



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 11:13 PM GMT

PDB ID : 2C13
Title : 5-HYDROXY-LEVULINIC ACID BOUND TO PORPHOBILINOGEN SYNTHASE FROM PSEUDOMONAS AERUGINOSA
Authors : Frere, F.; Nentwich, M.; Gacond, S.; Heinz, D.W.; Neier, R.; Frankenberg-Dinkel, N.
Deposited on : 2005-09-11
Resolution : 2.15 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>

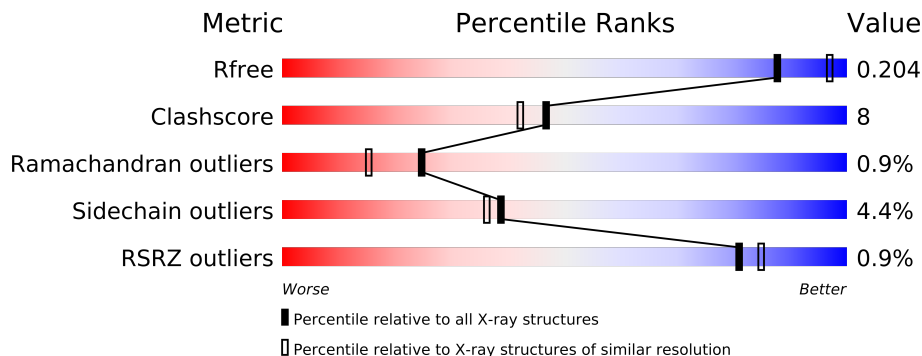
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
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) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.15 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	1094 (2.18-2.14)
Clashscore	79885	1299 (2.18-2.14)
Ramachandran outliers	78287	1272 (2.18-2.14)
Sidechain outliers	78261	1272 (2.18-2.14)
RSRZ outliers	66119	1094 (2.18-2.14)

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	337	
1	B	337	

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

Mol	Type	Chain	Res	Geometry	Electron density
6	PGE	B	1338	-	X
6	PGE	B	1339	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6020 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 DELTA-AMINOLEVULINIC ACID DEHYDRATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	334	Total	C	N	O	S	0	21	1
			2777	1740	492	533	12			
1	B	331	Total	C	N	O	S	0	22	1
			2753	1723	482	536	12			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	199	VAL	ILE	ENGINEERED MUTATION	UNP Q59643
B	199	VAL	ILE	ENGINEERED MUTATION	UNP Q59643

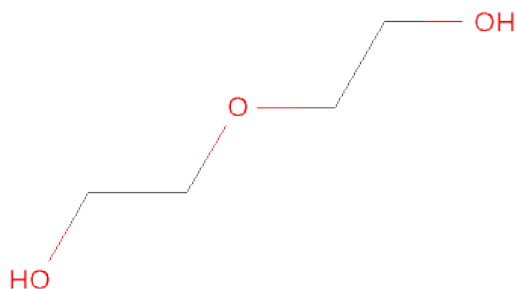
- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		

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

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

- Molecule 4 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).

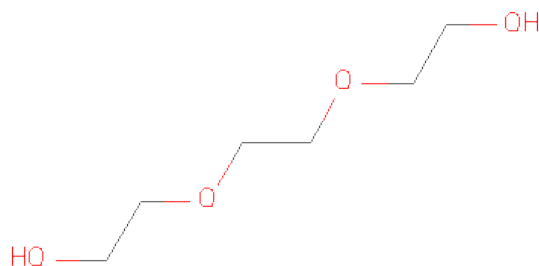


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 5 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	K	0	0
			1	1		
5	A	1	Total	K	0	0
			1	1		

- Molecule 6 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C₆H₁₄O₄).



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

- Molecule 7 is water.

