Full wwPDB X-ray Structure Validation Report (i)

Mar 13, 2023 – 02:18 PM EDT

PDB ID	:	4OO2
Title	:	Streptomyces globisporus C-1027 FAD dependent (S)-3-chloro-β-tyrosine-S-SgcC2 C-5 hydroxylase SgcC apo form
Authors	:	Cao, H.; Xu, W.; Bingman, C.A.; Lohman, J.R.; Yennamalli, R.; Shen, B.; Phillips Jr., G.N.; Enzyme Discovery for Natural Product Biosynthesis (NatPro)
Deposited on	:	2014-01-29
Resolution	:	2.63 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:

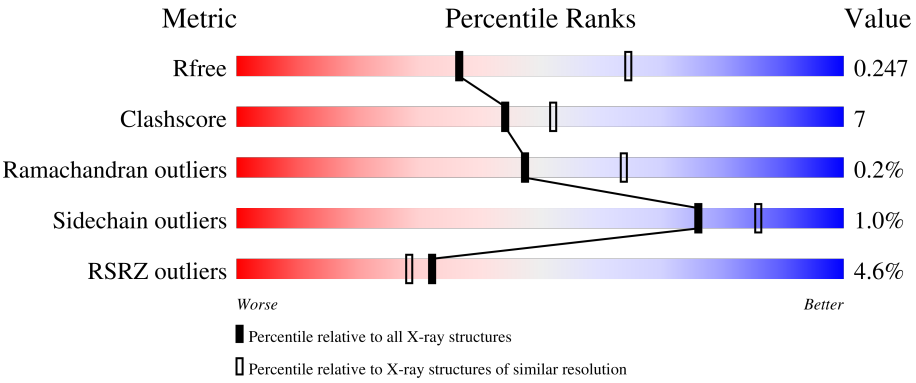
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.32.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.32.1

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 2.63 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	1426 (2.66-2.62)
Clashscore	141614	1472 (2.66-2.62)
Ramachandran outliers	138981	1446 (2.66-2.62)
Sidechain outliers	138945	1446 (2.66-2.62)
RSRZ outliers	127900	1408 (2.66-2.62)

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	527	7% 77% 19% ·
1	B	527	4% 78% 15% · 5%
1	C	527	5% 78% 18% ·
1	D	527	% 80% 17% ·

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 16163 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.

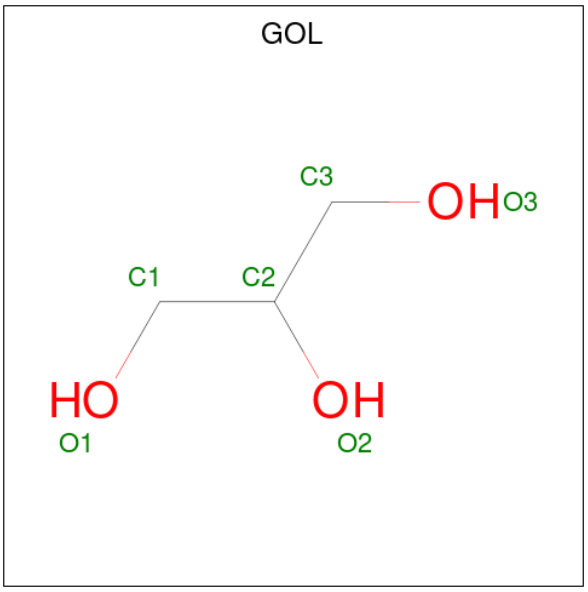
- Molecule 1 is a protein called Chlorophenol-4-monooxygenase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	506	Total	C	N	O	S	Se	0	2	0
			3982	2523	699	744	3	13			
1	B	499	Total	C	N	O	S	Se	0	4	0
			3942	2501	692	733	3	13			
1	C	510	Total	C	N	O	S	Se	0	2	0
			4006	2538	704	749	3	12			
1	D	508	Total	C	N	O	S	Se	0	4	0
			4008	2539	703	750	3	13			

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ca	0	0
			1	1		
2	C	1	Total	Ca	0	0
			1	1		

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



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

- Molecule 4 is water.

