



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:33 AM UTC

PDB ID : 2YPO / pdb_00002ypo
Title : 3-deoxy-D-arabino-heptulosonate 7-phosphate synthase with phenylalanine bound in only one site
Authors : Blackmore, N.J.; Reichau, S.; Jiao, W.; Hutton, R.D.; Baker, E.N.; Jameson, G.B.; Parker, E.J.
Deposited on : 2012-10-31
Resolution : 2.00 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

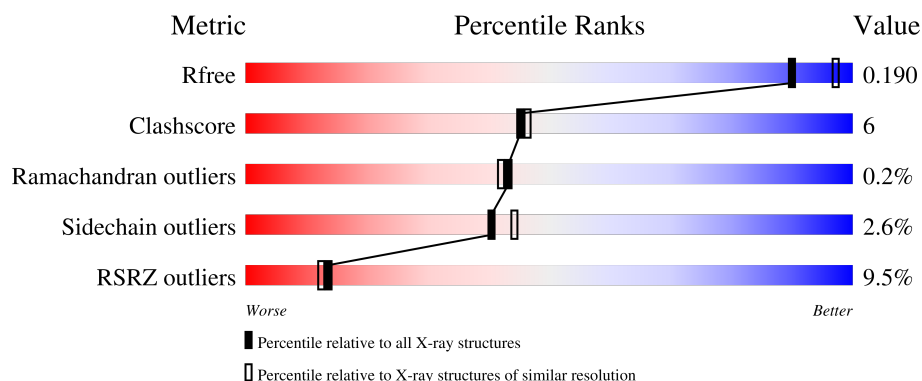
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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 $\leq 5\%$. 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	462	10% 84% 14% ..
1	B	462	9% 84% 13% ..

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	SO4	A	803	-	-	-	X

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 7568 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.

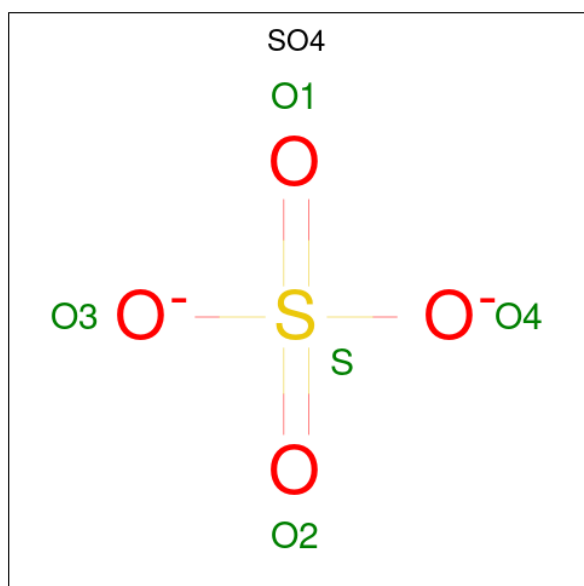
- Molecule 1 is a protein called PHOSPHO-2-DEHYDRO-3-DEOXYHEPTONATE ALDOLASE AROG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	458	Total	C	N	O	S	0	4	0
			3547	2217	644	668	18			
1	B	457	Total	C	N	O	S	0	5	0
			3547	2217	646	666	18			

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

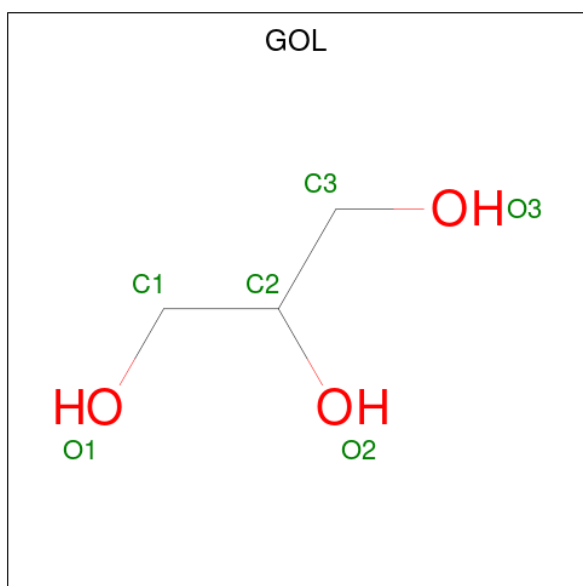
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mn	0	0
			1	1		
2	B	1	Total	Mn	0	0
			1	1		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



