Full wwPDB X-ray Structure Validation Report (i)

Dec 5, 2023 – 02:49 am GMT

PDB ID	:	2V4O
Title	:	Crystal structure of Salmonella typhimurium SurE at 2.75 angstrom resolution in monoclinic form
Authors	:	Anju, P.; Savithri, H.S.; Murthy, M.R.N.
Deposited on	:	2008-09-26
Resolution	:	2.71 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:

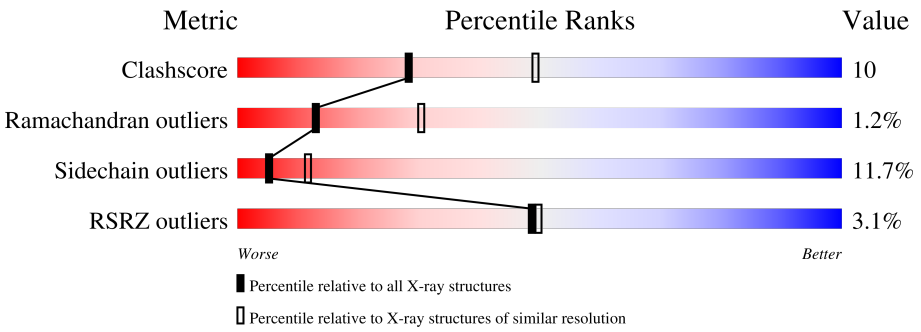
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.71 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))
Clashscore	141614	3686 (2.74-2.70)
Ramachandran outliers	138981	3622 (2.74-2.70)
Sidechain outliers	138945	3623 (2.74-2.70)
RSRZ outliers	127900	3276 (2.74-2.70)

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	267	3%; 68%, 22%, 5%, 5%
1	B	267	4%; 67%, 23%, 5%, 5%
1	C	267	4%; 74%, 19%, •, 5%
1	D	267	75%, 17%, 5%, •

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	PO4	D	1256	-	-	X	-

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 7636 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.

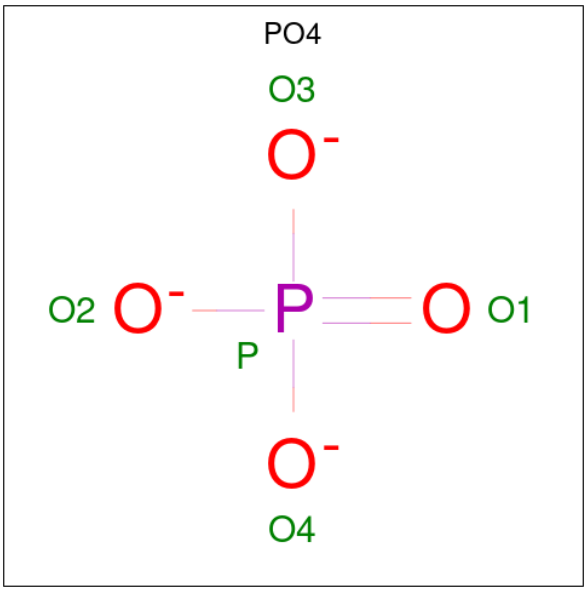
- Molecule 1 is a protein called MULTIFUNCTIONAL PROTEIN SUR E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	254	1883	1171	335	370	7	0	0	0
1	B	254	1864	1159	329	369	7	0	0	0
1	C	254	1885	1172	337	369	7	0	0	0
1	D	258	1910	1187	342	373	8	0	0	0

