Full wwPDB X-ray Structure Validation Report (i)

Feb 20, 2024 – 09:57 AM EST

PDB ID	:	4MM1
Title	:	GGGPS from Methanothermobacter thermautotrophicus
Authors	:	Rajendran, C.; Peterhoff, D.; Beer, B.; Kumpula, E.P.; Kapetaniou, E.; Guldan, H.; Wierenga, R.K.; Sterner, R.; Babinger, P.
Deposited on	:	2013-09-07
Resolution	:	2.80 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 (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

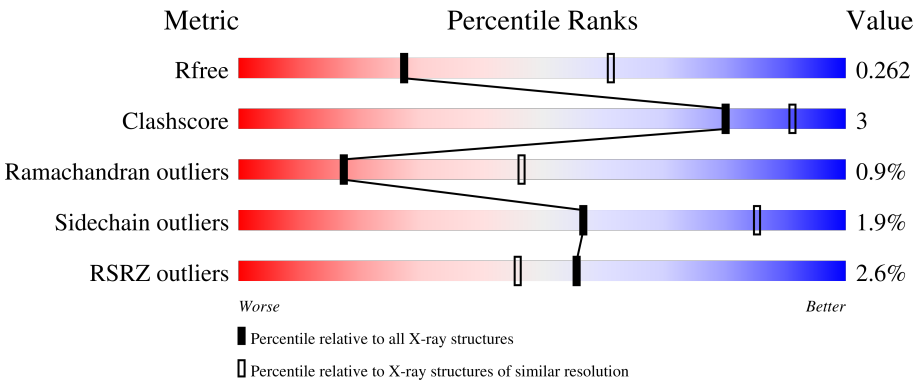
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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 <=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	250	91% • • 5%
1	B	250	4% 77% 8% • 14%
1	C	250	2% 88% 6% 6%
1	D	250	2% 83% • • 12%
1	E	250	3% 83% 7% 10%

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Mol	Chain	Length	Quality of chain
1	F	250	3% 77% 5% • 17%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 19944 atoms, of which 9883 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 Geranylgeranylglyceryl phosphate synthase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	238	3482	1114	1722	298	338	10	0	0	0
1	B	215	3187	1013	1588	266	312	8	0	0	0
1	C	234	3386	1086	1673	292	324	11	0	0	0
1	D	220	3216	1032	1592	278	305	9	0	0	0
1	E	225	3330	1057	1663	282	318	10	0	0	0
1	F	207	3029	967	1519	259	275	9	0	0	0

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP O26652
A	2	PHE	-	expression tag	UNP O26652
A	3	LYS	-	expression tag	UNP O26652
A	249	LEU	-	expression tag	UNP O26652
A	250	GLU	-	expression tag	UNP O26652
B	1	MET	-	expression tag	UNP O26652
B	2	PHE	-	expression tag	UNP O26652
B	3	LYS	-	expression tag	UNP O26652
B	249	LEU	-	expression tag	UNP O26652
B	250	GLU	-	expression tag	UNP O26652
C	1	MET	-	expression tag	UNP O26652
C	2	PHE	-	expression tag	UNP O26652
C	3	LYS	-	expression tag	UNP O26652
C	249	LEU	-	expression tag	UNP O26652
C	250	GLU	-	expression tag	UNP O26652
D	1	MET	-	expression tag	UNP O26652
D	2	PHE	-	expression tag	UNP O26652

