Full wwPDB X-ray Structure Validation Report (i)

Jun 15, 2024 – 02:28 PM EDT

PDB ID	:	1T8Q
Title	:	Structural genomics, Crystal structure of Glycerophosphoryl diester phosphodiesterase from E. coli
Authors	:	Zhang, R.; Kim, Y.; Dementieva, I.; Duke, N.; Stols, L.; Donnelly, M.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on	:	2004-05-13
Resolution	:	2.00 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:

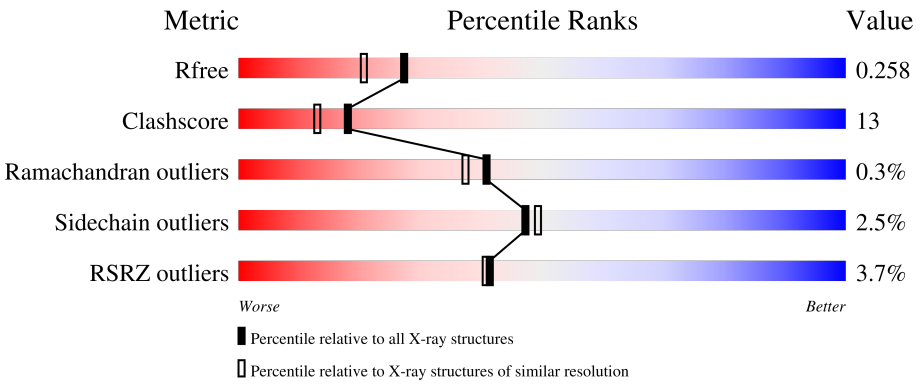
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	336	% 79% 18% • •
1	B	336	% 77% 20% • •
1	C	336	4% 70% 26% • •
1	D	336	8% 65% 31% • •

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
3	GOL	A	1609	-	-	-	X
3	GOL	C	1611	-	-	-	X

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 11867 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 Glycerophosphoryl diester phosphodiesterase, periplasmic.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	328	2659	1698	443	508	1	9	0	0	0
1	B	329	2668	1704	445	509	1	9	0	0	0
1	C	328	2659	1698	443	508	1	9	0	0	0
1	D	327	2651	1694	441	506	1	9	0	0	0

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	23	SER	-	CLONING ARTIFACT	UNP P09394
A	24	ASN	-	CLONING ARTIFACT	UNP P09394
A	53	MSE	MET	MODIFIED RESIDUE	UNP P09394
A	68	MSE	MET	MODIFIED RESIDUE	UNP P09394
A	138	MSE	MET	MODIFIED RESIDUE	UNP P09394
A	226	MSE	MET	MODIFIED RESIDUE	UNP P09394
A	228	MSE	MET	MODIFIED RESIDUE	UNP P09394
A	261	MSE	MET	MODIFIED RESIDUE	UNP P09394
A	267	MSE	MET	MODIFIED RESIDUE	UNP P09394
A	283	MSE	MET	MODIFIED RESIDUE	UNP P09394
A	299	MSE	MET	MODIFIED RESIDUE	UNP P09394
B	23	SER	-	CLONING ARTIFACT	UNP P09394
B	24	ASN	-	CLONING ARTIFACT	UNP P09394
B	53	MSE	MET	MODIFIED RESIDUE	UNP P09394
B	68	MSE	MET	MODIFIED RESIDUE	UNP P09394
B	138	MSE	MET	MODIFIED RESIDUE	UNP P09394
B	226	MSE	MET	MODIFIED RESIDUE	UNP P09394
B	228	MSE	MET	MODIFIED RESIDUE	UNP P09394
B	261	MSE	MET	MODIFIED RESIDUE	UNP P09394
B	267	MSE	MET	MODIFIED RESIDUE	UNP P09394
B	283	MSE	MET	MODIFIED RESIDUE	UNP P09394

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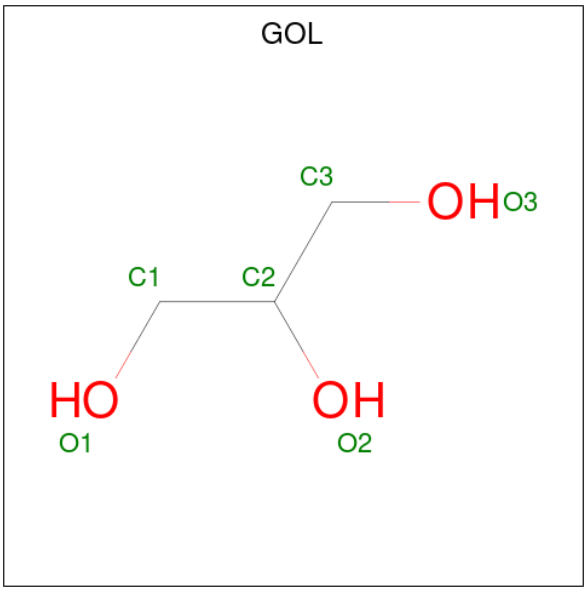
Chain	Residue	Modelled	Actual	Comment	Reference
B	299	MSE	MET	MODIFIED RESIDUE	UNP P09394
C	23	SER	-	CLONING ARTIFACT	UNP P09394
C	24	ASN	-	CLONING ARTIFACT	UNP P09394
C	53	MSE	MET	MODIFIED RESIDUE	UNP P09394
C	68	MSE	MET	MODIFIED RESIDUE	UNP P09394
C	138	MSE	MET	MODIFIED RESIDUE	UNP P09394
C	226	MSE	MET	MODIFIED RESIDUE	UNP P09394
C	228	MSE	MET	MODIFIED RESIDUE	UNP P09394
C	261	MSE	MET	MODIFIED RESIDUE	UNP P09394
C	267	MSE	MET	MODIFIED RESIDUE	UNP P09394
C	283	MSE	MET	MODIFIED RESIDUE	UNP P09394
C	299	MSE	MET	MODIFIED RESIDUE	UNP P09394
D	23	SER	-	CLONING ARTIFACT	UNP P09394
D	24	ASN	-	CLONING ARTIFACT	UNP P09394
D	53	MSE	MET	MODIFIED RESIDUE	UNP P09394
D	68	MSE	MET	MODIFIED RESIDUE	UNP P09394
D	138	MSE	MET	MODIFIED RESIDUE	UNP P09394
D	226	MSE	MET	MODIFIED RESIDUE	UNP P09394
D	228	MSE	MET	MODIFIED RESIDUE	UNP P09394
D	261	MSE	MET	MODIFIED RESIDUE	UNP P09394
D	267	MSE	MET	MODIFIED RESIDUE	UNP P09394
D	283	MSE	MET	MODIFIED RESIDUE	UNP P09394
D	299	MSE	MET	MODIFIED RESIDUE	UNP P09394

