Full wwPDB X-ray Structure Validation Report (i)

Jun 12, 2024 – 11:51 PM EDT

PDB ID	:	3WIC
Title	:	Structure of a substrate/cofactor-unbound glucose dehydrogenase
Authors	:	Sakuraba, H.; Kanoh, Y.; Yoneda, K.; Ohshima, T.
Deposited on	:	2013-09-10
Resolution	:	2.60 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:

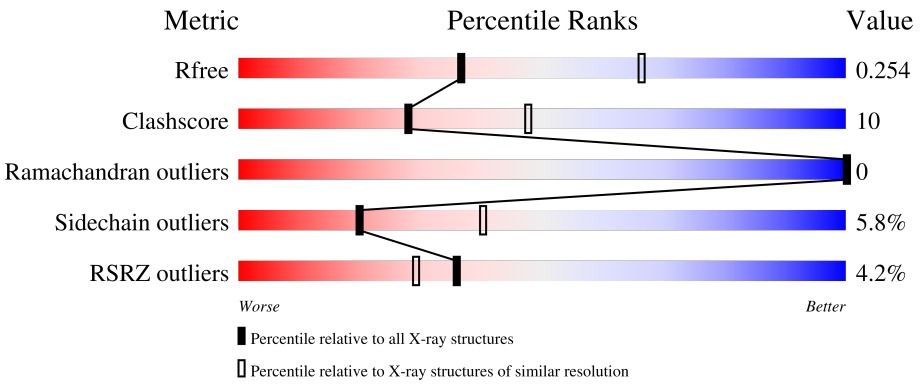
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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	361	2% 77% 20% •
1	B	361	6% 76% 20% •
1	C	361	4% 77% 21% •
1	D	361	5% 81% 17% •

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	PGO	B	1004	-	-	X	-
4	1PE	A	1004	-	-	-	X
4	1PE	B	1005	-	-	-	X
4	1PE	B	1007	-	-	-	X

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 11696 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.

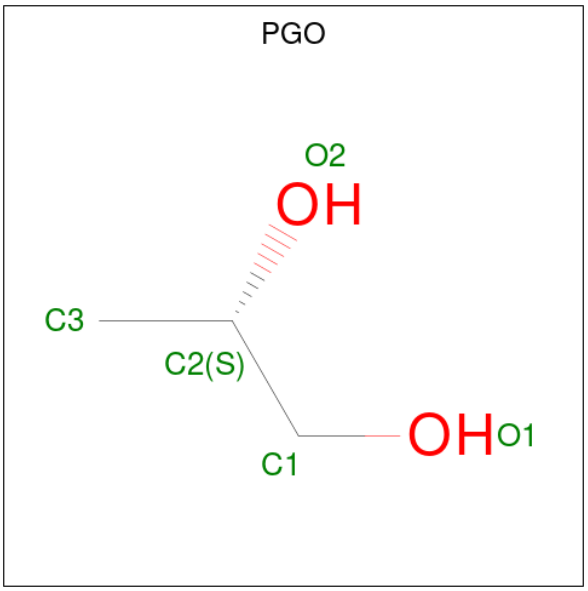
- Molecule 1 is a protein called Glucose 1-dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	360	2824	1799	473	539	13	0	0	0
1	B	360	2824	1799	473	539	13	0	0	0
1	C	360	2824	1799	473	539	13	0	0	0
1	D	360	2824	1799	473	539	13	0	0	0

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

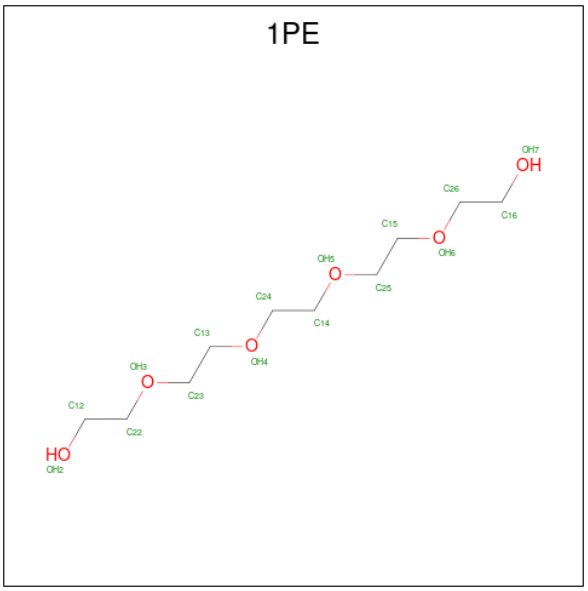
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
2	A	2	2	2	0	0
2	B	2	2	2	0	0
2	C	2	2	2	0	0
2	D	2	2	2	0	0

- Molecule 3 is S-1,2-PROPANEDIOL (three-letter code: PGO) (formula: $C_3H_8O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	5	3	2	0	0
3	B	1	5	3	2	0	0
3	B	1	5	3	2	0	0
3	C	1	5	3	2	0	0
3	D	1	5	3	2	0	0
3	D	1	5	3	2	0	0

- Molecule 4 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: $C_{10}H_{22}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
4	A	1	16	10	6	0	0
			Total	C	O		
4	A	1	16	10	6	0	0
			Total	C	O		
4	B	1	16	10	6	0	0
			Total	C	O		
4	B	1	16	10	6	0	0
			Total	C	O		
4	B	1	16	10	6	0	0
			Total	C	O		
4	C	1	16	10	6	0	0
			Total	C	O		
4	C	1	16	10	6	0	0
			Total	C	O		
4	D	1	16	10	6	0	0

- Molecule 5 is water.

