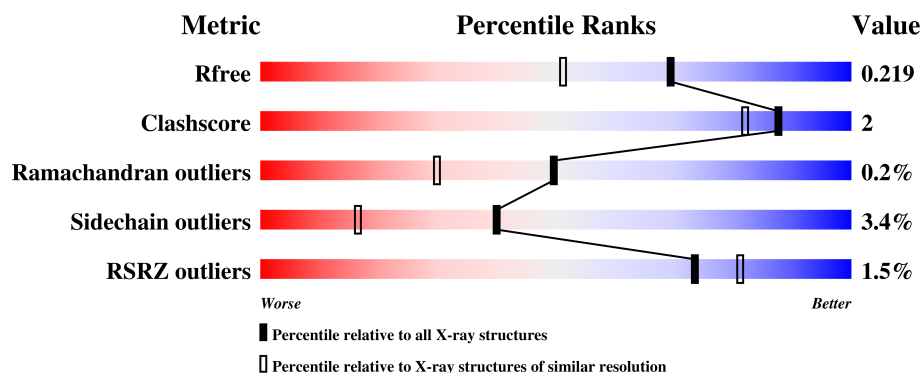
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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}	111664	1952 (1.76-1.76)
Clashscore	122126	2072 (1.76-1.76)
Ramachandran outliers	120053	2050 (1.76-1.76)
Sidechain outliers	120020	2050 (1.76-1.76)
RSRZ outliers	108989	1913 (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. 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	424	<div> <div>%</div> <div> <div></div> <div>90%</div> <div>5% . .</div> </div> </div>
1	B	424	<div> <div>%</div> <div> <div></div> <div>91%</div> <div>8% .</div> </div> </div>
1	C	424	<div> <div>2%</div> <div> <div></div> <div>90%</div> <div>5% . .</div> </div> </div>
1	D	424	<div> <div>3%</div> <div> <div></div> <div>91%</div> <div>7% .</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13702 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 Neamine transaminase NeoN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	409	Total	C	N	O	S	0	4	0
			3120	1972	566	574	8			
1	B	419	Total	C	N	O	S	0	3	0
			3206	2021	588	589	8			
1	C	410	Total	C	N	O	S	0	1	0
			3103	1958	561	576	8			
1	D	422	Total	C	N	O	S	0	5	0
			3233	2036	589	600	8			

There are 32 discrepancies between the modelled and reference sequences:

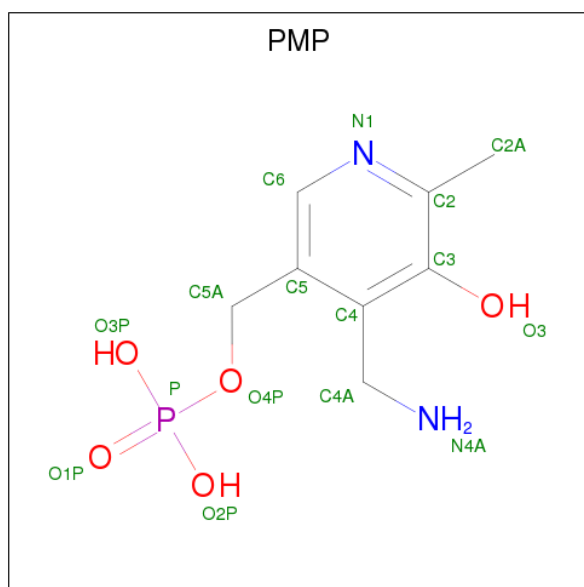
Chain	Residue	Modelled	Actual	Comment	Reference
A	417	LEU	-	expression tag	UNP Q53U08
A	418	GLU	-	expression tag	UNP Q53U08
A	419	HIS	-	expression tag	UNP Q53U08
A	420	HIS	-	expression tag	UNP Q53U08
A	421	HIS	-	expression tag	UNP Q53U08
A	422	HIS	-	expression tag	UNP Q53U08
A	423	HIS	-	expression tag	UNP Q53U08
A	424	HIS	-	expression tag	UNP Q53U08
B	417	LEU	-	expression tag	UNP Q53U08
B	418	GLU	-	expression tag	UNP Q53U08
B	419	HIS	-	expression tag	UNP Q53U08
B	420	HIS	-	expression tag	UNP Q53U08
B	421	HIS	-	expression tag	UNP Q53U08
B	422	HIS	-	expression tag	UNP Q53U08
B	423	HIS	-	expression tag	UNP Q53U08
B	424	HIS	-	expression tag	UNP Q53U08
C	417	LEU	-	expression tag	UNP Q53U08
C	418	GLU	-	expression tag	UNP Q53U08
C	419	HIS	-	expression tag	UNP Q53U08
C	420	HIS	-	expression tag	UNP Q53U08
C	421	HIS	-	expression tag	UNP Q53U08