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

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

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	357	357	357	0	0

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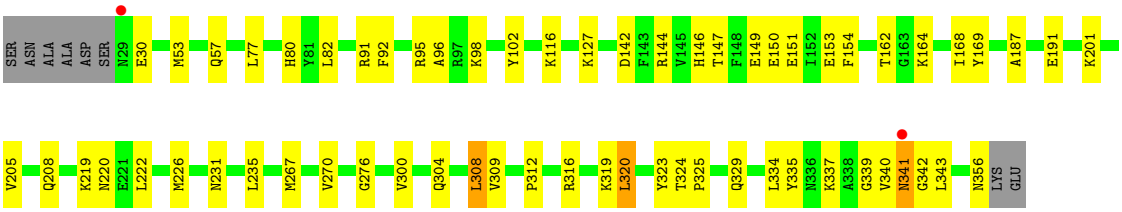
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	B	333	333	333	0	0
			Total	O		
4	C	255	255	255	0	0
			Total	O		
4	D	215	215	215	0	0

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.

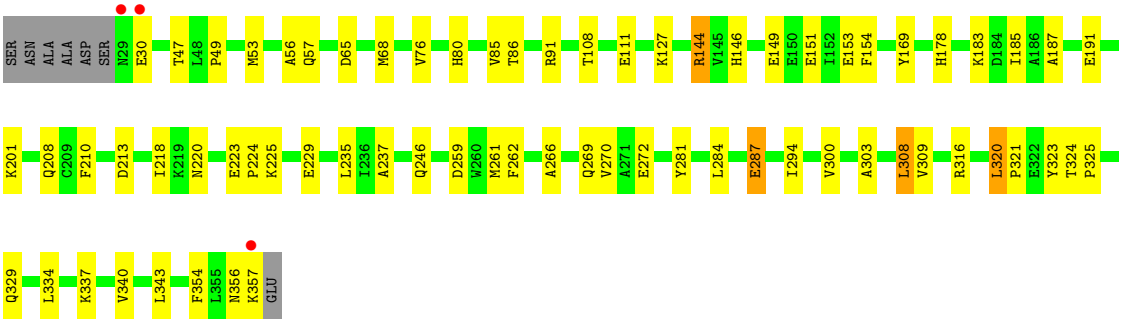
• Molecule 1: Glycerophosphoryl diester phosphodiesterase, periplasmic

Chain A:



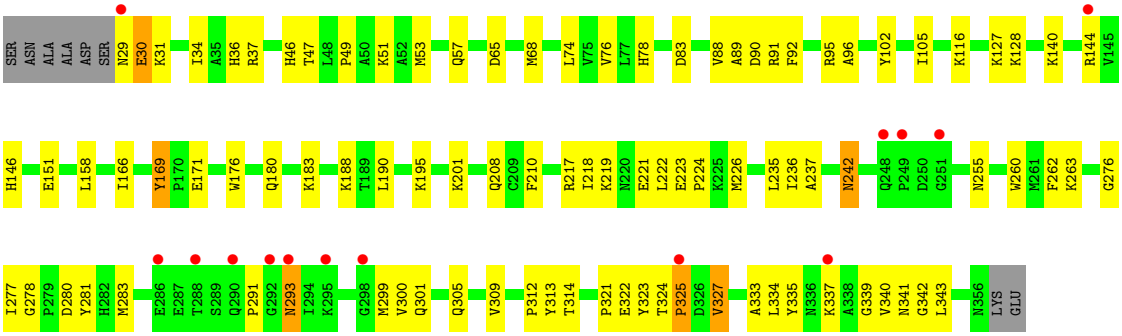
• Molecule 1: Glycerophosphoryl diester phosphodiesterase, periplasmic

Chain B:



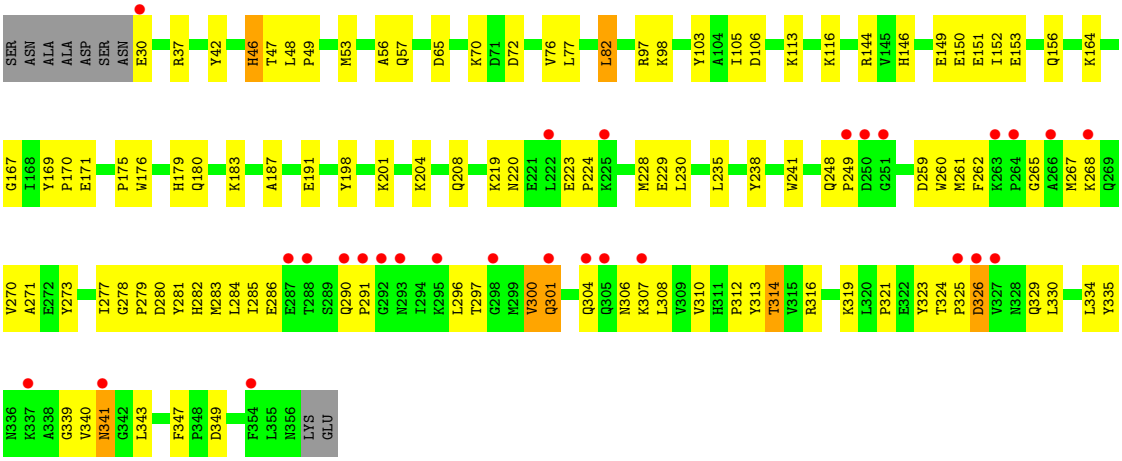
• Molecule 1: Glycerophosphoryl diester phosphodiesterase, periplasmic

