Full wwPDB X-ray Structure Validation Report

Jun 16, 2024 – 10:00 AM EDT

PDB ID	:	2PT5
Title	:	Crystal Structure Of Shikimate Kinase (aq_2177) From Aquifex Aeolicus vf5
Authors	:	Jeyakanthan, J.; Nithya, N.; Shimada, A.; Velmurugan, D.; Ebihara, A.; Shinkai, A.; Kuramitsu, S.; Shiro, Y.; Yokoyama, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on	:	2007-05-08
Resolution	:	2.10 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 (1)) were used in the production of this report:

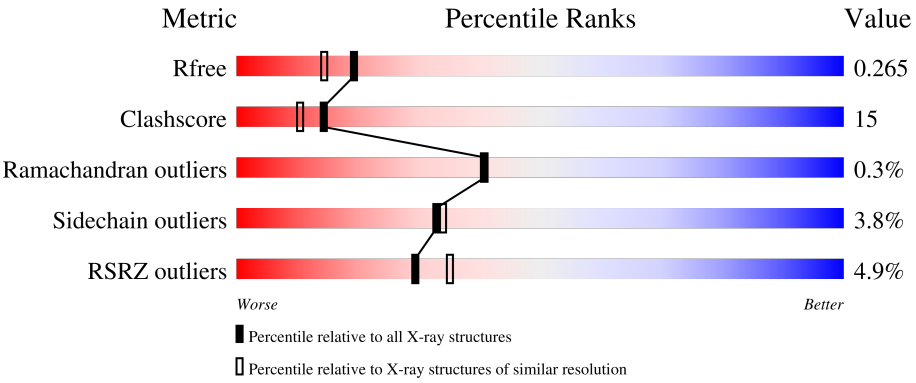
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, 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 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	168	2% 73% 22%
1	B	168	8% 67% 30%
1	C	168	4% 70% 22% 7%
1	D	168	5% 60% 33% 6%

2 Entry composition

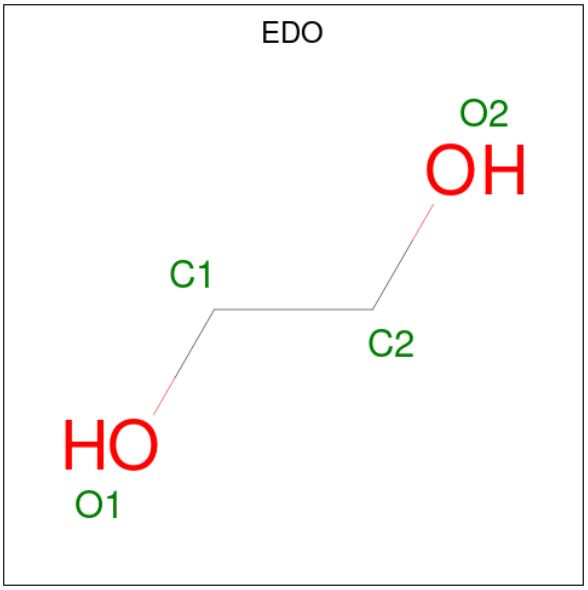
There are 4 unique types of molecules in this entry. The entry contains 5453 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Shikimate kinase.

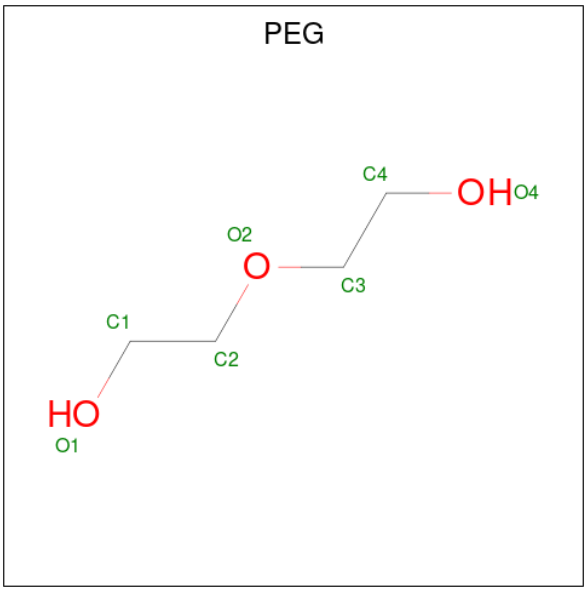
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	166	Total	C	N	O	S	Se	0	0	0
			1333	854	222	252	2	3			
1	B	168	Total	C	N	O	S	Se	0	0	0
			1341	858	224	254	2	3			
1	C	156	Total	C	N	O	S	Se	0	0	0
			1249	801	206	237	2	3			
1	D	158	Total	C	N	O	S	Se	0	0	0
			1257	805	208	239	2	3			

