



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 06:55 PM GMT

PDB ID : 4P7P
Title : Structure of the dehydratase domain of PpsC from Mycobacterium tuberculosis in complex with crotonyl-coenzyme A
Authors : Faille, A.; Mourey, L.; Pedelacq, J.D.
Deposited on : 2014-03-27
Resolution : 2.90 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 : unknown
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
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-20026982

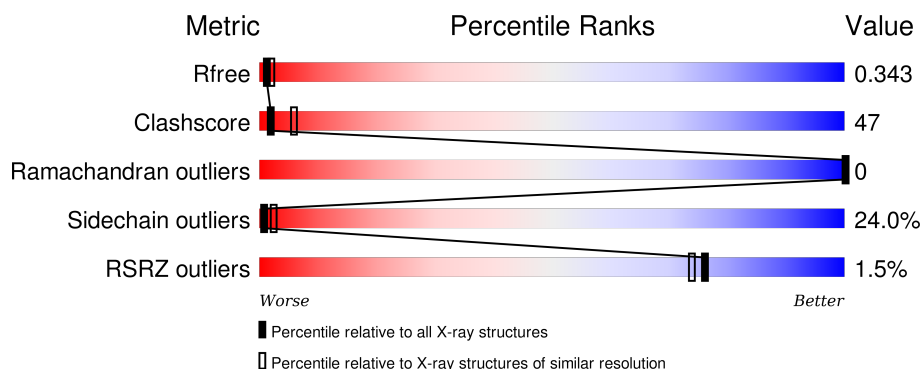
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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	325	

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	COO	A	1301	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 1967 atoms, of which 25 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.

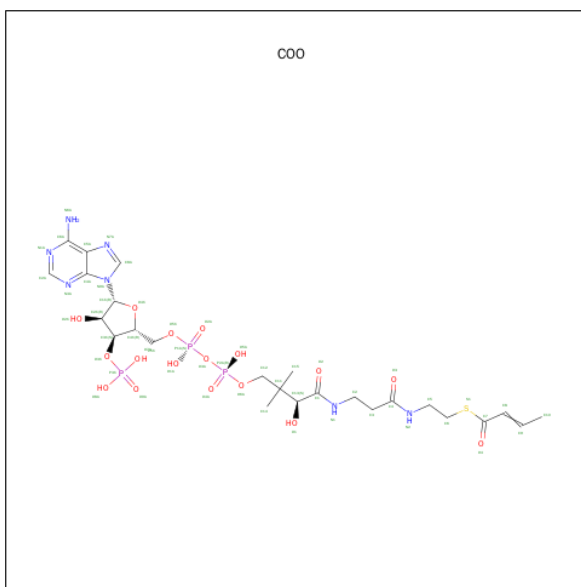
- Molecule 1 is a protein called Phthiocerol synthesis polyketide synthase type I PpsC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	259	Total	C	N	O	S	0	0	0
			1905	1193	332	373	7			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	900	MET	-	initiating methionine	UNP P96202
A	901	GLY	-	expression tag	UNP P96202
A	902	SER	-	expression tag	UNP P96202
A	903	SER	-	expression tag	UNP P96202
A	904	HIS	-	expression tag	UNP P96202
A	905	HIS	-	expression tag	UNP P96202
A	906	HIS	-	expression tag	UNP P96202
A	907	HIS	-	expression tag	UNP P96202
A	908	HIS	-	expression tag	UNP P96202
A	909	HIS	-	expression tag	UNP P96202
A	910	SER	-	expression tag	UNP P96202
A	911	SER	-	expression tag	UNP P96202
A	912	GLY	-	expression tag	UNP P96202
A	913	LEU	-	expression tag	UNP P96202
A	914	VAL	-	expression tag	UNP P96202
A	915	PRO	-	expression tag	UNP P96202
A	916	ARG	-	expression tag	UNP P96202
A	917	GLY	-	expression tag	UNP P96202
A	918	SER	-	expression tag	UNP P96202
A	919	HIS	-	expression tag	UNP P96202
A	920	MET	-	expression tag	UNP P96202
A	959	PHE	HIS	engineered mutation	UNP P96202
A	1223	GLY	-	expression tag	UNP P96202
A	1224	SER	-	expression tag	UNP P96202

- Molecule 2 is CROTONYL COENZYME A (three-letter code: COO) (formula: C₂₅H₄₀N₇O₁₇P₃S).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	P	S	0	0
			52	15	25	2	8	1	1		