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

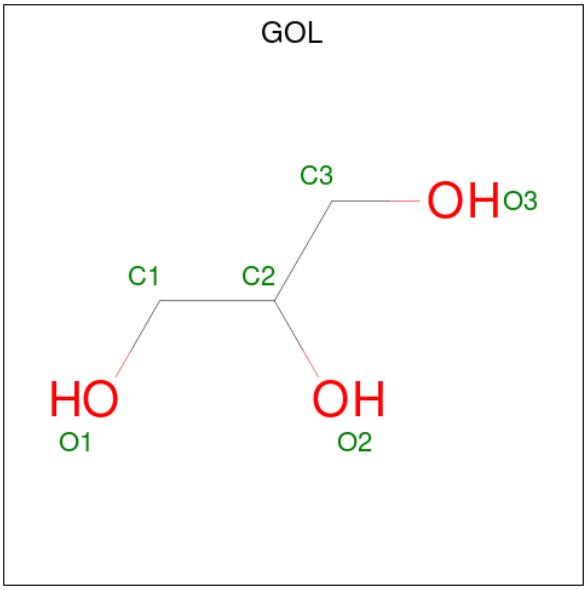
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
2	A	1	1	1	0	0
2	B	1	1	1	0	0
2	C	1	1	1	0	0
2	D	1	1	1	0	0

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	P		
3	A	1	5	4	1	0	0
			Total	O	P		
3	D	1	5	4	1	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 5 is water.

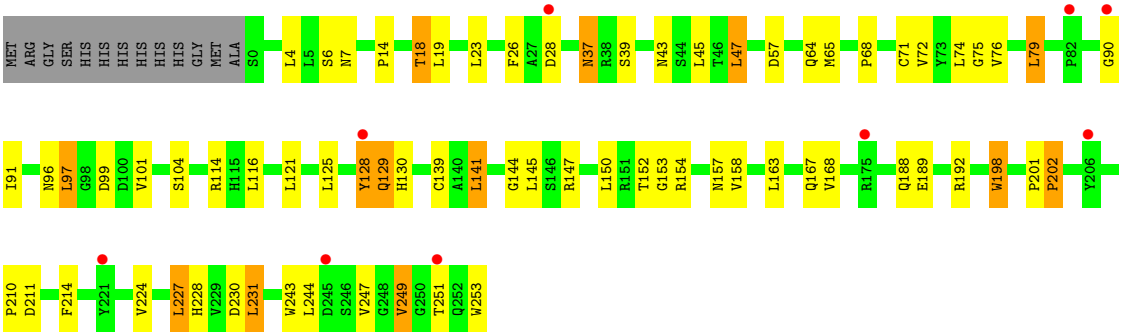
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	14	14	14	0	0
			Total	O		
5	B	11	11	11	0	0
			Total	O		
5	C	22	22	22	0	0
			Total	O		
5	D	27	27	27	0	0

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.

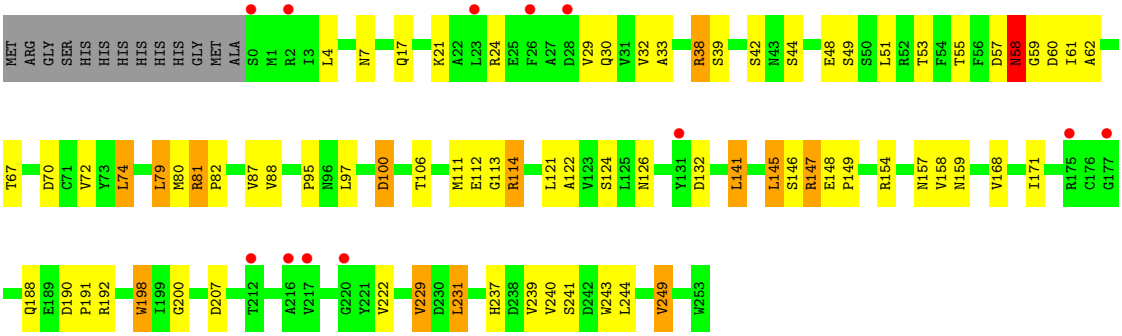
• Molecule 1: MULTIFUNCTIONAL PROTEIN SUR E

