Full wwPDB X-ray Structure Validation Report (i)

Jun 16, 2024 – 05:47 PM EDT

PDB ID	:	3KDR
Title	:	The Crystal Structure of a HK97 Family Phage Portal Protein from Corynebacterium diphtheriae to 2.9A
Authors	:	Nocek, B.; Stein, A.J.; Mulligan, R.; Duggan, E.; Abdullah, J.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on	:	2009-10-23
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 (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:

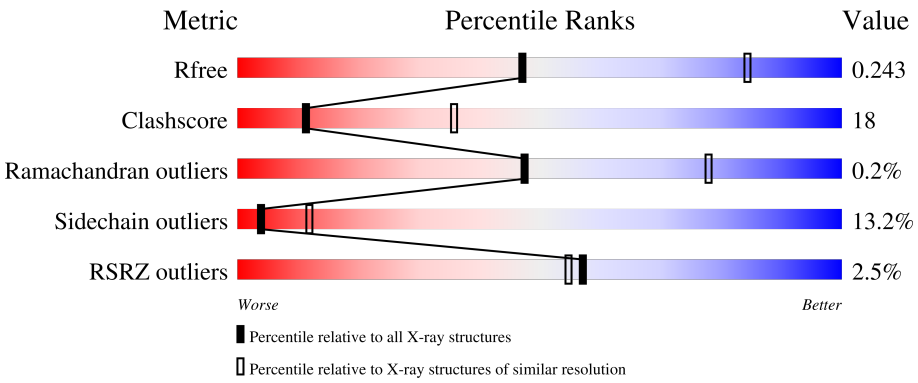
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, 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 2.90 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 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	310	2% 64% 22% 5% 8%
1	B	310	3% 57% 29% 5% 9%
1	C	310	2% 65% 23% • 7%

2 Entry composition (i)

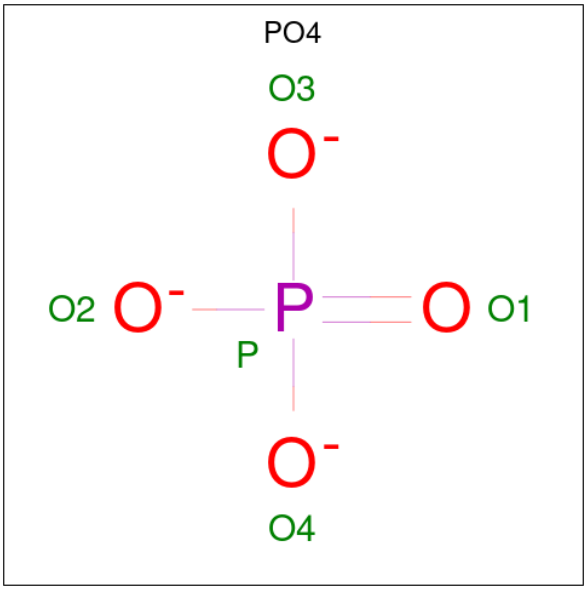
There are 5 unique types of molecules in this entry. The entry contains 6311 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 HK97 Family Phage Portal Protein.

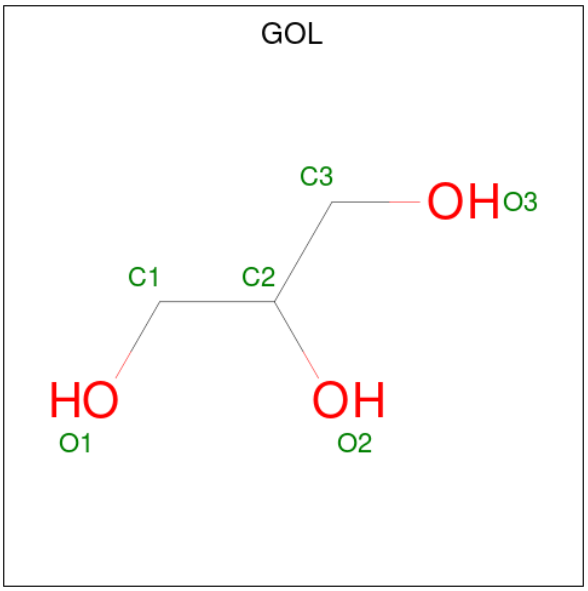
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	285	2087	1305	373	397	4	8	0	0	0
1	B	283	2085	1303	375	395	4	8	0	0	0
1	C	287	2106	1317	377	400	4	8	0	0	0