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Na	0	0
			1	1		

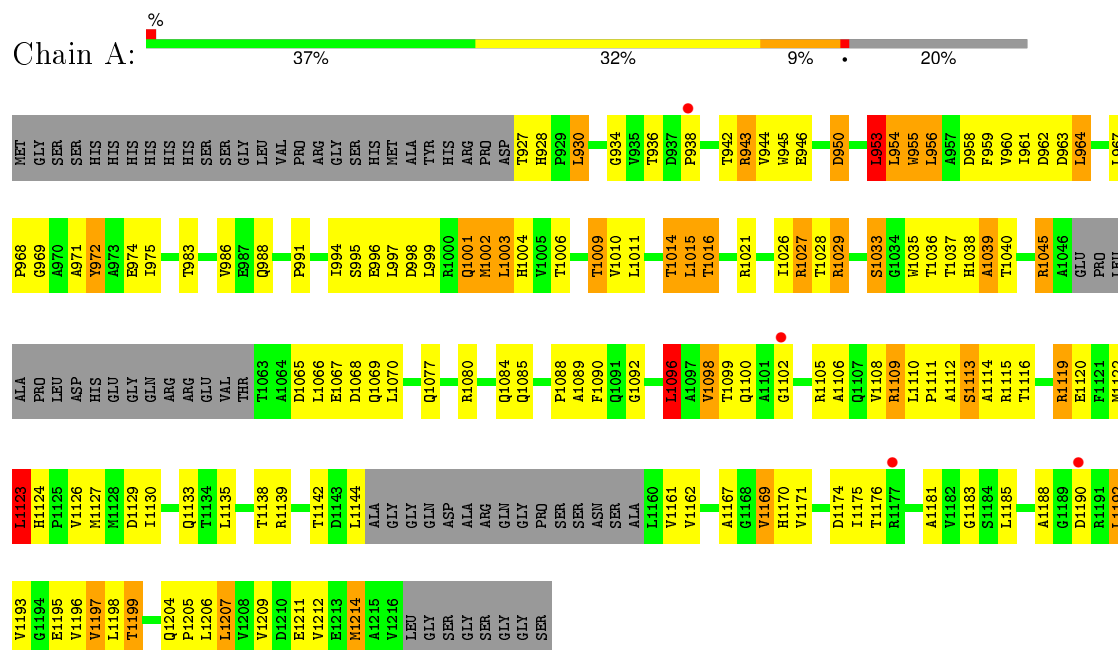
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	9	Total	O	0	0
			9	9		

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 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: Phthiocerol synthesis polyketide synthase type I PpsC



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	83.85Å 83.85Å 166.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.92 – 2.90 46.34 – 2.90	Depositor EDS
% Data completeness (in resolution range)	93.6 (41.92-2.90) 94.0 (46.34-2.90)	Depositor EDS
R_{merge}	0.28	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.03 (at 2.91Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.272 , 0.318 0.303 , 0.343	Depositor DCC
R_{free} test set	1300 reflections (11.11%)	DCC
Wilson B-factor (Å ²)	80.1	Xtriage
Anisotropy	0.600	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 25.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	3 of 13005 reflections (0.023%)	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	1967	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.57% 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, 0.375 respectively for untwinned datasets, and 0.333, 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, COO

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	0.96	1/1936 (0.1%)	1.13	8/2651 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1039	ALA	CA-CB	-5.14	1.41	1.52

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	954	LEU	CB-CA-C	-5.77	99.24	110.20
1	A	954	LEU	N-CA-C	5.54	125.96	111.00
1	A	964	LEU	CB-CG-CD1	-5.52	101.62	111.00
1	A	1001	GLN	CB-CA-C	-5.44	99.51	110.40
1	A	1096	LEU	CA-CB-CG	5.34	127.58	115.30
1	A	1188	ALA	N-CA-CB	-5.25	102.75	110.10
1	A	1123	LEU	CA-CB-CG	5.21	127.28	115.30
1	A	953	LEU	N-CA-C	-5.15	97.09	111.00

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	1905	0	1861	174	0
2	A	27	25	25	21	0
3	A	1	0	0	0	0
4	A	9	0	0	0	0
All	All	1942	25	1886	179	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 47.