Chain C:



• Molecule 1: Glycerophosphoryl diester phosphodiesterase, periplasmic

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	62.76Å 118.39Å 242.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.56 – 2.00 43.56 – 2.00	Depositor EDS
% Data completeness (in resolution range)	89.5 (43.56-2.00) 89.6 (43.56-2.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.59 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.233 , 0.265 0.225 , 0.258	Depositor DCC
R_{free} test set	21097 reflections (9.21%)	wwPDB-VP
Wilson B-factor (Å^2)	25.3	Xtriage
Anisotropy	0.666	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 44.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	11867	wwPDB-VP
Average B, all atoms (Å^2)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% 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: MG, GOL

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	0.43	0/2717	0.66	0/3663
1	B	0.45	0/2726	0.66	0/3674
1	C	0.41	0/2717	0.64	0/3663
1	D	0.41	0/2709	0.61	0/3652
All	All	0.43	0/10869	0.64	0/14652

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2659	0	2569	56	0
1	B	2668	0	2582	60	0
1	C	2659	0	2569	85	0
1	D	2651	0	2563	86	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	18	0	23	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	12	0	14	1	0
3	C	18	0	24	3	0
3	D	18	0	24	0	0
4	A	357	0	0	9	0
4	B	333	0	0	11	0
4	C	255	0	0	10	0
4	D	215	0	0	12	0
All	All	11867	0	10368	284	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 13.