Chain A: 3% 68% 22% 5% 5%



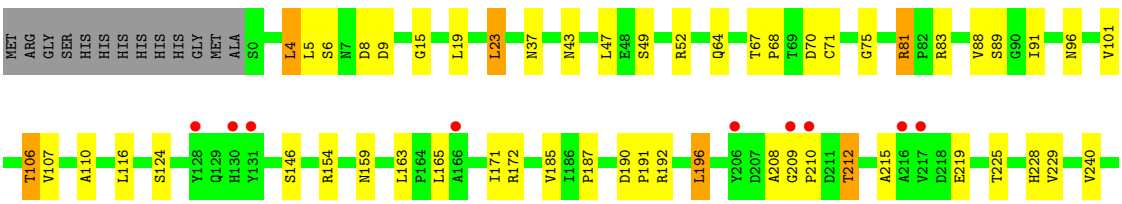
• Molecule 1: MULTIFUNCTIONAL PROTEIN SUR E

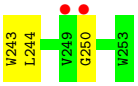
Chain B: 4% 67% 23% 5% 5%



• Molecule 1: MULTIFUNCTIONAL PROTEIN SUR E

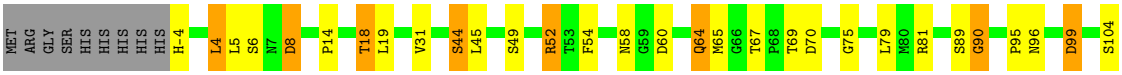
Chain C: 4% 74% 19% 5%

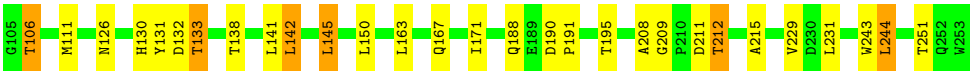




• Molecule 1: MULTIFUNCTIONAL PROTEIN SUR E

Chain D:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	161.08Å 95.30Å 94.57Å 90.00° 98.93° 90.00°	Depositor
Resolution (Å)	93.25 – 2.71 47.65 – 2.71	Depositor EDS
% Data completeness (in resolution range)	97.7 (93.25-2.71) 94.2 (47.65-2.71)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.96 (at 2.73Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.190 , 0.254 0.196 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	50.9	Xtriage
Anisotropy	0.300	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 56.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	7636	wwPDB-VP
Average B, all atoms (Å^2)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.09% 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, 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	0.79	3/1924 (0.2%)	0.87	1/2637 (0.0%)
1	B	0.75	0/1905	0.83	0/2614
1	C	0.79	0/1926	0.85	0/2638
1	D	0.82	1/1952 (0.1%)	0.96	5/2673 (0.2%)
All	All	0.79	4/7707 (0.1%)	0.88	6/10562 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	D	0	1
All	All	0	2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	253	TRP	CZ2-CH2	6.17	1.49	1.37
1	A	253	TRP	CD2-CE3	6.00	1.49	1.40
1	D	90	GLY	N-CA	5.44	1.54	1.46
1	A	253	TRP	NE1-CE2	5.31	1.44	1.37

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	81	ARG	NE-CZ-NH1	6.32	123.46	120.30
1	D	211	ASP	N-CA-CB	-5.78	100.20	110.60
1	D	60	ASP	CB-CG-OD1	5.53	123.27	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	253	TRP	CE3-CZ3-CH2	-5.47	115.19	121.20
1	D	4	LEU	CA-CB-CG	5.41	127.75	115.30
1	D	8	ASP	N-CA-CB	-5.09	101.44	110.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	200	GLY	Peptide
1	D	89	SER	Peptide