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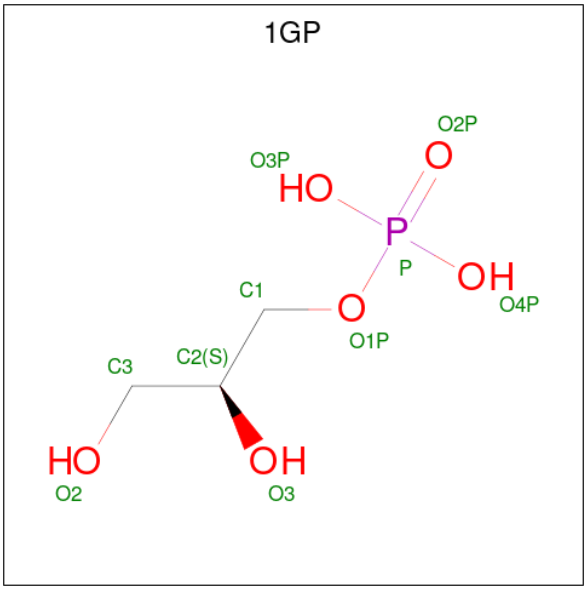
Chain	Residue	Modelled	Actual	Comment	Reference
D	3	LYS	-	expression tag	UNP O26652
D	249	LEU	-	expression tag	UNP O26652
D	250	GLU	-	expression tag	UNP O26652
E	1	MET	-	expression tag	UNP O26652
E	2	PHE	-	expression tag	UNP O26652
E	3	LYS	-	expression tag	UNP O26652
E	249	LEU	-	expression tag	UNP O26652
E	250	GLU	-	expression tag	UNP O26652
F	1	MET	-	expression tag	UNP O26652
F	2	PHE	-	expression tag	UNP O26652
F	3	LYS	-	expression tag	UNP O26652
F	249	LEU	-	expression tag	UNP O26652
F	250	GLU	-	expression tag	UNP O26652

- Molecule 2 is SN-GLYCEROL-1-PHOSPHATE (three-letter code: 1GP) (formula: $C_3H_9O_6P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	H	O	P		
2	A	1	17	3	7	6	1	0	0
			Total	C	H	O	P		
2	B	1	17	3	7	6	1	0	0
			Total	C	H	O	P		
2	C	1	17	3	7	6	1	0	0
			Total	C	H	O	P		
2	D	1	17	3	7	6	1	0	0

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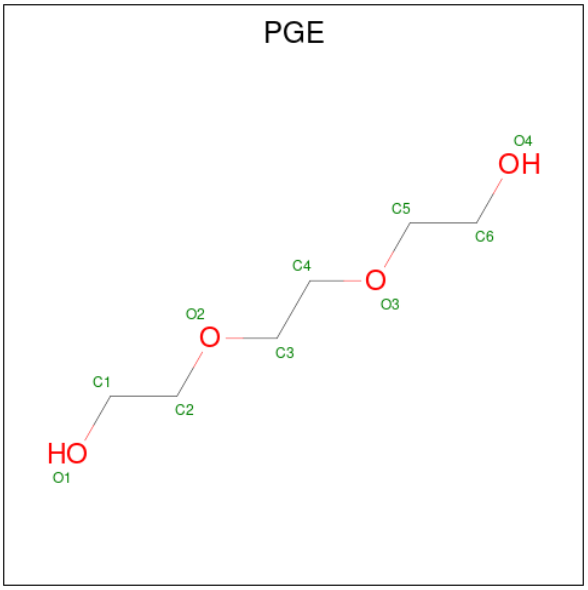
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	H	O	P		
2	E	1	17	3	7	6	1	0	0
			Total	C	H	O	P		
2	F	1	17	3	7	6	1	0	0

- Molecule 3 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
3	A	1	24	6	14	4	0	0
			Total	C	H	O		
3	B	1	24	6	14	4	0	0
			Total	C	H	O		
3	C	1	24	6	14	4	0	0
			Total	C	H	O		
3	D	1	24	6	14	4	0	0
			Total	C	H	O		
3	E	1	24	6	14	4	0	0
			Total	C	H	O		
3	F	1	24	6	14	4	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	19	19	19	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	B	10	10	10	0	0
			Total	O		
4	C	15	15	15	0	0
			Total	O		
4	D	9	9	9	0	0
			Total	O		
4	E	11	11	11	0	0
			Total	O		
4	F	4	4	4	0	0