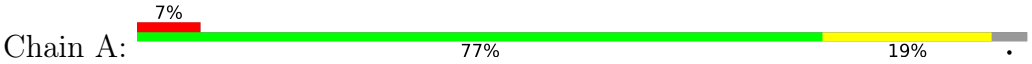
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	40	Total O 40 40	0	0
4	B	60	Total O 60 60	0	0
4	C	52	Total O 52 52	0	0
4	D	65	Total O 65 65	0	0

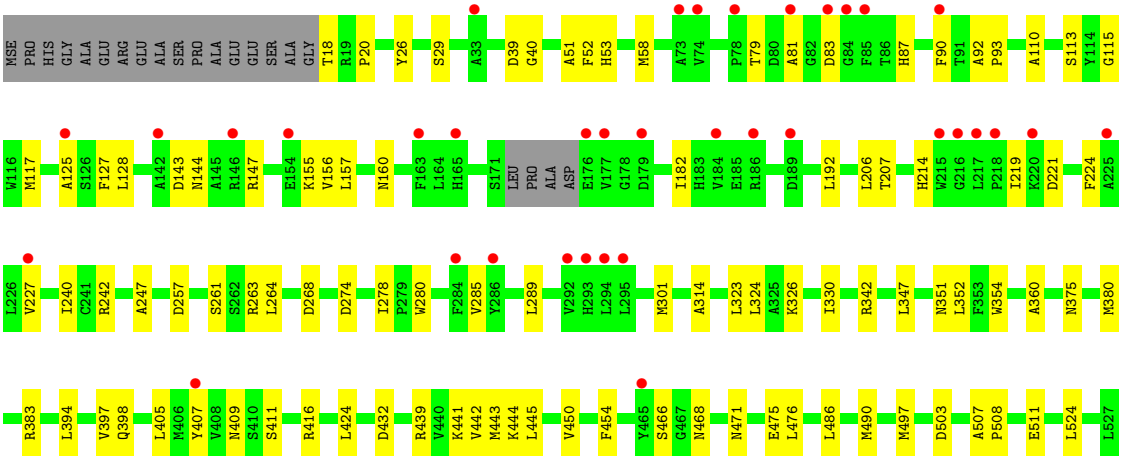
3 Residue-property plots (i)

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: Chlorophenol-4-monooxygenase

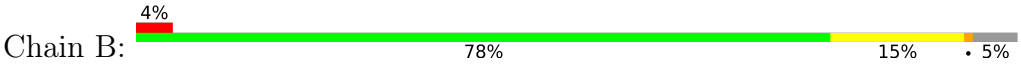
Chain A:

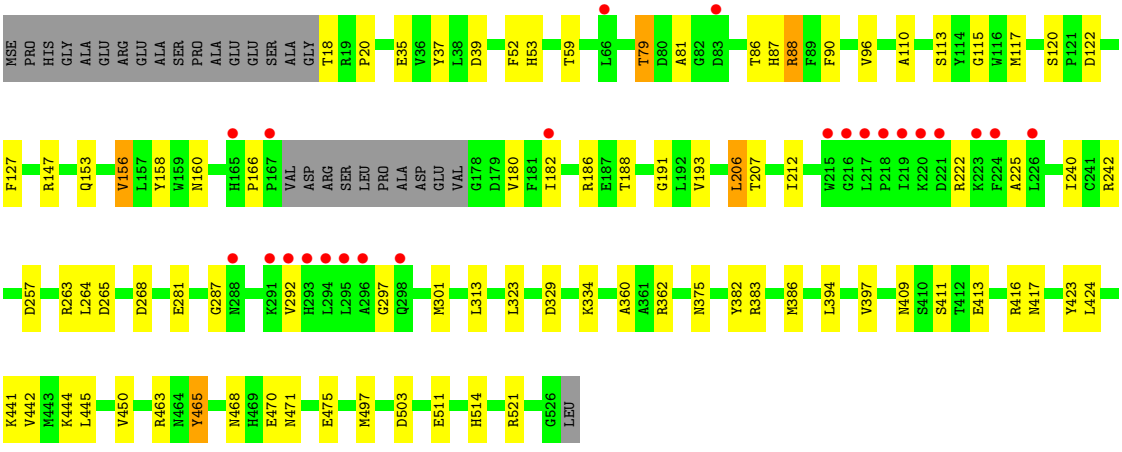




• Molecule 1: Chlorophenol-4-monooxygenase

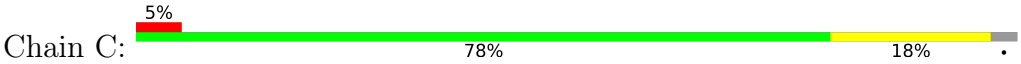
Chain B:

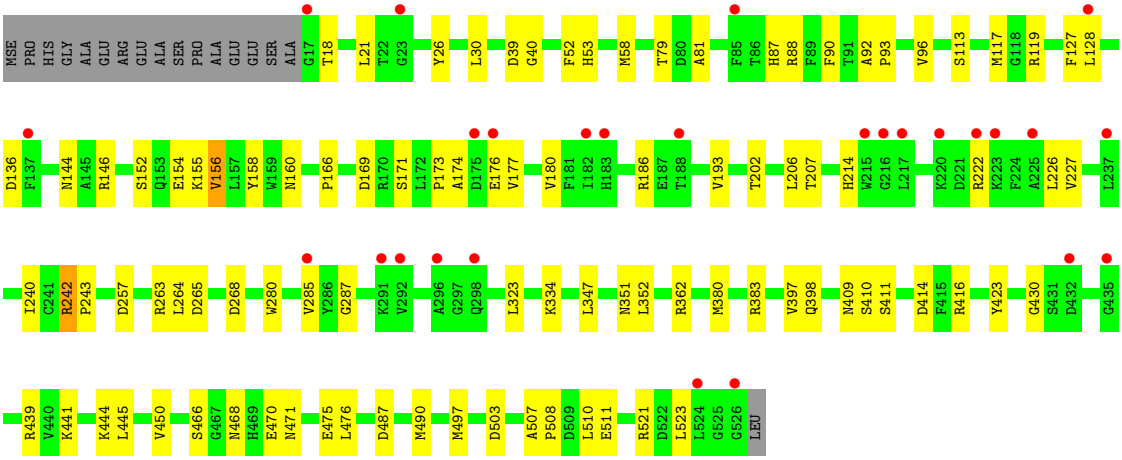




• Molecule 1: Chlorophenol-4-monooxygenase

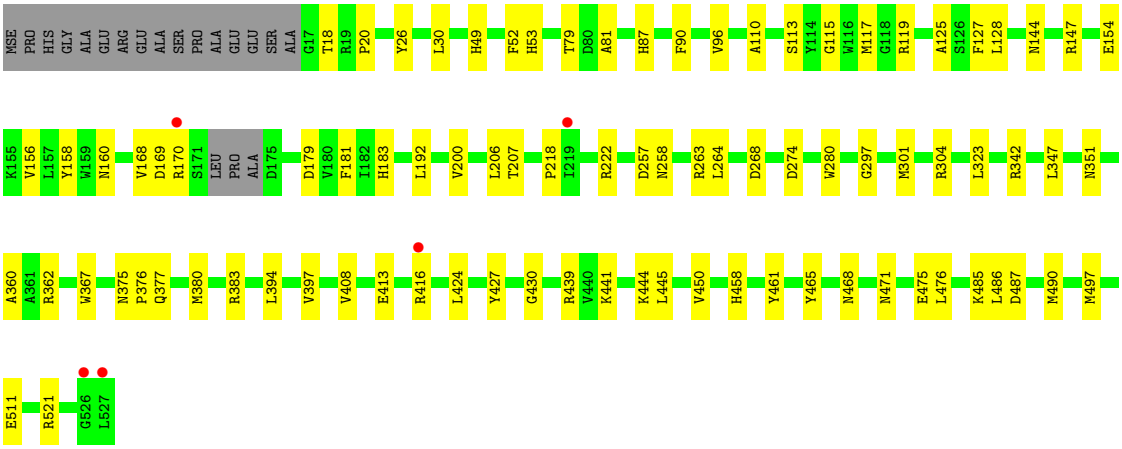
Chain C:





- Molecule 1: Chlorophenol-4-monooxygenase

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	99.16Å 173.73Å 113.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.67 – 2.63 47.67 – 2.63	Depositor EDS
% Data completeness (in resolution range)	94.4 (47.67-2.63) 79.8 (47.67-2.63)	Depositor EDS
R_{merge}	0.27	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	0.28 (at 2.61Å)	Xtriage
Refinement program	REFMAC, PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.217 , 0.246 0.222 , 0.247	Depositor DCC
R_{free} test set	1651 reflections (2.96%)	wwPDB-VP
Wilson B-factor (Å^2)	40.3	Xtriage
Anisotropy	0.413	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 45.0	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	0.92	EDS
Total number of atoms	16163	wwPDB-VP
Average B, all atoms (Å^2)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.21% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, 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	0.21	0/4072	0.40	0/5511
1	B	0.21	0/4032	0.39	0/5457
1	C	0.21	0/4098	0.38	0/5550
1	D	0.21	0/4099	0.39	0/5546
All	All	0.21	0/16301	0.39	0/22064

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	3982	0	3854	60	0
1	B	3942	0	3819	58	0
1	C	4006	0	3878	63	0
1	D	4008	0	3873	57	0
2	A	1	0	0	0	0
2	C	1	0	0	0	0
3	B	6	0	8	0	0
4	A	40	0	0	2	0
4	B	60	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	52	0	0	1	0
4	D	65	0	0	4	0
All	All	16163	0	15432	206	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 7.