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Chain	Residue	Modelled	Actual	Comment	Reference
C	422	HIS	-	expression tag	UNP Q53U08
C	423	HIS	-	expression tag	UNP Q53U08
C	424	HIS	-	expression tag	UNP Q53U08
D	417	LEU	-	expression tag	UNP Q53U08
D	418	GLU	-	expression tag	UNP Q53U08
D	419	HIS	-	expression tag	UNP Q53U08
D	420	HIS	-	expression tag	UNP Q53U08
D	421	HIS	-	expression tag	UNP Q53U08
D	422	HIS	-	expression tag	UNP Q53U08
D	423	HIS	-	expression tag	UNP Q53U08
D	424	HIS	-	expression tag	UNP Q53U08

- Molecule 2 is 4'-DEOXY-4'-AMINOPYRIDOXAL-5'-PHOSPHATE (three-letter code: PMP) (formula: C₈H₁₃N₂O₅P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			16	8	2	5	1		
2	B	1	Total	C	N	O	P	0	0
			16	8	2	5	1		
2	C	1	Total	C	N	O	P	0	0
			16	8	2	5	1		
2	D	1	Total	C	N	O	P	0	0
			16	8	2	5	1		

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	1	Total	Na	0	0
			1	1		

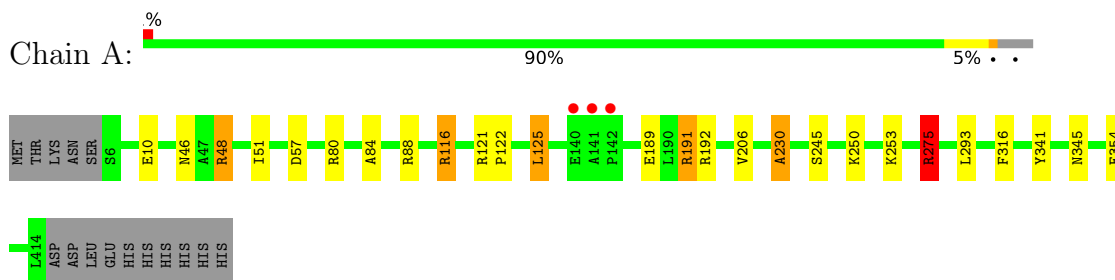
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	252	Total	O	0	0
			252	252		
5	B	240	Total	O	0	0
			240	240		
5	C	203	Total	O	0	0
			203	203		
5	D	276	Total	O	0	0
			276	276		

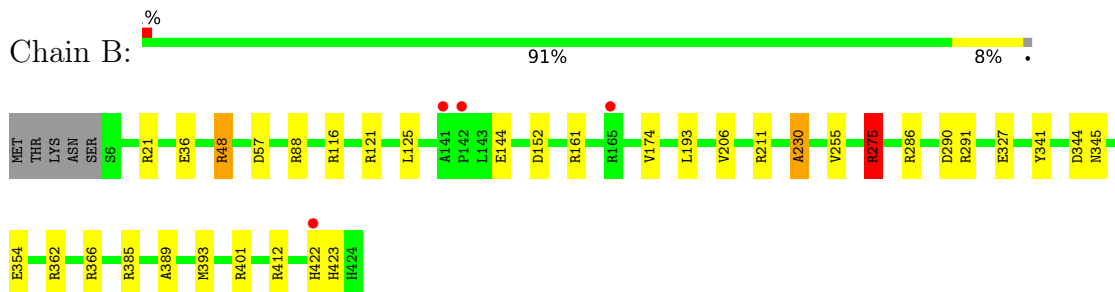
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