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



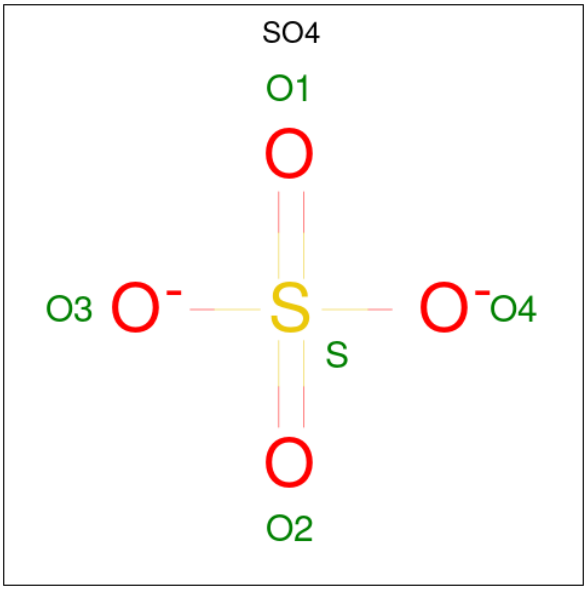
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	P		
2	A	1	5	4	1	0	0
2	C	1	5	4	1	0	0

- 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	B	1	6	3	3	0	0

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
4	B	1	5	4	1	0	0
			Total	O	S		
4	C	1	5	4	1	0	0

- Molecule 5 is water.

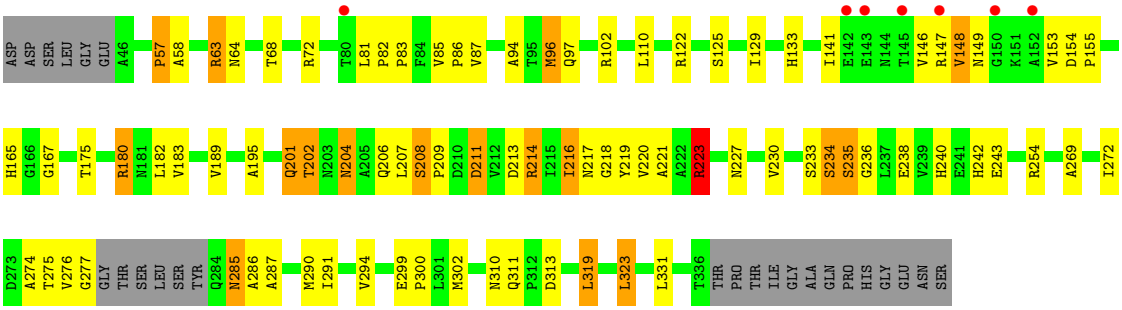
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	1	1	1	0	0

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.

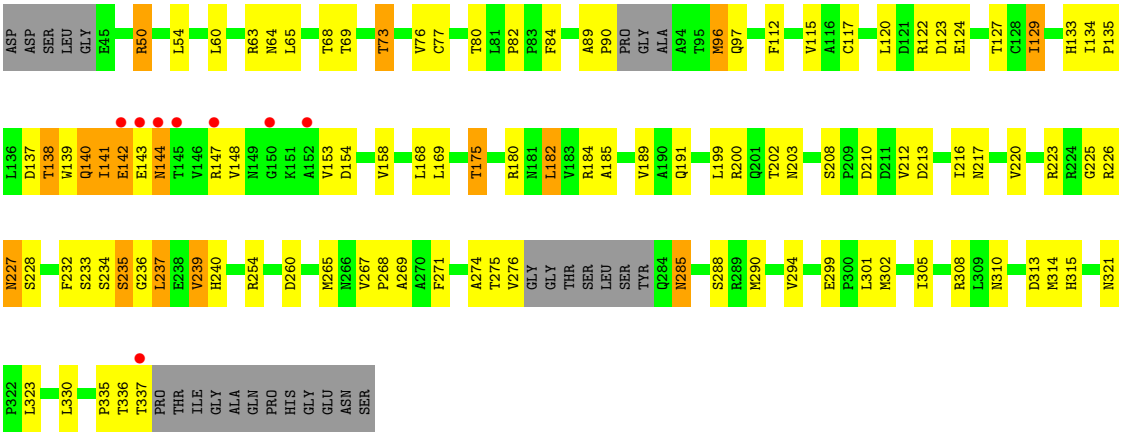
• Molecule 1: HK97 Family Phage Portal Protein

Chain A:



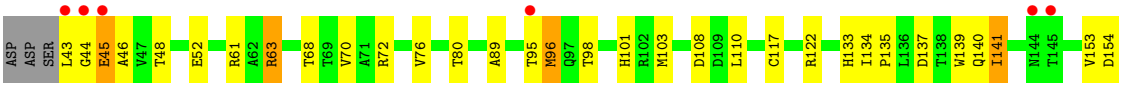
• Molecule 1: HK97 Family Phage Portal Protein