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	1883	0	1784	44	0
1	B	1864	0	1749	46	0
1	C	1885	0	1799	32	0
1	D	1910	0	1816	40	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	5	0	0	0	0
3	D	5	0	0	2	0
4	D	6	0	8	0	0
5	A	14	0	0	1	0
5	B	11	0	0	0	0
5	C	22	0	0	1	0
5	D	27	0	0	0	0
All	All	7636	0	7156	144	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 (144) 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:A:188:GLN:OE1	1:A:198:TRP:HZ3	1.31	1.11
1:D:130:HIS:HB3	1:D:133:THR:HG23	1.39	1.04
1:B:38:ARG:HH21	1:B:38:ARG:HG3	1.23	1.01
1:D:133:THR:HG21	1:D:163:LEU:O	1.63	0.98
1:A:188:GLN:OE1	1:A:198:TRP:CZ3	2.21	0.93
1:C:81:ARG:H	1:C:81:ARG:HH11	1.18	0.91
1:D:130:HIS:HB3	1:D:133:THR:CG2	2.03	0.89
1:A:14:PRO:O	1:A:18:THR:HG23	1.73	0.88
1:A:14:PRO:O	1:A:18:THR:CG2	2.23	0.87
1:D:209:GLY:O	1:D:212:THR:HB	1.74	0.86
1:A:243:TRP:CZ2	1:B:145:LEU:HD13	2.13	0.83
1:C:96:ASN:H	1:C:159:ASN:HD21	1.28	0.81
1:C:243:TRP:CZ2	1:D:145:LEU:HD13	2.16	0.80
1:D:54:PHE:HB2	1:D:64:GLN:HE22	1.48	0.79
1:B:171:ILE:HB	1:B:222:VAL:HB	1.67	0.76
1:C:81:ARG:H	1:C:81:ARG:NH1	1.85	0.74
1:C:209:GLY:O	1:C:212:THR:HB	1.88	0.74
1:A:243:TRP:O	1:A:247:VAL:HG22	1.89	0.71
1:D:208:ALA:HA	1:D:212:THR:HG21	1.73	0.70
1:D:212:THR:HG22	1:D:215:ALA:H	1.57	0.70
1:C:208:ALA:HA	1:C:212:THR:HG21	1.73	0.69
1:B:38:ARG:HG3	1:B:38:ARG:NH2	2.02	0.69
1:B:188:GLN:HG2	1:D:188:GLN:NE2	2.08	0.69
1:D:44:SER:HB3	1:D:67:THR:HG21	1.76	0.68
1:C:8:ASP:OD1	1:C:106:THR:HG22	1.94	0.67
1:B:38:ARG:HH21	1:B:38:ARG:CG	2.05	0.66
1:A:144:GLY:HA2	1:A:147:ARG:HB2	1.78	0.66
1:A:249:VAL:HG21	1:B:141:LEU:HD12	1.76	0.66
1:D:106:THR:CG2	3:D:1256:PO4:O1	2.45	0.65
1:A:14:PRO:O	1:A:18:THR:HG22	1.97	0.64
1:A:145:LEU:HD13	1:B:243:TRP:CZ2	2.34	0.63
1:B:42:SER:H	1:B:44:SER:HB2	1.66	0.61
1:C:52:ARG:HD2	1:C:64:GLN:OE1	2.01	0.61
1:B:33:ALA:O	1:B:62:ALA:HA	2.00	0.61
1:A:129:GLN:HB3	1:A:130:HIS:CD2	2.36	0.60
1:D:5:LEU:O	1:D:31:VAL:HA	2.01	0.60
1:D:96:ASN:ND2	1:D:104:SER:HB2	2.17	0.60
1:B:67:THR:O	1:B:70:ASP:HB2	2.03	0.59
1:A:101:VAL:HG12	1:B:231:LEU:HD23	1.83	0.59
1:D:14:PRO:O	1:D:18:THR:HG23	2.03	0.59
1:D:95:PRO:HG3	1:D:126:ASN:OD1	2.02	0.59
1:B:24:ARG:HG3	1:B:29:VAL:HG13	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:212:THR:HG22	1:C:215:ALA:H	1.67	0.59