All (179) 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:954:LEU:O	1:A:954:LEU:HD12	1.41	1.21
1:A:999:LEU:HD13	2:A:1301:COO:C6	1.79	1.11
1:A:999:LEU:HD13	2:A:1301:COO:H61	1.31	1.11
1:A:975:ILE:HG21	1:A:1026:ILE:HD13	1.34	1.07
1:A:955:TRP:HE3	1:A:955:TRP:O	1.44	0.99
1:A:959:PHE:HD2	1:A:968:PRO:HG3	1.24	0.99
1:A:1192:LEU:CD1	1:A:1192:LEU:N	2.30	0.95
1:A:954:LEU:O	1:A:954:LEU:CD1	2.15	0.94
1:A:967:LEU:HD23	1:A:972:TYR:CE1	2.05	0.91
1:A:1192:LEU:N	1:A:1192:LEU:HD12	1.85	0.90
1:A:943:ARG:HG2	1:A:945:TRP:CH2	2.07	0.90
1:A:959:PHE:CD2	1:A:968:PRO:HG3	2.06	0.90
1:A:961:ILE:HD11	2:A:1301:COO:C10	2.04	0.87
1:A:1124:HIS:CE1	1:A:1126:VAL:HG22	2.10	0.86
1:A:958:ASP:OD2	1:A:1088:PRO:HD2	1.74	0.86
1:A:955:TRP:O	1:A:955:TRP:CE3	2.30	0.85
1:A:961:ILE:HD11	2:A:1301:COO:H103	1.59	0.83
1:A:1092:GLY:HA3	1:A:1130:ILE:HD13	1.57	0.83
1:A:999:LEU:CD1	2:A:1301:COO:H61	2.09	0.82
1:A:943:ARG:HG2	1:A:945:TRP:CZ2	2.15	0.82
1:A:1123:LEU:O	1:A:1123:LEU:HD12	1.80	0.82
1:A:1192:LEU:H	1:A:1192:LEU:HD13	1.46	0.80
1:A:1080:ARG:NH1	1:A:1085:GLN:HG3	1.97	0.80
1:A:1130:ILE:HA	1:A:1133:GLN:HG3	1.65	0.79
1:A:961:ILE:CD1	2:A:1301:COO:C10	2.61	0.78
1:A:975:ILE:HG21	1:A:1026:ILE:CD1	2.15	0.76
2:A:1301:COO:O2	2:A:1301:COO:H153	1.84	0.75
1:A:975:ILE:CG2	1:A:1026:ILE:HD13	2.14	0.75
1:A:991:PRO:HB3	1:A:1045:ARG:NH2	2.02	0.75
1:A:1209:VAL:HG21	1:A:1212:VAL:CG2	2.18	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1092:GLY:HA3	1:A:1130:ILE:CD1	2.19	0.72
1:A:962:ASP:O	1:A:963:ASP:HB2	1.89	0.72
1:A:1080:ARG:HH12	1:A:1085:GLN:HG3	1.54	0.71
1:A:1192:LEU:H	1:A:1192:LEU:CD1	1.99	0.71
2:A:1301:COO:O1	2:A:1301:COO:P2A	2.49	0.71
1:A:954:LEU:HD13	1:A:955:TRP:CD1	2.26	0.70
1:A:999:LEU:HD13	2:A:1301:COO:C5	2.21	0.70
1:A:1138:THR:HG21	1:A:1185:LEU:HB3	1.73	0.70
1:A:956:LEU:HD12	1:A:956:LEU:O	1.91	0.70
1:A:1065:ASP:OD2	1:A:1065:ASP:O	2.10	0.70
1:A:972:TYR:HD2	1:A:1039:ALA:CB	2.05	0.69
1:A:991:PRO:HB3	1:A:1045:ARG:HH22	1.58	0.69
1:A:1070:LEU:CD1	1:A:1098:VAL:HG12	2.23	0.69
1:A:1209:VAL:CG2	1:A:1212:VAL:HG23	2.21	0.68
1:A:1002:MET:N	2:A:1301:COO:O3	2.26	0.68
1:A:1209:VAL:HG21	1:A:1212:VAL:HG23	1.77	0.67
1:A:975:ILE:CG2	1:A:1026:ILE:CD1	2.73	0.67
2:A:1301:COO:HO1	2:A:1301:COO:P2A	2.17	0.67
1:A:942:THR:CG2	1:A:1014:THR:HG22	2.24	0.66
1:A:960:VAL:HG23	1:A:961:ILE:N	2.09	0.66
1:A:1171:VAL:HG13	1:A:1175:ILE:HD11	1.76	0.66
1:A:1175:ILE:HD12	1:A:1175:ILE:H	1.59	0.66
1:A:961:ILE:CD1	2:A:1301:COO:H101	2.25	0.66
1:A:1014:THR:OG1	1:A:1027:ARG:NH1	2.27	0.66
1:A:1010:VAL:HG23	1:A:1029:ARG:HB3	1.78	0.66
1:A:1119:ARG:CD	1:A:1119:ARG:H	2.09	0.65
1:A:954:LEU:O	1:A:955:TRP:CG	2.50	0.65
1:A:1003:LEU:HG	1:A:1003:LEU:O	1.97	0.65
1:A:961:ILE:HD13	2:A:1301:COO:H101	1.78	0.64
1:A:953:LEU:O	1:A:956:LEU:HB2	1.97	0.63