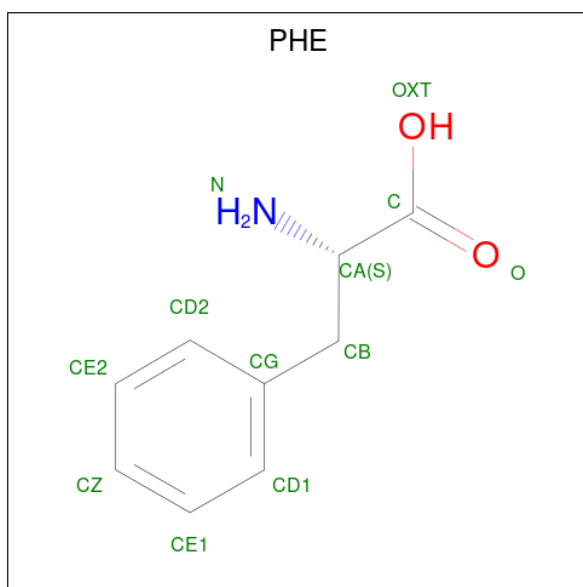
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 5 is PHENYLALANINE (CCD ID: PHE) (formula: $C_9H_{11}NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			12	9	1	2		
5	B	1	Total	C	N	O	0	0
			12	9	1	2		

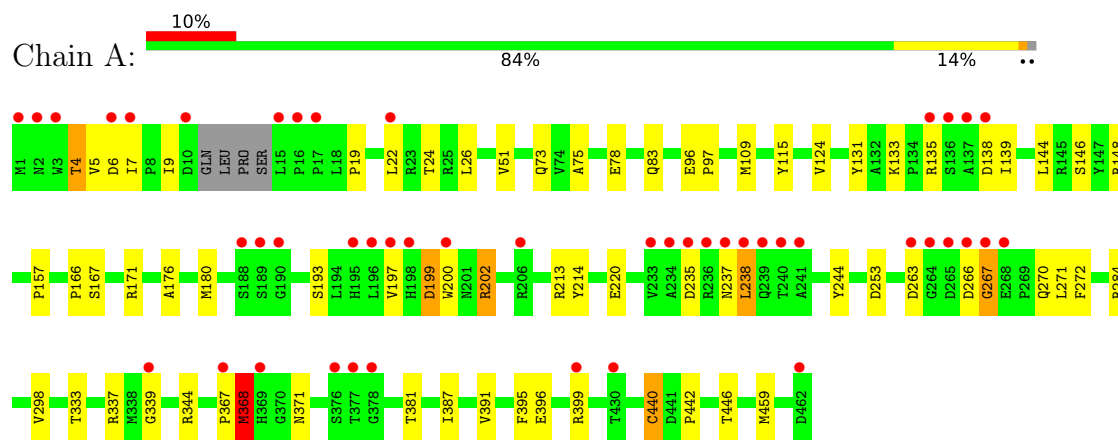
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	168	Total	O	0	0
			168	168		
6	B	219	Total	O	0	0
			219	219		

