



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 04:49 am BST

PDB ID : 4I8A
Title : Alanine-glyoxylate aminotransferase variant S187F
Authors : Fodor, K.; Oppici, E.; Williams, C.; Cellini, B.; Wilmanns, M.
Deposited on : 2012-12-03
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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

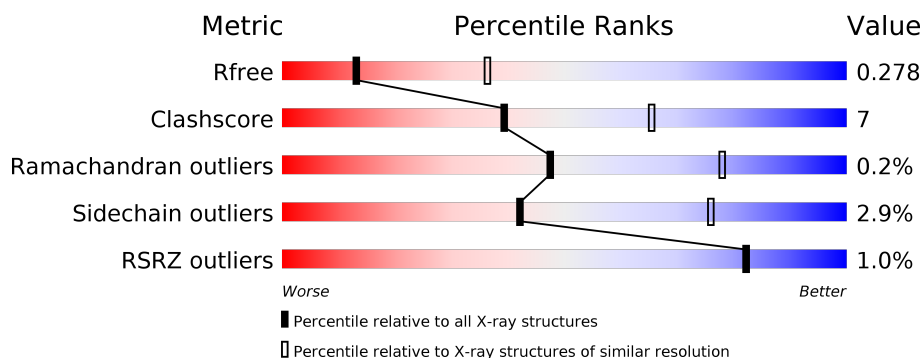
1 Overall quality at a glance

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}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (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 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 $\leq 5\%$. 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	394	<div> <div>%</div> <div> <div></div> <div>82%</div> <div>15%</div> <div>..</div> </div> </div>
1	B	394	<div> <div>%</div> <div> <div></div> <div>81%</div> <div>16%</div> <div>..</div> </div> </div>
1	C	394	<div> <div>%</div> <div> <div></div> <div>82%</div> <div>15%</div> <div>..</div> </div> </div>
1	D	394	<div> <div>%</div> <div> <div></div> <div>80%</div> <div>17%</div> <div>..</div> </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
2	GOL	D	402	-	-	-	X
2	GOL	D	403	-	-	X	X

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11929 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 Serine-pyruvate aminotransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	387	Total	C	N	O	P	S	0	1	0
			2958	1902	500	539	1	16			
1	B	386	Total	C	N	O	P	S	0	0	0
			2981	1911	512	541	1	16			
1	C	385	Total	C	N	O	P	S	0	0	0
			2960	1899	507	537	1	16			
1	D	386	Total	C	N	O	P	S	0	1	0
			2952	1896	499	540	1	16			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP P21549
A	0	ALA	-	EXPRESSION TAG	UNP P21549
A	187	PHE	SER	ENGINEERED MUTATION	UNP P21549
B	-1	GLY	-	EXPRESSION TAG	UNP P21549
B	0	ALA	-	EXPRESSION TAG	UNP P21549
B	187	PHE	SER	ENGINEERED MUTATION	UNP P21549
C	-1	GLY	-	EXPRESSION TAG	UNP P21549
C	0	ALA	-	EXPRESSION TAG	UNP P21549
C	187	PHE	SER	ENGINEERED MUTATION	UNP P21549
D	-1	GLY	-	EXPRESSION TAG	UNP P21549
D	0	ALA	-	EXPRESSION TAG	UNP P21549
D	187	PHE	SER	ENGINEERED MUTATION	UNP P21549

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).

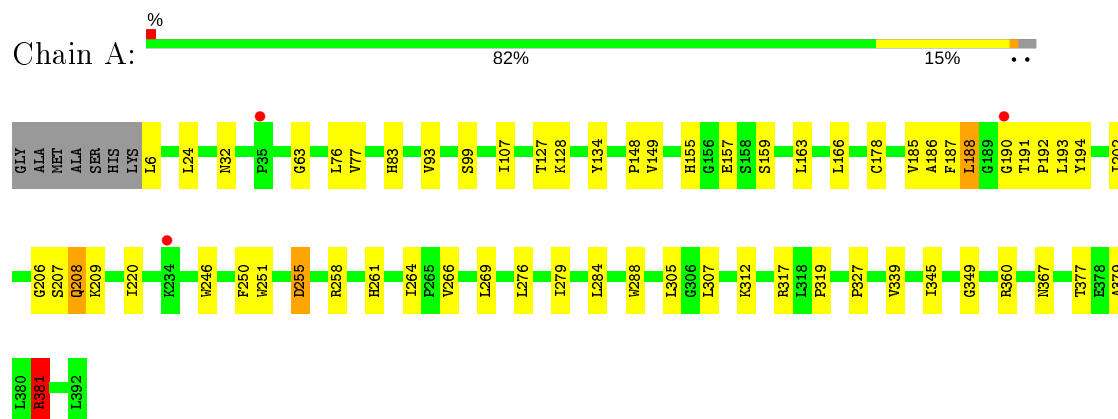


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total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