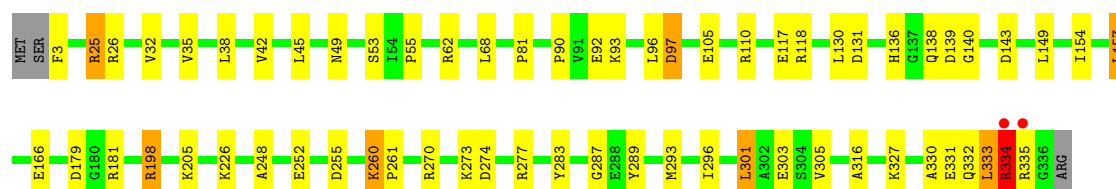
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	228	Total	O	0	0
			228	228		
7	B	229	Total	O	0	0
			229	229		

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.

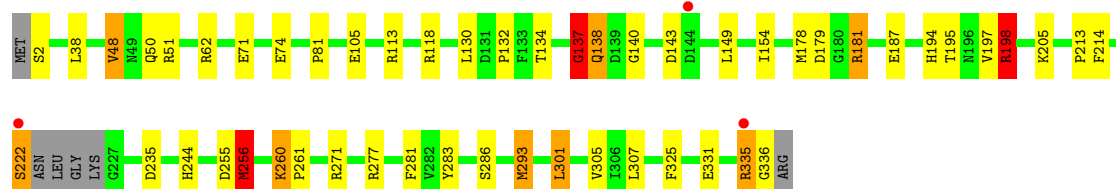
• Molecule 1: DELTA-AMINOLEVULINIC ACID DEHYDRATASE

Chain A: 



• Molecule 1: DELTA-AMINOLEVULINIC ACID DEHYDRATASE

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α , β , γ	126.31Å 126.31Å 85.48Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	89.80 – 2.15 39.94 – 2.15	Depositor EDS
% Data completeness (in resolution range)	99.2 (89.80-2.15) 99.2 (39.94-2.15)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.14 (at 2.16Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.152 , 0.204 0.153 , 0.204	Depositor DCC
R_{free} test set	1931 reflections (5.38%)	DCC
Wilson B-factor (Å ²)	19.5	Xtriage
Anisotropy	0.009	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 30.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 37848 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6020	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.88% 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: FHL, MG, PGE, CL, K, 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	0.81	0/2791	0.87	9/3782 (0.2%)
1	B	0.83	0/2766	0.88	10/3749 (0.3%)
All	All	0.82	0/5557	0.88	19/7531 (0.3%)

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	3
1	B	0	4
All	All	0	7

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	198	ARG	NE-CZ-NH2	-11.13	114.73	120.30
1	A	25[A]	ARG	NE-CZ-NH2	-8.98	115.81	120.30
1	A	25[B]	ARG	NE-CZ-NH2	-8.98	115.81	120.30
1	B	198	ARG	NE-CZ-NH1	7.57	124.08	120.30
1	A	97	ASP	CB-CG-OD1	6.80	124.42	118.30
1	B	113	ARG	NE-CZ-NH2	-6.38	117.11	120.30
1	A	25[A]	ARG	NE-CZ-NH1	6.15	123.38	120.30
1	A	25[B]	ARG	NE-CZ-NH1	6.15	123.38	120.30
1	B	138	GLN	N-CA-C	5.80	126.67	111.00
1	B	277	ARG	NE-CZ-NH1	5.76	123.18	120.30
1	A	333	LEU	N-CA-C	5.39	125.56	111.00
1	B	277	ARG	NE-CZ-NH2	-5.37	117.61	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	256[A]	MET	CG-SD-CE	5.34	108.74	100.20
1	B	256[B]	MET	CG-SD-CE	5.34	108.74	100.20
1	A	26	ARG	NE-CZ-NH2	-5.34	117.63	120.30
1	A	157	LEU	CA-CB-CG	5.29	127.48	115.30
1	B	271	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	A	274	ASP	CB-CG-OD1	5.14	122.93	118.30
1	B	235	ASP	CB-CG-OD1	5.02	122.81	118.30

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	260	FHL	Mainchain,Peptide
1	A	333	LEU	Peptide
1	B	137	GLY	Peptide
1	B	205	FHL	Mainchain
1	B	260	FHL	Mainchain,Peptide

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	2777	0	2717	43	6
1	B	2753	0	2669	48	6
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	7	0	10	2	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	B	20	0	28	1	0
7	A	228	0	0	16	1
7	B	229	0	0	16	1
All	All	6020	0	5424	90	7

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 8.