3 Residue-property plots (i)

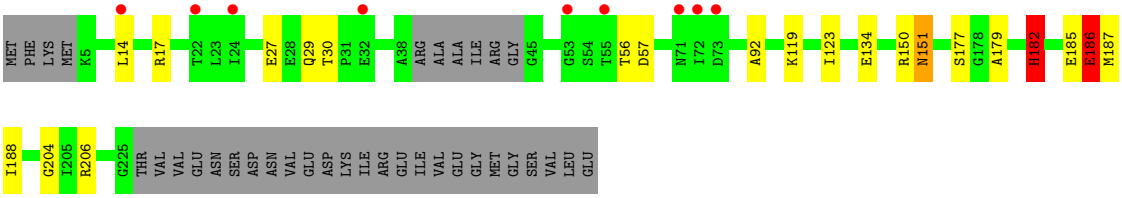
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Geranylgeranylglyceryl phosphate synthase

Chain A: 91% • • 5%

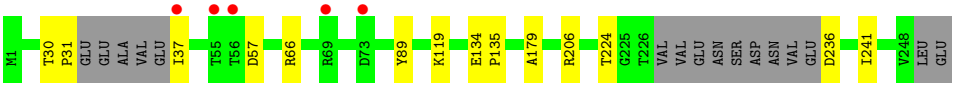
• Molecule 1: Geranylgeranylglyceryl phosphate synthase

Chain B: 4% 77% 8% • 14%



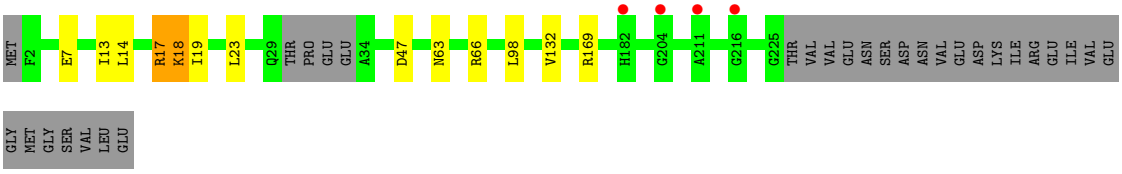
• Molecule 1: Geranylgeranylglyceryl phosphate synthase

Chain C: 2% 88% 6% 6%



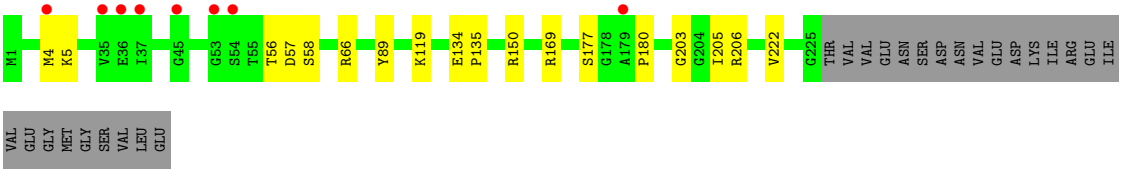
• Molecule 1: Geranylgeranylglyceryl phosphate synthase

Chain D: 2% 83% • • 12%



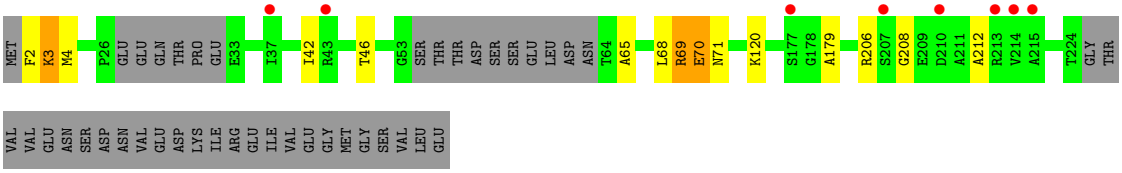
• Molecule 1: Geranylgeranylglyceryl phosphate synthase