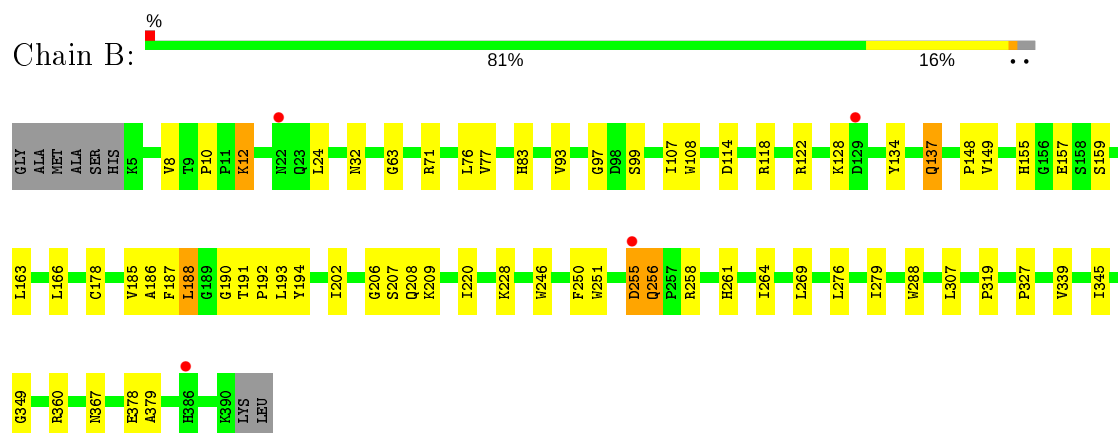
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

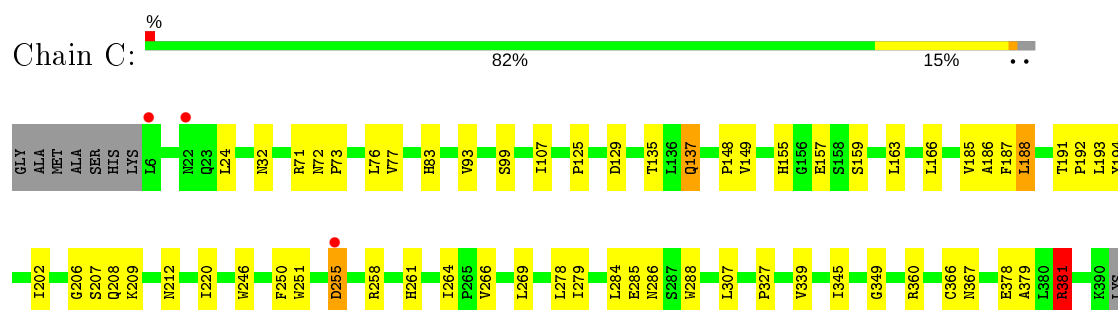
• Molecule 1: Serine-pyruvate aminotransferase



• Molecule 1: Serine-pyruvate aminotransferase

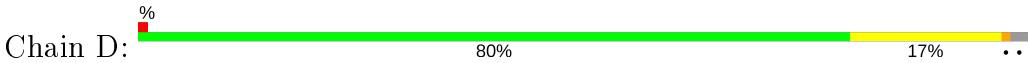


• Molecule 1: Serine-pyruvate aminotransferase



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● Molecule 1: Serine-pyruvate aminotransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	78.89Å 101.56Å 116.75Å 90.00° 90.64° 90.00°	Depositor
Resolution (Å)	18.07 – 2.90 18.07 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.8 (18.07-2.90) 99.3 (18.07-2.90)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.18 (at 2.92Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.251 , 0.278 0.255 , 0.278	Depositor DCC
R_{free} test set	2041 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	35.7	Xtriage
Anisotropy	0.532	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 28.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.048 for h,-k,-l	Xtriage
F_o, F_c correlation	0.82	EDS
Total number of atoms	11929	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.46% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, LLP

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	0.58	0/3003	0.72	3/4083 (0.1%)
1	B	0.58	0/3026	0.73	0/4109
1	C	0.58	0/3005	0.72	1/4084 (0.0%)
1	D	0.58	0/2997	0.71	0/4079
All	All	0.58	0/12031	0.72	4/16355 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	381	ARG	CB-CA-C	5.71	121.81	110.40
1	A	6	LEU	CB-CG-CD2	5.52	120.38	111.00
1	C	381	ARG	NE-CZ-NH1	5.20	122.90	120.30
1	A	381	ARG	CG-CD-NE	-5.14	101.00	111.80

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	2958	0	2930	40	0
1	B	2981	0	2990	46	0
1	C	2960	0	2953	43	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2952	0	2918	52	0
2	A	12	0	16	0	0
2	B	12	0	16	0	0
2	C	24	0	32	4	0
2	D	30	0	40	11	0
All	All	11929	0	11895	177	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 7.