1:A:1090:PHE:CZ	1:A:1126:VAL:HG12	2.33	0.63
1:A:986:VAL:HG21	1:A:991:PRO:HB3	1.80	0.63
1:A:1124:HIS:CE1	1:A:1126:VAL:CG2	2.81	0.62
1:A:967:LEU:HD23	1:A:972:TYR:HE1	1.63	0.62
1:A:1001:GLN:HG2	1:A:1001:GLN:O	2.00	0.62
1:A:1070:LEU:HD12	1:A:1098:VAL:HG12	1.81	0.61
1:A:986:VAL:HG21	1:A:991:PRO:CB	2.30	0.61
1:A:1001:GLN:O	1:A:1038:HIS:ND1	2.24	0.61
1:A:1015:LEU:HD23	1:A:1016:THR:N	2.16	0.61
1:A:1109:ARG:O	1:A:1111:PRO:HD3	2.00	0.61
1:A:1123:LEU:C	1:A:1123:LEU:HD12	2.21	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:999:LEU:HB3	2:A:1301:COO:H51	1.84	0.58
1:A:972:TYR:HD2	1:A:1039:ALA:HB2	1.67	0.58
1:A:1014:THR:HG1	1:A:1027:ARG:HH12	1.50	0.58
1:A:972:TYR:CD2	1:A:1039:ALA:HB2	2.38	0.58
1:A:1112:ALA:HA	1:A:1115:ARG:HG3	1.85	0.58
1:A:942:THR:HG21	1:A:1014:THR:HG22	1.86	0.58
1:A:983:THR:OG1	1:A:1015:LEU:HD13	2.04	0.58
1:A:1181:ALA:HB2	1:A:1198:LEU:HD12	1.86	0.57
1:A:1066:LEU:O	1:A:1100:GLN:NE2	2.29	0.56
1:A:1077:GLN:HA	1:A:1077:GLN:OE1	2.06	0.56
1:A:1175:ILE:HG23	1:A:1198:LEU:HD23	1.88	0.55
1:A:1119:ARG:HD2	1:A:1119:ARG:H	1.70	0.55
1:A:1070:LEU:HD11	1:A:1098:VAL:HG12	1.89	0.55
1:A:991:PRO:CB	1:A:1045:ARG:NH2	2.69	0.54
1:A:1110:LEU:HB2	1:A:1127:MET:HE1	1.90	0.54
1:A:1130:ILE:CA	1:A:1133:GLN:HG3	2.36	0.54
1:A:998:ASP:HB2	1:A:1040:THR:HG23	1.89	0.54
1:A:1161:VAL:HG22	1:A:1214:MET:HB3	1.90	0.53
1:A:1199:THR:HG22	1:A:1204:GLN:O	2.07	0.53
1:A:953:LEU:O	1:A:956:LEU:CB	2.56	0.53
1:A:1162:VAL:HG12	2:A:1301:COO:C10	2.39	0.53
1:A:928:HIS:HA	1:A:1120:GLU:HG2	1.89	0.53
1:A:1183:GLY:HA3	1:A:1196:VAL:HG22	1.91	0.53
1:A:1110:LEU:HB2	1:A:1127:MET:CE	2.38	0.53
1:A:1162:VAL:HG12	2:A:1301:COO:H102	1.89	0.53
1:A:1124:HIS:HE1	1:A:1126:VAL:HG22	1.68	0.53
1:A:1006:THR:O	1:A:1009:THR:OG1	2.27	0.52
1:A:1199:THR:HG22	1:A:1205:PRO:HA	1.90	0.52
1:A:1124:HIS:ND1	1:A:1126:VAL:HG22	2.23	0.52
1:A:1130:ILE:HA	1:A:1133:GLN:NE2	2.24	0.52
1:A:969:GLY:N	2:A:1301:COO:O4	2.43	0.51
1:A:997:LEU:HD23	1:A:997:LEU:C	2.30	0.51
1:A:972:TYR:CD2	1:A:1039:ALA:CB	2.91	0.51
1:A:1175:ILE:HG23	1:A:1198:LEU:CD2	2.41	0.51
1:A:1129:ASP:O	1:A:1133:GLN:HG3	2.11	0.51
1:A:1109:ARG:HB3	1:A:1109:ARG:CZ	2.41	0.51
1:A:960:VAL:CG2	1:A:961:ILE:N	2.75	0.50
1:A:1010:VAL:CG2	1:A:1029:ARG:HB3	2.42	0.50
1:A:1070:LEU:HD11	1:A:1098:VAL:CG1	2.41	0.50
1:A:959:PHE:CD2	1:A:968:PRO:CG	2.89	0.49
1:A:1066:LEU:O	1:A:1100:GLN:HG2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1301:COO:O1	2:A:1301:COO:O5A	2.30	0.49
1:A:1067:GLU:HA	1:A:1099:THR:HA	1.94	0.49
1:A:1089:ALA:HB2	1:A:1113:SER:HB2	1.95	0.49
1:A:1003:LEU:HB2	1:A:1038:HIS:CE1	2.47	0.49
1:A:934:GLY:HA2	1:A:944:VAL:O	2.12	0.49
1:A:972:TYR:CD1	1:A:972:TYR:N	2.81	0.49
1:A:994:ILE:HB	1:A:1169:VAL:HG13	1.95	0.48
1:A:1096:LEU:HD23	1:A:1106:ALA:HB2	1.95	0.48
1:A:1108:VAL:HG21	1:A:1198:LEU:HD11	1.94	0.48