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

- Molecule 3 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	C	1	Total C O 7 4 3	0	0

- Molecule 4 is water.

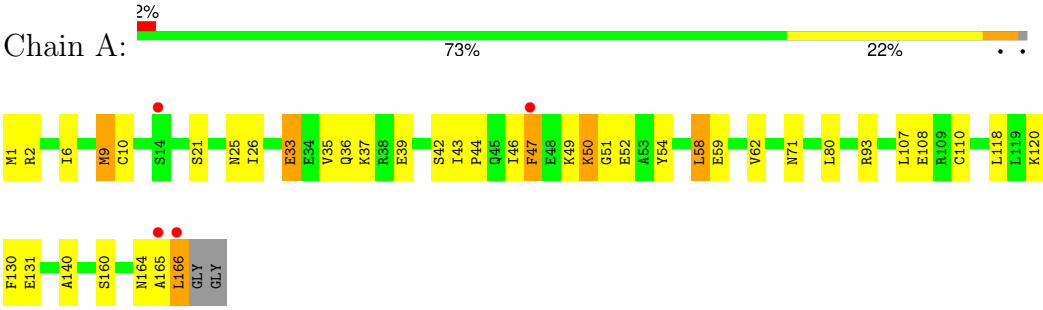
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	81	Total O 81 81	0	0
4	B	47	Total O 47 47	0	0
4	C	79	Total O 79 79	0	0
4	D	55	Total O 55 55	0	0

3 Residue-property plots

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.

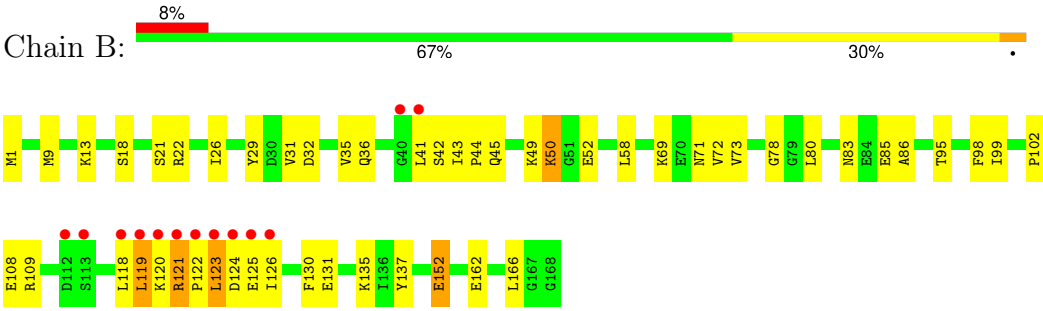
• Molecule 1: Shikimate kinase

Chain A:



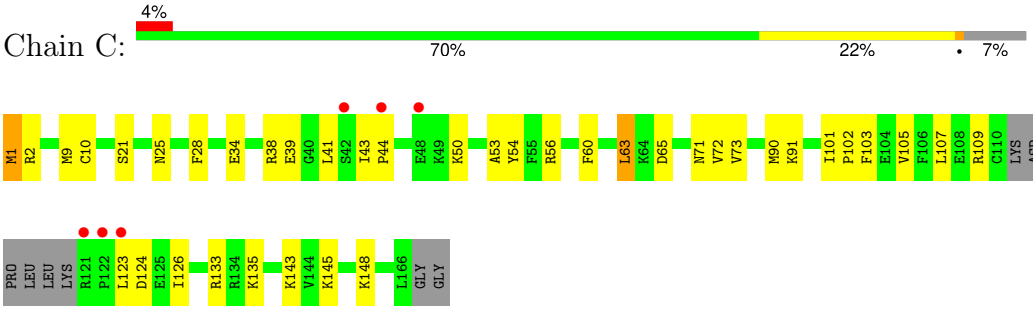
• Molecule 1: Shikimate kinase

Chain B:



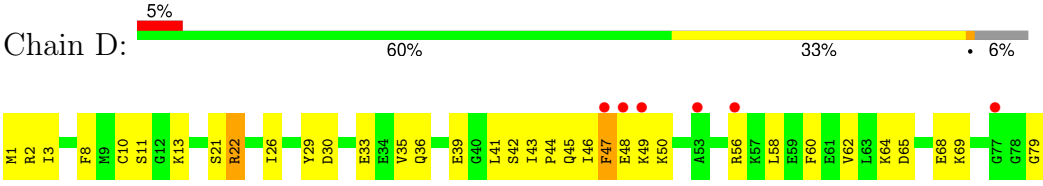
• Molecule 1: Shikimate kinase

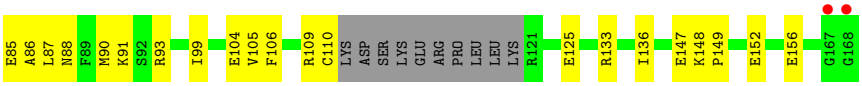
Chain C:



• Molecule 1: Shikimate kinase

Chain D:




E85
A86
L87
N88
F89
M90
K91
S92
R93
I99
E104
V105
F106
R109
C110
LYS
ASP
SER
LYS
GLU
ARG
PRO
LEU
LEU
LYS
R121
E125
R133
I136
E147
K148
P149
E152
E156
G167
G168

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	51.60Å 57.93Å 97.14Å 90.00° 90.59° 90.00°	Depositor
Resolution (Å)	38.53 – 2.10 49.76 – 2.10	Depositor EDS
% Data completeness (in resolution range)	96.8 (38.53-2.10) 96.9 (49.76-2.10)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	2.24 (at 2.10Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.214 , 0.269 0.208 , 0.265	Depositor DCC
R_{free} test set	1661 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å^2)	38.9	Xtriage
Anisotropy	0.321	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 56.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.027 for h,-k,-l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	5453	wwPDB-VP
Average B, all atoms (Å^2)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 37.64 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.1580e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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

5.1 Standard geometry

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

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	0.35	0/1351	0.58	0/1804
1	B	0.34	0/1359	0.64	1/1814 (0.1%)
1	C	0.37	0/1265	0.61	0/1689
1	D	0.36	0/1273	0.54	0/1699
All	All	0.36	0/5248	0.60	1/7006 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	123	LEU	N-CA-C	-6.90	92.36	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1333	0	1377	36	0
1	B	1341	0	1383	49	0
1	C	1249	0	1280	33	0
1	D	1257	0	1286	55	0
2	C	4	0	6	1	0
3	C	7	0	10	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	81	0	0	5	0
4	B	47	0	0	4	0
4	C	79	0	0	5	0
4	D	55	0	0	1	0
All	All	5453	0	5342	163	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 15.