All (177) 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:136[B]:LEU:HD13	1:D:136[B]:LEU:H	1.17	1.03
1:D:136[B]:LEU:N	1:D:136[B]:LEU:HD13	1.89	0.84
1:C:378:GLU:OE2	1:C:381:ARG:NH1	2.14	0.80
1:A:127:THR:OG1	1:C:129:ASP:OD1	2.03	0.76
1:D:136[B]:LEU:H	1:D:136[B]:LEU:CD1	1.86	0.72
1:C:286:ASN:ND2	2:C:404:GOL:O2	2.22	0.71
1:D:82:GLY:N	1:D:209:LLP:OP3	2.27	0.67
1:C:135:THR:HG22	2:C:401:GOL:H31	1.79	0.65
1:C:83:HIS:ND1	1:C:209:LLP:OP3	2.27	0.64
1:D:255:ASP:HB3	2:D:403:GOL:C3	2.29	0.63
1:A:83:HIS:ND1	1:A:209:LLP:OP3	2.25	0.62
1:D:83:HIS:ND1	1:D:209:LLP:OP1	2.29	0.62
1:D:255:ASP:HA	2:D:403:GOL:H2	1.82	0.61
1:D:76:LEU:O	1:D:220:ILE:HG22	2.01	0.60
1:C:76:LEU:O	1:C:220:ILE:HG22	2.02	0.60
1:A:76:LEU:O	1:A:220:ILE:HG22	2.02	0.60
1:B:76:LEU:O	1:B:220:ILE:HG22	2.03	0.59
1:B:83:HIS:ND1	1:B:209:LLP:OP3	2.29	0.58
1:B:157:GLU:OE2	1:B:159:SER:HB3	2.05	0.56
1:A:99:SER:HB3	1:A:148:PRO:HA	1.87	0.56
1:D:157:GLU:OE2	1:D:159:SER:HB3	2.05	0.56
1:D:99:SER:HB3	1:D:148:PRO:HA	1.88	0.56
1:C:157:GLU:OE2	1:C:159:SER:HB3	2.04	0.56
1:C:349:GLY:O	1:C:360:ARG:HD3	2.06	0.56
1:A:157:GLU:OE2	1:A:159:SER:HB3	2.05	0.56
1:B:192:PRO:HB3	1:B:288:TRP:CH2	2.42	0.55
1:C:192:PRO:HB3	1:C:288:TRP:CH2	2.42	0.55
1:D:209:LLP:NZ	1:D:209:LLP:O3	2.40	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:192:PRO:HB3	1:A:288:TRP:CH2	2.42	0.54
1:B:209:LLP:O3	1:B:209:LLP:NZ	2.40	0.54
1:B:255:ASP:OD2	1:D:231:TYR:CD2	2.61	0.54
1:D:349:GLY:O	1:D:360:ARG:HD3	2.08	0.53
1:B:349:GLY:O	1:B:360:ARG:HD3	2.09	0.53
1:A:349:GLY:O	1:A:360:ARG:HD3	2.09	0.53
1:C:99:SER:HB3	1:C:148:PRO:HA	1.90	0.53
1:B:255:ASP:N	1:B:255:ASP:OD1	2.31	0.52
1:A:192:PRO:HB2	1:A:194:TYR:CE1	2.45	0.52
1:B:192:PRO:HB2	1:B:194:TYR:CE1	2.45	0.52
1:A:188:LEU:HD22	1:A:279:ILE:CD1	2.40	0.52
1:B:255:ASP:OD2	1:D:231:TYR:HD2	1.92	0.52
1:D:108:TRP:CD1	1:D:209:LLP:H2'3	2.45	0.52
1:B:99:SER:HB3	1:B:148:PRO:HA	1.92	0.52
1:D:192:PRO:HB2	1:D:194:TYR:CE1	2.45	0.51
1:C:188:LEU:HD22	1:C:279:ILE:CD1	2.40	0.51
1:C:187:PHE:CD1	1:C:188:LEU:HD23	2.45	0.51
1:C:250:PHE:O	1:C:258:ARG:HD3	2.10	0.51
1:D:188:LEU:HD22	1:D:279:ILE:CD1	2.39	0.51
1:B:187:PHE:CD1	1:B:188:LEU:HD23	2.46	0.51
1:D:192:PRO:HB3	1:D:288:TRP:CH2	2.45	0.51
1:D:338:TYR:HH	1:D:386:HIS:HD1	1.58	0.51
1:A:250:PHE:O	1:A:258:ARG:HD3	2.10	0.51
1:B:228:LYS:CG	2:D:403:GOL:H11	2.41	0.51
1:C:107:ILE:HG22	1:C:159:SER:HB2	1.93	0.51
1:C:192:PRO:HB2	1:C:194:TYR:CE1	2.46	0.51
1:B:188:LEU:HD22	1:B:279:ILE:CD1	2.41	0.50
1:B:107:ILE:HG22	1:B:159:SER:HB2	1.93	0.50
1:D:255:ASP:CA	2:D:403:GOL:H2	2.41	0.50
1:B:250:PHE:O	1:B:258:ARG:HD3	2.12	0.50
1:A:255:ASP:N	1:A:255:ASP:OD1	2.31	0.49
1:D:187:PHE:CD1	1:D:188:LEU:HD23	2.47	0.49
1:C:255:ASP:N	1:C:255:ASP:OD1	2.30	0.49