1:A:1183:GLY:CA	1:A:1196:VAL:HG22	2.44	0.48
1:A:1167:ALA:HB3	1:A:1211:GLU:HB2	1.95	0.48
1:A:972:TYR:HD1	1:A:972:TYR:N	2.11	0.48
1:A:1033:SER:O	1:A:1035:TRP:CE3	2.67	0.47
1:A:1135:LEU:HD21	1:A:1196:VAL:HG23	1.97	0.47
1:A:986:VAL:HG21	1:A:991:PRO:CG	2.45	0.47
1:A:1130:ILE:HA	1:A:1133:GLN:CG	2.39	0.47
1:A:955:TRP:CE3	1:A:955:TRP:C	2.87	0.47
1:A:961:ILE:CD1	2:A:1301:COO:H103	2.35	0.47
1:A:953:LEU:HD12	1:A:954:LEU:N	2.30	0.46
1:A:961:ILE:HD11	2:A:1301:COO:C9	2.45	0.46
1:A:1102:GLY:O	1:A:1138:THR:HG23	2.15	0.46
1:A:1015:LEU:HD23	1:A:1015:LEU:C	2.34	0.46
1:A:1089:ALA:HB1	1:A:1114:ALA:HB2	1.97	0.46
2:A:1301:COO:O2	2:A:1301:COO:C15	2.56	0.46
1:A:1080:ARG:HA	1:A:1084:GLN:O	2.15	0.46
1:A:936:THR:O	1:A:938:PRO:HD3	2.16	0.46
1:A:1212:VAL:CG1	1:A:1214:MET:CE	2.94	0.46
1:A:996:GLU:HG3	1:A:997:LEU:N	2.30	0.46
1:A:959:PHE:HD2	1:A:968:PRO:CG	2.12	0.46
1:A:1027:ARG:HD3	1:A:1035:TRP:HB3	1.97	0.46
1:A:971:ALA:O	1:A:975:ILE:CD1	2.64	0.46
1:A:1127:MET:O	1:A:1130:ILE:HB	2.17	0.45
1:A:1212:VAL:HG11	1:A:1214:MET:CE	2.46	0.45
1:A:942:THR:HG22	1:A:943:ARG:N	2.32	0.45
1:A:1185:LEU:HD12	1:A:1193:VAL:O	2.17	0.45
1:A:1080:ARG:HG3	1:A:1080:ARG:NH1	2.32	0.45
1:A:967:LEU:HA	1:A:968:PRO:HD3	1.77	0.44
1:A:1033:SER:O	1:A:1035:TRP:CZ3	2.70	0.44
1:A:1138:THR:HG21	1:A:1185:LEU:H	1.82	0.43
1:A:1175:ILE:HG22	1:A:1175:ILE:O	2.17	0.43
1:A:1212:VAL:CG1	1:A:1214:MET:HE2	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1139:ARG:HA	1:A:1142:THR:OG1	2.18	0.43
1:A:1112:ALA:O	1:A:1115:ARG:HB2	2.18	0.43
1:A:1110:LEU:HD13	1:A:1176:THR:HG22	2.01	0.43
1:A:959:PHE:HE2	1:A:968:PRO:HA	1.84	0.43
1:A:1170:HIS:O	1:A:1207:LEU:HA	2.19	0.42
1:A:971:ALA:O	1:A:975:ILE:HD13	2.19	0.42
1:A:1174:ASP:OD2	1:A:1176:THR:OG1	2.34	0.42
1:A:942:THR:HG22	1:A:1014:THR:HG22	2.00	0.42
1:A:950:ASP:OD2	1:A:950:ASP:N	2.52	0.42
1:A:964:LEU:HA	1:A:964:LEU:HD23	1.87	0.42
1:A:1127:MET:HG2	1:A:1175:ILE:HG22	2.01	0.42
1:A:953:LEU:HD12	1:A:953:LEU:C	2.41	0.42
1:A:1126:VAL:HG23	1:A:1127:MET:N	2.35	0.42
1:A:1123:LEU:C	1:A:1123:LEU:CD1	2.86	0.42
1:A:991:PRO:CB	1:A:1045:ARG:HH21	2.32	0.41
1:A:975:ILE:CG2	1:A:1026:ILE:HD11	2.51	0.41
1:A:1068:ASP:OD1	1:A:1100:GLN:NE2	2.37	0.41
1:A:1196:VAL:CG1	1:A:1197:VAL:N	2.84	0.41
1:A:930:LEU:HD12	1:A:930:LEU:HA	1.80	0.41
1:A:1138:THR:HG21	1:A:1185:LEU:CB	2.48	0.41
1:A:1090:PHE:CE2	1:A:1126:VAL:HG12	2.56	0.41
1:A:1209:VAL:O	1:A:1209:VAL:CG2	2.68	0.40
1:A:974:GLU:OE1	1:A:974:GLU:HA	2.22	0.40
1:A:986:VAL:O	1:A:986:VAL:HG13	2.21	0.40
1:A:1138:THR:CG2	1:A:1185:LEU:HB3	2.48	0.40
1:A:998:ASP:HB2	1:A:1040:THR:CG2	2.52	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	253/325 (78%)	244 (96%)	9 (4%)	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	196/254 (77%)	149 (76%)	47 (24%)	1	2