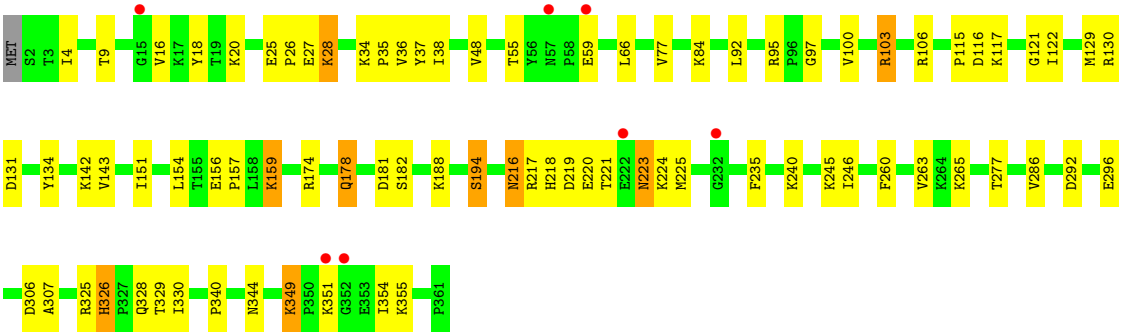
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	51	51	51	0	0
			Total	O		
5	B	53	53	53	0	0
			Total	O		
5	C	63	63	63	0	0
			Total	O		
5	D	67	67	67	0	0

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.

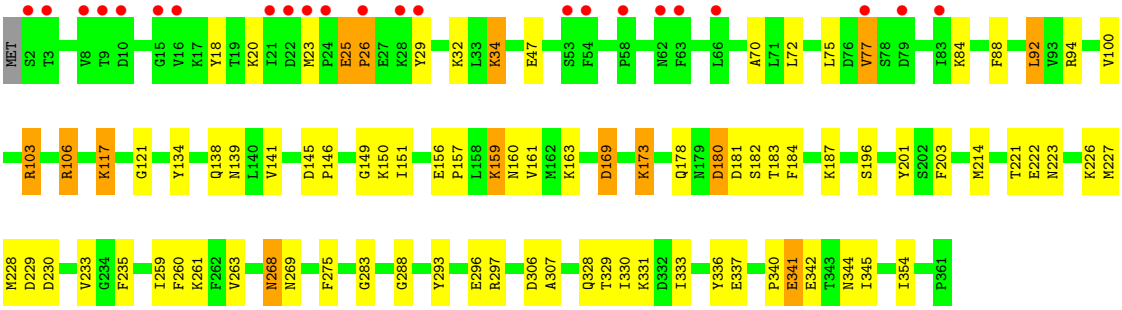
• Molecule 1: Glucose 1-dehydrogenase

Chain A:



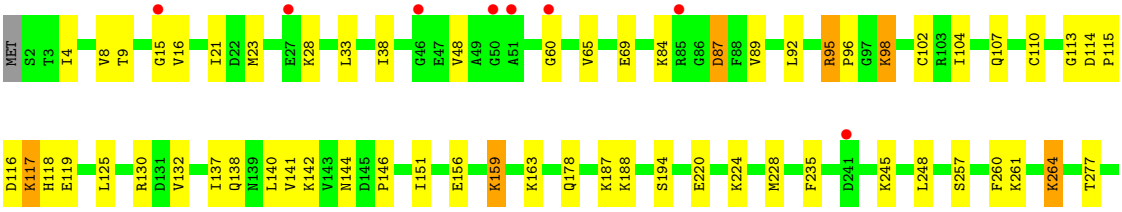
• Molecule 1: Glucose 1-dehydrogenase

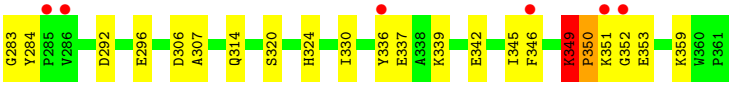
Chain B:



• Molecule 1: Glucose 1-dehydrogenase

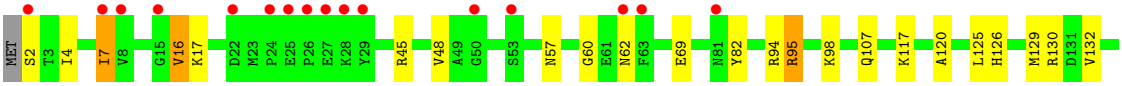
Chain C:

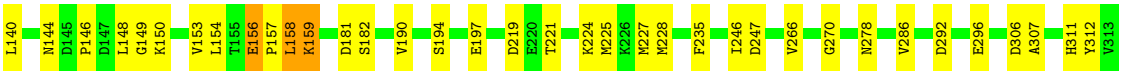


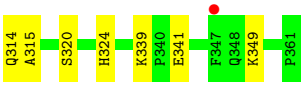


- Molecule 1: Glucose 1-dehydrogenase

Chain D:







4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	79.09Å 122.17Å 87.43Å 90.00° 109.92° 90.00°	Depositor
Resolution (Å)	49.08 – 2.60 49.03 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.08-2.60) 99.8 (49.03-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	7.62 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.205 , 0.257 0.211 , 0.254	Depositor DCC
R_{free} test set	2432 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å^2)	36.9	Xtriage
Anisotropy	0.284	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 54.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.52$, $< L^2 > = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	11696	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 4.36% 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: 1PE, PGO, ZN

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	0.75	0/2883	0.85	3/3898 (0.1%)
1	B	0.71	2/2883 (0.1%)	0.81	4/3898 (0.1%)
1	C	0.75	1/2883 (0.0%)	0.81	2/3898 (0.1%)
1	D	0.73	0/2883	0.82	2/3898 (0.1%)
All	All	0.73	3/11532 (0.0%)	0.82	11/15592 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	26	PRO	N-CD	5.59	1.55	1.47
1	B	180	ASP	CB-CG	-5.32	1.40	1.51
1	C	350	PRO	N-CD	5.30	1.55	1.47