Chain E: 3% 83% 7% 10%



• Molecule 1: Geranylgeranylglyceryl phosphate synthase

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	143.68Å 148.76Å 91.50Å 90.00° 109.50° 90.00°	Depositor
Resolution (Å)	46.56 – 2.80 48.86 – 2.80	Depositor EDS
% Data completeness (in resolution range)	95.0 (46.56-2.80) 95.1 (48.86-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.21 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1417)	Depositor
R, R_{free}	0.200 , 0.271 0.222 , 0.262	Depositor DCC
R_{free} test set	2120 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	42.2	Xtriage
Anisotropy	0.363	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 40.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.45$, $< L^2 > = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.89	EDS
Total number of atoms	19944	wwPDB-VP
Average B, all atoms (Å^2)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.38% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: 1GP, PGE

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	0.24	0/1790	0.52	1/2430 (0.0%)
1	B	0.24	0/1628	0.54	1/2213 (0.0%)
1	C	0.26	0/1742	0.51	0/2364
1	D	0.24	0/1653	0.48	0/2243
1	E	0.25	0/1698	0.47	0/2307
1	F	0.24	0/1538	0.51	0/2088
All	All	0.25	0/10049	0.50	2/13645 (0.0%)

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	B	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	44	GLY	N-CA-C	-6.45	96.97	113.10
1	B	186	GLU	N-CA-C	-5.20	96.96	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	182	HIS	Peptide

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	1760	1722	1744	6	0
1	B	1599	1588	1585	10	0
1	C	1713	1673	1694	4	0
1	D	1624	1592	1618	13	0
1	E	1667	1663	1660	13	0
1	F	1510	1519	1501	7	0
2	A	10	7	7	0	0
2	B	10	7	7	0	0
2	C	10	7	7	0	0
2	D	10	7	7	0	0
2	E	10	7	7	2	0
2	F	10	7	7	0	0
3	A	10	14	14	0	0
3	B	10	14	14	0	0
3	C	10	14	14	0	0
3	D	10	14	14	0	0
3	E	10	14	14	0	0
3	F	10	14	14	0	0
4	A	19	0	0	0	0
4	B	10	0	0	0	0
4	C	15	0	0	0	0
4	D	9	0	0	0	0
4	E	11	0	0	2	0
4	F	4	0	0	0	0
All	All	10061	9883	9928	53	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 3.