All (284) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:313:TYR:HB2	1:C:314:THR:HG22	1.39	1.01
1:B:68:MSE:HE2	1:B:185:ILE:HB	1.41	1.00
1:C:144:ARG:HB2	4:C:1839:HOH:O	1.65	0.96
1:C:301:GLN:O	1:C:305:GLN:HB2	1.70	0.91
1:C:333:ALA:HA	1:C:337:LYS:HG3	1.53	0.90
1:A:267:MSE:HE2	1:A:308:LEU:CD2	2.02	0.90
1:A:309:VAL:HG22	1:A:341:ASN:HD21	1.36	0.90
1:D:176:TRP:HD1	4:D:1767:HOH:O	1.55	0.88
1:A:276:GLY:HA3	1:A:309:VAL:HG12	1.55	0.85
1:A:267:MSE:HE3	1:A:267:MSE:HA	1.60	0.83
1:A:267:MSE:HE2	1:A:308:LEU:HD23	1.60	0.83
1:B:224:PRO:HB2	4:B:1919:HOH:O	1.85	0.76
1:A:300:VAL:O	1:A:304:GLN:HG3	1.86	0.75
1:D:179:HIS:HD2	4:D:1768:HOH:O	1.70	0.74
1:D:53:MSE:O	1:D:57:GLN:HG3	1.87	0.74
1:B:144:ARG:HD2	4:C:1826:HOH:O	1.87	0.73
1:D:105:ILE:HD11	1:D:183:LYS:HD2	1.71	0.73
1:D:187:ALA:O	1:D:191:GLU:HG2	1.89	0.72
1:C:222:LEU:HB3	1:C:226:MSE:CE	2.20	0.71
1:B:294:ILE:O	1:B:337:LYS:HE2	1.89	0.71
1:C:146:HIS:HE1	1:C:151:GLU:OE1	1.73	0.71
1:A:309:VAL:HG22	1:A:341:ASN:ND2	2.06	0.71
1:B:68:MSE:HE2	1:B:185:ILE:CB	2.21	0.70
1:A:153:GLU:OE2	4:A:1930:HOH:O	2.09	0.70
1:A:267:MSE:CE	1:A:270:VAL:HB	2.22	0.70
1:A:276:GLY:HA3	1:A:309:VAL:CG1	2.22	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:269:GLN:O	1:B:272:GLU:HG2	1.93	0.68
1:A:267:MSE:HE3	1:A:270:VAL:HB	1.73	0.68
1:A:309:VAL:CG2	1:A:341:ASN:ND2	2.55	0.68
1:B:287:GLU:HG2	4:B:1707:HOH:O	1.93	0.68
1:C:313:TYR:HB2	1:C:314:THR:CG2	2.19	0.68
1:B:325:PRO:HD2	1:B:329:GLN:HE22	1.60	0.67
1:D:179:HIS:CD2	4:D:1768:HOH:O	2.46	0.66
1:B:325:PRO:HD2	1:B:329:GLN:NE2	2.11	0.66
1:A:219:LYS:HG2	1:A:220:ASN:ND2	2.09	0.66
1:D:146:HIS:HE1	1:D:151:GLU:OE1	1.79	0.65
1:A:144:ARG:HG3	1:A:144:ARG:HH11	1.62	0.65
1:B:80:HIS:H	1:B:80:HIS:CD2	2.12	0.65
1:C:105:ILE:HD11	1:C:183:LYS:HD2	1.78	0.65
1:C:293:ASN:OD1	4:C:1804:HOH:O	2.15	0.64
1:D:334:LEU:O	1:D:340:VAL:HG22	1.96	0.64
1:D:175:PRO:HG2	4:D:1767:HOH:O	1.95	0.64
1:A:309:VAL:CG2	1:A:341:ASN:HD21	2.06	0.64
1:A:267:MSE:HE2	1:A:308:LEU:HD21	1.78	0.63
1:A:276:GLY:CA	1:A:309:VAL:HG12	2.29	0.63
1:B:218:ILE:O	1:B:223:GLU:HG3	2.00	0.62
1:B:225:LYS:HG3	4:B:1919:HOH:O	1.98	0.62
1:C:333:ALA:HA	1:C:337:LYS:CG	2.27	0.62
1:B:30:GLU:O	1:B:30:GLU:HG3	1.99	0.62
1:C:281:TYR:OH	1:C:321:PRO:HD3	2.00	0.62
1:B:201:LYS:HE3	1:B:229:GLU:HG2	1.80	0.62
1:D:235:LEU:HD23	1:D:278:GLY:HA3	1.82	0.62
1:D:267:MSE:HE2	1:D:267:MSE:HA	1.82	0.61
1:B:146:HIS:HE1	1:B:151:GLU:OE1	1.83	0.61
1:B:208:GLN:OE1	1:B:235:LEU:HD21	2.00	0.61
1:C:322:GLU:HA	1:C:322:GLU:OE2	2.00	0.61
1:B:47:THR:HB	1:B:49:PRO:HD2	1.83	0.61
1:B:91:ARG:HD3	4:B:1704:HOH:O	2.01	0.60
1:A:80:HIS:HE1	4:A:1699:HOH:O	1.84	0.60
4:B:1853:HOH:O	1:C:140:LYS:HE2	2.01	0.59
1:B:151:GLU:O	1:B:154:PHE:HB3	2.01	0.59
1:C:323:TYR:HD1	1:C:324:THR:HG23	1.68	0.59
1:C:195:LYS:HE2	3:C:1611:GOL:O3	2.03	0.58
1:C:236:ILE:HD12	1:C:299:MSE:HE1	1.85	0.58
1:D:306:ASN:O	1:D:308:LEU:HD13	2.04	0.58
1:C:260:TRP:HA	1:C:263:LYS:HD3	1.84	0.58
1:D:65:ASP:O	1:D:76:VAL:HA	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:MSE:O	1:A:57:GLN:HG3	2.04	0.57
1:D:77:LEU:HD21	1:D:82:LEU:CD1	2.33	0.57
1:C:46:HIS:HE1	1:C:65:ASP:OD2	1.87	0.57
1:A:80:HIS:H	1:A:80:HIS:CD2	2.21	0.57
1:C:327:VAL:HB	4:C:1859:HOH:O	2.02	0.57
1:C:335:TYR:O	1:C:339:GLY:HA2	2.05	0.57
1:A:168:ILE:HG13	1:A:205:VAL:HG23	1.85	0.57
1:A:149:GLU:O	1:A:153:GLU:HG3	2.05	0.56
1:C:222:LEU:HB3	1:C:226:MSE:HE3	1.86	0.56
1:D:116:LYS:HE3	1:D:144:ARG:HH12	1.71	0.56
1:A:335:TYR:O	1:A:339:GLY:HA2	2.06	0.56
1:D:280:ASP:OD2	1:D:282:HIS:ND1	2.39	0.56
1:D:149:GLU:O	1:D:153:GLU:HG3	2.05	0.56
1:C:218:ILE:O	1:C:223:GLU:HG3	2.06	0.55
1:D:47:THR:HB	1:D:49:PRO:HD2	1.88	0.55
1:A:144:ARG:NE	4:A:1923:HOH:O	2.38	0.55
1:B:144:ARG:NH2	4:B:1904:HOH:O	2.38	0.55
1:D:77:LEU:HD21	1:D:82:LEU:HD12	1.87	0.55
1:D:164:LYS:NZ	4:D:1765:HOH:O	2.39	0.55
1:B:68:MSE:HE1	1:B:183:LYS:HB3	1.88	0.55
1:D:259:ASP:HA	1:D:262:PHE:HD2	1.70	0.55
1:A:201:LYS:HD3	1:A:231:ASN:ND2	2.22	0.55
1:D:219:LYS:HG3	1:D:220:ASN:ND2	2.21	0.55
1:D:176:TRP:CD1	4:D:1767:HOH:O	2.41	0.55
1:C:34:ILE:HD13	1:C:169:TYR:CE1	2.43	0.55
1:D:116:LYS:CE	1:D:144:ARG:NH1	2.70	0.54
1:D:284:LEU:HA	1:D:297:THR:OG1	2.06	0.54
1:A:146:HIS:HE1	1:A:151:GLU:OE1	1.91	0.54
1:B:284:LEU:HD22	1:B:300:VAL:HG23	1.89	0.54