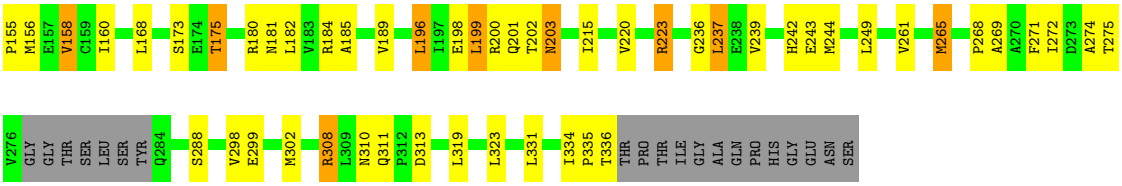
Chain B:



• Molecule 1: HK97 Family Phage Portal Protein

Chain C:




P155
M156
E157
V158
C159
I160
L168
S173
E174
T175
R180
N181
L182
V183
R184
A185
V189
L196
I197
E198
L199
R200
Q201
T202
N203
I215
V220
R223
G236
L237
E238
V239
H242
E243
M244
L249
V261
M265
P268
A269
A270
F271
I272
D273
A274
T275
V276
GLY
GLY
THR
SER
LEU
SER
TYR
Q284
S288
V298
E299
M302
R308
L309
N310
Q311
P312
D313
L319
L323
L331
I334
P335
T336
THR
PRO
THR
ILE
GLY
ALA
GLN
PRO
HIS
GLY
GLU
ASN
SER

4 Data and refinement statistics (i)

Property	Value	Source
Space group	F 4 3 2	Depositor
Cell constants a, b, c, α, β, γ	340.07Å 340.07Å 340.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.25 – 2.90 36.25 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.5 (36.25-2.90) 98.5 (36.25-2.90)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.79 (at 2.90Å)	Xtriage
Refinement program	REFMAC refmac_5.5.0102	Depositor
R, R_{free}	0.198 , 0.245 0.198 , 0.243	Depositor DCC
R_{free} test set	1854 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å^2)	48.9	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 28.6	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	0.93	EDS
Total number of atoms	6311	wwPDB-VP
Average B, all atoms (Å^2)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.29% 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: SO4, PO4, 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	0.97	0/2117	0.91	2/2886 (0.1%)
1	B	1.03	1/2113 (0.0%)	0.91	1/2878 (0.0%)
1	C	1.05	0/2136	0.91	1/2911 (0.0%)
All	All	1.02	1/6366 (0.0%)	0.91	4/8675 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	77	CYS	CB-SG	-8.47	1.67	1.82

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	223	ARG	NE-CZ-NH1	5.55	123.08	120.30
1	B	154	ASP	CB-CG-OD2	5.51	123.26	118.30
1	A	211	ASP	CB-CG-OD1	-5.28	113.55	118.30
1	C	141	ILE	CB-CA-C	-5.20	101.19	111.60

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	2087	0	2049	91	0
1	B	2085	0	2052	88	0
1	C	2106	0	2071	72	0
2	A	5	0	0	1	0
2	C	5	0	0	1	0
3	A	6	0	8	2	0
3	B	6	0	8	3	0
4	B	5	0	0	0	0
4	C	5	0	0	1	0
5	A	1	0	0	0	0
All	All	6311	0	6188	223	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 18.