1:B:97:LEU:O	1:B:100:ASP:HB2	2.04	0.58
1:C:67:THR:O	1:C:70:ASP:HB2	2.02	0.58
1:B:124:SER:HB2	1:B:159:ASN:OD1	2.02	0.58
1:C:81:ARG:O	1:C:81:ARG:HD2	2.04	0.58
1:A:45:LEU:HG	1:A:47:LEU:HD13	1.86	0.58
1:A:141:LEU:HD13	1:B:249:VAL:HG21	1.84	0.58
1:C:19:LEU:HG	1:C:23:LEU:HD22	1.86	0.57
1:D:6:SER:O	1:D:90:GLY:N	2.36	0.56
1:C:6:SER:O	1:C:89:SER:HA	2.07	0.55
1:D:243:TRP:CZ3	1:D:244:LEU:HD13	2.41	0.55
1:B:114:ARG:O	1:B:154:ARG:HD3	2.06	0.55
1:C:9:ASP:HA	1:C:37:ASN:ND2	2.22	0.55
1:D:106:THR:HG22	3:D:1256:PO4:O1	2.07	0.55
1:B:72:VAL:CG2	1:B:112:GLU:HB3	2.37	0.54
1:A:26:PHE:HE1	1:A:139:CYS:HG	1.55	0.54
1:A:19:LEU:HD22	1:A:91:ILE:CG1	2.38	0.54
1:C:81:ARG:HH11	1:C:81:ARG:N	1.99	0.53
1:A:97:LEU:HB3	1:A:214:PHE:CE1	2.44	0.53
1:D:243:TRP:CE3	1:D:244:LEU:HD13	2.44	0.53
1:A:4:LEU:HD11	1:A:75:GLY:HA3	1.91	0.52
1:B:42:SER:C	1:B:44:SER:H	2.13	0.52
1:D:58:ASN:C	1:D:58:ASN:OD1	2.48	0.52
1:A:157:ASN:HB2	1:A:227:LEU:HD22	1.93	0.51
1:B:111:MET:CE	1:B:229:VAL:HG23	2.40	0.51
1:A:150:LEU:HD21	1:B:240:VAL:HG22	1.92	0.51
1:C:190:ASP:HB2	1:C:191:PRO:CD	2.40	0.51
1:B:74:LEU:HB3	1:B:79:LEU:HD22	1.93	0.51
1:B:111:MET:HE2	1:B:229:VAL:HG23	1.93	0.50
1:C:15:GLY:HA3	1:C:91:ILE:O	2.12	0.50
1:C:107:VAL:O	1:C:110:ALA:HB3	2.12	0.50
1:C:4:LEU:HD11	1:C:75:GLY:HA3	1.92	0.49
1:A:6:SER:O	1:A:90:GLY:N	2.38	0.49
1:B:88:VAL:HG22	1:B:121:LEU:HD12	1.95	0.49
1:B:190:ASP:HB2	1:B:191:PRO:HD2	1.94	0.49
5:A:2003:HOH:O	1:B:48:GLU:HB2	2.13	0.48
1:B:58:ASN:HD22	1:B:60:ASP:H	1.61	0.48
1:B:81:ARG:HA	1:B:82:PRO:C	2.33	0.48
1:C:250:GLY:HA3	1:D:171:ILE:CG2	2.43	0.48
1:A:145:LEU:HA	1:B:243:TRP:HZ2	1.78	0.48
1:C:250:GLY:HA3	1:D:171:ILE:HG22	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:95:PRO:HD3	1:B:126:ASN:HA	1.94	0.48
1:C:101:VAL:HG12	1:D:231:LEU:HD12	1.96	0.47
1:A:247:VAL:HG23	1:A:249:VAL:HG13	1.96	0.47
1:B:72:VAL:HG22	1:B:112:GLU:HB3	1.96	0.47
1:C:49:SER:HA	5:C:2003:HOH:O	2.14	0.47
1:D:190:ASP:HB2	1:D:191:PRO:CD	2.45	0.47
1:B:87:VAL:HG13	1:B:113:GLY:HA3	1.97	0.47
1:A:163:LEU:HD22	1:A:167:GLN:HE21	1.80	0.47
1:B:147:ARG:HE	1:B:147:ARG:HB3	1.35	0.46
1:A:125:LEU:HD23	1:A:128:TYR:HA	1.96	0.46
1:A:37:ASN:ND2	1:A:39:SER:H	2.14	0.46
1:A:158:VAL:HG22	1:A:224:VAL:HG22	1.97	0.46
1:A:74:LEU:HB3	1:A:79:LEU:HD22	1.98	0.46
1:A:141:LEU:O	1:A:145:LEU:HB2	2.17	0.45