1:A:92:PHE:HB3	1:A:95:ARG:HG3	1.90	0.54
1:C:333:ALA:O	1:C:337:LYS:HB2	2.08	0.54
1:D:281:TYR:OH	1:D:321:PRO:HD3	2.08	0.54
1:D:259:ASP:HA	1:D:262:PHE:CD2	2.43	0.53
1:B:259:ASP:HA	1:B:262:PHE:CD2	2.43	0.53
1:B:337:LYS:HD2	4:B:1926:HOH:O	2.09	0.53
1:D:70:LYS:HE2	4:D:1683:HOH:O	2.09	0.53
1:A:267:MSE:HA	1:A:267:MSE:CE	2.36	0.52
1:A:146:HIS:HD2	4:A:1712:HOH:O	1.92	0.52
1:B:53:MSE:O	1:B:57:GLN:HG3	2.09	0.52
1:B:56:ALA:HB2	1:C:53:MSE:HG3	1.92	0.52
1:A:337:LYS:HD3	4:A:1813:HOH:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:72:ASP:OD2	1:D:183:LYS:HE3	2.11	0.51
1:D:290:GLN:OE1	1:D:290:GLN:HA	2.11	0.51
1:C:210:PHE:CE1	1:C:237:ALA:HB2	2.46	0.50
1:D:201:LYS:HG3	1:D:229:GLU:O	2.11	0.50
1:D:268:LYS:HA	1:D:306:ASN:OD1	2.12	0.50
1:D:296:LEU:HD12	1:D:300:VAL:HG21	1.92	0.50
1:A:316:ARG:O	1:A:320:LEU:HD13	2.11	0.50
1:C:313:TYR:HD1	1:C:314:THR:HG23	1.77	0.50
1:D:286:GLU:OE1	1:D:286:GLU:HA	2.11	0.50
1:C:171:GLU:HG3	1:C:208:GLN:HG3	1.93	0.50
1:C:208:GLN:OE1	1:C:235:LEU:HD21	2.11	0.50
1:B:85:VAL:HG23	1:B:86:THR:HG23	1.93	0.50
1:B:356:ASN:N	1:B:356:ASN:HD22	2.08	0.50
1:B:68:MSE:HE3	1:B:178:HIS:CD2	2.47	0.49
1:C:343:LEU:HD12	1:C:343:LEU:C	2.32	0.49
1:D:113:LYS:NZ	1:D:150:GLU:OE2	2.40	0.49
1:D:228:MSE:HE2	1:D:230:LEU:HD11	1.94	0.49
1:D:198:TYR:O	1:D:228:MSE:HE3	2.12	0.49
1:A:222:LEU:HB3	1:A:226:MSE:HE3	1.94	0.49
1:B:146:HIS:CE1	1:B:151:GLU:OE1	2.66	0.49
1:D:290:GLN:HB3	1:D:291:PRO:HD2	1.94	0.49
1:A:316:ARG:HB2	1:A:319:LYS:HB3	1.94	0.49
1:D:116:LYS:HE2	1:D:116:LYS:HA	1.95	0.49
1:C:47:THR:HB	1:C:49:PRO:HD2	1.93	0.49
1:A:53:MSE:HG3	1:D:56:ALA:HB2	1.95	0.49
1:A:337:LYS:HE3	4:A:1893:HOH:O	2.12	0.49
1:B:68:MSE:CE	1:B:183:LYS:HB3	2.43	0.48
1:C:74:LEU:HD22	1:C:188:LYS:HG3	1.96	0.48
1:C:92:PHE:HB3	1:C:95:ARG:HG3	1.95	0.48
1:C:280:ASP:HA	1:C:313:TYR:O	2.14	0.48
1:D:277:ILE:HD11	1:D:310:VAL:HG22	1.96	0.48
1:D:326:ASP:OD2	1:D:329:GLN:HG3	2.14	0.48
1:D:271:ALA:CB	1:D:308:LEU:HD11	2.44	0.48
1:B:144:ARG:HE	1:B:144:ARG:HB3	1.46	0.47
1:D:46:HIS:CG	1:D:76:VAL:HG12	2.49	0.47
1:D:70:LYS:HE3	1:D:106:ASP:O	2.13	0.47
1:B:281:TYR:OH	1:B:321:PRO:HD3	2.14	0.47
1:D:223:GLU:HB2	1:D:224:PRO:HD3	1.96	0.47
1:C:53:MSE:O	1:C:57:GLN:HG3	2.15	0.47
1:C:116:LYS:NZ	4:C:1773:HOH:O	2.46	0.47
1:C:146:HIS:CE1	1:C:151:GLU:OE1	2.61	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:262:PHE:HZ	1:C:283:MSE:HE2	1.80	0.47
1:C:334:LEU:O	1:C:340:VAL:HG22	2.15	0.47
1:A:151:GLU:O	1:A:154:PHE:HB3	2.15	0.47
1:C:222:LEU:HB3	1:C:226:MSE:HE1	1.97	0.47
1:C:65:ASP:CB	1:C:78:HIS:HA	2.45	0.47
1:D:98:LYS:HG3	4:D:1690:HOH:O	2.13	0.47
1:D:262:PHE:HZ	1:D:283:MSE:HE2	1.80	0.46
1:A:187:ALA:O	1:A:191:GLU:HG3	2.14	0.46
1:A:142:ASP:OD2	1:A:144:ARG:NH2	2.47	0.46
1:B:220:ASN:N	1:B:220:ASN:HD22	2.13	0.46
1:C:300:VAL:HG23	4:C:1866:HOH:O	2.15	0.46
1:D:116:LYS:CE	1:D:144:ARG:HH12	2.26	0.46
1:C:65:ASP:O	1:C:76:VAL:HA	2.15	0.46
1:B:146:HIS:CE1	1:B:151:GLU:HB2	2.50	0.46
1:B:146:HIS:HD2	4:B:1710:HOH:O	1.99	0.46
1:A:201:LYS:HD3	1:A:231:ASN:HD21	1.79	0.46
1:B:343:LEU:O	1:B:343:LEU:HD12	2.16	0.46
1:D:335:TYR:O	1:D:339:GLY:HA2	2.16	0.46
1:C:201:LYS:NZ	4:C:1785:HOH:O	2.44	0.46
1:D:249:PRO:HG2	4:D:1792:HOH:O	2.15	0.46
1:D:265:GLY:HA2	4:D:1756:HOH:O	2.16	0.46
1:A:334:LEU:O	1:A:340:VAL:HG22	2.15	0.46
1:B:213:ASP:OD2	1:B:246:GLN:NE2	2.40	0.46
1:C:37:ARG:HA	1:C:46:HIS:CD2	2.51	0.46
1:A:144:ARG:HG3	1:A:144:ARG:NH1	2.30	0.45
1:B:144:ARG:HD3	4:B:1878:HOH:O	2.16	0.45
1:B:354:PHE:O	1:B:357:LYS:HG3	2.15	0.45
1:D:48:LEU:N	1:D:49:PRO:CD	2.80	0.45
1:C:91:ARG:HD2	4:C:1777:HOH:O	2.16	0.45
1:D:37:ARG:HA	1:D:46:HIS:CE1	2.52	0.45
1:C:235:LEU:HD23	1:C:278:GLY:HA3	1.99	0.45
1:D:57:GLN:CD	1:D:349:ASP:HB3	2.37	0.45
1:A:77:LEU:HD21	1:A:82:LEU:HD23	1.98	0.45
1:B:149:GLU:O	1:B:153:GLU:HG3	2.16	0.45
1:B:334:LEU:O	1:B:340:VAL:HG22	2.17	0.45
1:C:223:GLU:HB2	1:C:224:PRO:HD3	1.98	0.45
1:D:171:GLU:HA	1:D:208:GLN:O	2.17	0.45
1:D:175:PRO:HG2	4:D:1664:HOH:O	2.16	0.45
1:A:98:LYS:HG3	4:A:1689:HOH:O	2.16	0.45
1:C:281:TYR:CD1	1:C:281:TYR:C	2.89	0.45
1:B:316:ARG:O	1:B:320:LEU:HD13	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:242:ASN:N	1:C:242:ASN:HD22	2.15	0.45
1:C:92:PHE:HB3	1:C:95:ARG:CG	2.47	0.45
1:B:303:ALA:O	1:B:308:LEU:HB2	2.16	0.44