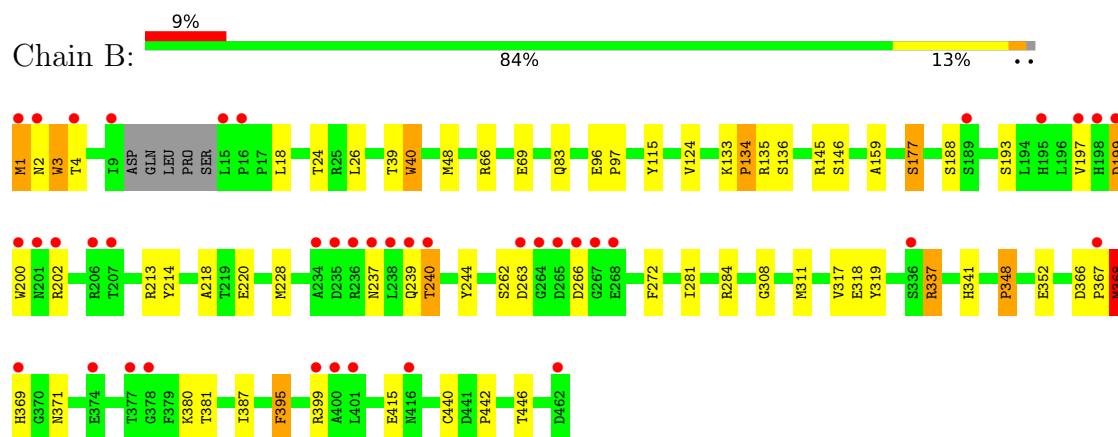
3 Residue-property plots [i](#)

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: PHOSPHO-2-DEHYDRO-3-DEOXYHEPTONATE ALDOLASE AROG



• Molecule 1: PHOSPHO-2-DEHYDRO-3-DEOXYHEPTONATE ALDOLASE AROG



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	203.97Å 203.97Å 66.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.16 – 2.00 44.16 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (44.16-2.00) 99.8 (44.16-2.00)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.89 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.138 , 0.159 0.171 , 0.190	Depositor DCC
R_{free} test set	5365 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	23.5	Xtriage
Anisotropy	0.056	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 34.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.000 for -h,-k,l	Xtriage
Reported twinning fraction	0.808 for H, K, L 0.192 for -h,-k,l	Depositor
Outliers	0 of 106995 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7568	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.71% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ 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: MN, GOL, SO4

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	1.47	15/3629 (0.4%)	1.17	7/4935 (0.1%)
1	B	1.49	22/3635 (0.6%)	1.19	6/4943 (0.1%)
All	All	1.48	37/7264 (0.5%)	1.18	13/9878 (0.1%)

All (37) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	139	ILE	C-O	-8.29	1.15	1.24
1	B	348	PRO	CA-C	-7.76	1.45	1.52
1	A	144	LEU	CA-C	-7.61	1.43	1.52
1	A	157	PRO	CA-C	-7.04	1.46	1.52
1	A	333	THR	N-CA	-6.88	1.37	1.46
1	B	134	PRO	N-CA	-6.80	1.40	1.46
1	A	199	ASP	N-CA	6.61	1.54	1.46
1	B	319	TYR	N-CA	6.17	1.53	1.46
1	B	199	ASP	N-CA	6.17	1.54	1.46
1	B	177	SER	N-CA	-6.13	1.38	1.46
1	B	188	SER	C-O	-6.12	1.15	1.23
1	A	399	ARG	CA-C	-6.10	1.44	1.52
1	B	40	TRP	C-O	-6.04	1.18	1.24
1	A	166	PRO	N-CA	-5.99	1.39	1.47
1	A	19	PRO	CA-C	5.98	1.59	1.52
1	B	40	TRP	CA-C	-5.88	1.46	1.52
1	B	317	VAL	CA-CB	5.85	1.61	1.54
1	B	159	ALA	N-CA	-5.81	1.39	1.46
1	A	176	ALA	CA-C	5.59	1.59	1.52
1	A	399	ARG	N-CA	-5.57	1.39	1.46
1	A	109	MET	CA-CB	5.55	1.61	1.53
1	A	171	ARG	C-O	5.49	1.30	1.24
1	A	180	MET	C-O	5.41	1.30	1.24
1	A	253	ASP	CA-C	5.31	1.59	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	318	GLU	CA-C	5.30	1.59	1.52
1	B	26	LEU	N-CA	5.27	1.52	1.46
1	B	218	ALA	CA-C	5.23	1.59	1.52
1	B	228	MET	CA-C	-5.20	1.46	1.52
1	A	75	ALA	CA-C	5.14	1.59	1.52
1	B	69	GLU	CA-C	-5.13	1.46	1.52
1	B	284	ARG	CA-C	-5.12	1.45	1.52
1	B	317	VAL	N-CA	-5.07	1.40	1.46
1	B	24	THR	CA-CB	5.07	1.61	1.53
1	B	281	ILE	CA-C	-5.02	1.46	1.52
1	B	395	PHE	N-CA	5.01	1.52	1.46
1	B	380	LYS	CA-C	5.01	1.59	1.52
1	B	311	MET	N-CA	-5.00	1.39	1.46