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	326	HIS	C-N-CD	6.00	140.99	128.40
1	D	95	ARG	NE-CZ-NH1	5.93	123.27	120.30
1	D	247	ASP	CB-CG-OD1	-5.93	112.96	118.30
1	B	25	GLU	C-N-CD	5.59	140.13	128.40
1	A	349	LYS	C-N-CD	5.54	140.03	128.40
1	B	169	ASP	CB-CG-OD1	5.50	123.25	118.30
1	B	180	ASP	CB-CA-C	-5.49	99.42	110.40
1	C	87	ASP	CB-CG-OD1	5.43	123.19	118.30
1	C	349	LYS	C-N-CD	5.28	139.49	128.40
1	A	174	ARG	NE-CZ-NH2	-5.25	117.67	120.30
1	B	106	ARG	NE-CZ-NH1	5.15	122.88	120.30

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	2824	0	2805	46	0
1	B	2824	0	2805	68	0
1	C	2824	0	2806	66	0
1	D	2824	0	2805	45	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	1	0
2	D	2	0	0	0	0
3	A	5	0	8	1	0
3	B	10	0	16	5	0
3	C	5	0	8	1	0
3	D	10	0	16	2	0
4	A	32	0	44	6	0
4	B	48	0	66	8	0
4	C	32	0	44	7	0
4	D	16	0	22	2	0
5	A	51	0	0	4	0
5	B	53	0	0	3	0
5	C	63	0	0	12	0
5	D	67	0	0	4	0
All	All	11696	0	11445	226	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 10.