All (163) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:56:ARG:HH22	1:D:79:GLY:HA3	1.01	1.08
1:A:166:LEU:HD22	1:D:56:ARG:HE	1.17	1.05
1:D:56:ARG:NH2	1:D:79:GLY:HA3	1.75	1.00
1:B:120:LYS:NZ	1:C:123:LEU:HB3	1.82	0.94
1:A:166:LEU:HD22	1:D:56:ARG:NE	1.87	0.89
1:A:166:LEU:CD2	1:D:56:ARG:HE	1.87	0.85
1:C:145:LYS:HB2	1:C:148:LYS:HE2	1.60	0.82
1:B:120:LYS:HZ2	1:C:123:LEU:HB3	1.46	0.80
1:D:56:ARG:HH22	1:D:79:GLY:CA	1.89	0.79
1:B:121:ARG:N	1:B:122:PRO:HD2	1.97	0.78
1:C:50:LYS:HD2	1:C:54:TYR:HB2	1.67	0.77
1:C:91:LYS:HD3	2:C:169:EDO:H12	1.67	0.77
1:A:166:LEU:HB2	1:D:133:ARG:HH22	1.51	0.76
1:A:43:ILE:HG22	1:A:47:PHE:CE1	2.23	0.73
1:A:9:MSE:HE2	1:A:10:CYS:HB3	1.71	0.71
1:C:25:ASN:ND2	1:D:125:GLU:HG3	2.06	0.70
1:D:82:ALA:O	1:D:136:ILE:HD13	1.92	0.70
1:B:83:ASN:HD22	1:B:86:ALA:H	1.40	0.69
1:B:123:LEU:H	1:B:125:GLU:H	1.40	0.69
1:C:65:ASP:HB3	4:C:431:HOH:O	1.93	0.68
1:D:35:VAL:HG13	1:D:58:LEU:HD23	1.74	0.68
1:B:121:ARG:N	1:B:122:PRO:CD	2.56	0.68
1:B:122:PRO:HB3	1:B:126:ILE:HG12	1.78	0.66
1:C:9:MSE:O	1:C:10:CYS:SG	2.53	0.66
1:D:82:ALA:O	1:D:136:ILE:HG21	1.96	0.66
1:D:42:SER:OG	1:D:44:PRO:HD2	1.96	0.65
1:C:133:ARG:HG2	1:C:133:ARG:HH11	1.62	0.64
1:A:43:ILE:HB	1:A:44:PRO:HD3	1.78	0.64
1:B:49:LYS:C	1:B:50:LYS:HD3	2.18	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:120:LYS:HZ1	1:C:123:LEU:HB3	1.61	0.64
1:D:22:ARG:NH2	1:D:22:ARG:HB2	2.12	0.64
1:B:122:PRO:HD3	1:B:126:ILE:HD11	1.80	0.64
1:D:1:MSE:HE2	1:D:3:ILE:HG12	1.80	0.64
1:D:43:ILE:HB	1:D:44:PRO:HD3	1.80	0.64
1:B:83:ASN:ND2	1:B:86:ALA:H	1.97	0.63
1:B:121:ARG:O	1:B:121:ARG:HG3	1.97	0.63
1:B:121:ARG:H	1:B:122:PRO:CD	2.11	0.62
1:D:42:SER:O	1:D:46:ILE:HG13	2.01	0.60
1:D:44:PRO:O	1:D:47:PHE:HB2	2.01	0.60
1:D:49:LYS:HG2	1:D:50:LYS:HD3	1.85	0.59
1:B:29:TYR:CE2	1:B:69:LYS:HE2	2.38	0.59
1:D:21:SER:HB2	1:D:26:ILE:O	2.02	0.59
1:A:165:ALA:HB3	4:A:415:HOH:O	2.02	0.59
1:B:152:GLU:HG3	4:B:279:HOH:O	2.01	0.58
1:A:42:SER:O	1:A:46:ILE:HG13	2.04	0.58
1:A:21:SER:HB2	1:A:26:ILE:O	2.02	0.58
4:C:257:HOH:O	1:D:125:GLU:HG3	2.03	0.58
1:A:33:GLU:O	1:A:37:LYS:HG3	2.03	0.58
1:A:9:MSE:HE1	1:A:110:CYS:HA	1.85	0.57
1:B:162:GLU:OE1	1:B:166:LEU:HD12	2.03	0.57
1:D:13:LYS:HB2	1:D:13:LYS:NZ	2.19	0.57
1:A:35:VAL:HG13	1:A:58:LEU:HD13	1.86	0.57
1:D:42:SER:H	1:D:45:GLN:HB2	1.70	0.57
1:C:39:GLU:O	1:C:41:LEU:HD22	2.04	0.57
1:C:124:ASP:HA	1:C:126:ILE:HG22	1.85	0.56
1:D:13:LYS:HG2	1:D:99:ILE:HD12	1.86	0.56