All (223) 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:202:THR:HG22	1:C:236:GLY:O	1.36	1.21
1:A:201:GLN:HG2	1:B:234:SER:O	1.42	1.14
1:B:97:GLN:HE22	1:B:133:HIS:N	1.51	1.07
1:B:97:GLN:NE2	1:B:133:HIS:H	1.56	1.03
1:C:43:LEU:HD12	1:C:52:GLU:HG2	1.49	0.94
1:A:97:GLN:HE22	1:A:133:HIS:H	0.95	0.92
1:B:310:ASN:HD21	1:B:323:LEU:H	1.17	0.90
1:B:285:ASN:HD22	1:B:285:ASN:H	1.21	0.87
1:A:201:GLN:CG	1:B:234:SER:O	2.22	0.86
1:B:69:THR:O	1:B:73:THR:HG23	1.75	0.84
1:A:227:ASN:O	1:A:230:VAL:HG23	1.77	0.83
1:C:43:LEU:HD12	1:C:52:GLU:CG	2.08	0.82
1:C:98:THR:OG1	1:C:101:HIS:HD2	1.62	0.82
1:B:185:ALA:O	1:B:189:VAL:HG13	1.81	0.81
1:C:215:ILE:HD12	1:C:237:LEU:HD22	1.62	0.80
1:C:199:LEU:HD23	1:C:237:LEU:HD21	1.63	0.79
1:A:97:GLN:NE2	1:A:133:HIS:H	1.77	0.79
1:A:310:ASN:HD21	1:A:323:LEU:H	1.32	0.77
1:B:64:ASN:O	1:B:68:THR:HB	1.86	0.76
1:A:97:GLN:HE22	1:A:133:HIS:N	1.80	0.76
1:A:235:SER:OG	1:C:201:GLN:HB3	1.85	0.76
1:C:269:ALA:HB1	1:C:274:ALA:HB3	1.68	0.75
1:B:285:ASN:HD22	1:B:285:ASN:N	1.83	0.75
1:C:223:ARG:HH11	1:C:223:ARG:HG2	1.50	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:ARG:HD2	1:C:243:GLU:OE2	1.87	0.74
1:B:315:HIS:HE1	1:B:321:ASN:O	1.69	0.74
1:C:95:THR:HG23	1:C:95:THR:O	1.88	0.73
1:A:165:HIS:HD2	1:A:167:GLY:H	1.36	0.73
1:B:299:GLU:OE2	1:B:302:MSE:HE3	1.89	0.72
1:C:43:LEU:HB3	1:C:44:GLY:HA3	1.72	0.72
1:B:299:GLU:CD	1:B:302:MSE:CE	2.58	0.72
1:A:180:ARG:O	1:A:183:VAL:HG12	1.90	0.72
1:B:299:GLU:OE2	1:B:302:MSE:CE	2.38	0.71
1:A:299:GLU:CD	1:A:302:MSE:CE	2.59	0.70
1:B:226:ARG:C	1:B:227:ASN:HD22	1.95	0.70
1:C:98:THR:OG1	1:C:101:HIS:CD2	2.45	0.69
1:A:201:GLN:NE2	1:B:234:SER:O	2.26	0.68
1:A:216:ILE:O	1:A:220:VAL:HG23	1.93	0.68
1:A:182:LEU:CD2	1:A:254:ARG:HA	2.22	0.68
1:A:240:HIS:HD2	3:A:350:GOL:H2	1.59	0.68
1:B:301:LEU:O	1:B:305:ILE:HG13	1.95	0.67
1:B:97:GLN:HE22	1:B:133:HIS:H	0.75	0.67
1:B:299:GLU:CD	1:B:302:MSE:HE2	2.14	0.67
1:C:43:LEU:O	1:C:184:ARG:NH2	2.28	0.67
1:A:202:THR:OG1	1:A:236:GLY:C	2.34	0.66
1:B:232:PHE:CE2	1:B:234:SER:HA	2.31	0.66
1:B:268:PRO:HG2	1:B:271:PHE:CD2	2.32	0.65
1:C:45:GLU:HG2	1:C:46:ALA:N	2.10	0.65
1:A:165:HIS:CD2	1:A:167:GLY:H	2.13	0.65
1:B:213:ASP:O	1:B:217:ASN:ND2	2.25	0.65
1:B:216:ILE:O	1:B:220:VAL:HG23	1.97	0.65
1:A:227:ASN:O	1:A:230:VAL:CG2	2.45	0.65
1:B:148:VAL:HG23	1:B:153:VAL:HG21	1.78	0.65
1:A:243:GLU:OE1	1:B:223:ARG:NH1	2.30	0.64
1:C:45:GLU:O	1:C:180:ARG:NH2	2.30	0.64
1:B:123:ASP:O	1:B:124:GLU:C	2.36	0.64
1:A:287:ALA:O	1:A:291:ILE:HG13	1.97	0.63
1:A:63:ARG:HH11	1:A:64:ASN:ND2	1.96	0.63
1:A:57:PRO:HG3	1:C:175:THR:HG21	1.80	0.62
1:A:299:GLU:HA	1:A:302:MSE:HE2	1.79	0.62
1:B:199:LEU:HB3	1:B:237:LEU:HD21	1.82	0.61
1:A:208:SER:H	1:A:211:ASP:HB2	1.64	0.61
1:C:261:VAL:O	1:C:265:MSE:HG3	2.01	0.60
1:A:201:GLN:HG2	1:B:234:SER:C	2.21	0.60
1:B:50:ARG:HD2	1:B:112:PHE:CD2	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:ASP:OD1	1:A:155:PRO:HD2	2.01	0.60
1:A:122:ARG:NH2	1:A:313:ASP:O	2.35	0.60
1:B:69:THR:O	1:B:73:THR:CG2	2.50	0.59
1:A:209:PRO:O	1:A:213:ASP:HB2	2.01	0.59
1:B:299:GLU:HA	1:B:302:MSE:HE2	1.85	0.58
1:B:285:ASN:H	1:B:285:ASN:ND2	1.99	0.58
1:A:290:MSE:O	1:A:294:VAL:HG23	2.04	0.58
1:B:310:ASN:HD21	1:B:323:LEU:N	1.94	0.58
1:B:168:LEU:CD1	1:B:265:MSE:HE3	2.34	0.58
1:A:63:ARG:HH11	1:A:64:ASN:HD22	1.51	0.58