All (206) 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:352:LEU:HD11	1:C:352:LEU:HD11	1.55	0.86
1:C:40:GLY:HA3	1:D:413:GLU:HB3	1.61	0.83
1:C:18:THR:HG21	1:C:154:GLU:HG3	1.64	0.79
1:B:503:ASP:OD1	1:D:416:ARG:NH2	2.15	0.78
1:A:503:ASP:OD1	1:C:416:ARG:NH2	2.20	0.75
1:B:88:ARG:NH1	4:B:737:HOH:O	2.22	0.71
1:B:188:THR:HG23	1:B:191:GLY:H	1.57	0.68
1:B:52:PHE:HB3	1:B:206:LEU:HD11	1.75	0.68
1:B:166:PRO:HG2	1:B:180:VAL:HG12	1.74	0.68
1:B:147:ARG:NH2	1:B:281:GLU:O	2.26	0.68
1:B:413:GLU:OE2	1:B:416:ARG:NH2	2.26	0.67
1:A:257:ASP:HA	1:A:444:LYS:HD3	1.76	0.66
1:A:416:ARG:NH2	1:C:503:ASP:OD1	2.28	0.66
1:C:323:LEU:HD11	1:C:445:LEU:HG	1.78	0.66
1:C:173:PRO:HG2	1:C:176:GLU:HG2	1.78	0.65
1:B:323:LEU:HD11	1:B:445:LEU:HG	1.80	0.64
1:C:160:ASN:HB3	1:C:207:THR:HG21	1.79	0.64
1:D:160:ASN:HB3	1:D:207:THR:HG21	1.80	0.64
1:A:342:ARG:NH2	1:B:470:GLU:OE2	2.31	0.64
1:D:486:LEU:HG	1:D:490:MSE:HE2	1.80	0.64
1:B:79:THR:HG22	1:B:153:GLN:HB3	1.81	0.63
1:D:468:ASN:OD1	1:D:471:ASN:ND2	2.32	0.63
1:B:360:ALA:HA	1:B:375:ASN:HB3	1.80	0.63
1:D:380[B]:MSE:HB3	1:D:476:LEU:HD13	1.81	0.62
1:B:160:ASN:HB3	1:B:207:THR:HG21	1.79	0.62
1:D:304:ARG:NH2	4:D:623:HOH:O	2.34	0.61
1:C:186:ARG:HH21	1:C:193:VAL:HG21	1.66	0.60
1:C:166:PRO:HG2	1:C:180:VAL:HG12	1.83	0.60
1:A:160:ASN:HB3	1:A:207:THR:HG21	1.84	0.60
1:B:257:ASP:HA	1:B:444:LYS:HD3	1.84	0.59
1:C:136:ASP:OD1	1:C:146:ARG:NH2	2.33	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:323:LEU:HD11	1:D:445:LEU:HG	1.85	0.59
1:C:409:ASN:ND2	1:C:411:SER:O	2.35	0.59
1:B:242:ARG:NH1	1:B:268:ASP:OD2	2.34	0.59
1:D:257:ASP:HA	1:D:444:LYS:HD3	1.85	0.59
1:B:222:ARG:HG3	1:B:287:GLY:HA2	1.85	0.58
1:D:362:ARG:NH1	4:D:630:HOH:O	2.36	0.58
1:B:383:ARG:NH2	1:B:475:GLU:OE1	2.31	0.58
1:D:383:ARG:NH2	1:D:475:GLU:OE1	2.33	0.58
1:A:383:ARG:NH2	1:A:475:GLU:OE1	2.32	0.57
1:C:257:ASP:HA	1:C:444:LYS:HD3	1.85	0.57
1:C:468:ASN:OD1	1:C:471:ASN:ND2	2.37	0.57
1:C:242:ARG:NH1	1:C:268:ASP:OD2	2.37	0.57
1:C:242:ARG:NH2	1:C:265:ASP:OD1	2.37	0.57
1:A:326:LYS:NZ	4:A:708:HOH:O	2.38	0.56
1:B:441:LYS:HG2	1:D:497:MSE:HE2	1.87	0.56
1:D:258:ASN:ND2	4:D:636:HOH:O	2.33	0.56
1:A:18:THR:N	1:A:83:ASP:O	2.38	0.56
1:A:330:ILE:HD12	1:A:442:VAL:HG11	1.88	0.56
1:C:398:GLN:NE2	4:C:725:HOH:O	2.39	0.56
1:D:144:ASN:ND2	1:D:280:TRP:O	2.37	0.56
1:A:398:GLN:OE1	1:B:463:ARG:NH2	2.38	0.56