All (53) 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:D:13:ILE:HG23	1:D:17:ARG:HE	1.08	1.09
1:D:14:LEU:HG	1:D:19:ILE:HD11	1.35	1.04
1:D:13:ILE:HG23	1:D:17:ARG:NE	1.80	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:186:GLU:O	1:B:188:ILE:N	2.05	0.90
1:D:14:LEU:HG	1:D:19:ILE:CD1	2.10	0.80
1:D:13:ILE:CG2	1:D:17:ARG:HE	1.95	0.76
1:D:13:ILE:CG2	1:D:17:ARG:NE	2.49	0.75
1:E:66:ARG:NH1	1:E:89:TYR:OH	2.23	0.71
1:B:150:ARG:NH1	1:B:182:HIS:O	2.28	0.67
1:A:226:THR:O	1:A:227:VAL:HG13	1.98	0.64
1:B:151:ASN:N	1:B:151:ASN:OD1	2.33	0.61
1:A:7:GLU:OE1	1:A:169:ARG:NH1	2.34	0.61
1:C:179:ALA:O	1:C:206:ARG:NH2	2.34	0.60
1:E:169:ARG:NH2	4:E:404:HOH:O	2.36	0.57
1:A:63:ASN:OD1	1:A:66:ARG:NH2	2.38	0.56
1:E:206:ARG:NH1	4:E:402:HOH:O	2.41	0.54
1:E:4:MET:HG2	1:E:5:LYS:H	1.73	0.53
1:D:14:LEU:CG	1:D:19:ILE:CD1	2.84	0.53
1:F:69:ARG:HA	1:F:71:ASN:N	2.22	0.53
1:C:30:THR:O	1:C:31:PRO:C	2.46	0.53
1:B:179:ALA:O	1:B:206:ARG:NH2	2.42	0.52
1:F:42:ILE:HA	1:F:46:THR:HG22	1.95	0.49
1:D:17:ARG:NH2	1:D:47:ASP:OD1	2.41	0.49
1:F:69:ARG:HA	1:F:71:ASN:H	1.77	0.49
1:D:63:ASN:OD1	1:D:66:ARG:NH2	2.46	0.49
1:E:180:PRO:O	1:E:206:ARG:NH2	2.41	0.48
1:B:134:GLU:HB3	1:B:150:ARG:HG3	1.96	0.48
1:B:27:GLU:HB2	1:B:56:THR:HG21	1.94	0.47
1:B:56:THR:HG22	1:B:57:ASP:N	2.29	0.47
1:A:226:THR:O	1:A:227:VAL:HG22	2.15	0.47
1:E:56:THR:HG22	1:E:57:ASP:N	2.30	0.47
1:B:14:LEU:HA	1:B:17:ARG:O	2.15	0.47
1:E:4:MET:HG2	1:E:5:LYS:N	2.30	0.47
1:B:185:GLU:O	1:B:186:GLU:CB	2.63	0.46
1:F:179:ALA:O	1:F:206:ARG:NH2	2.44	0.46
1:D:7:GLU:OE1	1:D:169:ARG:NH1	2.49	0.45
1:D:14:LEU:HD21	1:D:19:ILE:HD12	2.00	0.44
1:E:4:MET:CG	1:E:5:LYS:H	2.31	0.44
1:E:134:GLU:HG2	1:E:135:PRO:HA	2.00	0.44
1:E:205:ILE:HD13	1:E:222:VAL:HG13	2.00	0.44
1:E:134:GLU:HB3	1:E:150:ARG:HG3	1.99	0.43
1:C:66:ARG:NH1	1:C:89:TYR:OH	2.52	0.42
1:B:92:ALA:CA	1:B:123:ILE:HD11	2.49	0.42
1:F:3:LYS:HA	1:F:4:MET:HA	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:226:THR:C	1:A:227:VAL:HG22	2.39	0.42
1:D:18:LYS:HD3	1:D:18:LYS:HA	1.76	0.42
1:E:177:SER:HA	2:E:301:1GP:H12	2.01	0.42
1:C:224:THR:HG21	1:C:241:ILE:HD13	2.02	0.41
1:A:42:ILE:HA	1:A:46:THR:HG22	2.02	0.41
1:D:98:LEU:HD11	1:D:132:VAL:HG23	2.02	0.41
1:F:208:GLY:O	1:F:212:ALA:N	2.52	0.41
1:E:203:GLY:HA2	2:E:301:1GP:H11	2.02	0.40
1:F:69:ARG:N	1:F:70:GLU:CB	2.84	0.40

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	234/250 (94%)	226 (97%)	7 (3%)	1 (0%)	34	66
1	B	211/250 (84%)	201 (95%)	6 (3%)	4 (2%)	8	26
1	C	228/250 (91%)	217 (95%)	9 (4%)	2 (1%)	17	46
1	D	216/250 (86%)	209 (97%)	6 (3%)	1 (0%)	29	61
1	E	223/250 (89%)	216 (97%)	7 (3%)	0	100	100
1	F	201/250 (80%)	193 (96%)	4 (2%)	4 (2%)	7	24
All	All	1313/1500 (88%)	1262 (96%)	39 (3%)	12 (1%)	17	46

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	182	HIS
1	B	186	GLU
1	B	187	MET

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Mol	Chain	Res	Type
1	C	135	PRO
1	F	65	ALA
1	F	69	ARG
1	D	18	LYS
1	F	68	LEU
1	A	226	THR
1	B	204	GLY
1	F	70	GLU
1	C	134	GLU