1:A:107:ILE:HG22	1:A:159:SER:HB2	1.94	0.49
1:A:187:PHE:CD1	1:A:188:LEU:HD23	2.48	0.49
1:A:77:VAL:HB	1:A:251:TRP:CZ2	2.48	0.49
1:D:136[B]:LEU:N	1:D:136[B]:LEU:HD22	2.27	0.49
1:D:250:PHE:O	1:D:258:ARG:HD3	2.12	0.49
1:A:305:LEU:HD22	1:A:381:ARG:HB2	1.93	0.48
1:D:339:VAL:HG13	1:D:379:ALA:HB1	1.95	0.48
1:C:155:HIS:CE1	1:C:186:ALA:HB3	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:77:VAL:HB	1:B:251:TRP:CZ2	2.49	0.48
1:C:185:VAL:O	1:C:185:VAL:HG12	2.13	0.47
1:D:107:ILE:HG22	1:D:159:SER:HB2	1.94	0.47
1:C:77:VAL:HB	1:C:251:TRP:CZ2	2.49	0.47
1:D:163:LEU:HD21	1:D:191:THR:HG23	1.97	0.47
1:C:93:VAL:HG21	1:C:202:ILE:HD11	1.97	0.47
1:B:228:LYS:HG3	2:D:403:GOL:C3	2.44	0.47
1:A:339:VAL:HG13	1:A:379:ALA:HB1	1.97	0.47
1:B:185:VAL:O	1:B:185:VAL:HG12	2.14	0.47
1:A:185:VAL:O	1:A:185:VAL:HG12	2.14	0.47
1:C:135:THR:HB	1:C:137:GLN:NE2	2.30	0.47
1:A:155:HIS:CE1	1:A:186:ALA:HB3	2.50	0.47
1:A:377:THR:O	1:A:381:ARG:HB3	2.15	0.47
1:D:155:HIS:CE1	1:D:186:ALA:HB3	2.50	0.47
1:D:122:ARG:HD3	2:D:401:GOL:H11	1.97	0.46
1:D:77:VAL:HB	1:D:251:TRP:CZ2	2.50	0.46
1:B:155:HIS:CE1	1:B:186:ALA:HB3	2.51	0.46
1:B:93:VAL:HG21	1:B:202:ILE:HD11	1.98	0.46
1:C:339:VAL:HG13	1:C:379:ALA:HB1	1.98	0.46
1:D:185:VAL:O	1:D:185:VAL:HG12	2.15	0.46
1:C:206:GLY:O	1:C:207:SER:C	2.55	0.46
1:C:307:LEU:CD2	1:C:327:PRO:HG3	2.46	0.46
1:A:93:VAL:HG21	1:A:202:ILE:HD11	1.98	0.45
1:B:228:LYS:HG2	2:D:403:GOL:H11	1.98	0.45
1:C:163:LEU:HD21	1:C:191:THR:HG23	1.98	0.45
1:B:97:GLY:O	1:B:122:ARG:CZ	2.64	0.45
1:A:206:GLY:O	1:A:207:SER:C	2.55	0.45
1:C:286:ASN:ND2	2:C:404:GOL:C2	2.80	0.45
1:B:228:LYS:HG3	2:D:403:GOL:C1	2.47	0.45
1:A:163:LEU:HD21	1:A:191:THR:HG23	1.98	0.45
1:D:10:PRO:O	1:D:12:LYS:HE3	2.17	0.45
1:A:155:HIS:CE1	1:A:193:LEU:HD13	2.52	0.45
1:A:307:LEU:CD2	1:A:327:PRO:HG3	2.47	0.45
1:B:206:GLY:O	1:B:207:SER:C	2.56	0.45
1:C:125:PRO:O	2:C:402:GOL:O2	2.35	0.45
1:D:255:ASP:HB3	2:D:403:GOL:H32	1.99	0.45
1:D:155:HIS:CE1	1:D:193:LEU:HD13	2.53	0.44
1:D:307:LEU:CD2	1:D:327:PRO:HG3	2.47	0.44
1:D:155:HIS:CG	1:D:166:LEU:HD11	2.53	0.44
1:D:246:TRP:HB3	1:D:261:HIS:CD2	2.52	0.44
1:B:339:VAL:HG13	1:B:379:ALA:HB1	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:163:LEU:HD21	1:B:191:THR:HG23	2.00	0.44
1:C:209:LLP:OP1	1:D:260:TYR:OH	2.34	0.43
1:C:77:VAL:HB	1:C:251:TRP:CH2	2.53	0.43
1:B:137:GLN:N	1:B:137:GLN:OE1	2.51	0.43
1:B:155:HIS:CG	1:B:166:LEU:HD11	2.53	0.43
1:B:155:HIS:CE1	1:B:193:LEU:HD13	2.54	0.43
1:C:155:HIS:CE1	1:C:193:LEU:HD13	2.54	0.43
1:D:206:GLY:O	1:D:207:SER:C	2.56	0.43
1:B:10:PRO:O	1:B:12:LYS:HE3	2.18	0.43
1:D:93:VAL:HG21	1:D:202:ILE:HD11	1.99	0.43
1:C:155:HIS:CG	1:C:166:LEU:HD11	2.54	0.43
1:D:255:ASP:HB3	2:D:403:GOL:C2	2.48	0.43