1:C:201:GLN:HG2	1:C:237:LEU:HD12	1.85	0.58
1:C:154:ASP:OD2	1:C:155:PRO:HD2	2.05	0.56
1:A:214:ARG:O	1:A:217:ASN:HB2	2.05	0.56
1:C:133:HIS:ND1	4:C:3:SO4:O1	2.34	0.56
1:A:299:GLU:CD	1:A:302:MSE:HE2	2.25	0.56
2:A:1:PO4:O1	2:C:2:PO4:O4	2.23	0.56
1:C:63:ARG:NH2	1:C:108:ASP:OD2	2.39	0.56
1:A:182:LEU:HD21	1:A:254:ARG:N	2.21	0.55
1:B:240:HIS:HD2	3:B:2:GOL:H32	1.71	0.55
1:B:240:HIS:HD2	3:B:2:GOL:C3	2.19	0.55
1:A:240:HIS:CD2	3:A:350:GOL:H2	2.42	0.55
1:C:223:ARG:HG2	1:C:223:ARG:NH1	2.17	0.55
1:C:202:THR:CG2	1:C:236:GLY:O	2.31	0.55
1:A:234:SER:O	1:A:236:GLY:N	2.25	0.55
1:A:195:ALA:HB2	1:B:191:GLN:O	2.08	0.54
1:B:73:THR:HG22	1:B:330:LEU:HD11	1.89	0.54
1:C:199:LEU:CD2	1:C:237:LEU:HD21	2.34	0.54
1:A:189:VAL:HG12	1:C:249:LEU:CD1	2.37	0.54
1:A:311:GLN:NE2	1:B:96:MSE:HE2	2.21	0.54
1:B:202:THR:OG1	1:B:236:GLY:O	2.23	0.54
1:A:97:GLN:HE21	1:A:102:ARG:HB2	1.73	0.54
1:B:180:ARG:HD3	1:B:184:ARG:CZ	2.38	0.54
1:A:57:PRO:CG	1:C:175:THR:HG21	2.38	0.53
1:C:168:LEU:HD11	1:C:265:MSE:HE3	1.90	0.53
1:A:235:SER:OG	1:C:201:GLN:CB	2.55	0.53
1:B:134:ILE:HD12	1:B:138:THR:HG22	1.89	0.53
1:B:285:ASN:N	1:B:285:ASN:ND2	2.55	0.53
1:C:117:CYS:SG	1:C:158:VAL:HG22	2.48	0.53
1:A:331:LEU:CD1	1:B:335:PRO:HB2	2.39	0.53
1:A:148:VAL:O	1:A:148:VAL:CG2	2.56	0.52
1:A:96:MSE:HE1	1:C:311:GLN:HG3	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:313:ASP:OD2	1:B:313:ASP:N	2.42	0.52
1:A:83:PRO:HB2	1:A:129:ILE:HA	1.92	0.52
1:A:319:LEU:HD12	1:A:319:LEU:H	1.75	0.52
1:A:234:SER:C	1:A:236:GLY:H	2.08	0.52
1:A:182:LEU:HD23	1:A:254:ARG:HA	1.90	0.52
1:B:315:HIS:CE1	1:B:321:ASN:O	2.59	0.52
1:B:189:VAL:HG23	1:B:189:VAL:O	2.09	0.51
1:C:299:GLU:CD	1:C:302:MSE:CE	2.79	0.51
1:B:63:ARG:HD3	1:B:63:ARG:C	2.30	0.51
1:B:144:ASN:C	1:B:144:ASN:OD1	2.50	0.51
1:C:117:CYS:HB2	1:C:139:TRP:CE2	2.46	0.50
1:A:242:HIS:NE2	1:B:239:VAL:HG22	2.26	0.50
1:B:82:PRO:HG3	1:B:315:HIS:CD2	2.47	0.50
1:A:202:THR:N	1:A:236:GLY:O	2.31	0.50
1:B:137:ASP:C	1:B:137:ASP:OD1	2.50	0.50
1:B:227:ASN:N	1:B:227:ASN:ND2	2.59	0.50
1:A:202:THR:OG1	1:A:236:GLY:HA2	2.11	0.50
1:B:235:SER:O	1:B:235:SER:OG	2.30	0.50
1:A:242:HIS:CE1	1:B:239:VAL:HG22	2.47	0.50
1:C:185:ALA:O	1:C:189:VAL:HG23	2.12	0.50
1:B:240:HIS:CD2	3:B:2:GOL:H32	2.47	0.49
1:B:73:THR:CG2	1:B:330:LEU:HD11	2.42	0.49
1:B:168:LEU:HD11	1:B:265:MSE:HE3	1.93	0.49
1:A:146:VAL:O	1:A:153:VAL:HG12	2.11	0.49
1:C:308:ARG:NH2	1:C:313:ASP:OD2	2.38	0.49
1:B:141:ILE:HG22	1:B:142:GLU:O	2.12	0.49
1:C:45:GLU:HG2	1:C:46:ALA:H	1.76	0.49
1:A:276:VAL:O	1:A:277:GLY:C	2.52	0.49
1:A:85:VAL:HB	1:A:86:PRO:HD3	1.95	0.49
1:A:218:GLY:O	1:A:221:ALA:HB3	2.13	0.49
1:B:175:THR:HG21	1:B:260:ASP:HB3	1.95	0.48
1:B:233:SER:HB3	1:B:237:LEU:HB3	1.94	0.48
1:C:45:GLU:CG	1:C:46:ALA:H	2.26	0.48
1:C:48:THR:N	1:C:52:GLU:OE1	2.38	0.48
1:C:334:ILE:HG22	1:C:335:PRO:O	2.13	0.48
1:C:299:GLU:HA	1:C:302:MSE:HE2	1.96	0.48
1:A:68:THR:O	1:A:72:ARG:HG3	2.14	0.47
1:A:202:THR:OG1	1:A:236:GLY:CA	2.62	0.47
1:B:203:ASN:OD1	1:B:203:ASN:C	2.51	0.47
1:A:299:GLU:N	1:A:300:PRO:HD2	2.29	0.47
1:B:310:ASN:ND2	1:B:323:LEU:H	2.00	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:GLN:HA	1:A:236:GLY:O	2.15	0.47
1:C:95:THR:O	1:C:95:THR:CG2	2.57	0.47
1:C:43:LEU:CD1	1:C:52:GLU:OE2	2.63	0.47
1:A:233:SER:HB2	1:C:200:ARG:HA	1.97	0.47
1:B:308:ARG:HD2	1:B:308:ARG:HA	1.73	0.47