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	368	MET	N-CA-C	7.94	119.94	111.28
1	B	368	MET	N-CA-C	6.94	118.84	111.28
1	B	3	TRP	N-CA-C	-6.92	99.70	110.42
1	B	348	PRO	N-CA-C	6.61	118.77	110.70
1	B	240	THR	N-CA-CB	6.16	119.17	110.36
1	A	138	ASP	N-CA-CB	6.01	120.19	110.40
1	B	177	SER	CA-CB-OG	-5.75	99.60	111.10
1	A	339	GLY	N-CA-C	-5.54	106.09	112.29
1	A	138	ASP	N-CA-C	-5.43	106.15	112.89
1	A	399	ARG	N-CA-C	5.16	117.30	111.11
1	A	131	TYR	OH-CZ-CE2	-5.12	104.54	119.90
1	A	399	ARG	CA-C-O	5.08	126.34	120.90
1	B	341	HIS	N-CA-C	-5.07	107.05	113.18

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	3547	0	3508	51	0
1	B	3547	0	3515	38	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	10	0	0	0	0
3	B	15	0	0	1	0
4	A	18	0	23	2	0
4	B	18	0	23	3	0
5	A	12	0	8	0	0
5	B	12	0	8	0	0
6	A	168	0	0	2	0
6	B	219	0	0	3	0
All	All	7568	0	7085	84	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 6.