All (90) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:25[B]:ARG:NH1	7:A:2037:HOH:O	1.71	1.22
1:A:25[A]:ARG:HD2	7:A:2037:HOH:O	1.57	1.04
1:B:335:ARG:HG3	1:B:336:GLY:N	1.80	0.96
1:B:187[B]:GLU:OE2	7:B:2152:HOH:O	1.84	0.96
1:A:166[B]:GLU:OE2	7:A:2138:HOH:O	1.88	0.89
1:A:32[A]:VAL:CG2	7:A:2045:HOH:O	2.21	0.87
1:B:293[B]:MET:HE2	1:B:325:PHE:HE1	1.42	0.84
1:B:118[A]:ARG:NE	7:B:2112:HOH:O	2.11	0.83
1:B:118[A]:ARG:NH2	7:B:2112:HOH:O	2.13	0.80
1:A:330:ALA:O	1:A:334:ARG:HB2	1.83	0.78
1:B:293[B]:MET:CE	1:B:325:PHE:HE1	1.97	0.76
1:B:222[A]:SER:C	7:B:2173:HOH:O	2.23	0.75
7:A:2002:HOH:O	1:B:244[B]:HIS:HD2	1.72	0.73
1:B:198:ARG:N	1:B:198:ARG:HD3	2.04	0.73
1:B:71[A]:GLU:OE2	7:B:2078:HOH:O	2.06	0.72
1:B:331[B]:GLU:OE2	7:B:2224:HOH:O	2.10	0.70
1:B:222[A]:SER:O	7:B:2171:HOH:O	2.12	0.68
1:B:335:ARG:CG	1:B:336:GLY:N	2.53	0.67
1:A:32[B]:VAL:HG13	7:A:2045:HOH:O	1.94	0.67
1:A:166[A]:GLU:OE2	7:A:2139:HOH:O	2.12	0.67
1:B:194:HIS:HB3	1:B:197[B]:VAL:CG1	2.24	0.67
1:B:154[B]:ILE:HD11	1:B:178:MET:HB3	1.78	0.66
1:B:154[B]:ILE:HD11	1:B:178:MET:CB	2.26	0.65
1:B:222[B]:SER:HA	7:B:2173:HOH:O	1.97	0.65
1:A:32[A]:VAL:HG23	7:A:2045:HOH:O	1.91	0.63
1:B:118[A]:ARG:CZ	7:B:2112:HOH:O	2.40	0.62
1:B:213:PRO:HB2	1:B:286:SER:HB2	1.81	0.62
1:A:136[A]:HIS:HD2	1:A:138:GLN:H	1.48	0.62
1:B:2:SER:HB2	7:B:2001:HOH:O	1.98	0.61
1:B:293[B]:MET:HE2	1:B:325:PHE:CE1	2.30	0.60
1:A:301:LEU:HG	1:A:305:VAL:HG11	1.83	0.60
1:B:105[B]:GLU:OE1	7:B:2103:HOH:O	2.17	0.60
1:B:143:ASP:HB3	1:B:149:LEU:HD21	1.84	0.59
1:A:327:LYS:HE2	1:A:331:GLU:OE2	2.03	0.58
1:A:252:GLU:HA	4:A:1338:PEG:H42	1.84	0.58
1:A:273[B]:LYS:NZ	7:A:2196:HOH:O	2.37	0.57
1:B:301:LEU:HG	1:B:305:VAL:HG11	1.87	0.56
1:A:252:GLU:HA	4:A:1338:PEG:C4	2.37	0.55
1:B:38:LEU:O	1:B:81:PRO:HD2	2.06	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:293[B]:MET:CE	1:B:325:PHE:CE1	2.86	0.55
1:A:289:TYR:CZ	1:A:293[A]:MET:HG3	2.42	0.54
1:A:45:LEU:HD21	1:A:62[A]:ARG:HD3	1.89	0.54
1:B:130:LEU:HD12	1:B:140:GLY:HA2	1.88	0.54