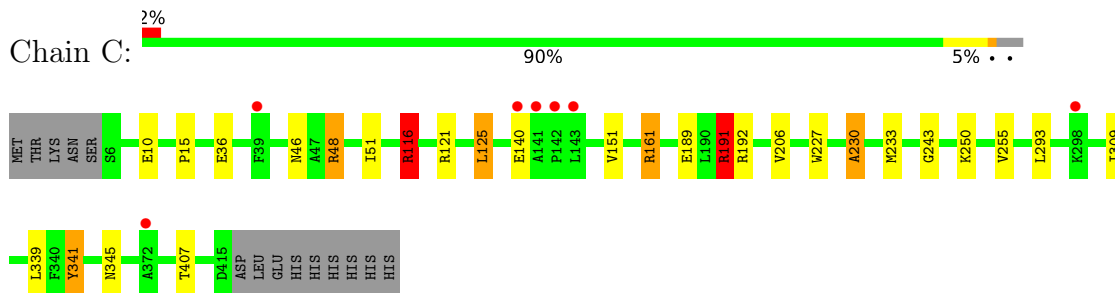
- Molecule 1: Neamine transaminase NeoN



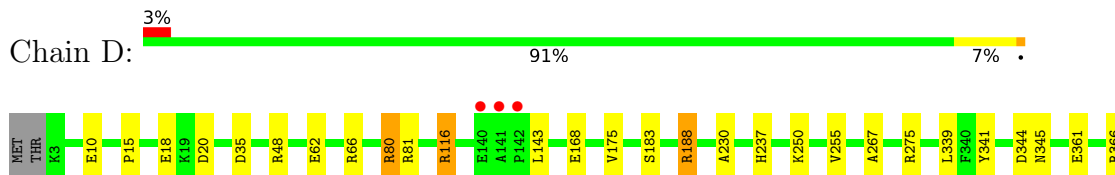
- Molecule 1: Neamine transaminase NeoN

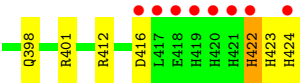


- Molecule 1: Neamine transaminase NeoN



- Molecule 1: Neamine transaminase NeoN





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	62.73Å 82.70Å 100.86Å 75.59° 88.85° 70.60°	Depositor
Resolution (Å)	97.46 – 1.75 29.52 – 1.75	Depositor EDS
% Data completeness (in resolution range)	95.9 (97.46-1.75) 95.9 (29.52-1.75)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.89 (at 1.75Å)	Xtriage
Refinement program	REFMAC 5.8.0124	Depositor
R, R_{free}	0.175 , 0.208 0.188 , 0.219	Depositor DCC
R_{free} test set	8819 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	22.2	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 43.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13702	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: NA, PMP, 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	0.56	0/3201	0.86	5/4354 (0.1%)
1	B	0.56	0/3290	0.87	10/4474 (0.2%)
1	C	0.54	0/3175	0.83	4/4320 (0.1%)
1	D	0.57	0/3323	0.92	10/4519 (0.2%)
All	All	0.56	0/12989	0.87	29/17667 (0.2%)

There are no bond length outliers.

