



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 1, 2016 – 02:07 PM EDT

PDB ID : 5D05
Title : Neisseria meningitidis 3 deoxy-D-arabino-heptulosonate 7-phosphate synthase
Lys107Ala variant regulated
Authors : Heyes, L.C.; Parker, E.J.
Deposited on : 2015-08-02
Resolution : 1.75 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027939
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027939

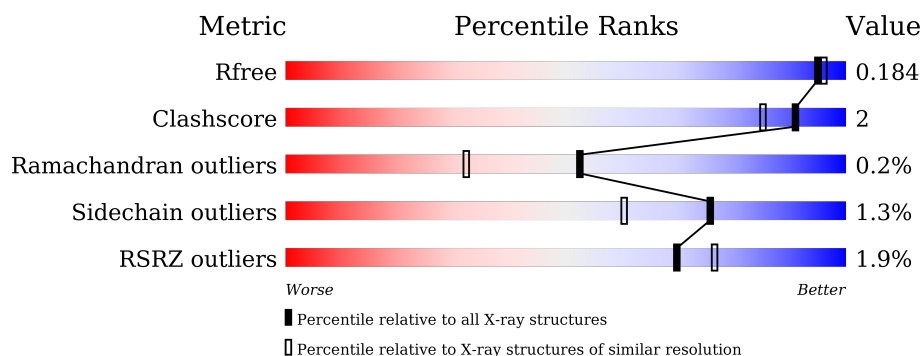
1 Overall quality at a glance

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}	91344	1609 (1.76-1.76)
Clashscore	102246	1730 (1.76-1.76)
Ramachandran outliers	100387	1711 (1.76-1.76)
Sidechain outliers	100360	1711 (1.76-1.76)
RSRZ outliers	91569	1610 (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	351	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 92% 6% .. </div> </div>
1	B	351	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 93% 5% . </div> </div>
1	C	351	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 93% 5% . </div> </div>
1	D	351	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 4% 95% . . </div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
8	PEG	B	403	-	-	-	X

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 11952 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 Phospho-2-dehydro-3-deoxyheptonate aldolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	346	Total	C	N	O	S	0	3	0
			2689	1683	487	505	14			
1	B	346	Total	C	N	O	S	0	1	0
			2663	1668	480	501	14			
1	C	346	Total	C	N	O	S	0	1	0
			2664	1668	480	502	14			
1	D	347	Total	C	N	O	S	0	2	0
			2658	1663	480	501	14			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	107	ALA	LYS	engineered mutation	UNP Q9K169
B	107	ALA	LYS	engineered mutation	UNP Q9K169
C	107	ALA	LYS	engineered mutation	UNP Q9K169
D	107	ALA	LYS	engineered mutation	UNP Q9K169

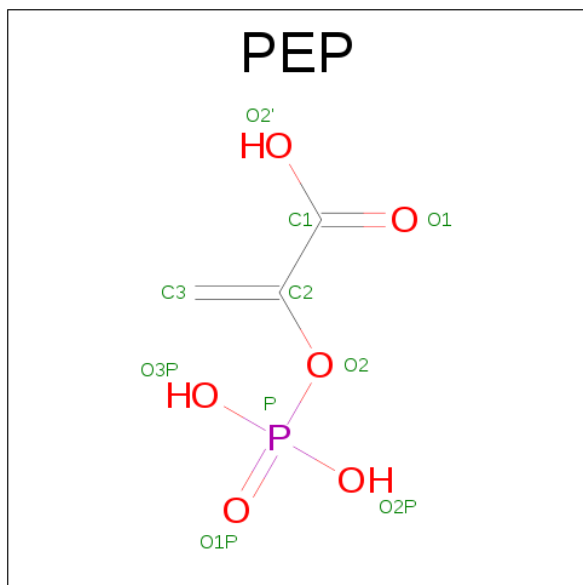
- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mn	0	0
			1	1		
2	A	1	Total	Mn	0	0
			1	1		
2	D	1	Total	Mn	0	0
			1	1		
2	C	1	Total	Mn	0	0
			1	1		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Cl 1 1	0	0
3	D	2	Total Cl 2 2	0	0

- Molecule 4 is PHOSPHOENOLPYRUVATE (three-letter code: PEP) (formula: $C_3H_5O_6P$).