1:A:524:LEU:HD13	1:D:49:HIS:CE1	2.41	0.55
1:A:497:MSE:HE2	1:C:441:LYS:HG2	1.87	0.55
1:B:468:ASN:OD1	1:B:471:ASN:ND2	2.40	0.55
1:C:470:GLU:OE2	1:D:342:ARG:NH2	2.39	0.55
1:B:497:MSE:HE2	1:D:445:LEU:HD23	1.89	0.54
1:C:383:ARG:NH2	1:C:475:GLU:OE1	2.36	0.54
1:D:377:GLN:HE21	1:D:485:LYS:HE3	1.72	0.54
1:A:323:LEU:HD11	1:A:445:LEU:HG	1.89	0.54
1:A:263:ARG:HG3	1:A:264:LEU:HG	1.90	0.54
1:B:409:ASN:ND2	1:B:411:SER:O	2.41	0.54
1:A:182:ILE:HB	1:A:214:HIS:HE1	1.73	0.53
1:A:143:ASP:O	1:A:147:ARG:HG2	2.08	0.53
1:A:445:LEU:HD23	1:C:497:MSE:HE2	1.89	0.53
1:C:96:VAL:HG13	1:C:362:ARG:HA	1.91	0.53
1:C:347:LEU:O	1:C:351:ASN:ND2	2.41	0.53
1:A:221:ASP:HB3	1:A:224:PHE:HD2	1.74	0.53
1:A:468:ASN:OD1	1:A:471:ASN:ND2	2.42	0.53
1:D:52:PHE:HB3	1:D:206:LEU:HD21	1.91	0.53
1:A:40:GLY:O	1:B:417:ASN:ND2	2.42	0.52
1:A:486:LEU:HG	1:A:490:MSE:HE2	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:263:ARG:HG3	1:C:264:LEU:HG	1.91	0.52
1:D:119:ARG:NH1	1:D:465:TYR:HE1	2.08	0.52
1:B:386:MSE:HE1	1:B:465:TYR:HD1	1.74	0.52
1:A:360:ALA:HA	1:A:375:ASN:HB3	1.92	0.52
1:B:79:THR:O	1:B:81:ALA:N	2.37	0.51
1:C:174:ALA:HA	1:C:177:VAL:HG22	1.90	0.51
1:B:394:LEU:HD22	1:B:450:VAL:HG22	1.93	0.51
1:A:144:ASN:ND2	1:A:280:TRP:O	2.44	0.51
1:A:221:ASP:HB3	1:A:224:PHE:CD2	2.45	0.51
1:B:263:ARG:HG3	1:B:264:LEU:HG	1.93	0.51
1:B:88:ARG:NH2	4:B:711:HOH:O	2.44	0.51
1:A:398:GLN:NE2	4:A:701:HOH:O	2.44	0.51
1:B:113:SER:HB2	1:B:117[A]:MSE:HG3	1.93	0.51
1:D:263:ARG:HG3	1:D:264:LEU:HG	1.93	0.51
1:A:394:LEU:HD22	1:A:450:VAL:HG22	1.92	0.50
1:C:127:PHE:HD2	1:C:128:LEU:HD12	1.76	0.50
1:C:144:ASN:ND2	1:C:280:TRP:O	2.44	0.50
1:B:503:ASP:O	1:B:514:HIS:NE2	2.31	0.50
1:D:458:HIS:HA	1:D:461:TYR:HB3	1.93	0.50
1:D:168:VAL:HG21	1:D:181:PHE:CZ	2.47	0.50
1:A:113:SER:HB2	1:A:117[A]:MSE:HG3	1.94	0.49
1:C:410:SER:OG	1:C:414:ASP:OD2	2.27	0.49
1:C:113:SER:HB2	1:C:117[A]:MSE:HG3	1.94	0.49
1:C:430:GLY:HA2	1:C:439:ARG:HH11	1.77	0.49
1:A:79:THR:O	1:A:81:ALA:N	2.45	0.48
1:C:240:ILE:HG21	1:D:424:LEU:HD21	1.94	0.48
1:D:18:THR:HG21	1:D:154:GLU:HG3	1.96	0.48
1:B:441:LYS:HE2	1:D:497:MSE:HB3	1.94	0.48
1:A:424:LEU:HD21	1:B:240:ILE:HG21	1.96	0.48
1:A:511:GLU:HB2	1:C:53:HIS:CE1	2.49	0.48
1:A:192:LEU:N	1:A:278:ILE:O	2.44	0.48
1:C:52:PHE:HB3	1:C:206:LEU:HD21	1.93	0.48
1:D:79:THR:O	1:D:81:ALA:N	2.44	0.48
1:D:347:LEU:O	1:D:351:ASN:ND2	2.46	0.48