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	927	THR
1	A	930	LEU
1	A	943	ARG
1	A	946	GLU
1	A	950	ASP
1	A	953	LEU
1	A	955	TRP
1	A	956	LEU
1	A	972	TYR
1	A	988	GLN
1	A	995	SER
1	A	1002	MET
1	A	1003	LEU
1	A	1004	HIS
1	A	1009	THR
1	A	1011	LEU
1	A	1014	THR
1	A	1015	LEU
1	A	1016	THR
1	A	1021	ARG
1	A	1027	ARG
1	A	1028	THR
1	A	1029	ARG

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Mol	Chain	Res	Type
1	A	1033	SER
1	A	1036	THR
1	A	1037	THR
1	A	1045	ARG
1	A	1069	GLN
1	A	1096	LEU
1	A	1098	VAL
1	A	1105	ARG
1	A	1109	ARG
1	A	1113	SER
1	A	1116	THR
1	A	1119	ARG
1	A	1122	MET
1	A	1123	LEU
1	A	1144	LEU
1	A	1169	VAL
1	A	1190	ASP
1	A	1192	LEU
1	A	1195	GLU
1	A	1197	VAL
1	A	1199	THR
1	A	1206	LEU
1	A	1207	LEU
1	A	1214	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1170	HIS

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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	COO	A	1301	-	21,26,55	0.66	0	28,35,81	1.03	2 (7%)

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	COO	A	1301	-	-	0/32/33/70	0/0/0/3

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1301	COO	C10-C9-C8	-2.47	119.97	125.42
2	A	1301	COO	O5A-P2A-O4A	2.86	119.95	110.63

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1301	COO	21	0

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 [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	259/325 (79%)	0.30	4 (1%) 76 74	17, 33, 60, 94	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1190	ASP	3.0
1	A	1102	GLY	2.8
1	A	938	PRO	2.2
1	A	1177	ARG	2.1

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
2	COO	A	1301	27/53	0.83	0.26	1.36	28,49,68,87	52
3	NA	A	1302	1/1	0.74	0.57	-	21,21,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.