1:D:316:ARG:HB2	1:D:319:LYS:HB3	1.99	0.44
1:C:105:ILE:HD11	1:C:183:LYS:CD	2.46	0.44
1:C:242:ASN:HA	1:C:255:ASN:OD1	2.17	0.44
1:B:187:ALA:O	1:B:191:GLU:HG3	2.17	0.44
1:B:108:THR:OG1	1:B:111:GLU:HG3	2.18	0.44
1:D:325:PRO:HD2	1:D:329:GLN:CD	2.38	0.44
1:D:325:PRO:HD2	1:D:329:GLN:OE1	2.17	0.44
1:C:323:TYR:CD1	1:C:324:THR:HG23	2.50	0.44
1:D:228:MSE:CE	1:D:230:LEU:HD11	2.48	0.44
1:D:304:GLN:O	1:D:307:LYS:HD2	2.18	0.44
1:C:190:LEU:HD11	1:C:218:ILE:HG23	1.99	0.44
1:C:236:ILE:CD1	1:C:299:MSE:HE1	2.46	0.44
1:C:277:ILE:HD12	1:C:299:MSE:HE2	1.99	0.44
1:A:309:VAL:HG23	1:A:341:ASN:ND2	2.32	0.43
1:B:68:MSE:HE1	1:B:183:LYS:CB	2.48	0.43
1:D:176:TRP:O	1:D:180:GLN:HG3	2.18	0.43
1:A:80:HIS:H	1:A:80:HIS:HD2	1.66	0.43
1:A:96:ALA:HB2	1:A:102:TYR:CZ	2.53	0.43
1:A:343:LEU:HD12	1:A:343:LEU:C	2.39	0.43
1:C:83:ASP:HB3	1:C:102:TYR:HD2	1.83	0.43
1:D:324:THR:HG21	1:D:330:LEU:HD13	2.00	0.43
1:C:29:ASN:O	1:C:30:GLU:CB	2.67	0.43
1:C:176:TRP:O	1:C:180:GLN:HG3	2.18	0.43
1:A:208:GLN:OE1	1:A:235:LEU:HD21	2.19	0.43
1:C:46:HIS:HA	1:C:51:LYS:HZ1	1.84	0.43
1:C:46:HIS:CE1	1:C:65:ASP:OD2	2.71	0.43
1:D:301:GLN:HA	1:D:301:GLN:OE1	2.18	0.43
1:B:127:LYS:NZ	4:B:1906:HOH:O	2.48	0.43
1:B:323:TYR:HD1	1:B:324:THR:HG23	1.84	0.43
1:C:96:ALA:HB2	1:C:102:TYR:CE1	2.54	0.43
1:C:277:ILE:CD1	1:C:299:MSE:HE2	2.49	0.43
1:D:248:GLN:HB3	1:D:249:PRO:HD2	2.00	0.43
1:B:261:MSE:HA	1:B:266:ALA:HB1	2.01	0.43
1:C:31:LYS:HG2	1:C:335:TYR:CD1	2.54	0.43
1:D:30:GLU:O	1:D:341:ASN:ND2	2.52	0.43
1:C:334:LEU:C	1:C:340:VAL:HG22	2.39	0.42
1:D:270:VAL:HA	1:D:273:TYR:CE2	2.54	0.42
1:A:312:PRO:HD2	1:A:342:GLY:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:162:THR:HG22	1:D:42:TYR:CD2	2.54	0.42
1:B:30:GLU:O	1:B:30:GLU:CG	2.68	0.42
1:C:219:LYS:NZ	4:C:1850:HOH:O	2.33	0.42
1:D:335:TYR:CZ	1:D:343:LEU:HG	2.54	0.42
1:A:323:TYR:HD1	1:A:324:THR:HG23	1.84	0.42
1:D:97:ARG:HD2	1:D:103:TYR:CE2	2.54	0.42
1:C:127:LYS:HG2	1:C:128:LYS:N	2.34	0.42
1:D:277:ILE:CG1	1:D:310:VAL:HG22	2.49	0.42
1:B:65:ASP:O	1:B:76:VAL:HA	2.20	0.42
1:B:210:PHE:HE2	3:B:1606:GOL:H32	1.83	0.42
1:C:313:TYR:HA	1:C:314:THR:HA	1.78	0.42
1:D:310:VAL:HG12	1:D:312:PRO:HD3	2.01	0.42
1:B:68:MSE:HE3	1:B:178:HIS:HD2	1.84	0.42
1:C:276:GLY:HA3	1:C:309:VAL:HG12	2.02	0.42
1:C:323:TYR:O	1:C:325:PRO:HD3	2.20	0.42
1:D:167:GLY:HA3	1:D:204:LYS:HB3	2.01	0.42
1:A:325:PRO:HD2	1:A:329:GLN:OE1	2.20	0.41
1:B:223:GLU:HB2	1:B:224:PRO:HD3	2.01	0.41
1:B:284:LEU:HD13	1:B:300:VAL:CG2	2.50	0.41
1:D:152:ILE:O	1:D:156:GLN:HG3	2.20	0.41
1:D:313:TYR:HA	1:D:314:THR:HA	1.87	0.41
1:C:322:GLU:OE2	1:C:322:GLU:CA	2.67	0.41
1:D:280:ASP:HA	1:D:313:TYR:O	2.20	0.41
1:D:343:LEU:HD12	1:D:343:LEU:O	2.20	0.41
1:A:147:THR:OG1	1:A:150:GLU:HG3	2.21	0.41
1:C:68:MSE:HE3	1:C:188:LYS:HG2	2.01	0.41
1:C:312:PRO:HD2	1:C:342:GLY:O	2.19	0.41
1:D:285:ILE:HG21	1:D:323:TYR:CE1	2.55	0.41
1:D:347:PHE:CD1	1:D:347:PHE:N	2.88	0.41
1:B:210:PHE:CE1	1:B:237:ALA:HB2	2.56	0.41
1:A:127:LYS:NZ	4:A:1875:HOH:O	2.53	0.41
1:D:260:TRP:CZ2	1:D:261:MSE:HE2	2.56	0.41
1:B:91:ARG:HE	1:B:91:ARG:HB3	1.54	0.41
1:C:217:ARG:O	1:C:221:GLU:HB2	2.20	0.41
1:D:170:PRO:O	1:D:208:GLN:HG2	2.21	0.41
1:D:262:PHE:CZ	1:D:283:MSE:HE2	2.55	0.41
1:A:116:LYS:HD2	4:A:1896:HOH:O	2.21	0.41
1:C:36:HIS:CE1	1:C:314:THR:HG21	2.56	0.41
1:B:266:ALA:O	1:B:269:GLN:HB2	2.20	0.41
1:C:88:VAL:HG13	1:C:89:ALA:N	2.35	0.41
1:C:90:ASP:HB3	1:C:91:ARG:HH21	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:48:LEU:HD23	1:D:48:LEU:HA	1.87	0.41
1:D:238:TYR:HD2	1:D:241:TRP:CZ2	2.39	0.40
1:B:80:HIS:HD2	4:B:1726:HOH:O	2.05	0.40
1:C:221:GLU:OE1	4:C:1722:HOH:O	2.21	0.40
1:C:291:PRO:C	1:C:293:ASN:H	2.24	0.40
1:C:210:PHE:HE2	3:C:1607:GOL:H32	1.86	0.40
1:D:341:ASN:HB2	4:D:1742:HOH:O	2.22	0.40
1:A:267:MSE:CE	1:A:270:VAL:CB	2.96	0.40
1:C:195:LYS:CE	3:C:1611:GOL:O3	2.67	0.40
1:C:222:LEU:C	1:C:226:MSE:HE3	2.42	0.40
1:C:276:GLY:HA3	1:C:309:VAL:CG1	2.52	0.40
1:C:166:ILE:N	1:C:166:ILE:HD13	2.36	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	326/336 (97%)	309 (95%)	17 (5%)	0	100	100
1	B	327/336 (97%)	314 (96%)	13 (4%)	0	100	100
1	C	326/336 (97%)	304 (93%)	21 (6%)	1 (0%)	41	37
1	D	325/336 (97%)	304 (94%)	18 (6%)	3 (1%)	17	11
All	All	1304/1344 (97%)	1231 (94%)	69 (5%)	4 (0%)	41	37