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	48	ARG	NE-CZ-NH2	10.36	125.48	120.30
1	B	275[A]	ARG	NE-CZ-NH1	-8.43	116.09	120.30
1	B	275[B]	ARG	NE-CZ-NH1	-8.43	116.09	120.30
1	A	116	ARG	NE-CZ-NH1	-8.40	116.10	120.30
1	A	116	ARG	NE-CZ-NH2	7.96	124.28	120.30
1	D	412	ARG	NE-CZ-NH2	-7.80	116.40	120.30
1	A	275[A]	ARG	NE-CZ-NH1	-7.63	116.49	120.30
1	A	275[B]	ARG	NE-CZ-NH1	-7.63	116.49	120.30
1	C	191	ARG	NE-CZ-NH1	-7.09	116.75	120.30
1	C	116	ARG	NE-CZ-NH2	-6.73	116.93	120.30
1	D	116	ARG	NE-CZ-NH2	-6.60	117.00	120.30
1	D	188	ARG	NE-CZ-NH2	-6.60	117.00	120.30
1	B	211	ARG	NE-CZ-NH1	-6.28	117.16	120.30
1	D	35	ASP	CB-CG-OD1	6.17	123.86	118.30
1	D	188	ARG	NE-CZ-NH1	6.13	123.36	120.30
1	D	412	ARG	NE-CZ-NH1	5.99	123.29	120.30
1	D	48	ARG	NE-CZ-NH1	-5.98	117.31	120.30
1	C	191	ARG	NE-CZ-NH2	5.88	123.24	120.30
1	D	366	ARG	NE-CZ-NH2	5.84	123.22	120.30
1	B	412	ARG	NE-CZ-NH2	-5.62	117.49	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	412	ARG	NE-CZ-NH1	5.51	123.06	120.30
1	B	366	ARG	NE-CZ-NH2	5.49	123.04	120.30
1	B	385	ARG	NE-CZ-NH1	5.35	122.97	120.30
1	B	362	ARG	NE-CZ-NH1	5.28	122.94	120.30
1	C	161	ARG	NE-CZ-NH1	5.24	122.92	120.30
1	B	290	ASP	CB-CG-OD1	5.16	122.94	118.30
1	D	35	ASP	CB-CG-OD2	-5.14	113.67	118.30
1	A	125	LEU	CB-CG-CD1	5.07	119.62	111.00
1	B	152	ASP	CB-CG-OD1	5.05	122.85	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	3120	0	3087	15	0
1	B	3206	0	3141	14	0
1	C	3103	0	3048	16	0
1	D	3233	0	3160	13	0
2	A	16	0	10	0	0
2	B	16	0	10	0	0
2	C	16	0	10	1	0
2	D	16	0	10	0	0
3	A	4	0	6	0	0
4	C	1	0	0	0	0
5	A	252	0	0	1	0
5	B	240	0	0	2	0
5	C	203	0	0	1	0
5	D	276	0	0	4	0
All	All	13702	0	12482	55	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 2.