All (226) 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:351:LYS:HD2	1:C:352:GLY:N	1.59	1.16
1:C:21:ILE:HD11	1:C:23:MET:HE2	1.47	0.97
1:A:325:ARG:O	1:A:326:HIS:ND1	2.00	0.93
1:C:141:VAL:HA	4:C:1005:1PE:H242	1.51	0.92
1:C:359:LYS:HD2	5:C:1162:HOH:O	1.72	0.89
1:C:351:LYS:HD2	1:C:352:GLY:H	1.39	0.87
1:A:181:ASP:O	1:A:182:SER:OG	1.92	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:VAL:HA	1:B:103:ARG:HD3	1.59	0.84
1:D:181:ASP:O	1:D:182:SER:OG	1.96	0.84
1:B:184:PHE:O	1:B:187:LYS:HB2	1.76	0.83
1:B:25:GLU:HG2	1:B:26:PRO:HD2	1.61	0.82
1:A:159:LYS:HD3	1:A:307:ALA:HB3	1.64	0.79
1:A:220:GLU:OE1	1:A:224:LYS:HD2	1.84	0.77
1:D:69:GLU:HG2	1:D:156:GLU:HG2	1.65	0.77
1:A:100:VAL:HA	1:A:103:ARG:HD3	1.66	0.75
1:C:4:ILE:HG21	1:C:132:VAL:HG12	1.69	0.75
1:A:328:GLN:HA	1:A:328:GLN:NE2	2.02	0.75
1:B:26:PRO:CG	1:B:29:TYR:HB3	2.18	0.74
1:C:337:GLU:HB3	5:C:1162:HOH:O	1.87	0.74
1:A:9:THR:HB	1:A:66:LEU:HD11	1.71	0.72
5:A:1143:HOH:O	4:B:1005:1PE:H142	1.89	0.72
5:A:1143:HOH:O	4:B:1005:1PE:C14	2.38	0.72
1:B:145:ASP:O	3:B:1004:PGO:H12	1.94	0.68
1:A:221:THR:O	1:A:225:MET:HG2	1.95	0.67
1:D:4:ILE:HG21	1:D:132:VAL:HG12	1.77	0.65
1:B:214:MET:HE3	1:B:233:VAL:HG11	1.77	0.65
1:D:69:GLU:HG2	1:D:156:GLU:CG	2.25	0.65
1:B:26:PRO:HG2	1:B:29:TYR:HB3	1.77	0.64
1:B:156:GLU:HB3	1:B:157:PRO:HD3	1.80	0.64
1:C:98:LYS:NZ	5:C:1158:HOH:O	2.31	0.64
1:D:7:ILE:HG23	1:D:129:MET:HB3	1.80	0.64
1:C:21:ILE:CD1	1:C:23:MET:HE2	2.26	0.64
1:A:194:SER:OG	1:A:216:ASN:ND2	2.24	0.63
1:C:351:LYS:CD	1:C:352:GLY:N	2.51	0.63
1:B:26:PRO:HG2	1:B:29:TYR:CD2	2.34	0.63
1:C:110:CYS:SG	2:C:1002:ZN:ZN	1.88	0.62
1:C:359:LYS:HB2	5:C:1162:HOH:O	2.00	0.62
4:D:1005:1PE:C25	5:D:1155:HOH:O	2.48	0.62
1:C:351:LYS:HD2	1:C:351:LYS:C	2.18	0.61
1:B:34:LYS:HG3	1:B:75:LEU:HD11	1.81	0.61
1:B:26:PRO:HG2	1:B:29:TYR:HD2	1.66	0.61
1:C:23:MET:HE3	1:C:130:ARG:HD3	1.83	0.60
1:B:25:GLU:CG	1:B:26:PRO:HD2	2.31	0.59
4:B:1007:1PE:H262	1:C:178:GLN:OE1	2.01	0.59
1:A:218:HIS:HD2	1:A:219:ASP:O	1.85	0.58
1:B:228:MET:SD	1:B:235:PHE:HB2	2.44	0.58
1:D:159:LYS:HG2	1:D:307:ALA:HB3	1.84	0.58
1:B:25:GLU:HG2	1:B:26:PRO:CD	2.31	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:145:ASP:O	3:B:1004:PGO:C1	2.52	0.57
1:D:306:ASP:OD1	1:D:307:ALA:N	2.33	0.57
1:C:314:GLN:CD	5:C:1130:HOH:O	2.42	0.57
1:C:23:MET:CE	1:C:130:ARG:HD3	2.34	0.56
1:B:18:TYR:CZ	1:B:340:PRO:HB2	2.41	0.56
1:C:260:PHE:O	1:C:264:LYS:HD3	2.04	0.56
1:D:156:GLU:O	1:D:159:LYS:HD3	2.05	0.56
1:C:110:CYS:SG	1:C:116:ASP:HB3	2.45	0.56
1:D:126:HIS:O	3:D:1003:PGO:O2	2.20	0.56
1:B:149:GLY:N	3:B:1004:PGO:O1	2.31	0.56
1:B:221:THR:OG1	1:B:223:ASN:ND2	2.39	0.56
1:C:138:GLN:HA	4:C:1005:1PE:H152	1.87	0.55
1:B:328:GLN:O	1:B:331:LYS:HB2	2.05	0.55
1:C:117:LYS:HD3	1:C:119:GLU:OE2	2.07	0.55
1:D:82:TYR:HB2	5:D:1155:HOH:O	2.05	0.55
1:B:25:GLU:CG	1:B:26:PRO:CD	2.85	0.55
5:A:1143:HOH:O	4:B:1005:1PE:H141	2.03	0.55
1:B:293:TYR:CE1	1:B:297:ARG:CZ	2.90	0.55
1:B:306:ASP:OD1	1:B:307:ALA:N	2.33	0.54
1:B:342:GLU:O	1:B:345:ILE:HG12	2.07	0.54
1:D:190:VAL:HG21	1:D:246:ILE:HG12	1.90	0.54
1:D:144:ASN:O	1:D:146:PRO:HD3	2.07	0.54
4:A:1004:1PE:C26	1:B:178:GLN:NE2	2.71	0.54
1:D:16:VAL:HG13	1:D:48:VAL:HG22	1.89	0.53
1:C:339:LYS:NZ	5:C:1141:HOH:O	2.41	0.53
1:D:311:HIS:HA	1:D:314:GLN:HG2	1.89	0.53
1:B:92:LEU:HD22	1:B:139:ASN:HA	1.90	0.53
1:D:130:ARG:HD2	1:D:132:VAL:O	2.08	0.53
1:B:203:PHE:HE2	1:B:333:ILE:HD11	1.74	0.53
1:C:137:ILE:HG22	4:C:1005:1PE:H251	1.89	0.53
1:D:292:ASP:O	1:D:296:GLU:HG2	2.08	0.53
1:A:221:THR:O	1:A:225:MET:CG	2.57	0.52
1:C:60:GLY:CA	5:C:1142:HOH:O	2.57	0.52
1:D:221:THR:O	1:D:225:MET:HG2	2.09	0.52
1:C:349:LYS:HB2	1:C:353:GLU:OE1	2.10	0.52
1:C:104:ILE:HD12	4:C:1004:1PE:H161	1.91	0.52
1:B:106:ARG:NH1	4:B:1005:1PE:H241	2.24	0.51
1:B:283:GLY:HA2	1:D:286:VAL:O	2.11	0.51
1:C:60:GLY:N	5:C:1142:HOH:O	2.42	0.51
1:B:18:TYR:CD2	1:B:341:GLU:HB3	2.46	0.51
1:C:69:GLU:HG2	1:C:156:GLU:OE2	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:156:GLU:O	1:C:159:LYS:HE3	2.11	0.51
1:C:350:PRO:O	1:C:353:GLU:HB2	2.11	0.51
1:B:141:VAL:HA	4:B:1006:1PE:H132	1.93	0.50
1:D:149:GLY:HA2	3:D:1004:PGO:H2	1.93	0.50
1:A:97:GLY:HA3	1:A:116:ASP:O	2.12	0.50
1:C:264:LYS:NZ	1:C:284:TYR:OH	2.43	0.50
1:B:26:PRO:HG3	1:B:29:TYR:HB3	1.91	0.50
4:C:1004:1PE:H162	1:D:182:SER:HA	1.94	0.50
1:A:156:GLU:O	1:A:159:LYS:HE3	2.12	0.50
1:C:188:LYS:HD2	1:C:245:LYS:O	2.12	0.49
1:B:181:ASP:O	1:B:182:SER:OG	2.19	0.49
1:A:188:LYS:HD2	1:A:245:LYS:O	2.13	0.49
1:A:292:ASP:OD2	1:C:277:THR:CG2	2.61	0.49
1:B:151:ILE:HD11	1:B:331:LYS:HA	1.94	0.49
1:D:150:LYS:O	1:D:153:VAL:HG12	2.13	0.49
4:D:1005:1PE:C14	5:D:1155:HOH:O	2.61	0.49
1:A:220:GLU:HG2	1:A:235:PHE:CZ	2.48	0.49
1:B:18:TYR:CE1	1:B:340:PRO:HB2	2.48	0.49
1:B:156:GLU:O	1:B:159:LYS:HG3	2.13	0.48
1:A:246:ILE:HD12	1:A:265:LYS:HB2	1.96	0.48
1:C:159:LYS:HD3	1:C:307:ALA:HB3	1.95	0.48
1:D:224:LYS:HA	1:D:227:MET:CE	2.43	0.48
1:D:7:ILE:HD11	1:D:16:VAL:CG1	2.43	0.48
1:C:320:SER:O	1:C:324:HIS:HD2	1.96	0.48
1:D:7:ILE:HD11	1:D:16:VAL:HG12	1.94	0.48
1:A:221:THR:OG1	1:A:223:ASN:ND2	2.46	0.47
1:B:259:ILE:O	1:B:263:VAL:HG23	2.14	0.47
1:B:47:GLU:OE1	1:B:121:GLY:O	2.32	0.47
1:A:34:LYS:HG3	1:A:35:PRO:HD2	1.96	0.47
1:B:261:LYS:NZ	5:B:1133:HOH:O	2.45	0.47