1:D:24:LEU:HA	1:D:32:ASN:HD21	1.84	0.43
1:A:307:LEU:HD23	1:A:327:PRO:HG3	2.01	0.43
1:C:24:LEU:HA	1:C:32:ASN:HD21	1.84	0.43
1:B:97:GLY:O	1:B:122:ARG:NH2	2.51	0.43
1:C:137:GLN:H	1:C:137:GLN:HG3	1.49	0.43
1:A:246:TRP:HB3	1:A:261:HIS:CD2	2.54	0.43
1:A:63:GLY:HA3	1:A:276:LEU:HD13	2.00	0.43
1:B:307:LEU:CD2	1:B:327:PRO:HG3	2.49	0.43
1:C:246:TRP:HB3	1:C:261:HIS:CD2	2.54	0.43
1:B:108:TRP:CD1	1:B:209:LLP:H2'3	2.54	0.42
1:A:305:LEU:CD2	1:A:381:ARG:HB2	2.50	0.42
1:B:128:LYS:HD2	1:B:134:TYR:CE2	2.54	0.42
1:B:190:GLY:HA2	1:B:319:PRO:HG2	2.02	0.42
1:D:338:TYR:OH	1:D:386:HIS:ND1	2.42	0.42
1:A:279:ILE:HD11	1:A:284:LEU:HD23	2.01	0.42
1:B:114:ASP:OD1	1:B:118:ARG:NE	2.49	0.42
1:B:256:GLN:CA	1:B:256:GLN:HE21	2.32	0.42
1:D:317:ARG:O	1:D:319:PRO:HD3	2.20	0.42
1:A:317:ARG:O	1:A:319:PRO:HD3	2.19	0.42
1:A:77:VAL:HB	1:A:251:TRP:CH2	2.55	0.42
1:C:32:ASN:HB2	1:C:367:ASN:HD21	1.85	0.42
1:D:32:ASN:HB2	1:D:367:ASN:HD21	1.85	0.42
1:A:32:ASN:HB2	1:A:367:ASN:HD21	1.85	0.42
1:B:339:VAL:HG12	1:B:345:ILE:HB	2.02	0.42
1:A:190:GLY:HA2	1:A:319:PRO:HG2	2.02	0.41
1:A:264:ILE:HD13	1:A:269:LEU:HD12	2.02	0.41
1:D:279:ILE:HD11	1:D:284:LEU:HD23	2.02	0.41
1:D:307:LEU:HD23	1:D:327:PRO:HG3	2.02	0.41
1:B:228:LYS:CG	2:D:403:GOL:C1	2.99	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:24:LEU:HA	1:B:32:ASN:HD21	1.84	0.41
1:B:264:ILE:HD13	1:B:269:LEU:HD12	2.01	0.41
1:B:63:GLY:HA3	1:B:276:LEU:HD13	2.03	0.41
1:D:77:VAL:HB	1:D:251:TRP:CH2	2.55	0.41
1:A:128:LYS:HD2	1:A:134:TYR:CE2	2.55	0.41
1:D:339:VAL:HG12	1:D:345:ILE:HB	2.02	0.41
1:B:246:TRP:HB3	1:B:261:HIS:CD2	2.56	0.41
1:B:77:VAL:HB	1:B:251:TRP:CH2	2.56	0.41
1:C:307:LEU:HD23	1:C:327:PRO:HG3	2.02	0.41
1:A:206:GLY:O	1:A:208:GLN:N	2.54	0.41
1:C:279:ILE:HD11	1:C:284:LEU:HD23	2.01	0.41
1:C:212:ASN:HD21	1:C:366:CYS:H	1.68	0.41
1:D:184:SER:O	1:D:205:SER:HB2	2.21	0.41
1:C:278:LEU:HD23	1:C:278:LEU:HA	1.97	0.41
1:A:155:HIS:CG	1:A:166:LEU:HD11	2.56	0.41
1:C:264:ILE:HD13	1:C:269:LEU:HD12	2.03	0.40
1:A:339:VAL:HG12	1:A:345:ILE:HB	2.03	0.40
1:C:285:GLU:HA	1:C:288:TRP:CE3	2.56	0.40
1:C:72:ASN:HA	1:C:73:PRO:HD3	1.97	0.40
1:A:188:LEU:CD2	1:A:279:ILE:CD1	3.00	0.40
1:B:32:ASN:HB2	1:B:367:ASN:HD21	1.87	0.40
1:D:188:LEU:CD2	1:D:279:ILE:CD1	2.99	0.40
1:A:24:LEU:HA	1:A:32:ASN:HD21	1.86	0.40
1:C:339:VAL:HG12	1:C:345:ILE:HB	2.02	0.40
1:D:128:LYS:HD2	1:D:134:TYR:CE2	2.57	0.40
1:D:264:ILE:HD13	1:D:269:LEU:HD12	2.02	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	385/394 (98%)	375 (97%)	9 (2%)	1 (0%)	41	71
1	B	383/394 (97%)	373 (97%)	10 (3%)	0	100	100
1	C	382/394 (97%)	373 (98%)	8 (2%)	1 (0%)	41	71
1	D	384/394 (98%)	372 (97%)	11 (3%)	1 (0%)	41	71
All	All	1534/1576 (97%)	1493 (97%)	38 (2%)	3 (0%)	47	78