All (55) 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:291:ARG:NH2	1:B:354:GLU:OE2	2.12	0.83
1:B:116:ARG:HG3	1:B:121:ARG:O	1.85	0.77
1:C:161:ARG:NH2	1:C:189:GLU:OE1	2.21	0.73
1:B:36:GLU:OE1	5:B:601:HOH:O	2.07	0.71
1:A:48:ARG:HG3	1:A:341:TYR:CE2	2.32	0.64
1:C:191:ARG:HD3	1:C:191:ARG:O	2.01	0.61
1:B:48:ARG:HG3	1:B:341:TYR:CE2	2.35	0.61
1:C:36:GLU:OE2	1:D:80:ARG:NH2	2.34	0.60
1:A:116:ARG:HG3	1:A:121:ARG:O	2.02	0.59
1:C:191:ARG:HD3	1:C:191:ARG:C	2.24	0.56
1:C:48:ARG:HG3	1:C:341:TYR:CE2	2.41	0.55
1:B:57:ASP:OD2	1:B:275[A]:ARG:NH1	2.32	0.54
1:D:361:GLU:HG3	5:D:739:HOH:O	2.07	0.53
1:B:48:ARG:HG3	1:B:341:TYR:CZ	2.44	0.53
1:C:116:ARG:HG3	1:C:121:ARG:O	2.09	0.53
1:C:192:ARG:NH1	5:C:601:HOH:O	2.37	0.53
1:D:422:HIS:HB2	1:D:424:HIS:CD2	2.44	0.52
1:B:344:ASP:OD2	1:B:423:HIS:CD2	2.63	0.52
1:C:10:GLU:HG2	1:D:255:VAL:HB	1.91	0.52
1:A:46:ASN:HB3	1:A:51:ILE:HB	1.94	0.49
1:D:188:ARG:NH2	5:D:604:HOH:O	2.45	0.49
1:B:286:ARG:HD3	5:B:739:HOH:O	2.12	0.49
1:C:255:VAL:HB	1:D:10:GLU:HG2	1.95	0.48
1:B:174:VAL:HG13	1:B:174:VAL:O	2.15	0.47
1:D:183:SER:H	1:D:398:GLN:NE2	2.11	0.47
1:B:206:VAL:HG13	1:B:230:ALA:HB3	1.97	0.46
1:D:18:GLU:HB2	5:D:786:HOH:O	2.14	0.46
1:A:206:VAL:HG13	1:A:230:ALA:HB3	1.97	0.46
1:D:237:HIS:HD2	5:D:741:HOH:O	1.98	0.46
1:A:57:ASP:OD2	1:A:275[A]:ARG:NH1	2.32	0.45
1:C:206:VAL:HG13	1:C:230:ALA:HB3	1.98	0.45
1:C:230:ALA:O	1:C:233:MET:HG2	2.16	0.45
2:C:501:PMP:N4A	2:C:501:PMP:O3	2.50	0.45
1:A:191:ARG:HD3	1:A:191:ARG:C	2.37	0.45
1:A:84:ALA:O	1:A:88:ARG:HG3	2.17	0.45
1:A:116:ARG:HD2	1:A:122:PRO:O	2.16	0.44
1:A:10:GLU:HG2	1:B:255:VAL:HB	1.99	0.44
1:A:191:ARG:HD2	5:A:605:HOH:O	2.16	0.44
1:A:48:ARG:HG3	1:A:341:TYR:CZ	2.52	0.44
1:D:81:ARG:HD3	1:D:267:ALA:HA	2.00	0.44
1:B:161:ARG:CZ	1:B:193:LEU:HD21	2.48	0.44
1:C:293:LEU:CD2	1:C:309:ILE:HG21	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:46:ASN:HB3	1:C:51:ILE:HB	2.01	0.43
1:A:191:ARG:HD3	1:A:191:ARG:O	2.19	0.42
1:C:227:TRP:CZ2	1:C:243:GLY:HA3	2.55	0.42
1:C:125:LEU:HD12	1:C:151:VAL:HB	2.01	0.42
1:B:389:ALA:O	1:B:393:MET:HG2	2.20	0.42
1:C:15:PRO:HB3	1:C:339:LEU:HD11	2.02	0.41
1:B:423:HIS:O	1:B:423:HIS:CD2	2.73	0.41
1:A:189:GLU:OE1	1:A:192:ARG:NH2	2.50	0.41
1:A:293:LEU:HD21	1:A:316:PHE:CG	2.55	0.41
1:A:80:ARG:HA	1:A:80:ARG:HD3	1.91	0.40
1:D:15:PRO:HB3	1:D:339:LEU:HD11	2.03	0.40
1:D:344:ASP:OD2	1:D:423:HIS:CD2	2.74	0.40

There are no symmetry-related clashes.