1:A:286:VAL:O	1:C:283:GLY:HA2	2.15	0.47
1:B:138:GLN:HG2	4:B:1006:1PE:H251	1.97	0.47
1:B:94:ARG:HD3	1:B:117:LYS:HG3	1.97	0.47
1:B:88:PHE:CG	3:B:1004:PGO:H31	2.50	0.46
1:B:268:ASN:N	1:B:268:ASN:HD22	2.12	0.46
1:A:292:ASP:O	1:A:296:GLU:HG2	2.15	0.46
1:C:84:LYS:HG2	1:C:87:ASP:OD2	2.15	0.46
1:A:36:VAL:HG12	1:A:37:TYR:CD1	2.50	0.46
1:D:94:ARG:HD3	1:D:107:GLN:O	2.15	0.46
5:C:1160:HOH:O	1:D:182:SER:HB2	2.16	0.46
1:A:55:THR:HG22	1:A:122:ILE:HG22	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:159:LYS:HD3	1:B:307:ALA:HB3	1.97	0.46
1:A:4:ILE:HB	1:A:131:ASP:HB2	1.98	0.46
1:A:142:LYS:O	4:A:1005:1PE:H222	2.16	0.46
1:C:89:VAL:HG12	1:C:142:LYS:HA	1.99	0.46
1:B:77:VAL:HG21	1:B:84:LYS:HA	1.98	0.45
1:B:228:MET:HB3	1:B:233:VAL:O	2.16	0.45
1:C:342:GLU:O	1:C:345:ILE:HG12	2.16	0.45
1:C:114:ASP:HA	1:C:115:PRO:HA	1.60	0.45
1:B:159:LYS:HD2	1:B:160:ASN:N	2.32	0.45
1:C:4:ILE:HG21	1:C:132:VAL:CG1	2.43	0.45
1:C:351:LYS:CD	1:C:352:GLY:H	2.19	0.45
1:C:151:ILE:HD12	1:C:330:ILE:HG22	1.99	0.45
1:A:151:ILE:HD12	1:A:330:ILE:HG22	1.98	0.45
1:B:146:PRO:C	3:B:1004:PGO:H12	2.37	0.45
1:C:23:MET:HE1	1:C:65:VAL:HG21	1.97	0.45
1:C:292:ASP:O	1:C:296:GLU:HG2	2.17	0.45
1:D:339:LYS:HB3	1:D:341:GLU:OE1	2.16	0.45
1:D:158:LEU:HD13	1:D:315:ALA:HA	1.99	0.45
1:B:26:PRO:CG	1:B:29:TYR:CD2	2.99	0.45
1:D:120:ALA:HA	1:D:125:LEU:HB3	1.99	0.45
1:B:330:ILE:HD13	1:B:333:ILE:HD12	1.98	0.45
1:A:328:GLN:HA	1:A:328:GLN:HE21	1.79	0.44
1:B:32:LYS:HG3	1:B:134:TYR:CE1	2.52	0.44
4:C:1005:1PE:H161	5:C:1151:HOH:O	2.17	0.44
1:D:154:LEU:O	1:D:157:PRO:HD2	2.18	0.44
1:D:224:LYS:HA	1:D:227:MET:HE3	1.98	0.44
1:C:349:LYS:HB3	1:C:349:LYS:HE2	1.81	0.44
1:A:355:LYS:HE3	5:A:1108:HOH:O	2.17	0.44
1:C:95:ARG:NH1	3:C:1003:PGO:H32	2.32	0.44
1:D:292:ASP:OD1	1:D:296:GLU:OE2	2.36	0.44
1:A:121:GLY:C	1:A:122:ILE:HG13	2.38	0.43
1:A:219:ASP:OD2	1:A:240:LYS:NZ	2.51	0.43
1:B:268:ASN:ND2	1:B:297:ARG:HG2	2.33	0.43
1:D:159:LYS:HG3	1:D:312:TYR:CE1	2.52	0.43
1:B:181:ASP:O	1:B:182:SER:CB	2.67	0.43
1:D:156:GLU:HG3	1:D:157:PRO:HD3	2.00	0.43
1:A:26:PRO:HD2	1:A:134:TYR:CZ	2.53	0.43
1:B:268:ASN:HD22	1:B:268:ASN:H	1.66	0.43
1:D:45:ARG:NH1	1:D:349:LYS:HD2	2.34	0.43
1:A:38:ILE:CD1	1:A:129:MET:HA	2.49	0.43
4:A:1004:1PE:H261	1:B:178:GLN:NE2	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:169:ASP:O	1:B:173:LYS:HG2	2.19	0.43
1:C:314:GLN:NE2	5:C:1130:HOH:O	2.50	0.43
1:B:181:ASP:OD1	1:B:183:THR:OG1	2.25	0.43
1:C:118:HIS:CB	1:C:125:LEU:HD22	2.48	0.43
1:A:292:ASP:OD2	1:C:277:THR:HG21	2.19	0.43
1:C:8:VAL:O	1:C:16:VAL:HA	2.19	0.43
1:C:144:ASN:O	1:C:146:PRO:HD3	2.18	0.43
1:A:106:ARG:HH11	4:A:1004:1PE:C16	2.31	0.43
1:D:45:ARG:NH1	1:D:349:LYS:CD	2.81	0.43
1:D:307:ALA:HB1	1:D:311:HIS:HB2	2.01	0.43
1:C:38:ILE:HD12	1:C:38:ILE:N	2.34	0.42
1:C:228:MET:SD	1:C:235:PHE:HB2	2.59	0.42
1:C:28:LYS:HA	5:C:1154:HOH:O	2.19	0.42
1:A:260:PHE:HA	1:A:263:VAL:HB	2.01	0.42
1:B:106:ARG:NH1	4:B:1005:1PE:C24	2.82	0.42
1:B:260:PHE:HA	1:B:263:VAL:HB	2.01	0.42
1:D:320:SER:O	1:D:324:HIS:HD2	2.01	0.42
1:A:16:VAL:HB	1:A:48:VAL:HG22	2.00	0.42
1:B:150:LYS:NZ	1:B:337:GLU:OE2	2.43	0.42
1:C:118:HIS:HB2	1:C:125:LEU:HD22	2.02	0.42
1:A:106:ARG:HH11	4:A:1004:1PE:H162	1.85	0.42
1:C:9:THR:HA	1:C:15:GLY:O	2.19	0.42
1:B:230:ASP:OD2	1:B:329:THR:HG21	2.19	0.42
1:C:257:SER:OG	1:C:284:TYR:HB2	2.20	0.41
1:A:154:LEU:C	1:A:157:PRO:HD2	2.41	0.41
1:C:102:CYS:SG	1:C:110:CYS:SG	3.15	0.41
1:A:4:ILE:HD12	1:A:130:ARG:HB2	2.02	0.41
1:A:292:ASP:OD2	1:C:277:THR:HG23	2.20	0.41
1:B:269:ASN:ND2	5:B:1147:HOH:O	2.50	0.41
1:C:96:PRO:HD3	1:C:107:GLN:NE2	2.36	0.41
1:C:113:GLY:O	1:C:116:ASP:N	2.37	0.41
1:D:190:VAL:CG2	1:D:246:ILE:HG12	2.50	0.41
1:A:18:TYR:CE1	1:A:340:PRO:HB2	2.54	0.41
1:A:224:LYS:HE2	1:A:354:ILE:HG12	2.03	0.41
1:D:60:GLY:N	5:D:1163:HOH:O	2.46	0.41
1:A:27:GLU:HG3	1:A:28:LYS:H	1.86	0.41
1:A:95:ARG:CZ	3:A:1003:PGO:H32	2.51	0.41
1:C:220:GLU:OE2	1:C:224:LYS:HD2	2.20	0.41
4:C:1005:1PE:H131	4:C:1005:1PE:H222	1.87	0.41
1:D:228:MET:SD	1:D:235:PHE:HB2	2.60	0.41
1:B:161:VAL:HG13	1:B:201:TYR:CD1	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:48:VAL:HG21	1:C:346:PHE:HB3	2.01	0.41
1:D:266:VAL:CG1	1:D:270:GLY:HA3	2.51	0.41
1:B:160:ASN:HB3	1:B:275:PHE:HE1	1.85	0.40
1:B:288:GLY:HA3	1:D:278:ASN:O	2.21	0.40
1:B:70:ALA:HB1	1:B:72:LEU:HD21	2.03	0.40
1:B:336:TYR:OH	1:B:342:GLU:OE2	2.28	0.40
1:A:178:GLN:H	1:A:178:GLN:NE2	2.19	0.40
4:A:1004:1PE:H262	1:B:178:GLN:NE2	2.36	0.40
1:B:329:THR:HG22	5:B:1140:HOH:O	2.21	0.40
1:C:187:LYS:HA	1:C:187:LYS:HD3	1.93	0.40
1:D:159:LYS:CG	1:D:307:ALA:HB3	2.49	0.40
1:A:26:PRO:HD2	1:A:134:TYR:CE2	2.56	0.40
1:C:336:TYR:OH	1:C:342:GLU:OE2	2.32	0.40
1:D:57:ASN:ND2	1:D:62:ASN:O	2.55	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	358/361 (99%)	341 (95%)	17 (5%)	0	100	100
1	B	358/361 (99%)	340 (95%)	18 (5%)	0	100	100
1	C	358/361 (99%)	342 (96%)	16 (4%)	0	100	100
1	D	358/361 (99%)	345 (96%)	13 (4%)	0	100	100
All	All	1432/1444 (99%)	1368 (96%)	64 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	313/314 (100%)	290 (93%)	23 (7%)	14	28
1	B	313/314 (100%)	292 (93%)	21 (7%)	16	33
1	C	313/314 (100%)	299 (96%)	14 (4%)	27	52
1	D	313/314 (100%)	298 (95%)	15 (5%)	25	49
All	All	1252/1256 (100%)	1179 (94%)	73 (6%)	20	40