1:B:134:THR:OG1	1:B:137:GLY:HA2	2.06	0.54
1:A:35[A]:VAL:HG11	7:A:2074:HOH:O	2.09	0.53
1:B:48:VAL:O	1:B:50:GLN:HG2	2.09	0.52
1:A:136[A]:HIS:CD2	1:A:138:GLN:HB2	2.44	0.52
1:A:143:ASP:HB3	1:A:149:LEU:HD21	1.90	0.52
1:B:132:PRO:HB3	7:B:2084:HOH:O	2.10	0.52
1:A:62[B]:ARG:NH2	7:A:2064:HOH:O	2.39	0.52
1:B:195[A]:THR:HG21	7:B:2157:HOH:O	2.11	0.51
1:A:130:LEU:HD12	1:A:140:GLY:HA2	1.92	0.51
1:A:105:GLU:O	1:A:110[A]:ARG:HD3	2.10	0.51
1:B:154[A]:ILE:HD11	1:B:179:ASP:O	2.11	0.50
1:B:2:SER:CB	7:B:2001:HOH:O	2.58	0.50
1:A:332[B]:GLN:HG2	1:A:335:ARG:NH2	2.26	0.50
1:B:222[A]:SER:HA	7:B:2173:HOH:O	2.11	0.50
1:A:136[A]:HIS:CD2	1:A:138:GLN:H	2.30	0.49
1:B:198:ARG:HA	1:B:255:ASP:OD2	2.12	0.49
1:A:38:LEU:O	1:A:81:PRO:HD2	2.12	0.49
1:A:45:LEU:HG	1:A:62[A]:ARG:HB3	1.95	0.48
1:A:55:PRO:HB3	6:B:1338:PGE:H52	1.96	0.48
1:B:154[B]:ILE:HD11	1:B:178:MET:HB2	1.95	0.47
1:A:198:ARG:HA	1:A:255:ASP:OD2	2.15	0.47
1:A:273[A]:LYS:NZ	7:A:2196:HOH:O	2.36	0.47
1:A:131:ASP:HB3	1:A:139:ASP:OD1	2.16	0.46
1:A:136[A]:HIS:HE1	7:A:2084:HOH:O	1.98	0.46
1:A:90:PRO:HD2	1:A:93:LYS:HD2	1.98	0.46
1:B:214:PHE:CG	1:B:260:FHL:H2C1	2.52	0.45
1:A:42[B]:VAL:HG21	1:A:68:LEU:HD13	1.99	0.45
1:B:260:FHL:H3C1	1:B:260:FHL:HEC2	1.81	0.45
1:A:97:ASP:HB3	7:A:2088:HOH:O	2.16	0.44
1:B:154[B]:ILE:CD1	1:B:178:MET:CB	2.93	0.44
1:B:256[A]:MET:HB2	1:B:281:PHE:CD1	2.53	0.43
1:A:42[B]:VAL:HG21	1:A:68:LEU:CD1	2.48	0.43
1:B:198:ARG:N	1:B:198:ARG:CD	2.79	0.42
1:A:205:FHL:H3C2	1:A:260:FHL:NZ	2.35	0.42
1:B:154[B]:ILE:HD12	1:B:181:ARG:CB	2.50	0.41
1:A:154:ILE:HD11	1:A:179:ASP:O	2.19	0.41
1:A:296:ILE:HG21	1:A:303[B]:GLU:HG2	2.01	0.41
1:A:248:ALA:O	1:A:252:GLU:HG3	2.21	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:92[A]:GLU:HG3	7:A:2081:HOH:O	2.20	0.41
1:A:270:ARG:HA	1:A:316:ALA:O	2.19	0.41
1:A:287:GLY:HA3	1:B:301:LEU:HD11	2.03	0.41
1:A:32[A]:VAL:HG21	7:A:2045:HOH:O	2.05	0.41
1:B:222[A]:SER:CA	7:B:2173:HOH:O	2.62	0.41
1:B:154[B]:ILE:CD1	1:B:178:MET:HB2	2.51	0.41
1:B:51:ARG:HA	1:B:62:ARG:O	2.21	0.40