1:C:196:LEU:HD12	1:D:52:ARG:HB2	1.98	0.45
1:D:75:GLY:HA2	1:D:79:LEU:HB2	1.97	0.45
1:A:37:ASN:HD22	1:A:37:ASN:C	2.20	0.45
1:D:8:ASP:HB2	1:D:106:THR:HB	1.98	0.45
1:B:4:LEU:HD11	1:B:32:VAL:HG13	1.98	0.44
1:C:116:LEU:O	1:C:154:ARG:NH2	2.47	0.44
1:C:124:SER:HB2	1:C:159:ASN:ND2	2.32	0.44
1:D:138:THR:O	1:D:142:LEU:HD22	2.17	0.44
1:A:4:LEU:HD23	1:A:4:LEU:C	2.38	0.44
1:C:172:ARG:NH2	1:C:219:GLU:OE1	2.40	0.44
1:D:45:LEU:HD13	1:D:69:THR:CG2	2.48	0.43
1:D:111:MET:HE3	1:D:229:VAL:HG22	1.99	0.43
1:A:7:ASN:OD1	1:A:7:ASN:C	2.56	0.43
1:B:95:PRO:HD3	1:B:126:ASN:HD22	1.83	0.43
1:A:72:VAL:O	1:A:76:VAL:HG13	2.18	0.43
1:A:231:LEU:CD1	1:B:231:LEU:HD11	2.49	0.43
1:B:188:GLN:CD	1:B:198:TRP:CZ3	2.92	0.43
1:B:148:GLU:HA	1:B:149:PRO:HD3	1.82	0.42
1:A:228:HIS:CD2	1:A:230:ASP:H	2.38	0.42
1:C:185:VAL:O	1:C:187:PRO:HD3	2.19	0.42
1:D:99:ASP:OD2	1:D:99:ASP:N	2.45	0.42
1:B:72:VAL:HG21	1:B:112:GLU:HB3	2.00	0.42
1:D:14:PRO:O	1:D:18:THR:CG2	2.68	0.42
1:A:228:HIS:CD2	1:A:228:HIS:C	2.92	0.42
1:D:132:ASP:OD2	1:D:132:ASP:N	2.52	0.42
1:A:96:ASN:OD1	1:A:104:SER:HB2	2.20	0.42
1:B:55:THR:HG22	1:B:61:ILE:CG1	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:THR:HB	1:B:239:VAL:HG11	2.00	0.42
1:C:5:LEU:HD13	1:C:88:VAL:HG22	2.02	0.42
1:D:45:LEU:HD13	1:D:69:THR:HG22	2.00	0.41
1:B:17:GLN:NE2	1:B:17:GLN:HA	2.34	0.41
1:D:45:LEU:HD12	1:D:70:ASP:OD1	2.20	0.41
1:A:152:THR:HG22	1:A:153:GLY:N	2.36	0.41
1:D:190:ASP:HB2	1:D:191:PRO:HD2	2.02	0.41
1:C:68:PRO:O	1:C:71:CYS:N	2.53	0.41
1:C:171:ILE:HG22	1:C:172:ARG:N	2.35	0.41
1:D:243:TRP:CZ3	1:D:244:LEU:CD1	3.04	0.41
1:A:145:LEU:HD13	1:B:243:TRP:CH2	2.55	0.41
1:A:68:PRO:O	1:A:71:CYS:HB2	2.21	0.40
1:A:201:PRO:HA	1:A:202:PRO:HD2	1.91	0.40
1:D:19:LEU:HD23	1:D:90:GLY:HA2	2.02	0.40
1:A:145:LEU:CA	1:B:243:TRP:HZ2	2.34	0.40
1:C:240:VAL:HG22	1:D:150:LEU:HD21	2.03	0.40
1:B:30:GLN:HG2	1:B:80:MET:HG3	2.04	0.40
1:A:37:ASN:ND2	1:A:37:ASN:C	2.75	0.40
1:B:122:ALA:HB3	1:B:157:ASN:HD22	1.86	0.40
1:D:96:ASN:HD22	1:D:104:SER:HB2	1.86	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	252/267 (94%)	233 (92%)	14 (6%)	5 (2%)	7 17
1	B	252/267 (94%)	228 (90%)	20 (8%)	4 (2%)	9 23
1	C	252/267 (94%)	238 (94%)	12 (5%)	2 (1%)	19 41
1	D	256/267 (96%)	240 (94%)	15 (6%)	1 (0%)	34 58
All	All	1012/1068 (95%)	939 (93%)	61 (6%)	12 (1%)	13 30