1:C:45:GLU:CG	1:C:46:ALA:N	2.77	0.47
1:C:202:THR:HG22	1:C:236:GLY:C	2.25	0.47
1:A:195:ALA:HB3	1:A:243:GLU:HB2	1.96	0.46
1:A:219:TYR:OH	1:C:196:LEU:HD22	2.15	0.46
1:A:269:ALA:HB1	1:A:274:ALA:HB3	1.97	0.46
1:A:204:ASN:OD1	1:A:204:ASN:N	2.47	0.46
1:B:267:VAL:HG23	1:B:268:PRO:HD2	1.98	0.46
1:B:182:LEU:HD22	1:B:254:ARG:HA	1.97	0.46
1:A:96:MSE:HE3	1:C:311:GLN:CD	2.36	0.46
1:A:299:GLU:OE1	1:A:302:MSE:CE	2.62	0.46
1:C:68:THR:O	1:C:72:ARG:HG3	2.15	0.46
1:C:298:VAL:HG12	1:C:302:MSE:HG3	1.98	0.46
1:A:238:GLU:OE1	1:A:240:HIS:HE1	1.99	0.46
1:A:82:PRO:HA	1:A:83:PRO:HD3	1.86	0.45
1:A:201:GLN:CD	1:B:234:SER:O	2.55	0.45
1:B:139:TRP:O	1:B:140:GLN:HG2	2.16	0.45
1:C:122:ARG:NH1	1:C:156:MSE:HG3	2.30	0.45
1:B:84:PHE:CE2	1:B:314:MSE:HE3	2.51	0.45
1:C:201:GLN:HE22	1:C:203:ASN:HB2	1.82	0.45
1:B:135:PRO:O	1:B:138:THR:HB	2.16	0.45
1:B:123:ASP:CB	1:B:129:ILE:HD11	2.46	0.45
1:B:139:TRP:C	1:B:140:GLN:HG2	2.37	0.45
1:A:148:VAL:HG23	1:A:149:ASN:ND2	2.32	0.45
1:A:195:ALA:HB3	1:A:243:GLU:CB	2.47	0.45
1:A:219:TYR:CE2	1:C:196:LEU:HD22	2.51	0.45
1:A:299:GLU:CD	1:A:302:MSE:HE1	2.37	0.45
1:C:70:VAL:HG12	1:C:103:MSE:HG3	1.99	0.45
1:A:182:LEU:HD21	1:A:254:ARG:CA	2.47	0.44
1:B:223:ARG:C	1:B:225:GLY:H	2.21	0.44
1:C:61:ARG:HG3	1:C:272:ILE:HG22	1.99	0.44
1:A:96:MSE:HE3	1:C:311:GLN:OE1	2.16	0.44
1:A:311:GLN:HE21	1:B:96:MSE:HE2	1.82	0.44
1:A:331:LEU:HD11	1:B:335:PRO:HB2	1.99	0.44
1:A:141:ILE:O	1:A:141:ILE:HG23	2.17	0.44
1:B:54:LEU:HD22	1:B:60:LEU:HD13	1.99	0.44
1:A:96:MSE:CE	1:C:311:GLN:CD	2.86	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:TYR:CZ	1:C:196:LEU:HD22	2.53	0.44
1:B:227:ASN:HD22	1:B:227:ASN:N	2.14	0.44
1:C:96:MSE:HE2	1:C:96:MSE:HB3	1.85	0.43
1:C:198:GLU:OE1	1:C:242:HIS:HE1	2.02	0.43
1:A:219:TYR:CE2	1:C:196:LEU:CD2	3.02	0.43
1:C:265:MSE:HE2	1:C:265:MSE:HB3	1.85	0.43
1:B:97:GLN:NE2	1:B:133:HIS:N	2.35	0.43
1:B:294:VAL:O	1:B:299:GLU:HB2	2.18	0.43
1:B:50:ARG:HD2	1:B:112:PHE:CG	2.54	0.42
1:A:110:LEU:HD23	1:A:110:LEU:HA	1.93	0.42
1:C:140:GLN:HG3	1:C:141:ILE:N	2.34	0.42
1:B:117:CYS:HB2	1:B:139:TRP:CE2	2.54	0.42
1:B:267:VAL:HG22	1:B:268:PRO:N	2.35	0.42
1:C:122:ARG:NH1	1:C:156:MSE:O	2.47	0.42
1:C:268:PRO:O	1:C:271:PHE:HB2	2.19	0.42
1:A:122:ARG:HH22	1:A:313:ASP:HB2	1.85	0.42
1:A:58:ALA:HB1	1:A:272:ILE:HD12	2.02	0.42
1:C:160:ILE:O	1:C:308:ARG:HD3	2.19	0.42
1:C:201:GLN:NE2	1:C:203:ASN:HB2	2.35	0.42
1:C:299:GLU:CD	1:C:302:MSE:HE2	2.41	0.42
1:B:169:LEU:HD23	1:B:169:LEU:HA	1.86	0.41
1:B:290:MSE:O	1:B:294:VAL:HG23	2.19	0.41
1:C:310:ASN:HD21	1:C:323:LEU:H	1.67	0.41
1:B:134:ILE:O	1:B:134:ILE:HG13	2.19	0.41
1:C:89:ALA:CB	1:C:95:THR:O	2.68	0.41
1:A:285:ASN:O	1:A:286:ALA:C	2.60	0.41
1:C:181:ASN:HD22	1:C:181:ASN:HA	1.67	0.41
1:A:238:GLU:OE1	1:A:240:HIS:CE1	2.74	0.40
1:B:142:GLU:O	1:B:143:GLU:C	2.57	0.40
1:B:269:ALA:HB1	1:B:274:ALA:HB3	2.03	0.40
1:A:94:ALA:O	1:C:308:ARG:NH1	2.54	0.40
1:A:285:ASN:ND2	1:A:285:ASN:H	2.19	0.40
1:B:89:ALA:HA	1:B:90:PRO:HD3	1.85	0.40
1:C:134:ILE:HA	1:C:135:PRO:HD3	1.90	0.40
1:A:299:GLU:OE1	1:A:302:MSE:HE2	2.22	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	281/310 (91%)	264 (94%)	16 (6%)	1 (0%)	34	66
1	B	277/310 (89%)	263 (95%)	13 (5%)	1 (0%)	34	66
1	C	283/310 (91%)	274 (97%)	9 (3%)	0	100	100
All	All	841/930 (90%)	801 (95%)	38 (4%)	2 (0%)	47	78