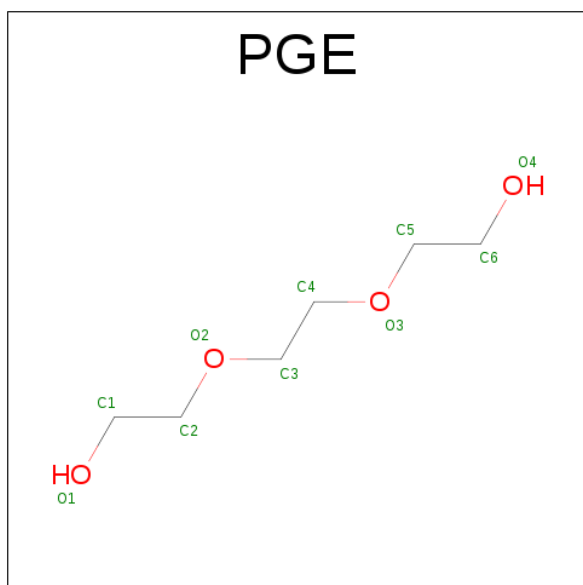
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O P 10 3 6 1	0	0
4	B	1	Total C O P 10 3 6 1	0	0
4	C	1	Total C O P 10 3 6 1	0	0
4	D	1	Total C O P 10 3 6 1	0	0

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



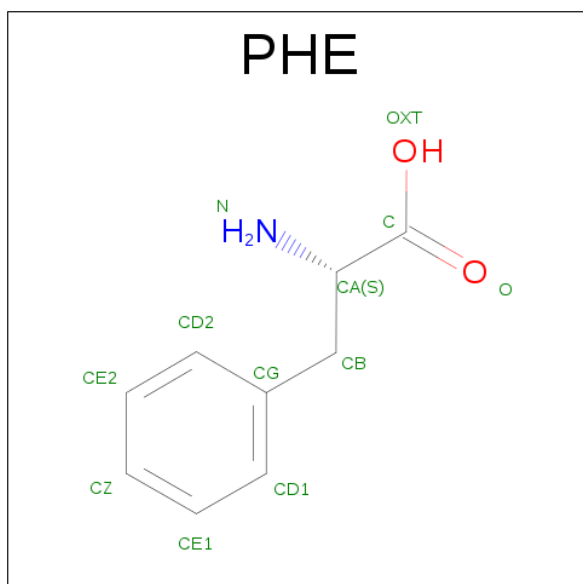
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).



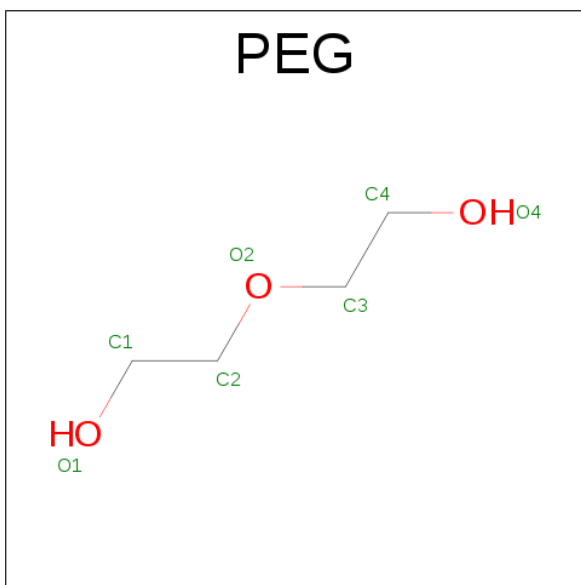
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			10	6	4		

- Molecule 7 is PHENYLALANINE (three-letter code: PHE) (formula: $C_9H_{11}NO_2$).



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

- Molecule 8 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	B	1	Total	C	O	0	0
			7	4	3		
8	B	1	Total	C	O	0	0
			7	4	3		
8	C	1	Total	C	O	0	0
			7	4	3		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	300	Total	O	0	0
			300	300		
9	B	304	Total	O	0	0
			304	304		
9	C	327	Total	O	0	0
			327	327		
9	D	201	Total	O	0	0
			201	201		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	75.33Å 141.03Å 76.03Å 90.00° 97.02° 90.00°	Depositor
Resolution (Å)	74.46 – 1.75 44.18 – 1.75	Depositor EDS
% Data completeness (in resolution range)	100.0 (74.46-1.75) 100.0 (44.18-1.75)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 1.75Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.150 , 0.176 0.162 , 0.184	Depositor DCC
R_{free} test set	9914 reflections (6.69%)	DCC
Wilson B-factor (Å ²)	20.7	Xtriage
Anisotropy	0.048	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.024 for l,-k,h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	11952	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.72% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PGE, CL, MN, PEP, SO4, PEG