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	266	VAL
1	A	266	VAL
1	D	266	VAL

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	303/325 (93%)	296 (98%)	7 (2%)	50	80
1	B	313/325 (96%)	302 (96%)	11 (4%)	36	70
1	C	308/325 (95%)	301 (98%)	7 (2%)	50	80
1	D	306/325 (94%)	296 (97%)	10 (3%)	38	72
All	All	1230/1300 (95%)	1195 (97%)	35 (3%)	42	76

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	149	VAL
1	A	178	CYS
1	A	188	LEU
1	A	208	GLN
1	A	255	ASP
1	A	312	LYS
1	A	381	ARG
1	B	8	VAL

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Mol	Chain	Res	Type
1	B	12	LYS
1	B	71	ARG
1	B	137	GLN
1	B	149	VAL
1	B	178	CYS
1	B	188	LEU
1	B	208	GLN
1	B	255	ASP
1	B	256	GLN
1	B	378	GLU
1	C	71	ARG
1	C	137	GLN
1	C	149	VAL
1	C	188	LEU
1	C	208	GLN
1	C	255	ASP
1	C	381	ARG
1	D	9	THR
1	D	12	LYS
1	D	71	ARG
1	D	137	GLN
1	D	149	VAL
1	D	178	CYS
1	D	188	LEU
1	D	208	GLN
1	D	255	ASP
1	D	256	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	ASN
1	A	55	GLN
1	A	72	ASN
1	A	208	GLN
1	A	212	ASN
1	A	299	HIS
1	B	32	ASN
1	B	55	GLN
1	B	72	ASN
1	B	208	GLN
1	B	212	ASN

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Mol	Chain	Res	Type
1	B	256	GLN
1	B	299	HIS
1	C	32	ASN
1	C	55	GLN
1	C	72	ASN
1	C	208	GLN
1	C	212	ASN
1	C	286	ASN
1	C	299	HIS
1	D	32	ASN
1	D	72	ASN
1	D	208	GLN
1	D	212	ASN
1	D	256	GLN
1	D	299	HIS

5.3.3 RNA ⓘ

There are no RNA molecules 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	LLP	D	209	1	23,24,25	1.82	7 (30%)	25,32,34	1.97	4 (16%)
1	LLP	C	209	1	23,24,25	1.92	5 (21%)	25,32,34	1.92	4 (16%)
1	LLP	B	209	1	23,24,25	1.86	5 (21%)	25,32,34	1.81	6 (24%)
1	LLP	A	209	1	23,24,25	1.73	4 (17%)	25,32,34	2.04	5 (20%)

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	LLP	D	209	1	-	4/16/17/19	0/1/1/1
1	LLP	C	209	1	-	8/16/17/19	0/1/1/1
1	LLP	B	209	1	-	4/16/17/19	0/1/1/1
1	LLP	A	209	1	-	8/16/17/19	0/1/1/1

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	209	LLP	O3-C3	-5.38	1.24	1.37
1	B	209	LLP	O3-C3	-5.32	1.24	1.37
1	A	209	LLP	O3-C3	-5.03	1.25	1.37
1	D	209	LLP	O3-C3	-4.65	1.26	1.37
1	C	209	LLP	C6-N1	3.65	1.42	1.34
1	C	209	LLP	C2-N1	3.42	1.40	1.33
1	B	209	LLP	CE-NZ	3.30	1.54	1.46
1	D	209	LLP	CE-NZ	3.08	1.53	1.46
1	A	209	LLP	C2-N1	3.00	1.39	1.33
1	B	209	LLP	C4'-NZ	3.00	1.37	1.27
1	B	209	LLP	C4-C4'	2.90	1.52	1.46
1	D	209	LLP	C4-C4'	2.81	1.52	1.46
1	D	209	LLP	C4'-NZ	2.59	1.35	1.27
1	D	209	LLP	P-OP3	-2.53	1.45	1.54
1	D	209	LLP	P-OP2	-2.42	1.45	1.54
1	C	209	LLP	C4-C4'	2.39	1.51	1.46
1	C	209	LLP	P-OP2	-2.17	1.46	1.54
1	D	209	LLP	C2-N1	2.15	1.37	1.33
1	A	209	LLP	P-OP4	2.08	1.66	1.60
1	B	209	LLP	C2-N1	2.02	1.37	1.33
1	A	209	LLP	C4-C4'	2.01	1.50	1.46