All (7) 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)
1:A:277[B]:ARG:NH1	1:B:2:SER:O[3_645]	0.99	1.21
1:A:277[B]:ARG:CZ	1:B:2:SER:O[3_645]	1.02	1.18
1:A:277[B]:ARG:NH2	1:B:2:SER:O[3_645]	1.72	0.48
1:A:277[B]:ARG:NH1	1:B:2:SER:C[3_645]	1.78	0.42
1:A:277[B]:ARG:NE	1:B:2:SER:O[3_645]	2.12	0.08
7:A:2082:HOH:O	7:B:2058:HOH:O[8_665]	2.12	0.08
1:A:277[B]:ARG:CZ	1:B:2:SER:C[3_645]	2.16	0.04

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	351/337 (104%)	340 (97%)	9 (3%)	2 (1%)	33	25
1	B	346/337 (103%)	331 (96%)	11 (3%)	4 (1%)	19	10
All	All	697/674 (103%)	671 (96%)	20 (3%)	6 (1%)	25	15

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	261	PRO
1	B	137	GLY
1	B	261	PRO

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Mol	Chain	Res	Type
1	A	334	ARG
1	B	138	GLN
1	B	335	ARG

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	285/267 (107%)	272 (95%)	13 (5%)	37	34
1	B	284/267 (106%)	270 (95%)	14 (5%)	35	31
All	All	569/534 (107%)	542 (95%)	27 (5%)	39	33

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	PHE
1	A	49	ASN
1	A	53	SER
1	A	96	LEU
1	A	117	GLU
1	A	118	ARG
1	A	157	LEU
1	A	181	ARG
1	A	198	ARG
1	A	226	LYS
1	A	283	TYR
1	A	301	LEU
1	A	334	ARG
1	B	48	VAL
1	B	74[A]	GLU
1	B	74[B]	GLU
1	B	181	ARG
1	B	198	ARG
1	B	222[A]	SER
1	B	222[B]	SER
1	B	256[A]	MET
1	B	256[B]	MET

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Mol	Chain	Res	Type
1	B	283	TYR
1	B	293[A]	MET
1	B	293[B]	MET
1	B	301	LEU
1	B	307	LEU

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 ⓘ

4 non-standard protein/DNA/RNA residues 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$
1	FHL	A	205	1	16,16,17	4.56	3 (18%)	16,18,20	1.94	2 (12%)
1	FHL	A	260	1	16,16,17	5.00	2 (12%)	16,18,20	2.94	6 (37%)
1	FHL	B	205	1	16,16,17	4.29	3 (18%)	16,18,20	2.12	3 (18%)
1	FHL	B	260	1	16,16,17	5.13	3 (18%)	16,18,20	2.20	4 (25%)

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
1	FHL	A	205	1	-	0/14/17/19	0/0/0/0
1	FHL	A	260	1	-	0/14/17/19	0/0/0/0
1	FHL	B	205	1	-	0/14/17/19	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	FHL	B	260	1	-	0/14/17/19	0/0/0/0