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	0.57	0/2740	0.69	0/3706
1	B	0.55	1/2714 (0.0%)	0.69	1/3674 (0.0%)
1	C	0.57	1/2716 (0.0%)	0.72	3/3678 (0.1%)
1	D	0.49	0/2710	0.65	0/3675
All	All	0.54	2/10880 (0.0%)	0.69	4/14733 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	145	GLU	CD-OE1	6.25	1.32	1.25
1	B	145	GLU	CD-OE2	5.86	1.32	1.25

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	52	ARG	NE-CZ-NH2	6.05	123.33	120.30
1	C	40	ARG	NE-CZ-NH2	5.42	123.01	120.30
1	C	40	ARG	NE-CZ-NH1	-5.39	117.61	120.30
1	B	267	ASP	CB-CG-OD1	5.29	123.06	118.30

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	2689	0	2667	20	0
1	B	2663	0	2636	10	0
1	C	2664	0	2632	14	0
1	D	2658	0	2596	9	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	D	2	0	0	1	0
4	A	10	0	2	0	0
4	B	10	0	2	0	0
4	C	10	0	2	0	0
4	D	10	0	2	0	0
5	A	5	0	0	0	0
5	B	10	0	0	0	0
5	D	5	0	0	0	0
6	A	10	0	14	1	0
7	A	12	0	8	0	0
7	B	12	0	8	0	0
7	C	12	0	8	0	0
7	D	12	0	8	0	0
8	B	14	0	20	1	0
8	C	7	0	10	0	0
9	A	300	0	0	9	0
9	B	304	0	0	2	0
9	C	327	0	0	5	0
9	D	201	0	0	4	0
All	All	11952	0	10615	48	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.

The worst 5 of 48 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:C:40:ARG:NH1	9:C:501:HOH:O	1.89	1.05
1:D:219:HIS:CD2	9:D:501:HOH:O	2.32	0.81
1:A:334:GLU:HG3	9:A:520:HOH:O	1.82	0.80
1:A:219:HIS:CD2	9:A:501:HOH:O	2.35	0.78
1:D:219:HIS:NE2	9:D:501:HOH:O	2.17	0.77

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	347/351 (99%)	338 (97%)	8 (2%)	1 (0%)	46	25
1	B	345/351 (98%)	336 (97%)	8 (2%)	1 (0%)	46	25
1	C	345/351 (98%)	338 (98%)	7 (2%)	0	100	100
1	D	347/351 (99%)	339 (98%)	7 (2%)	1 (0%)	46	25
All	All	1384/1404 (99%)	1351 (98%)	30 (2%)	3 (0%)	52	32

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	294	ASP
1	A	294	ASP
1	D	294	ASP

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	283/288 (98%)	276 (98%)	7 (2%)	55	30
1	B	279/288 (97%)	277 (99%)	2 (1%)	88	79
1	C	280/288 (97%)	277 (99%)	3 (1%)	80	66
1	D	275/288 (96%)	272 (99%)	3 (1%)	80	66

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1117/1152 (97%)	1102 (99%)	15 (1%)	76 60

5 of 15 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	334	GLU
1	B	189	ASN
1	D	189	ASN
1	A	239	LYS
1	C	236	ARG

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

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 23 ligands modelled in this entry, 7 are monoatomic - leaving 16 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
4	PEP	A	403	-	6,9,9	1.00	0	6,13,13	1.11	0
5	SO4	A	404	-	4,4,4	0.41	0	6,6,6	0.22	0
6	PGE	A	405	-	9,9,9	0.52	0	8,8,8	0.57	0
7	PHE	A	406	-	9,12,12	0.53	0	10,15,15	0.32	0
8	PEG	B	402	-	6,6,6	0.41	0	5,5,5	0.29	0
8	PEG	B	403	-	6,6,6	0.94	0	5,5,5	1.26	1 (20%)
4	PEP	B	404	-	6,9,9	1.36	1 (16%)	6,13,13	1.51	1 (16%)
5	SO4	B	405	-	4,4,4	0.40	0	6,6,6	0.22	0
5	SO4	B	406	-	4,4,4	0.35	0	6,6,6	0.15	0
7	PHE	B	407	-	9,12,12	0.56	0	10,15,15	0.32	0
8	PEG	C	402	-	6,6,6	0.48	0	5,5,5	0.61	0
4	PEP	C	403	-	6,9,9	0.93	0	6,13,13	1.63	2 (33%)
7	PHE	C	404	-	9,12,12	0.43	0	10,15,15	0.55	0
4	PEP	D	404	-	6,9,9	1.47	1 (16%)	6,13,13	1.96	2 (33%)
5	SO4	D	405	-	4,4,4	0.37	0	6,6,6	0.09	0
7	PHE	D	406	-	9,12,12	0.33	0	10,15,15	0.48	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
4	PEP	A	403	-	-	0/5/9/9	0/0/0/0
5	SO4	A	404	-	-	0/0/0/0	0/0/0/0
6	PGE	A	405	-	-	0/7/7/7	0/0/0/0
7	PHE	A	406	-	-	0/4/8/8	0/1/1/1
8	PEG	B	402	-	-	0/4/4/4	0/0/0/0
8	PEG	B	403	-	-	0/4/4/4	0/0/0/0
4	PEP	B	404	-	-	0/5/9/9	0/0/0/0
5	SO4	B	405	-	-	0/0/0/0	0/0/0/0
5	SO4	B	406	-	-	0/0/0/0	0/0/0/0
7	PHE	B	407	-	-	0/4/8/8	0/1/1/1
8	PEG	C	402	-	-	0/4/4/4	0/0/0/0
4	PEP	C	403	-	-	0/5/9/9	0/0/0/0
7	PHE	C	404	-	-	0/4/8/8	0/1/1/1
4	PEP	D	404	-	-	0/5/9/9	0/0/0/0
5	SO4	D	405	-	-	0/0/0/0	0/0/0/0
7	PHE	D	406	-	-	0/4/8/8	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	404	PEP	P-O2P	-2.37	1.46	1.54
4	D	404	PEP	P-O2	2.31	1.63	1.59