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	235	SER
1	B	142	GLU

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	214/235 (91%)	190 (89%)	24 (11%)	6	18
1	B	215/235 (92%)	181 (84%)	34 (16%)	2	8
1	C	216/235 (92%)	189 (88%)	27 (12%)	4	14
All	All	645/705 (92%)	560 (87%)	85 (13%)	4	12

All (85) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	PRO

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Mol	Chain	Res	Type
1	A	63	ARG
1	A	81	LEU
1	A	87	VAL
1	A	96	MSE
1	A	125	SER
1	A	147	ARG
1	A	148	VAL
1	A	175	THR
1	A	180	ARG
1	A	201	GLN
1	A	202	THR
1	A	204	ASN
1	A	206	GLN
1	A	207	LEU
1	A	208	SER
1	A	214	ARG
1	A	216	ILE
1	A	223	ARG
1	A	234	SER
1	A	275	THR
1	A	285	ASN
1	A	319	LEU
1	A	323	LEU
1	B	50	ARG
1	B	65	LEU
1	B	73	THR
1	B	76	VAL
1	B	80	THR
1	B	96	MSE
1	B	115	VAL
1	B	120	LEU
1	B	122	ARG
1	B	127	THR
1	B	129	ILE
1	B	138	THR
1	B	140	GLN
1	B	141	ILE
1	B	144	ASN
1	B	147	ARG
1	B	158	VAL
1	B	175	THR
1	B	182	LEU