1:A:160:SER:O	1:A:164:ASN:ND2	2.34	0.56
1:B:107:LEU:O	1:B:119:LEU:HD13	2.06	0.55
1:C:50:LYS:O	1:C:50:LYS:HD3	2.07	0.55
1:B:120:LYS:NZ	1:C:123:LEU:O	2.39	0.55
1:B:123:LEU:C	1:B:125:GLU:H	2.03	0.55
1:A:1:MSE:O	1:A:71:ASN:HA	2.07	0.55
1:C:50:LYS:HD2	1:C:54:TYR:CB	2.37	0.55
1:B:43:ILE:N	1:B:44:PRO:HD2	2.22	0.54
1:A:36:GLN:HE21	1:A:42:SER:HA	1.71	0.54
1:A:118:LEU:HD13	1:A:126:ILE:HD12	1.90	0.54
1:B:131:GLU:O	1:B:135:LYS:HG3	2.08	0.54
1:D:83:ASN:ND2	1:D:86:ALA:H	2.06	0.54
1:D:87:LEU:O	1:D:91:LYS:HG3	2.08	0.54
1:B:42:SER:H	1:B:45:GLN:HE21	1.56	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:123:LEU:O	1:B:124:ASP:HB3	2.07	0.54
1:A:47:PHE:CD1	1:A:47:PHE:N	2.75	0.53
1:A:49:LYS:HA	4:A:380:HOH:O	2.08	0.53
1:D:39:GLU:O	1:D:41:LEU:HG	2.09	0.53
1:A:43:ILE:HG22	1:A:47:PHE:CZ	2.43	0.53
1:A:47:PHE:HA	1:A:50:LYS:O	2.09	0.53
1:C:143:LYS:HE3	4:C:399:HOH:O	2.08	0.53
1:A:2:ARG:HD2	1:A:93:ARG:HB3	1.90	0.52
1:A:47:PHE:H	1:A:47:PHE:HD1	1.52	0.52
1:C:1:MSE:SE	1:C:1:MSE:N	2.92	0.52
1:A:54:TYR:O	1:A:58:LEU:HB2	2.10	0.52
1:A:59:GLU:HG3	1:A:80:LEU:HD13	1.90	0.52
1:B:121:ARG:H	1:B:122:PRO:HD2	1.68	0.52
1:B:105:VAL:O	1:B:108:GLU:HG2	2.10	0.52
1:D:10:CYS:HB2	1:D:109:ARG:HB3	1.91	0.52
1:A:127:LYS:O	1:A:131:GLU:HG3	2.09	0.52
1:B:106:PHE:CE2	1:B:130:PHE:HB2	2.45	0.51
1:D:106:PHE:O	1:D:110:CYS:HB2	2.11	0.51
1:B:122:PRO:HD3	1:B:126:ILE:CD1	2.39	0.51
1:B:118:LEU:O	1:B:121:ARG:HB3	2.10	0.51
1:B:32:ASP:O	1:B:36:GLN:HG3	2.10	0.51
1:B:123:LEU:HB2	1:B:125:GLU:HB2	1.93	0.51
1:C:1:MSE:O	1:C:71:ASN:HA	2.10	0.50
1:D:152:GLU:O	1:D:156:GLU:HG3	2.12	0.50
1:C:53:ALA:HB3	4:C:271:HOH:O	2.11	0.50
1:D:8:PHE:O	1:D:11:SER:HB2	2.12	0.49
1:B:13:LYS:HG2	1:B:99:ILE:HD12	1.93	0.49
1:B:9:MSE:HE2	1:B:78:GLY:H	1.77	0.49
1:C:50:LYS:O	1:C:50:LYS:CD	2.60	0.49
1:C:102:PRO:HG2	1:C:105:VAL:HG23	1.95	0.49
1:B:109:ARG:NH1	4:B:367:HOH:O	2.45	0.49
1:A:43:ILE:HG22	1:A:47:PHE:HE1	1.75	0.48
1:C:1:MSE:HG2	1:C:2:ARG:N	2.29	0.48
1:D:60:PHE:CZ	1:D:64:LYS:HE3	2.49	0.48
1:A:6:ILE:HD11	1:A:140:ALA:HB2	1.95	0.48
1:D:8:PHE:HD1	4:D:245:HOH:O	1.97	0.48
1:B:41:LEU:HD22	1:B:45:GLN:NE2	2.30	0.47
1:C:34:GLU:O	1:C:38:ARG:HG3	2.14	0.47
1:C:72:VAL:HG22	1:C:73:VAL:N	2.30	0.47
1:A:9:MSE:CE	1:A:110:CYS:HA	2.43	0.47
1:B:31:VAL:O	1:B:35:VAL:HG12	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:36:GLN:NE2	1:D:43:ILE:HG13	2.30	0.47
1:B:123:LEU:C	1:B:125:GLU:N	2.62	0.47