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	178/196 (91%)	175 (98%)	3 (2%)	60	87
1	B	165/196 (84%)	160 (97%)	5 (3%)	41	75
1	C	171/196 (87%)	167 (98%)	4 (2%)	50	82
1	D	162/196 (83%)	160 (99%)	2 (1%)	71	92
1	E	169/196 (86%)	167 (99%)	2 (1%)	71	92
1	F	146/196 (74%)	143 (98%)	3 (2%)	53	84
All	All	991/1176 (84%)	972 (98%)	19 (2%)	57	85

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	GLU
1	A	57	ASP
1	A	227	VAL
1	B	29	GLN
1	B	30	THR
1	B	119	LYS
1	B	151	ASN
1	B	177	SER
1	C	37	ILE

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Mol	Chain	Res	Type
1	C	57	ASP
1	C	119	LYS
1	C	236	ASP
1	D	17	ARG
1	D	23	LEU
1	E	58	SER
1	E	119	LYS
1	F	2	PHE
1	F	3	LYS
1	F	120	LYS

Sometimes 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 monosaccharides in this entry.

5.6 Ligand geometry (i)

12 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	PGE	B	302	-	9,9,9	0.67	0	8,8,8	0.82	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	1GP	E	301	-	9,9,9	1.08	0	11,12,12	0.68	0
2	1GP	D	301	-	9,9,9	1.06	0	11,12,12	0.61	0
2	1GP	C	301	-	9,9,9	1.10	0	11,12,12	0.73	0
3	PGE	F	302	-	9,9,9	0.68	0	8,8,8	0.72	0
2	1GP	F	301	-	9,9,9	1.08	0	11,12,12	0.65	0
3	PGE	A	302	-	9,9,9	0.65	0	8,8,8	0.78	0
2	1GP	A	301	-	9,9,9	1.13	0	11,12,12	0.65	0
3	PGE	C	302	-	9,9,9	0.64	0	8,8,8	0.82	0
3	PGE	E	302	-	9,9,9	0.67	0	8,8,8	0.79	0
2	1GP	B	301	-	9,9,9	1.10	0	11,12,12	0.60	0
3	PGE	D	302	-	9,9,9	0.66	0	8,8,8	0.91	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
3	PGE	B	302	-	-	3/7/7/7	-
2	1GP	E	301	-	-	2/8/8/8	-
2	1GP	D	301	-	-	4/8/8/8	-
2	1GP	C	301	-	-	3/8/8/8	-
3	PGE	F	302	-	-	2/7/7/7	-
2	1GP	F	301	-	-	0/8/8/8	-
3	PGE	A	302	-	-	6/7/7/7	-
2	1GP	A	301	-	-	2/8/8/8	-
3	PGE	C	302	-	-	4/7/7/7	-
3	PGE	E	302	-	-	3/7/7/7	-
2	1GP	B	301	-	-	5/8/8/8	-
3	PGE	D	302	-	-	2/7/7/7	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (36) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	1GP	C1-C2-C3-O2

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Mol	Chain	Res	Type	Atoms
2	B	301	1GP	O1P-C1-C2-O3
2	B	301	1GP	C1-O1P-P-O3P
2	B	301	1GP	C1-O1P-P-O4P
2	C	301	1GP	C1-O1P-P-O3P
2	C	301	1GP	C1-O1P-P-O4P
2	E	301	1GP	O1P-C1-C2-C3
2	E	301	1GP	O1P-C1-C2-O3
3	D	302	PGE	O2-C3-C4-O3
2	B	301	1GP	O1P-C1-C2-C3
3	E	302	PGE	O2-C3-C4-O3
3	C	302	PGE	O3-C5-C6-O4
2	A	301	1GP	O3-C2-C3-O2
2	D	301	1GP	O1P-C1-C2-O3
2	D	301	1GP	O1P-C1-C2-C3
3	B	302	PGE	O1-C1-C2-O2
3	C	302	PGE	O2-C3-C4-O3
2	C	301	1GP	C1-O1P-P-O2P
3	C	302	PGE	C3-C4-O3-C5
3	A	302	PGE	O3-C5-C6-O4
3	D	302	PGE	O1-C1-C2-O2
3	C	302	PGE	C4-C3-O2-C2
3	E	302	PGE	O1-C1-C2-O2
3	E	302	PGE	C3-C4-O3-C5
3	A	302	PGE	C4-C3-O2-C2
3	A	302	PGE	O1-C1-C2-O2
3	A	302	PGE	C3-C4-O3-C5
3	F	302	PGE	O3-C5-C6-O4
2	D	301	1GP	C1-C2-C3-O2
3	A	302	PGE	C6-C5-O3-C4
3	B	302	PGE	O3-C5-C6-O4
2	D	301	1GP	O3-C2-C3-O2
2	B	301	1GP	C1-O1P-P-O2P
3	F	302	PGE	O2-C3-C4-O3
3	A	302	PGE	O2-C3-C4-O3
3	B	302	PGE	O2-C3-C4-O3