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	128	TYR
1	B	58	ASN
1	B	100	ASP
1	A	43	ASN
1	A	114	ARG
1	D	131	TYR
1	A	202	PRO
1	B	237	HIS
1	C	210	PRO
1	C	83	ARG
1	A	210	PRO
1	B	59	GLY

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	191/215 (89%)	165 (86%)	26 (14%)	3 8
1	B	188/215 (87%)	159 (85%)	29 (15%)	2 6
1	C	193/215 (90%)	177 (92%)	16 (8%)	11 25
1	D	194/215 (90%)	175 (90%)	19 (10%)	8 18
All	All	766/860 (89%)	676 (88%)	90 (12%)	5 12

All (90) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	THR
1	A	23	LEU
1	A	28	ASP
1	A	37	ASN
1	A	47	LEU
1	A	57	ASP
1	A	64	GLN
1	A	65	MET

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Mol	Chain	Res	Type
1	A	79	LEU
1	A	97	LEU
1	A	99	ASP
1	A	116	LEU
1	A	121	LEU
1	A	129	GLN
1	A	141	LEU
1	A	154	ARG
1	A	168	VAL
1	A	189	GLU
1	A	192	ARG
1	A	198	TRP
1	A	211	ASP
1	A	227	LEU
1	A	231	LEU
1	A	244	LEU
1	A	249	VAL
1	A	251	THR
1	B	7	ASN
1	B	21	LYS
1	B	38	ARG
1	B	39	SER
1	B	49	SER
1	B	51	LEU
1	B	53	THR
1	B	57	ASP
1	B	58	ASN
1	B	74	LEU
1	B	79	LEU
1	B	81	ARG
1	B	106	THR
1	B	114	ARG
1	B	132	ASP
1	B	141	LEU
1	B	145	LEU
1	B	146	SER
1	B	147	ARG
1	B	158	VAL
1	B	168	VAL
1	B	192	ARG
1	B	198	TRP
1	B	207	ASP