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	30	GLU
1	D	46	HIS
1	D	301	GLN

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Mol	Chain	Res	Type
1	D	300	VAL

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	283/280 (101%)	275 (97%)	8 (3%)	43	44
1	B	284/280 (101%)	277 (98%)	7 (2%)	47	49
1	C	283/280 (101%)	276 (98%)	7 (2%)	47	49
1	D	282/280 (101%)	276 (98%)	6 (2%)	53	57
All	All	1132/1120 (101%)	1104 (98%)	28 (2%)	47	49

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	GLU
1	A	91	ARG
1	A	164	LYS
1	A	169	TYR
1	A	308	LEU
1	A	320	LEU
1	A	341	ASN
1	A	356	ASN
1	B	144	ARG
1	B	169	TYR
1	B	270	VAL
1	B	287	GLU
1	B	308	LEU
1	B	309	VAL
1	B	320	LEU
1	C	158	LEU
1	C	169	TYR
1	C	242	ASN
1	C	293	ASN
1	C	325	PRO

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Mol	Chain	Res	Type
1	C	327	VAL
1	C	341	ASN
1	D	82	LEU
1	D	169	TYR
1	D	279	PRO
1	D	314	THR
1	D	326	ASP
1	D	341	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	HIS
1	A	146	HIS
1	A	220	ASN
1	A	231	ASN
1	A	242	ASN
1	A	290	GLN
1	A	305	GLN
1	A	336	ASN
1	A	341	ASN
1	A	356	ASN
1	B	80	HIS
1	B	146	HIS
1	B	220	ASN
1	B	248	GLN
1	B	304	GLN
1	B	329	GLN
1	B	336	ASN
1	B	356	ASN
1	C	46	HIS
1	C	146	HIS
1	C	165	ASN
1	C	242	ASN
1	C	336	ASN
1	D	146	HIS
1	D	220	ASN
1	D	242	ASN
1	D	305	GLN
1	D	336	ASN
1	D	341	ASN