1:D:430:GLY:HA2	1:D:439:ARG:HH11	1.78	0.48
1:B:53:HIS:CE1	1:D:511:GLU:HB2	2.48	0.48
1:B:127:PHE:HB2	1:B:301:MSE:HG3	1.96	0.48
1:D:181:PHE:CZ	1:D:218:PRO:HG2	2.49	0.48
1:A:20:PRO:HB3	1:A:156:VAL:HG23	1.96	0.47
1:C:58:MSE:HE2	1:C:264:LEU:HD21	1.95	0.47
1:C:334[B]:LYS:HB3	1:C:334[B]:LYS:HE3	1.62	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:LYS:HE3	1:A:157:LEU:HD21	1.96	0.47
1:B:242:ARG:NH2	1:B:265:ASP:OD1	2.44	0.47
1:B:59:THR:OG1	1:B:117[B]:MSE:HE1	2.15	0.47
1:C:119:ARG:HG3	1:C:202:THR:HB	1.97	0.47
1:C:186:ARG:NH2	1:C:193:VAL:HG21	2.29	0.46
1:C:169:ASP:OD2	1:C:171:SER:OG	2.22	0.46
1:C:180:VAL:HG13	1:D:427:TYR:HD1	1.81	0.46
1:A:432:ASP:N	1:A:432:ASP:OD1	2.49	0.46
1:C:397:VAL:HG11	1:C:450:VAL:HG21	1.97	0.46
1:B:497:MSE:HB3	1:D:441:LYS:HE2	1.98	0.46
1:D:179:ASP:HB2	1:D:183:HIS:CE1	2.51	0.46
1:A:51:ALA:HB2	1:A:247:ALA:HB2	1.98	0.46
1:A:53:HIS:CE1	1:C:511:GLU:HB2	2.51	0.46
1:A:39:ASP:OD1	1:D:521:ARG:NH1	2.49	0.45
1:A:441:LYS:HE2	1:C:497:MSE:HB3	1.98	0.45
1:C:423:TYR:OH	1:D:274:ASP:OD2	2.26	0.45
1:B:521:ARG:NH1	1:C:39:ASP:OD1	2.49	0.45
1:D:20:PRO:HB3	1:D:156:VAL:HG23	1.99	0.45
1:B:20:PRO:HB3	1:B:156:VAL:HG23	1.99	0.45
1:A:26:TYR:O	1:A:29:SER:OG	2.29	0.45
1:D:487:ASP:HA	1:D:490:MSE:HE3	1.97	0.45
1:A:240:ILE:HG21	1:B:424:LEU:HD21	1.98	0.45
1:A:274:ASP:OD2	1:B:423:TYR:OH	2.25	0.45
1:B:313[B]:LEU:HD12	1:B:382:TYR:HE1	1.82	0.45
1:A:127:PHE:HB2	1:A:301:MSE:SE	2.67	0.45
1:A:347:LEU:O	1:A:351:ASN:ND2	2.49	0.45
1:D:360:ALA:HA	1:D:375:ASN:HB3	1.99	0.44
1:A:110:ALA:O	1:A:115:GLY:N	2.48	0.44
1:A:409:ASN:ND2	1:A:411:SER:O	2.49	0.44
1:D:376:PRO:HB2	1:D:380[B]:MSE:HE3	1.99	0.44
1:D:125:ALA:HA	1:D:128:LEU:HB2	2.00	0.44
1:A:242:ARG:NE	1:A:268:ASP:H	2.16	0.44
1:D:380[A]:MSE:HB3	1:D:476:LEU:HD13	1.98	0.44
1:C:383:ARG:HD2	1:C:466:SER:HB2	2.00	0.44
1:D:110:ALA:O	1:D:115:GLY:N	2.49	0.44
1:B:360:ALA:HA	1:B:375:ASN:CB	2.48	0.44
1:D:376:PRO:O	1:D:380[B]:MSE:HG3	2.18	0.44
1:A:407:TYR:HD1	1:A:439:ARG:HH21	1.65	0.44
1:C:380:MSE:HB3	1:C:476:LEU:HD13	1.99	0.44
1:A:227:VAL:HB	1:A:285:VAL:HB	1.99	0.44
1:A:324:LEU:HD13	1:A:397:VAL:HG22	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:35:GLU:HG2	1:B:37:TYR:CE2	2.52	0.43
1:B:329:ASP:HA	1:B:334[B]:LYS:HG2	1.98	0.43
1:D:87:HIS:HB3	1:D:90:PHE:CD2	2.53	0.43
1:A:507:ALA:HA	1:A:508:PRO:HD3	1.86	0.43
1:D:169:ASP:OD1	1:D:170:ARG:N	2.50	0.43