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	LYS
1	A	25	GLU
1	A	28	LYS
1	A	59	GLU
1	A	77	VAL
1	A	84	LYS
1	A	92	LEU
1	A	103	ARG
1	A	115	PRO
1	A	117	LYS
1	A	143	VAL
1	A	159	LYS
1	A	178	GLN
1	A	194	SER
1	A	216	ASN
1	A	217	ARG
1	A	223	ASN
1	A	277	THR
1	A	306	ASP
1	A	329	THR
1	A	344	ASN
1	A	349	LYS
1	A	351	LYS
1	B	20	LYS

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Mol	Chain	Res	Type
1	B	23	MET
1	B	34	LYS
1	B	77	VAL
1	B	92	LEU
1	B	103	ARG
1	B	117	LYS
1	B	159	LYS
1	B	163	LYS
1	B	173	LYS
1	B	180	ASP
1	B	196	SER
1	B	222	GLU
1	B	226	LYS
1	B	227	MET
1	B	229	ASP
1	B	268	ASN
1	B	296	GLU
1	B	341	GLU
1	B	344	ASN
1	B	354	ILE
1	C	33	LEU
1	C	92	LEU
1	C	95	ARG
1	C	98	LYS
1	C	117	LYS
1	C	140	LEU
1	C	159	LYS
1	C	163	LYS
1	C	194	SER
1	C	248	LEU
1	C	261	LYS
1	C	264	LYS
1	C	306	ASP
1	C	349	LYS
1	D	2	SER
1	D	7	ILE
1	D	16	VAL
1	D	17	LYS
1	D	95	ARG
1	D	98	LYS
1	D	117	LYS
1	D	140	LEU