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Mol	Chain	Res	Type
1	B	229	VAL
1	B	231	LEU
1	B	241	SER
1	B	244	LEU
1	B	249	VAL
1	C	4	LEU
1	C	23	LEU
1	C	43	ASN
1	C	47	LEU
1	C	81	ARG
1	C	106	THR
1	C	146	SER
1	C	163	LEU
1	C	165	LEU
1	C	192	ARG
1	C	196	LEU
1	C	212	THR
1	C	225	THR
1	C	228	HIS
1	C	229	VAL
1	C	244	LEU
1	D	-4	HIS
1	D	4	LEU
1	D	18	THR
1	D	44	SER
1	D	49	SER
1	D	52	ARG
1	D	64	GLN
1	D	65	MET
1	D	99	ASP
1	D	106	THR
1	D	133	THR
1	D	141	LEU
1	D	142	LEU
1	D	145	LEU
1	D	167	GLN
1	D	195	THR
1	D	212	THR
1	D	244	LEU
1	D	251	THR

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	17	GLN
1	A	37	ASN
1	A	92	ASN
1	A	130	HIS
1	A	167	GLN
1	A	228	HIS
1	A	252	GLN
1	B	7	ASN
1	B	17	GLN
1	B	58	ASN
1	B	126	ASN
1	B	157	ASN
1	B	228	HIS
1	B	237	HIS
1	C	37	ASN
1	C	43	ASN
1	C	92	ASN
1	C	126	ASN
1	C	159	ASN
1	C	188	GLN
1	C	234	HIS
1	D	64	GLN
1	D	96	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 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 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	PO4	D	1256	2	4,4,4	0.45	0	6,6,6	0.92	0
3	PO4	A	1255	2	4,4,4	0.72	0	6,6,6	0.78	0
4	GOL	D	1255	-	5,5,5	0.57	0	5,5,5	1.40	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
4	GOL	D	1255	-	-	1/4/4/4	-

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
4	D	1255	GOL	C1-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	1256	PO4	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	254/267 (95%)	0.26	9 (3%)	44	44	22, 47, 58, 65	1 (0%)
1	B	254/267 (95%)	0.37	12 (4%)	31	30	30, 53, 66, 75	0
1	C	254/267 (95%)	0.27	11 (4%)	35	34	34, 44, 55, 61	0
1	D	258/267 (96%)	0.07	0	100	100	34, 40, 49, 58	0
All	All	1020/1068 (95%)	0.24	32 (3%)	49	50	22, 46, 61, 75	1 (0%)

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	216	ALA	4.2
1	C	209	GLY	4.0
1	B	28	ASP	3.3
1	C	217	VAL	3.2
1	A	175	ARG	3.0
1	B	0	SER	2.8
1	C	249	VAL	2.8
1	C	210	PRO	2.7
1	C	250	GLY	2.7
1	B	216	ALA	2.7
1	A	128	TYR	2.6
1	B	23	LEU	2.6
1	B	220	GLY	2.5
1	B	175	ARG	2.5
1	C	130	HIS	2.5
1	A	221	TYR	2.4
1	C	206	TYR	2.4
1	A	82	PRO	2.4
1	B	177	GLY	2.3
1	C	128	TYR	2.3
1	B	217	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	90	GLY	2.2
1	A	251	THR	2.2
1	A	245	ASP	2.2
1	A	206	TYR	2.2
1	C	131	TYR	2.2
1	B	131	TYR	2.1
1	B	212	THR	2.1
1	A	28	ASP	2.1
1	B	26	PHE	2.1
1	C	166	ALA	2.0
1	B	2	ARG	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
2	MG	B	1254	1/1	0.57	0.13	59,59,59,59	0
2	MG	C	1254	1/1	0.92	0.12	44,44,44,44	0
3	PO4	A	1255	5/5	0.92	0.17	37,37,41,41	5
4	GOL	D	1255	6/6	0.92	0.16	49,54,55,57	0
3	PO4	D	1256	5/5	0.93	0.21	31,32,35,36	5
2	MG	A	1254	1/1	0.96	0.07	35,35,35,35	0
2	MG	D	1254	1/1	0.98	0.07	29,29,29,29	0

6.5 Other polymers (i)

There are no such residues in this entry.