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	209	LLP	OP4-C5'-C5	6.93	122.56	109.35
1	A	209	LLP	C4-C4'-NZ	-5.98	96.87	124.31
1	A	209	LLP	OP4-C5'-C5	5.71	120.22	109.35
1	B	209	LLP	OP4-C5'-C5	5.57	119.97	109.35
1	C	209	LLP	C4-C4'-NZ	-5.39	99.58	124.31
1	C	209	LLP	OP4-C5'-C5	5.09	119.04	109.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	209	LLP	OP3-P-OP2	3.72	121.84	107.64
1	C	209	LLP	C3-C4-C5	3.53	120.97	118.26
1	B	209	LLP	C4-C4'-NZ	-3.50	108.25	124.31
1	D	209	LLP	OP3-P-OP4	-3.07	98.56	106.73
1	D	209	LLP	C4-C4'-NZ	-3.03	110.39	124.31
1	B	209	LLP	OP3-P-OP2	2.65	117.75	107.64
1	B	209	LLP	OP4-P-OP1	-2.56	99.28	106.47
1	A	209	LLP	OP2-P-OP4	-2.51	100.06	106.73
1	A	209	LLP	C3-C4-C5	2.30	120.03	118.26
1	C	209	LLP	C3-C2-N1	-2.24	117.87	120.77
1	B	209	LLP	C5'-C5-C6	-2.11	115.90	119.37
1	A	209	LLP	C5-C6-N1	-2.10	120.31	123.82
1	B	209	LLP	C5-C6-N1	-2.01	120.47	123.82

There are no chirality outliers.

All (24) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	D	209	LLP	C5'-OP4-P-OP1
1	D	209	LLP	C5'-OP4-P-OP2
1	D	209	LLP	C5'-OP4-P-OP3
1	C	209	LLP	C5'-OP4-P-OP1
1	C	209	LLP	C5'-OP4-P-OP2
1	B	209	LLP	C5'-OP4-P-OP1
1	B	209	LLP	C5'-OP4-P-OP2
1	B	209	LLP	C5'-OP4-P-OP3
1	A	209	LLP	C5'-OP4-P-OP2
1	D	209	LLP	C4-C4'-NZ-CE
1	B	209	LLP	C4-C4'-NZ-CE
1	C	209	LLP	CG-CD-CE-NZ
1	A	209	LLP	CG-CD-CE-NZ
1	C	209	LLP	C4-C4'-NZ-CE
1	C	209	LLP	C3-C4-C4'-NZ
1	A	209	LLP	C3-C4-C4'-NZ
1	A	209	LLP	C4-C4'-NZ-CE
1	A	209	LLP	C5'-OP4-P-OP1
1	A	209	LLP	CD-CE-NZ-C4'
1	C	209	LLP	CD-CE-NZ-C4'
1	C	209	LLP	C5-C4-C4'-NZ
1	C	209	LLP	C5'-OP4-P-OP3
1	A	209	LLP	C5'-OP4-P-OP3
1	A	209	LLP	C5-C4-C4'-NZ

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	209	LLP	4	0
1	C	209	LLP	2	0
1	B	209	LLP	3	0
1	A	209	LLP	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