1:A:125:ALA:HA	1:A:128:LEU:HB2	1.99	0.43
1:C:92:ALA:HA	1:C:93:PRO:HD3	1.81	0.43
1:D:26:TYR:CZ	1:D:30:LEU:HD11	2.53	0.43
1:A:87:HIS:HB3	1:A:90:PHE:CD2	2.53	0.43
1:B:442:VAL:HG22	1:D:497:MSE:HE1	2.01	0.43
1:C:21:LEU:HG	1:C:155:LYS:HB3	2.01	0.43
1:A:52:PHE:HB3	1:A:206:LEU:HD21	2.01	0.43
1:A:58:MSE:HE1	1:C:510:LEU:HD21	2.01	0.43
1:B:182:ILE:HG21	1:B:212:ILE:HG21	2.01	0.43
1:D:408:VAL:HB	1:D:444:LYS:NZ	2.34	0.43
1:B:186:ARG:HB3	1:B:193:VAL:HB	2.01	0.43
1:C:523:LEU:HD23	1:C:523:LEU:HA	1.89	0.43
1:B:110:ALA:O	1:B:115:GLY:N	2.47	0.42
1:A:261:SER:HB3	1:A:454:PHE:HB2	2.00	0.42
1:A:92:ALA:HA	1:A:93:PRO:HD3	1.89	0.42
1:B:511:GLU:HB2	1:D:53:HIS:CE1	2.54	0.42
1:C:79:THR:O	1:C:81:ALA:N	2.46	0.42
1:B:18:THR:O	1:B:18:THR:OG1	2.35	0.42
1:D:200:VAL:HG13	1:D:268:ASP:HB3	2.01	0.42
1:A:383:ARG:HH11	1:A:466:SER:HB2	1.85	0.42
1:C:113:SER:HB3	1:C:117[A]:MSE:HE2	2.00	0.42
1:B:79:THR:HG23	1:B:86:THR:HB	2.01	0.42
1:B:39:ASP:OD1	1:C:521:ARG:NH1	2.53	0.41
1:C:507:ALA:HA	1:C:508:PRO:HD3	1.83	0.41
1:D:397:VAL:HG11	1:D:450:VAL:HG21	2.02	0.41
1:D:96:VAL:HG13	1:D:362:ARG:HA	2.01	0.41
1:C:214:HIS:NE2	1:C:226:LEU:O	2.40	0.41
1:C:227:VAL:HB	1:C:285:VAL:HB	2.01	0.41
1:C:242:ARG:HG2	1:C:243:PRO:N	2.35	0.41
1:D:113:SER:HB3	1:D:117[B]:MSE:HE2	2.01	0.41
1:D:222:ARG:NH1	4:D:642:HOH:O	2.39	0.41
1:B:87:HIS:HB3	1:B:90:PHE:CD2	2.55	0.41
1:B:96:VAL:HG13	1:B:362:ARG:HA	2.02	0.41
1:A:405:LEU:HA	1:A:443:MSE:HG3	2.03	0.41
1:B:120:SER:OG	1:B:122:ASP:OD1	2.31	0.41
1:C:128:LEU:HD23	1:C:152:SER:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:394:LEU:HD22	1:D:450:VAL:HG22	2.03	0.41
1:C:487:ASP:HA	1:C:490:MSE:HE2	2.02	0.40
1:C:222:ARG:HG3	1:C:287:GLY:HA2	2.03	0.40
1:A:314:ALA:HB1	1:A:354:TRP:CD1	2.57	0.40
1:B:147:ARG:NH2	1:B:281:GLU:HB2	2.37	0.40
1:C:90:PHE:CE2	1:C:156:VAL:HG13	2.57	0.40
1:D:127:PHE:HB2	1:D:301:MSE:HG3	2.03	0.40
1:B:225:ALA:HB1	1:B:292:VAL:HG21	2.04	0.40
1:B:397:VAL:HG11	1:B:450:VAL:HG21	2.03	0.40
1:C:26:TYR:CZ	1:C:30:LEU:HD11	2.57	0.40
1:C:87:HIS:HB3	1:C:90:PHE:CD2	2.57	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	504/527 (96%)	482 (96%)	22 (4%)	0	100	100
1	B	499/527 (95%)	477 (96%)	20 (4%)	2 (0%)	34	48
1	C	510/527 (97%)	486 (95%)	23 (4%)	1 (0%)	47	64
1	D	508/527 (96%)	483 (95%)	24 (5%)	1 (0%)	47	64
All	All	2021/2108 (96%)	1928 (95%)	89 (4%)	4 (0%)	47	64