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	301	1GP	2	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	238/250 (95%)	0.03	1 (0%)	92	91	16, 29, 56, 130	0
1	B	215/250 (86%)	0.32	9 (4%)	36	26	20, 47, 90, 120	0
1	C	234/250 (93%)	0.10	5 (2%)	63	54	17, 30, 73, 128	0
1	D	220/250 (88%)	0.13	4 (1%)	68	61	17, 32, 63, 96	0
1	E	225/250 (90%)	0.40	8 (3%)	42	32	18, 43, 81, 138	0
1	F	207/250 (82%)	0.26	8 (3%)	39	29	22, 44, 79, 128	0
All	All	1339/1500 (89%)	0.20	35 (2%)	56	46	16, 36, 79, 138	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	54	SER	4.1
1	A	55	THR	3.8
1	F	207	SER	3.4
1	E	53	GLY	3.4
1	E	37	ILE	3.3
1	C	55	THR	3.2
1	C	73	ASP	3.1
1	E	36	GLU	2.9
1	B	53	GLY	2.9
1	E	179	ALA	2.8
1	F	37	ILE	2.8
1	F	210	ASP	2.8
1	D	211	ALA	2.8
1	E	45	GLY	2.7
1	B	14	LEU	2.7
1	F	43	ARG	2.4
1	E	4	MET	2.3
1	B	22	THR	2.3
1	B	73	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	F	213	ARG	2.3
1	C	56	THR	2.3
1	D	204	GLY	2.3
1	F	177	SER	2.2
1	B	71	ASN	2.2
1	E	35	VAL	2.2
1	F	215	ALA	2.1
1	D	182	HIS	2.1
1	B	32	GLU	2.1
1	B	72	ILE	2.1
1	D	216	GLY	2.1
1	C	69	ARG	2.1
1	B	55	THR	2.1
1	F	214	VAL	2.1
1	B	24	ILE	2.0
1	C	37	ILE	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 monosaccharides 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(Å^2)	Q<0.9
3	PGE	C	302	10/10	0.84	0.26	42,57,70,77	0
3	PGE	F	302	10/10	0.93	0.18	42,53,63,67	0
3	PGE	A	302	10/10	0.94	0.20	35,47,58,58	0
3	PGE	B	302	10/10	0.94	0.18	36,45,52,56	0
2	1GP	E	301	10/10	0.95	0.17	29,40,55,62	0
3	PGE	D	302	10/10	0.95	0.19	29,43,64,81	0
3	PGE	E	302	10/10	0.95	0.19	30,48,60,63	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	1GP	F	301	10/10	0.95	0.21	31,53,66,74	0
2	1GP	B	301	10/10	0.96	0.13	41,49,61,71	0
2	1GP	D	301	10/10	0.97	0.13	23,33,43,46	0
2	1GP	C	301	10/10	0.98	0.14	16,28,34,40	0
2	1GP	A	301	10/10	0.98	0.13	11,18,30,36	0

6.5 Other polymers (i)

There are no such residues in this entry.