13 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	$\# Z > 2$
2	GOL	C	404	-	5,5,5	0.42	0	5,5,5	0.68	0
2	GOL	C	403	-	5,5,5	0.45	0	5,5,5	0.50	0
2	GOL	D	402	-	5,5,5	0.47	0	5,5,5	0.59	0
2	GOL	B	402	-	5,5,5	0.45	0	5,5,5	0.47	0
2	GOL	D	401	-	5,5,5	0.41	0	5,5,5	0.41	0
2	GOL	B	401	-	5,5,5	0.40	0	5,5,5	0.27	0
2	GOL	D	405	-	5,5,5	0.49	0	5,5,5	0.23	0
2	GOL	C	401	-	5,5,5	0.36	0	5,5,5	0.23	0
2	GOL	A	401	-	5,5,5	0.39	0	5,5,5	0.37	0
2	GOL	A	402	-	5,5,5	0.49	0	5,5,5	0.76	0
2	GOL	C	402	-	5,5,5	0.43	0	5,5,5	0.67	0
2	GOL	D	403	-	5,5,5	0.55	0	5,5,5	0.74	0
2	GOL	D	404	-	5,5,5	0.45	0	5,5,5	0.39	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
2	GOL	C	404	-	-	4/4/4/4	-
2	GOL	C	403	-	-	4/4/4/4	-
2	GOL	D	402	-	-	2/4/4/4	-
2	GOL	B	402	-	-	4/4/4/4	-
2	GOL	D	401	-	-	2/4/4/4	-
2	GOL	B	401	-	-	2/4/4/4	-
2	GOL	D	405	-	-	2/4/4/4	-
2	GOL	C	401	-	-	2/4/4/4	-
2	GOL	A	401	-	-	0/4/4/4	-
2	GOL	A	402	-	-	1/4/4/4	-
2	GOL	C	402	-	-	3/4/4/4	-
2	GOL	D	403	-	-	3/4/4/4	-
2	GOL	D	404	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (31) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	404	GOL	C1-C2-C3-O3
2	C	403	GOL	O1-C1-C2-C3
2	C	403	GOL	C1-C2-C3-O3
2	D	402	GOL	O1-C1-C2-C3
2	B	402	GOL	O1-C1-C2-C3
2	C	401	GOL	C1-C2-C3-O3
2	C	402	GOL	C1-C2-C3-O3
2	D	403	GOL	C1-C2-C3-O3
2	C	404	GOL	O2-C2-C3-O3
2	C	402	GOL	O2-C2-C3-O3
2	C	404	GOL	O1-C1-C2-C3
2	B	402	GOL	C1-C2-C3-O3
2	D	401	GOL	C1-C2-C3-O3
2	B	401	GOL	O1-C1-C2-C3
2	C	402	GOL	O1-C1-C2-C3
2	D	404	GOL	O1-C1-C2-C3
2	C	404	GOL	O1-C1-C2-O2
2	C	403	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
2	C	403	GOL	O2-C2-C3-O3
2	B	402	GOL	O1-C1-C2-O2
2	B	402	GOL	O2-C2-C3-O3
2	D	401	GOL	O2-C2-C3-O3
2	B	401	GOL	O1-C1-C2-O2
2	D	404	GOL	O1-C1-C2-O2
2	D	402	GOL	O1-C1-C2-O2
2	C	401	GOL	O2-C2-C3-O3
2	D	405	GOL	O2-C2-C3-O3
2	A	402	GOL	O2-C2-C3-O3
2	D	405	GOL	C1-C2-C3-O3
2	D	403	GOL	O1-C1-C2-C3
2	D	403	GOL	O2-C2-C3-O3

There are no ring outliers.

5 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	404	GOL	2	0
2	D	401	GOL	1	0
2	C	401	GOL	1	0
2	C	402	GOL	1	0
2	D	403	GOL	10	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	386/394 (97%)	0.03	3 (0%) 86 86	19, 29, 43, 65	1 (0%)
1	B	385/394 (97%)	0.05	4 (1%) 82 82	17, 30, 45, 57	2 (0%)
1	C	384/394 (97%)	-0.01	3 (0%) 86 86	18, 28, 41, 57	1 (0%)
1	D	385/394 (97%)	0.15	5 (1%) 77 77	19, 31, 46, 70	2 (0%)
All	All	1540/1576 (97%)	0.05	15 (0%) 82 82	17, 30, 45, 70	6 (0%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	386	HIS	5.2
1	B	386	HIS	4.8
1	A	234	LYS	3.8
1	D	315	ALA	3.1
1	C	22	ASN	3.1
1	D	286	ASN	3.1
1	C	255	ASP	3.0
1	C	6	LEU	2.9
1	B	22	ASN	2.8
1	D	141	GLU	2.3
1	A	190	GLY	2.3
1	A	35	PRO	2.2
1	B	129	ASP	2.1
1	B	255	ASP	2.1
1	D	146	HIS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column

labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
1	LLP	C	209	24/25	0.93	0.18	25,28,29,30	0
1	LLP	B	209	24/25	0.93	0.17	25,33,34,36	0
1	LLP	D	209	24/25	0.94	0.18	26,31,34,36	0
1	LLP	A	209	24/25	0.94	0.19	26,30,33,34	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	GOL	D	403	6/6	0.60	0.52	49,52,53,65	0
2	GOL	B	401	6/6	0.68	0.39	47,50,50,50	0
2	GOL	D	402	6/6	0.69	0.51	44,47,49,51	0
2	GOL	A	402	6/6	0.79	0.33	42,48,51,51	0
2	GOL	B	402	6/6	0.79	0.27	34,38,40,40	0
2	GOL	C	401	6/6	0.81	0.29	26,27,28,30	0
2	GOL	C	404	6/6	0.82	0.35	38,45,47,51	0
2	GOL	D	404	6/6	0.84	0.28	45,47,48,48	0
2	GOL	C	403	6/6	0.86	0.24	32,34,34,34	0
2	GOL	D	405	6/6	0.90	0.34	34,37,37,38	0
2	GOL	D	401	6/6	0.90	0.29	30,33,33,35	0
2	GOL	A	401	6/6	0.90	0.20	50,52,53,57	0
2	GOL	C	402	6/6	0.91	0.18	33,34,39,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.