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	403	PEP	O3P-P-O2	-2.35	98.28	105.49
4	D	404	PEP	O3P-P-O2	-2.31	98.40	105.49
8	B	403	PEG	C3-O2-C2	2.15	122.51	113.31
4	B	404	PEP	O2P-P-O1P	2.19	117.77	110.63
4	C	403	PEP	O3P-P-O2P	2.55	116.82	107.44

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
6	A	405	PGE	1	0
8	B	403	PEG	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	346/351 (98%)	-0.07	4 (1%) 81 86	14, 21, 32, 47	0
1	B	346/351 (98%)	-0.24	4 (1%) 81 86	14, 20, 31, 52	0
1	C	346/351 (98%)	-0.21	4 (1%) 81 86	13, 19, 31, 52	0
1	D	347/351 (98%)	0.21	14 (4%) 42 48	16, 29, 43, 65	0
All	All	1385/1404 (98%)	-0.08	26 (1%) 70 76	13, 21, 38, 65	0

The worst 5 of 26 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	348	ALA	5.6
1	D	294	ASP	3.9
1	C	5	TYR	3.7
1	B	5	TYR	3.5
1	D	3	HIS	3.5

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. 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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
8	PEG	B	403	7/7	0.72	0.27	28.20	33,35,40,43	0
8	PEG	C	402	7/7	0.85	0.16	1.97	44,46,51,53	0
6	PGE	A	405	10/10	0.87	0.16	0.69	38,39,51,54	0
7	PHE	D	406	12/12	0.93	0.09	0.08	18,19,23,24	0
7	PHE	C	404	12/12	0.97	0.08	-0.10	15,15,16,16	0
7	PHE	A	406	12/12	0.96	0.10	-0.13	15,17,17,17	0
7	PHE	B	407	12/12	0.97	0.08	-0.40	16,17,17,17	0
2	MN	C	401	1/1	1.00	0.08	-0.62	16,16,16,16	0
2	MN	A	401	1/1	1.00	0.08	-0.73	18,18,18,18	0
4	PEP	A	403	10/10	0.99	0.09	-0.90	17,19,21,22	0
4	PEP	D	404	10/10	0.98	0.08	-1.00	25,27,29,29	0
4	PEP	C	403	10/10	0.98	0.07	-1.40	16,18,21,21	0
4	PEP	B	404	10/10	0.99	0.05	-1.55	17,19,20,20	0
2	MN	D	401	1/1	0.97	0.06	-1.65	27,27,27,27	0
2	MN	B	401	1/1	1.00	0.05	-1.94	17,17,17,17	0
3	CL	D	403	1/1	0.98	0.05	-2.71	37,37,37,37	0
5	SO4	D	405	5/5	0.73	0.32	-	92,94,97,98	0
5	SO4	B	406	5/5	0.82	0.40	-	89,92,99,100	0
3	CL	D	402	1/1	0.99	0.04	-	25,25,25,25	0
5	SO4	B	405	5/5	0.82	0.34	-	68,75,80,84	0
8	PEG	B	402	7/7	0.93	0.14	-	44,46,50,50	0
5	SO4	A	404	5/5	0.88	0.24	-	88,89,92,95	0
3	CL	A	402	1/1	1.00	0.06	-	23,23,23,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.