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	297	GLY
1	B	297	GLY
1	C	156	VAL
1	B	156	VAL

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	412/412 (100%)	407 (99%)	5 (1%)	71	83
1	B	407/412 (99%)	402 (99%)	5 (1%)	71	83
1	C	414/412 (100%)	411 (99%)	3 (1%)	84	91
1	D	415/412 (101%)	411 (99%)	4 (1%)	76	86
All	All	1648/1648 (100%)	1631 (99%)	17 (1%)	76	86

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	219	ILE
1	A	289	LEU
1	A	380[A]	MSE
1	A	380[B]	MSE
1	A	476	LEU
1	B	79	THR
1	B	88	ARG
1	B	158	TYR
1	B	206	LEU
1	B	465	TYR
1	C	88	ARG
1	C	158	TYR
1	C	242	ARG
1	D	147	ARG
1	D	158	TYR
1	D	192	LEU
1	D	367	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	468	ASN
1	A	471	ASN

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Mol	Chain	Res	Type
1	C	377	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

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

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	GOL	B	601	-	5,5,5	0.37	0	5,5,5	0.24	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	B	601	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	601	GOL	O1-C1-C2-C3
3	B	601	GOL	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

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	495/527 (93%)	0.49	36 (7%)	15	12	41, 59, 95, 117	0
1	B	488/527 (92%)	0.33	23 (4%)	31	27	40, 53, 82, 114	0
1	C	499/527 (94%)	0.40	27 (5%)	25	22	38, 56, 85, 103	0
1	D	497/527 (94%)	0.10	5 (1%)	82	81	34, 47, 68, 114	0
All	All	1979/2108 (93%)	0.33	91 (4%)	32	28	34, 53, 87, 117	0

All (91) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	220	LYS	6.0
1	C	292	VAL	5.8
1	B	217	LEU	5.7
1	C	17	GLY	5.1
1	C	296	ALA	4.0
1	B	295	LEU	3.9
1	A	292	VAL	3.8
1	A	215	TRP	3.7
1	C	137	PHE	3.7
1	A	83	ASP	3.7
1	A	85	PHE	3.6
1	C	526	GLY	3.6
1	A	284	PHE	3.5
1	C	216	GLY	3.4
1	C	217	LEU	3.3
1	A	179	ASP	3.3
1	A	293	HIS	3.3
1	B	224	PHE	3.2
1	A	407	TYR	3.2
1	B	215	TRP	3.1
1	B	288	ASN	3.1

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Mol	Chain	Res	Type	RSRZ
1	C	220	LYS	3.1
1	B	293	HIS	3.0
1	B	182	ILE	3.0
1	A	227	VAL	3.0
1	A	217	LEU	2.9
1	A	81	ALA	2.9
1	B	219	ILE	2.9
1	B	218	PRO	2.9
1	B	220	LYS	2.8
1	D	170	ARG	2.8
1	D	527	LEU	2.7
1	A	154	GLU	2.7
1	C	182	ILE	2.7
1	C	85	PHE	2.7
1	C	128	LEU	2.7
1	A	73	ALA	2.7
1	B	83	ASP	2.7
1	B	221	ASP	2.7
1	B	216	GLY	2.6
1	B	292	VAL	2.6
1	C	285	VAL	2.6
1	B	223	LYS	2.5
1	B	294	LEU	2.5
1	D	416	ARG	2.5
1	A	84	GLY	2.5
1	A	225	ALA	2.4
1	A	189	ASP	2.4
1	A	33	ALA	2.4
1	C	237	LEU	2.4
1	A	90	PHE	2.4
1	C	435	GLY	2.4
1	B	165	HIS	2.3
1	D	219	ILE	2.3
1	A	184	VAL	2.3
1	A	177	VAL	2.3
1	A	295	LEU	2.3
1	C	524	LEU	2.3
1	C	222	ARG	2.3
1	A	125	ALA	2.3
1	A	465	TYR	2.3
1	A	142	ALA	2.3
1	C	291	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	175	ASP	2.2
1	C	23	GLY	2.2
1	C	176	GLU	2.2
1	B	66	LEU	2.2
1	C	215	TRP	2.2
1	A	216	GLY	2.2
1	A	78	PRO	2.2
1	C	188	THR	2.2
1	A	286	TYR	2.2
1	A	186	ARG	2.2
1	A	218	PRO	2.2
1	C	183	HIS	2.2
1	A	163	PHE	2.2
1	A	176	GLU	2.2
1	B	226	LEU	2.2
1	C	432	ASP	2.2
1	A	146	ARG	2.1
1	C	223	LYS	2.1
1	C	225	ALA	2.1
1	C	298	GLN	2.1
1	A	74	VAL	2.1
1	B	296	ALA	2.1
1	B	167	PRO	2.1
1	B	298	GLN	2.1
1	A	294	LEU	2.0
1	B	291	LYS	2.0
1	A	165	HIS	2.0
1	D	526	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
2	CA	C	601	1/1	0.78	0.22	59,59,59,59	0
3	GOL	B	601	6/6	0.78	0.27	54,63,64,68	0
2	CA	A	601	1/1	0.86	0.22	60,60,60,60	0

6.5 Other polymers (i)

There are no such residues in this entry.