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Mol	Chain	Res	Type
1	D	148	LEU
1	D	156	GLU
1	D	158	LEU
1	D	159	LYS
1	D	194	SER
1	D	197	GLU
1	D	219	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	118	HIS
1	A	178	GLN
1	A	218	HIS
1	A	223	ASN
1	A	269	ASN
1	A	328	GLN
1	B	144	ASN
1	B	178	GLN
1	B	223	ASN
1	B	268	ASN
1	B	269	ASN
1	B	311	HIS
1	B	321	ASN
1	B	328	GLN
1	B	348	GLN
1	C	267	ASN
1	C	269	ASN
1	C	321	ASN
1	D	138	GLN
1	D	139	ASN
1	D	321	ASN
1	D	324	HIS
1	D	348	GLN

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 22 ligands modelled in this entry, 8 are monoatomic - leaving 14 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	PGO	B	1004	-	4,4,4	0.54	0	4,4,4	1.33	0
4	1PE	C	1004	-	15,15,15	0.69	0	14,14,14	0.55	0
3	PGO	D	1004	-	4,4,4	0.72	0	4,4,4	0.79	0
4	1PE	A	1005	-	15,15,15	0.54	0	14,14,14	0.46	0
4	1PE	A	1004	-	15,15,15	0.58	0	14,14,14	0.65	0
4	1PE	B	1005	-	15,15,15	0.62	0	14,14,14	0.51	0
4	1PE	B	1006	-	15,15,15	0.62	0	14,14,14	0.58	0
3	PGO	A	1003	-	4,4,4	0.31	0	4,4,4	1.20	0
4	1PE	B	1007	-	15,15,15	0.71	0	14,14,14	0.49	0
3	PGO	B	1003	-	4,4,4	0.58	0	4,4,4	0.65	0
3	PGO	D	1003	-	4,4,4	0.66	0	4,4,4	0.70	0
3	PGO	C	1003	-	4,4,4	0.50	0	4,4,4	0.67	0
4	1PE	D	1005	-	15,15,15	0.61	0	14,14,14	0.46	0
4	1PE	C	1005	-	15,15,15	0.47	0	14,14,14	0.65	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	PGO	B	1004	-	-	0/2/2/2	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	1PE	C	1004	-	-	8/13/13/13	-
3	PGO	D	1004	-	-	0/2/2/2	-
4	1PE	A	1005	-	-	9/13/13/13	-
4	1PE	A	1004	-	-	4/13/13/13	-
4	1PE	B	1005	-	-	4/13/13/13	-
4	1PE	B	1006	-	-	8/13/13/13	-
3	PGO	A	1003	-	-	0/2/2/2	-
4	1PE	B	1007	-	-	6/13/13/13	-
3	PGO	B	1003	-	-	0/2/2/2	-
3	PGO	D	1003	-	-	0/2/2/2	-
3	PGO	C	1003	-	-	0/2/2/2	-
4	1PE	D	1005	-	-	9/13/13/13	-
4	1PE	C	1005	-	-	7/13/13/13	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (55) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	1006	1PE	OH6-C15-C25-OH5
4	C	1004	1PE	OH6-C15-C25-OH5
4	A	1005	1PE	OH4-C13-C23-OH3
4	B	1007	1PE	OH6-C15-C25-OH5
4	C	1005	1PE	OH4-C13-C23-OH3
4	B	1005	1PE	OH2-C12-C22-OH3
4	B	1007	1PE	OH7-C16-C26-OH6
4	D	1005	1PE	OH2-C12-C22-OH3
4	D	1005	1PE	OH4-C13-C23-OH3
4	B	1006	1PE	OH4-C13-C23-OH3
4	A	1005	1PE	OH2-C12-C22-OH3
4	D	1005	1PE	OH7-C16-C26-OH6
4	B	1006	1PE	OH5-C14-C24-OH4
4	C	1005	1PE	OH7-C16-C26-OH6
4	D	1005	1PE	OH6-C15-C25-OH5
4	C	1005	1PE	OH6-C15-C25-OH5
4	C	1004	1PE	OH5-C14-C24-OH4
4	A	1005	1PE	OH5-C14-C24-OH4
4	A	1004	1PE	OH6-C15-C25-OH5

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Mol	Chain	Res	Type	Atoms
4	B	1007	1PE	OH5-C14-C24-OH4
4	B	1007	1PE	OH4-C13-C23-OH3
4	B	1005	1PE	OH6-C15-C25-OH5
4	C	1005	1PE	C15-C25-OH5-C14
4	A	1005	1PE	C14-C24-OH4-C13
4	B	1006	1PE	C16-C26-OH6-C15
4	B	1006	1PE	C12-C22-OH3-C23
4	C	1004	1PE	C24-C14-OH5-C25
4	A	1005	1PE	C16-C26-OH6-C15
4	A	1005	1PE	C23-C13-OH4-C24
4	D	1005	1PE	C15-C25-OH5-C14
4	D	1005	1PE	C24-C14-OH5-C25
4	A	1004	1PE	OH4-C13-C23-OH3
4	C	1004	1PE	C16-C26-OH6-C15
4	C	1005	1PE	C14-C24-OH4-C13
4	B	1006	1PE	C25-C15-OH6-C26
4	A	1005	1PE	C24-C14-OH5-C25
4	D	1005	1PE	C16-C26-OH6-C15
4	C	1005	1PE	C13-C23-OH3-C22
4	D	1005	1PE	C12-C22-OH3-C23
4	C	1004	1PE	C23-C13-OH4-C24
4	A	1005	1PE	C15-C25-OH5-C14
4	A	1004	1PE	C14-C24-OH4-C13
4	A	1005	1PE	C12-C22-OH3-C23
4	C	1004	1PE	C25-C15-OH6-C26
4	B	1005	1PE	C24-C14-OH5-C25
4	D	1005	1PE	C25-C15-OH6-C26
4	B	1007	1PE	C15-C25-OH5-C14
4	B	1005	1PE	OH5-C14-C24-OH4
4	B	1007	1PE	OH2-C12-C22-OH3
4	C	1004	1PE	OH4-C13-C23-OH3
4	B	1006	1PE	C13-C23-OH3-C22
4	A	1004	1PE	C12-C22-OH3-C23
4	C	1004	1PE	C15-C25-OH5-C14
4	C	1005	1PE	OH5-C14-C24-OH4
4	B	1006	1PE	OH2-C12-C22-OH3

There are no ring outliers.