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	411/424 (97%)	402 (98%)	8 (2%)	1 (0%)	49	30
1	B	420/424 (99%)	410 (98%)	9 (2%)	1 (0%)	49	30
1	C	409/424 (96%)	401 (98%)	7 (2%)	1 (0%)	49	30
1	D	425/424 (100%)	415 (98%)	9 (2%)	1 (0%)	49	30
All	All	1665/1696 (98%)	1628 (98%)	33 (2%)	4 (0%)	49	30

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	230	ALA
1	B	230	ALA
1	C	230	ALA

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Mol	Chain	Res	Type
1	D	230	ALA

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	307/318 (96%)	297 (97%)	10 (3%)	41	17
1	B	315/318 (99%)	304 (96%)	11 (4%)	39	15
1	C	304/318 (96%)	295 (97%)	9 (3%)	44	20
1	D	320/318 (101%)	305 (95%)	15 (5%)	29	9
All	All	1246/1272 (98%)	1201 (96%)	45 (4%)	40	14

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	ARG
1	A	125	LEU
1	A	191	ARG
1	A	245	SER
1	A	250	LYS
1	A	253	LYS
1	A	275[A]	ARG
1	A	275[B]	ARG
1	A	345	ASN
1	A	354	GLU
1	B	21	ARG
1	B	48	ARG
1	B	88	ARG
1	B	125	LEU
1	B	144	GLU
1	B	275[A]	ARG
1	B	275[B]	ARG
1	B	327	GLU
1	B	345	ASN
1	B	401	ARG

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Mol	Chain	Res	Type
1	B	422	HIS
1	C	48	ARG
1	C	116	ARG
1	C	125	LEU
1	C	140	GLU
1	C	191	ARG
1	C	250	LYS
1	C	341	TYR
1	C	345	ASN
1	C	407	THR
1	D	20	ASP
1	D	62	GLU
1	D	66	ARG
1	D	80	ARG
1	D	116	ARG
1	D	143	LEU
1	D	168	GLU
1	D	250	LYS
1	D	275[A]	ARG
1	D	275[B]	ARG
1	D	341	TYR
1	D	345	ASN
1	D	401	ARG
1	D	416	ASP
1	D	422	HIS

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	334	ASN
1	B	90	HIS
1	B	420	HIS
1	B	423	HIS
1	C	237	HIS
1	C	398	GLN
1	D	90	HIS
1	D	237	HIS
1	D	398	GLN
1	D	423	HIS
1	D	424	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 1 is monoatomic - leaving 5 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$
2	PMP	A	501	-	16,16,16	2.93	3 (18%)	20,23,23	1.64	5 (25%)
3	EDO	A	502	-	3,3,3	0.61	0	2,2,2	0.49	0
2	PMP	B	501	-	16,16,16	3.09	4 (25%)	20,23,23	2.21	9 (45%)
2	PMP	C	501	-	16,16,16	3.61	3 (18%)	20,23,23	1.69	6 (30%)
2	PMP	D	501	-	16,16,16	3.20	3 (18%)	20,23,23	2.41	8 (40%)

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	PMP	A	501	-	-	0/8/8/8	0/1/1/1
3	EDO	A	502	-	-	0/1/1/1	0/0/0/0
2	PMP	B	501	-	-	0/8/8/8	0/1/1/1
2	PMP	C	501	-	-	0/8/8/8	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PMP	D	501	-	-	0/8/8/8	0/1/1/1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	PMP	P-O4P	2.47	1.68	1.60
2	B	501	PMP	C3-C4	3.87	1.46	1.40
2	B	501	PMP	C5-C4	4.56	1.46	1.40
2	A	501	PMP	C5-C4	4.73	1.47	1.40
2	D	501	PMP	C3-C4	4.80	1.47	1.40
2	C	501	PMP	C3-C4	4.82	1.47	1.40
2	A	501	PMP	C3-C4	5.33	1.48	1.40
2	D	501	PMP	C5-C4	5.36	1.48	1.40
2	C	501	PMP	C5-C4	5.61	1.48	1.40
2	A	501	PMP	C3-C2	9.00	1.47	1.40
2	D	501	PMP	C3-C2	10.11	1.47	1.40
2	B	501	PMP	C3-C2	10.30	1.47	1.40
2	C	501	PMP	C3-C2	12.04	1.49	1.40