1:A:166:LEU:HB2	1:D:133:ARG:NH2	2.25	0.47
1:A:35:VAL:HG22	1:A:62:VAL:HG21	1.97	0.47
1:D:87:LEU:HA	1:D:90:MSE:HE3	1.96	0.47
1:D:148:LYS:HB2	1:D:149:PRO:CD	2.45	0.47
1:C:43:ILE:HB	1:C:44:PRO:HD3	1.97	0.47
1:D:1:MSE:HE2	1:D:3:ILE:CG1	2.45	0.46
1:D:2:ARG:NH2	1:D:93:ARG:HH21	2.13	0.46
1:A:39:GLU:CD	1:A:58:LEU:HD11	2.36	0.46
1:C:10:CYS:HB2	1:C:109:ARG:CB	2.46	0.46
1:D:13:LYS:HB2	1:D:13:LYS:HZ3	1.81	0.45
1:B:22:ARG:HD2	4:B:338:HOH:O	2.15	0.45
1:D:58:LEU:O	1:D:62:VAL:HG23	2.17	0.45
1:B:83:ASN:HD21	1:B:85:GLU:HB3	1.81	0.45
1:D:35:VAL:HG13	1:D:58:LEU:CD2	2.44	0.45
1:B:102:PRO:HG2	1:B:105:VAL:HG23	1.97	0.45
1:B:52:GLU:H	1:B:52:GLU:CD	2.21	0.44
1:A:120:LYS:HB2	4:A:270:HOH:O	2.17	0.44
1:C:101:ILE:O	1:C:101:ILE:HG13	2.16	0.44
1:D:148:LYS:HB2	1:D:149:PRO:HD2	1.99	0.44
1:C:135:LYS:HE2	4:C:276:HOH:O	2.18	0.44
1:B:98:PHE:CD1	1:B:137:TYR:HB3	2.52	0.44
1:D:65:ASP:O	1:D:68:GLU:HB2	2.18	0.43
1:B:41:LEU:HB3	1:B:45:GLN:HE21	1.82	0.43
1:B:72:VAL:HG22	1:B:73:VAL:N	2.32	0.43
1:C:63:LEU:HD11	1:C:90:MSE:HE2	2.01	0.43
1:D:2:ARG:NH2	1:D:93:ARG:NH2	2.65	0.43
4:A:373:HOH:O	1:C:1:MSE:HE3	2.17	0.43
1:B:21:SER:HB2	1:B:26:ILE:O	2.17	0.43
1:C:10:CYS:HB2	1:C:109:ARG:HB3	1.99	0.43
1:D:22:ARG:HB2	1:D:22:ARG:CZ	2.47	0.43
1:D:148:LYS:HE2	1:D:148:LYS:HB3	1.91	0.43
1:D:104:GLU:HG2	1:D:105:VAL:N	2.33	0.43
1:D:29:TYR:CE2	1:D:69:LYS:HE3	2.54	0.43
1:D:82:ALA:C	1:D:136:ILE:HG21	2.39	0.43
1:D:30:ASP:HB3	1:D:33:GLU:HB2	2.01	0.42
1:A:51:GLY:O	1:A:54:TYR:HB3	2.19	0.42
1:D:88:ASN:HD22	1:D:88:ASN:HA	1.69	0.42
1:B:1:MSE:O	1:B:71:ASN:HA	2.19	0.42
1:D:35:VAL:HG22	1:D:62:VAL:HG21	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:83:ASN:HD21	1:D:85:GLU:HB3	1.85	0.42
1:A:108:GLU:HG3	4:A:269:HOH:O	2.18	0.42
1:A:130:PHE:CD1	1:A:130:PHE:C	2.94	0.42
1:C:103:PHE:CE1	1:C:107:LEU:HD13	2.55	0.41
1:B:41:LEU:CD2	1:B:45:GLN:NE2	2.83	0.41
1:B:42:SER:H	1:B:45:GLN:NE2	2.18	0.41
1:C:21:SER:CB	1:C:28:PHE:HB2	2.51	0.41
1:D:83:ASN:HD22	1:D:86:ALA:HB2	1.85	0.41
1:B:121:ARG:HB2	4:B:453:HOH:O	2.20	0.41
1:B:18:SER:O	1:B:22:ARG:HG3	2.21	0.41
1:D:49:LYS:CG	1:D:50:LYS:HD3	2.51	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	164/168 (98%)	161 (98%)	3 (2%)	0	100	100
1	B	166/168 (99%)	159 (96%)	6 (4%)	1 (1%)	25	21
1	C	152/168 (90%)	145 (95%)	7 (5%)	0	100	100
1	D	154/168 (92%)	146 (95%)	7 (4%)	1 (1%)	25	21
All	All	636/672 (95%)	611 (96%)	23 (4%)	2 (0%)	41	41