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	260	FHL	O-C	19.39	1.24	1.11
1	A	260	FHL	O-C	19.08	1.24	1.11
1	A	205	FHL	O-C	17.58	1.23	1.11
1	B	205	FHL	O-C	16.53	1.22	1.11
1	B	260	FHL	C5-C4	4.96	1.54	1.50
1	A	260	FHL	C5-C4	4.94	1.54	1.50
1	A	205	FHL	CA-C	3.21	1.54	1.48
1	B	205	FHL	C5-C4	2.93	1.52	1.50
1	B	260	FHL	CA-C	2.68	1.53	1.48
1	A	205	FHL	C5-C4	2.12	1.52	1.50
1	B	205	FHL	CA-C	2.04	1.52	1.48

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	205	FHL	C-CA-N	-6.69	107.15	113.83
1	A	260	FHL	C3-C4-NZ	-6.50	115.90	119.55
1	A	260	FHL	C-CA-N	-6.10	107.73	113.83
1	B	260	FHL	C-CA-N	-5.50	108.34	113.83
1	B	205	FHL	C3-C4-NZ	-5.02	116.73	119.55
1	B	205	FHL	C-CA-N	-4.77	109.06	113.83
1	B	260	FHL	C3-C4-NZ	-4.52	117.01	119.55
1	A	260	FHL	O3-C5-C4	4.42	122.46	110.83
1	B	260	FHL	O3-C5-C4	3.58	120.25	110.83
1	A	260	FHL	CD-CE-NZ	-3.51	105.80	110.75
1	A	260	FHL	CE-NZ-C4	3.22	128.72	119.90
1	A	260	FHL	C2-C3-C4	-2.85	109.99	114.67
1	B	205	FHL	CE-NZ-C4	2.40	126.49	119.90
1	A	205	FHL	CE-NZ-C4	2.33	126.30	119.90
1	B	260	FHL	CD-CE-NZ	-2.16	107.71	110.75

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 9 ligands modelled in this entry, 6 are monoatomic - leaving 3 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	PEG	A	1338	-	6,6,6	0.44	0	5,5,5	0.77	0
6	PGE	B	1338	-	9,9,9	0.63	0	8,8,8	0.40	0
6	PGE	B	1339	-	9,9,9	0.67	0	8,8,8	0.32	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	PEG	A	1338	-	-	0/4/4/4	0/0/0/0
6	PGE	B	1338	-	-	0/7/7/7	0/0/0/0
6	PGE	B	1339	-	-	0/7/7/7	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

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	334/337 (99%)	-0.59	2 (0%) 86 90	11, 16, 27, 58	0
1	B	331/337 (98%)	-0.63	3 (0%) 81 85	9, 15, 29, 50	3 (0%)
All	All	665/674 (98%)	-0.61	5 (0%) 81 87	9, 15, 28, 58	3 (0%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	335	ARG	3.6
1	A	334	ARG	3.0
1	B	144	ASP	2.3
1	B	335	ARG	2.2
1	B	222[A]	SER	2.0

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

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
1	FHL	A	260	17/18	0.13	2.42	2,12,14,15	0
1	FHL	B	205	17/18	0.13	2.13	15,20,29,33	0
1	FHL	A	205	17/18	0.11	1.80	10,13,19,22	0
1	FHL	B	260	17/18	0.10	1.47	5,12,14,16	0

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
6	PGE	B	1338	10/10	0.34	10.10	23,30,33,34	10
6	PGE	B	1339	10/10	0.21	6.97	36,43,47,47	0
3	CL	B	1337	1/1	0.08	0.61	23,23,23,23	0
4	PEG	A	1338	7/7	0.11	-0.09	29,31,35,36	0
3	CL	A	1337	1/1	0.08	-0.12	23,23,23,23	0
5	K	B	1340	1/1	0.03	-2.25	22,22,22,22	0
5	K	A	1339	1/1	0.02	-3.36	20,20,20,20	0
2	MG	A	1336	1/1	0.06	-3.41	10,10,10,10	0
2	MG	B	1336	1/1	0.04	-3.57	10,10,10,10	0

6.5 Other polymers ⓘ

There are no such residues in this entry.