13 monomers are involved in 32 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1004	PGO	5	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	1004	1PE	2	0
3	D	1004	PGO	1	0
4	A	1005	1PE	1	0
4	A	1004	1PE	5	0
4	B	1005	1PE	5	0
4	B	1006	1PE	2	0
3	A	1003	PGO	1	0
4	B	1007	1PE	1	0
3	D	1003	PGO	1	0
3	C	1003	PGO	1	0
4	D	1005	1PE	2	0
4	C	1005	1PE	5	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	360/361 (99%)	0.04	7 (1%)	66	62	18, 32, 56, 86	0
1	B	360/361 (99%)	0.34	23 (6%)	19	14	17, 37, 62, 108	0
1	C	360/361 (99%)	0.13	14 (3%)	39	32	18, 31, 53, 85	0
1	D	360/361 (99%)	0.28	17 (4%)	31	25	16, 31, 54, 80	0
All	All	1440/1444 (99%)	0.20	61 (4%)	36	29	16, 33, 58, 108	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	352	GLY	6.0
1	B	53	SER	5.9
1	B	26	PRO	5.7
1	D	62	ASN	5.6
1	B	16	VAL	5.0
1	D	28	LYS	4.9
1	D	347	PHE	4.4
1	D	25	GLU	4.3
1	B	28	LYS	4.2
1	D	63	PHE	4.2
1	B	63	PHE	3.7
1	D	7	ILE	3.7
1	B	29	TYR	3.5
1	B	58	PRO	3.5
1	B	54	PHE	3.3
1	D	8	VAL	3.2
1	B	23	MET	3.2
1	A	351	LYS	3.2
1	B	10	ASP	3.1
1	B	9	THR	3.1
1	D	29	TYR	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	26	PRO	3.0
1	B	8	VAL	2.9
1	C	51	ALA	2.9
1	B	15	GLY	2.9
1	C	351	LYS	2.9
1	A	232	GLY	2.8
1	B	79	ASP	2.8
1	B	66	LEU	2.8
1	D	2	SER	2.7
1	C	60	GLY	2.7
1	B	77	VAL	2.5
1	D	27	GLU	2.5
1	B	62	ASN	2.5
1	B	24	PRO	2.5
1	C	46	GLY	2.4
1	D	50	GLY	2.4
1	C	336	TYR	2.4
1	C	27	GLU	2.4
1	D	22	ASP	2.3
1	C	50	GLY	2.3
1	D	24	PRO	2.3
1	C	352	GLY	2.2
1	B	21	ILE	2.2
1	C	15	GLY	2.2
1	C	346	PHE	2.2
1	B	83	ILE	2.2
1	A	59	GLU	2.1
1	D	81	ASN	2.1
1	B	3	THR	2.1
1	A	57	ASN	2.1
1	B	2	SER	2.1
1	B	22	ASP	2.1
1	A	222	GLU	2.1
1	C	286	VAL	2.1
1	C	285	PRO	2.1
1	C	241	ASP	2.1
1	D	15	GLY	2.0
1	C	85	ARG	2.0
1	D	53	SER	2.0
1	A	15	GLY	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
4	1PE	A	1004	16/16	0.65	0.42	56,66,73,74	0
4	1PE	B	1005	16/16	0.65	0.43	66,73,86,87	0
4	1PE	C	1004	16/16	0.72	0.40	42,66,72,72	0
4	1PE	B	1007	16/16	0.74	0.42	49,64,69,69	0
4	1PE	D	1005	16/16	0.78	0.32	43,55,59,63	0
4	1PE	C	1005	16/16	0.80	0.30	50,58,62,64	0
4	1PE	B	1006	16/16	0.83	0.24	44,49,73,74	0
3	PGO	A	1003	5/5	0.85	0.31	42,43,48,48	0
4	1PE	A	1005	16/16	0.85	0.26	49,57,62,63	0
3	PGO	B	1004	5/5	0.86	0.25	39,40,44,46	0
3	PGO	C	1003	5/5	0.93	0.32	45,45,48,49	0
3	PGO	B	1003	5/5	0.95	0.24	45,46,46,49	0
3	PGO	D	1004	5/5	0.95	0.18	34,36,38,39	0
3	PGO	D	1003	5/5	0.96	0.18	40,43,48,53	0
2	ZN	B	1001	1/1	0.96	0.08	34,34,34,34	0
2	ZN	A	1002	1/1	0.99	0.07	25,25,25,25	0
2	ZN	A	1001	1/1	0.99	0.09	25,25,25,25	0
2	ZN	B	1002	1/1	0.99	0.07	31,31,31,31	0
2	ZN	C	1001	1/1	0.99	0.14	26,26,26,26	0
2	ZN	C	1002	1/1	0.99	0.05	33,33,33,33	0
2	ZN	D	1001	1/1	0.99	0.14	31,31,31,31	0
2	ZN	D	1002	1/1	0.99	0.08	32,32,32,32	0

6.5 Other polymers (i)

There are no such residues in this entry.