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	PMP	C2A-C2-C3	-6.58	113.12	120.96
2	B	501	PMP	C2A-C2-C3	-4.87	115.15	120.96
2	C	501	PMP	O4P-P-O1P	-3.41	96.91	106.47
2	A	501	PMP	C2A-C2-C3	-3.16	117.20	120.96
2	B	501	PMP	C4-C3-C2	-2.90	115.55	120.01
2	B	501	PMP	C5-C6-N1	-2.71	119.25	123.83
2	D	501	PMP	C4-C3-C2	-2.38	116.34	120.01
2	D	501	PMP	C5-C6-N1	-2.37	119.83	123.83
2	A	501	PMP	C4-C3-C2	-2.21	116.61	120.01
2	D	501	PMP	O3P-P-O4P	-2.13	101.06	106.73
2	C	501	PMP	C3-C2-N1	-2.13	117.96	120.75
2	B	501	PMP	C4A-C4-C3	-2.05	117.17	120.38
2	D	501	PMP	O3P-P-O2P	2.08	115.80	107.59
2	C	501	PMP	O3-C3-C2	2.23	122.44	117.78
2	C	501	PMP	C5A-C5-C6	2.26	123.14	119.34
2	A	501	PMP	C2A-C2-N1	2.27	122.35	117.86
2	A	501	PMP	C6-N1-C2	2.36	123.73	119.19
2	B	501	PMP	O3P-P-O2P	2.39	117.06	107.59
2	A	501	PMP	O3P-P-O2P	2.43	117.21	107.59
2	D	501	PMP	C6-N1-C2	2.54	124.08	119.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	PMP	C6-N1-C2	2.75	124.49	119.19
2	C	501	PMP	C2A-C2-N1	2.80	123.41	117.86
2	B	501	PMP	O3-C3-C2	2.87	123.79	117.78
2	B	501	PMP	C6-N1-C2	3.00	124.97	119.19
2	B	501	PMP	C6-C5-C4	3.20	120.47	118.14
2	B	501	PMP	C2A-C2-N1	3.30	124.41	117.86
2	D	501	PMP	C2A-C2-N1	3.94	125.66	117.86
2	D	501	PMP	C6-C5-C4	4.39	121.34	118.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	501	PMP	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	409/424 (96%)	-0.29	3 (0%) 87 92	14, 23, 40, 69	0
1	B	419/424 (98%)	-0.11	4 (0%) 82 88	16, 23, 45, 79	0
1	C	410/424 (96%)	-0.08	7 (1%) 70 78	14, 27, 46, 68	0
1	D	422/424 (99%)	-0.13	11 (2%) 56 62	15, 22, 47, 94	0
All	All	1660/1696 (97%)	-0.15	25 (1%) 73 81	14, 24, 45, 94	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	141	ALA	5.8
1	D	422	HIS	5.4
1	D	417	LEU	5.1
1	B	142	PRO	4.9
1	D	142	PRO	4.9
1	C	142	PRO	4.8
1	B	422	HIS	3.9
1	C	141	ALA	3.6
1	D	418	GLU	3.5
1	D	140	GLU	3.5
1	D	416	ASP	3.2
1	D	419	HIS	3.2
1	D	421	HIS	3.1
1	C	298	LYS	2.8
1	D	420	HIS	2.6
1	C	140	GLU	2.5
1	A	142	PRO	2.5
1	A	140	GLU	2.4
1	D	424	HIS	2.2
1	C	372	ALA	2.2
1	A	141	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	143	LEU	2.1
1	B	141	ALA	2.1
1	C	39	PHE	2.0
1	B	165	ARG	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. 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(\AA^2)	Q<0.9
3	EDO	A	502	4/4	0.91	0.11	34,36,40,42	0
2	PMP	D	501	16/16	0.98	0.07	16,21,27,42	0
2	PMP	C	501	16/16	0.98	0.08	18,23,27,40	0
2	PMP	A	501	16/16	0.98	0.10	16,19,24,42	0
2	PMP	B	501	16/16	0.98	0.09	16,23,26,37	0
4	NA	C	502	1/1	0.99	0.03	20,20,20,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.