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Mol	Chain	Res	Type
1	B	200	ARG
1	B	208	SER
1	B	210	ASP
1	B	212	VAL
1	B	227	ASN
1	B	228	SER
1	B	235	SER
1	B	237	LEU
1	B	239	VAL
1	B	275	THR
1	B	276	VAL
1	B	285	ASN
1	B	288	SER
1	B	336	THR
1	B	337	THR
1	C	45	GLU
1	C	63	ARG
1	C	76	VAL
1	C	80	THR
1	C	96	MSE
1	C	110	LEU
1	C	137	ASP
1	C	153	VAL
1	C	158	VAL
1	C	173	SER
1	C	175	THR
1	C	182	LEU
1	C	196	LEU
1	C	199	LEU
1	C	203	ASN
1	C	220	VAL
1	C	223	ARG
1	C	237	LEU
1	C	239	VAL
1	C	244	MSE
1	C	265	MSE
1	C	275	THR
1	C	288	SER
1	C	308	ARG
1	C	319	LEU
1	C	331	LEU
1	C	336	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	ASN
1	A	97	GLN
1	A	149	ASN
1	A	165	HIS
1	A	171	HIS
1	A	206	GLN
1	A	240	HIS
1	A	285	ASN
1	A	310	ASN
1	A	311	GLN
1	A	321	ASN
1	B	97	GLN
1	B	149	ASN
1	B	192	ASN
1	B	201	GLN
1	B	227	ASN
1	B	240	HIS
1	B	285	ASN
1	B	310	ASN
1	B	315	HIS
1	C	64	ASN
1	C	101	HIS
1	C	181	ASN
1	C	242	HIS
1	C	310	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)

6 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	GOL	B	2	-	5,5,5	0.59	0	5,5,5	1.52	0
4	SO4	C	3	-	4,4,4	0.30	0	6,6,6	0.50	0
2	PO4	A	1	-	4,4,4	0.94	0	6,6,6	1.17	1 (16%)
2	PO4	C	2	-	4,4,4	0.84	0	6,6,6	1.04	0
4	SO4	B	4	-	4,4,4	0.34	0	6,6,6	0.38	0
3	GOL	A	350	-	5,5,5	0.57	0	5,5,5	1.12	1 (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
3	GOL	B	2	-	-	2/4/4/4	-
3	GOL	A	350	-	-	2/4/4/4	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1	PO4	O4-P-O2	2.32	115.13	107.91
3	A	350	GOL	O2-C2-C3	2.02	117.53	109.18

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	2	GOL	C1-C2-C3-O3
3	A	350	GOL	O2-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
3	A	350	GOL	C1-C2-C3-O3
3	B	2	GOL	O2-C2-C3-O3

There are no ring outliers.

5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	2	GOL	3	0
4	C	3	SO4	1	0
2	A	1	PO4	1	0
2	C	2	PO4	1	0
3	A	350	GOL	2	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	277/310 (89%)	-0.35	7 (2%)	57	55	2, 16, 27, 32	0
1	B	275/310 (88%)	-0.28	8 (2%)	51	47	2, 14, 25, 34	0
1	C	279/310 (90%)	-0.46	6 (2%)	62	59	2, 16, 24, 34	0
All	All	831/930 (89%)	-0.36	21 (2%)	57	55	2, 16, 26, 34	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	44	GLY	5.4
1	B	142	GLU	4.4
1	B	152	ALA	4.1
1	B	337	THR	3.9
1	B	145	THR	3.7
1	C	43	LEU	3.4
1	C	45	GLU	3.3
1	A	150	GLY	3.2
1	C	145	THR	3.1
1	B	144	ASN	3.1
1	B	147	ARG	3.0
1	A	145	THR	2.9
1	B	150	GLY	2.7
1	A	147	ARG	2.4
1	A	142	GLU	2.4
1	A	80	THR	2.4
1	A	143	GLU	2.4
1	C	95	THR	2.4
1	B	143	GLU	2.2
1	C	144	ASN	2.2
1	A	152	ALA	2.1

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates (i)

There are no 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
2	PO4	A	1	5/5	0.82	0.23	78,79,82,83	0
4	SO4	B	4	5/5	0.83	0.29	101,101,102,102	0
3	GOL	A	350	6/6	0.85	0.19	58,61,62,62	0
2	PO4	C	2	5/5	0.86	0.25	77,77,78,81	0
3	GOL	B	2	6/6	0.91	0.16	45,47,51,55	0
4	SO4	C	3	5/5	0.93	0.19	76,78,80,80	0

6.5 Other polymers (i)

There are no such residues in this entry.