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	121	ARG
1	D	47	PHE

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	149/146 (102%)	140 (94%)	9 (6%)	19	16
1	B	149/146 (102%)	143 (96%)	6 (4%)	31	32
1	C	139/146 (95%)	135 (97%)	4 (3%)	42	46
1	D	139/146 (95%)	136 (98%)	3 (2%)	52	57
All	All	576/584 (99%)	554 (96%)	22 (4%)	33	34

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	MSE
1	A	25	ASN
1	A	33	GLU
1	A	47	PHE
1	A	50	LYS
1	A	52	GLU
1	A	58	LEU
1	A	107	LEU
1	A	166	LEU
1	B	50	LYS
1	B	58	LEU
1	B	80	LEU
1	B	95	THR
1	B	119	LEU
1	B	152	GLU
1	C	1	MSE
1	C	56	ARG
1	C	60	PHE
1	C	63	LEU
1	D	22	ARG
1	D	48	GLU
1	D	147	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	36	GLN
1	B	45	GLN
1	B	83	ASN
1	C	25	ASN
1	C	88	ASN
1	C	164	ASN
1	D	36	GLN
1	D	45	GLN
1	D	83	ASN
1	D	88	ASN
1	D	128	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)

2 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	PEG	C	170	-	6,6,6	0.80	0	5,5,5	0.50	0
2	EDO	C	169	-	3,3,3	0.62	0	2,2,2	0.49	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	PEG	C	170	-	-	1/4/4/4	-
2	EDO	C	169	-	-	0/1/1/1	-

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	C	170	PEG	O2-C3-C4-O4

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	169	EDO	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

6.1 Protein, DNA and RNA chains

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, 95^{th} 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	163/168 (97%)	-0.01	4 (2%)	57	62	28, 40, 77, 100	0
1	B	165/168 (98%)	0.40	13 (7%)	12	16	29, 49, 93, 99	0
1	C	153/168 (91%)	0.16	6 (3%)	39	45	26, 42, 78, 100	0
1	D	155/168 (92%)	0.26	8 (5%)	27	32	26, 44, 88, 101	0
All	All	636/672 (94%)	0.20	31 (4%)	29	35	26, 44, 87, 101	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	123	LEU	7.6
1	A	165	ALA	5.5
1	D	47	PHE	5.2
1	D	56	ARG	5.0
1	D	48	GLU	4.8
1	B	124	ASP	4.3
1	C	123	LEU	4.3
1	B	125	GLU	4.3
1	B	126	ILE	4.3
1	B	122	PRO	4.2
1	C	122	PRO	3.3
1	B	119	LEU	3.2
1	C	48	GLU	3.2
1	B	118	LEU	3.1
1	D	49	LYS	3.1
1	A	166	LEU	2.9
1	D	77	GLY	2.9
1	D	168	GLY	2.7
1	B	113	SER	2.6
1	B	41	LEU	2.6
1	B	120	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	47	PHE	2.5
1	B	112	ASP	2.5
1	C	121	ARG	2.4
1	C	42	SER	2.4
1	C	44	PRO	2.2
1	B	121	ARG	2.1
1	D	53	ALA	2.1
1	A	14	SER	2.0
1	D	167	GLY	2.0
1	B	40	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, 95^{th} 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	PEG	C	170	7/7	0.84	0.23	50,55,60,60	0
2	EDO	C	169	4/4	0.89	0.13	47,47,50,54	0

6.5 Other polymers (i)

There are no such residues in this entry.