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)

Of 15 ligands modelled in this entry, 4 are monoatomic - leaving 11 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	$\#\|Z\| > 2$
3	GOL	C	1611	-	5,5,5	0.46	0	5,5,5	0.54	0
3	GOL	A	1605	2	5,5,5	0.35	0	5,5,5	0.37	0
3	GOL	C	1613	-	5,5,5	0.44	0	5,5,5	0.50	0
3	GOL	D	1612	-	5,5,5	0.46	0	5,5,5	0.52	0
3	GOL	C	1607	2	5,5,5	0.33	0	5,5,5	0.40	0
3	GOL	A	1609	-	5,5,5	0.37	0	5,5,5	0.28	0
3	GOL	D	1608	2	5,5,5	0.34	0	5,5,5	0.19	0
3	GOL	B	1606	2	5,5,5	0.44	0	5,5,5	0.48	0
3	GOL	B	1610	-	5,5,5	0.43	0	5,5,5	0.52	0
3	GOL	D	1615	-	5,5,5	0.35	0	5,5,5	0.75	0
3	GOL	A	1614	-	5,5,5	0.37	0	5,5,5	0.35	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	GOL	C	1611	-	-	0/4/4/4	-
3	GOL	A	1605	2	-	1/4/4/4	-
3	GOL	C	1613	-	-	0/4/4/4	-
3	GOL	D	1612	-	-	0/4/4/4	-
3	GOL	C	1607	2	-	0/4/4/4	-
3	GOL	A	1609	-	-	0/4/4/4	-
3	GOL	D	1608	2	-	0/4/4/4	-
3	GOL	B	1606	2	-	0/4/4/4	-
3	GOL	B	1610	-	-	0/4/4/4	-
3	GOL	D	1615	-	-	0/4/4/4	-
3	GOL	A	1614	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1605	GOL	C1-C2-C3-O3

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1611	GOL	2	0
3	C	1607	GOL	1	0
3	B	1606	GOL	1	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	319/336 (94%)	-0.08	2 (0%)	89	88	18, 28, 46, 75	0
1	B	320/336 (95%)	-0.02	3 (0%)	84	83	17, 26, 47, 78	0
1	C	319/336 (94%)	0.44	14 (4%)	34	33	17, 37, 56, 67	0
1	D	318/336 (94%)	0.61	28 (8%)	10	9	26, 41, 60, 73	0
All	All	1276/1344 (94%)	0.24	47 (3%)	41	41	17, 33, 55, 78	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	29	ASN	10.9
1	A	29	ASN	6.2
1	D	292	GLY	4.4
1	D	222	LEU	3.5
1	B	30	GLU	3.5
1	D	291	PRO	3.4
1	C	251	GLY	3.2
1	D	354	PHE	3.2
1	C	292	GLY	3.2
1	D	293	ASN	3.1
1	C	298	GLY	3.0
1	D	298	GLY	3.0
1	D	301	GLN	3.0
1	B	357	LYS	2.9
1	D	287	GLU	2.9
1	C	293	ASN	2.8
1	D	307	LYS	2.7
1	D	268	LYS	2.7
1	D	341	ASN	2.7
1	D	326	ASP	2.7
1	C	249	PRO	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	325	PRO	2.7
1	D	263	LYS	2.7
1	D	304	GLN	2.6
1	D	251	GLY	2.6
1	D	266	ALA	2.5
1	D	325	PRO	2.5
1	D	250	ASP	2.4
1	D	249	PRO	2.4
1	C	337	LYS	2.4
1	D	290	GLN	2.4
1	C	288	THR	2.3
1	D	288	THR	2.3
1	D	305	GLN	2.3
1	D	264	PRO	2.2
1	C	286	GLU	2.2
1	D	225	LYS	2.2
1	D	30	GLU	2.2
1	D	327	VAL	2.2
1	C	144	ARG	2.2
1	C	295	LYS	2.1
1	D	337	LYS	2.1
1	D	295	LYS	2.1
1	A	341	ASN	2.1
1	C	248	GLN	2.0
1	C	290	GLN	2.0
1	C	29	ASN	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	GOL	D	1612	6/6	0.59	0.33	50,53,54,54	0
3	GOL	C	1611	6/6	0.63	0.50	50,52,54,54	0
3	GOL	B	1610	6/6	0.64	0.37	45,51,53,55	0
3	GOL	A	1609	6/6	0.73	0.44	56,56,56,57	0
3	GOL	C	1613	6/6	0.76	0.31	45,48,48,48	0
3	GOL	D	1615	6/6	0.79	0.35	45,47,48,49	0
2	MG	D	1604	1/1	0.88	0.20	35,35,35,35	0
3	GOL	D	1608	6/6	0.89	0.27	38,41,42,42	0
3	GOL	A	1614	6/6	0.89	0.30	42,45,49,52	0
2	MG	C	1603	1/1	0.89	0.14	29,29,29,29	0
3	GOL	C	1607	6/6	0.91	0.23	33,35,37,38	0
3	GOL	A	1605	6/6	0.94	0.20	28,32,33,33	0
2	MG	A	1601	1/1	0.95	0.18	27,27,27,27	0
3	GOL	B	1606	6/6	0.96	0.20	23,26,28,29	0
2	MG	B	1602	1/1	0.99	0.19	22,22,22,22	0

6.5 Other polymers (i)

There are no such residues in this entry.