All (84) 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:5:VAL:HG11	1:B:48:MET:HE1	1.17	1.10
1:A:244:TYR:CE2	1:A:244:TYR:CE1	2.36	1.05
1:A:244:TYR:CE1	1:A:244:TYR:OH	2.17	0.97
1:A:395:PHE:CD1	1:A:459:MET:CE	2.50	0.94
1:A:22:LEU:HD11	1:A:271:LEU:HD22	1.47	0.93
1:A:395:PHE:CD1	1:A:459:MET:HE2	2.07	0.88
1:A:244:TYR:CE2	1:A:244:TYR:OH	2.28	0.87
1:A:24:THR:OG1	4:A:801:GOL:H32	1.79	0.82
1:A:5:VAL:CG1	1:B:48:MET:HE1	2.07	0.82
1:B:133:LYS:NZ	1:B:440:CYS:SG	2.53	0.80
1:A:344:ARG:NH1	1:A:396:GLU:OE1	2.21	0.74
1:A:9:ILE:HD12	1:B:2:ASN:HA	1.70	0.73
1:B:197:VAL:HA	1:B:200:TRP:CE3	2.26	0.71
1:A:395:PHE:CD1	1:A:459:MET:HE1	2.25	0.70
1:B:136:SER:N	3:B:806:SO4:O1	2.23	0.68
1:A:24:THR:OG1	4:A:801:GOL:C3	2.42	0.67
1:A:197:VAL:HA	1:A:200:TRP:CE3	2.30	0.66
1:A:5:VAL:HG11	1:B:48:MET:CE	2.11	0.64
1:A:7:ILE:O	1:B:2:ASN:HB2	2.00	0.61
1:A:135:ARG:HG3	1:A:146:SER:HB3	1.80	0.61
1:A:9:ILE:HG21	1:A:51:VAL:HG21	1.84	0.59
1:A:266:ASP:O	1:A:267:GLY:C	2.46	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:THR:HG23	1:B:4:THR:HG23	1.86	0.57
1:A:391:VAL:HG12	1:A:459:MET:HE1	1.86	0.56
1:A:22:LEU:HD21	1:A:271:LEU:HD13	1.86	0.56
1:B:135:ARG:HH11	1:B:146:SER:HB3	1.70	0.56
1:A:395:PHE:HD1	1:A:459:MET:HE1	1.70	0.55
1:A:235:ASP:HA	1:A:238:LEU:HD22	1.89	0.55
1:A:135:ARG:HG3	1:A:146:SER:CB	2.36	0.55
1:A:6:ASP:C	1:A:7:ILE:HD12	2.32	0.55
1:B:308:GLY:HA2	1:B:337:ARG:O	2.06	0.54
1:A:367:PRO:HB2	1:A:387:ILE:HG23	1.89	0.53
1:A:367:PRO:O	1:A:371:ASN:ND2	2.42	0.52
1:B:367:PRO:HB2	1:B:387:ILE:HG23	1.91	0.52
4:B:803:GOL:H11	6:B:2156:HOH:O	2.09	0.52
1:B:96:GLU:HB3	1:B:97:PRO:HD3	1.91	0.52
1:A:133:LYS:NZ	1:A:440:CYS:SG	2.72	0.50
1:B:395:PHE:O	1:B:399:ARG:HG3	2.12	0.50
1:B:83:GLN:HA	1:B:124:VAL:O	2.13	0.49
1:A:83:GLN:HA	1:A:124:VAL:O	2.13	0.49
1:B:262:SER:HB3	1:B:272:PHE:CE1	2.48	0.48
1:A:395:PHE:CG	1:A:459:MET:CE	2.96	0.48
1:B:145:ARG:NH1	6:B:2112:HOH:O	2.47	0.47
1:B:368:MET:H	1:B:368:MET:HG2	1.49	0.47
1:A:199:ASP:HA	1:A:202:ARG:HG3	1.96	0.47
1:B:1:MET:HE2	1:B:2:ASN:H	1.79	0.47
1:A:135:ARG:HG2	1:A:148:ARG:HD2	1.97	0.47
1:A:368:MET:HE3	1:A:368:MET:HB3	1.67	0.46
1:B:96:GLU:N	1:B:97:PRO:CD	2.79	0.46
1:A:135:ARG:HA	1:A:135:ARG:HD2	1.71	0.46
1:A:167:SER:HB2	1:B:3:TRP:HA	1.97	0.46
1:B:367:PRO:O	1:B:371:ASN:ND2	2.47	0.46
1:B:381:THR:HA	1:B:442:PRO:HG3	1.97	0.46
1:B:213:ARG:HD3	1:B:214:TYR:CZ	2.51	0.46
1:A:73:GLN:NE2	1:A:78:GLU:OE1	2.48	0.46
1:B:368:MET:HB3	1:B:368:MET:HE3	1.74	0.46
1:A:368:MET:H	1:A:368:MET:HG2	1.51	0.45
1:A:133:LYS:HE3	6:A:2082:HOH:O	2.15	0.45
1:A:22:LEU:CD1	1:A:271:LEU:HD22	2.33	0.45
1:A:270:GLN:HB2	1:A:272:PHE:CE1	2.52	0.45
1:A:395:PHE:CG	1:A:459:MET:HE2	2.49	0.45
1:A:263:ASP:OD1	1:A:263:ASP:C	2.60	0.44
6:A:2109:HOH:O	4:B:804:GOL:H12	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:39:THR:O	1:B:39:THR:OG1	2.33	0.44
1:B:115:TYR:OH	1:B:220:GLU:HG2	2.18	0.43
1:A:381:THR:HA	1:A:442:PRO:HG3	2.00	0.43
1:B:134:PRO:O	1:B:135:ARG:HD2	2.17	0.43
4:B:804:GOL:H32	6:B:2136:HOH:O	2.18	0.43
1:B:133:LYS:HZ2	1:B:440:CYS:CB	2.30	0.43
1:A:96[B]:GLU:HB3	1:A:97:PRO:HD3	2.00	0.43
1:B:199:ASP:HA	1:B:202:ARG:HE	1.84	0.43
1:A:213:ARG:HD3	1:A:214:TYR:CZ	2.54	0.43
1:A:96[A]:GLU:HB3	1:A:97:PRO:HD3	2.01	0.42
1:B:48:MET:N	1:B:48:MET:HE2	2.33	0.42
1:B:366:ASP:OD2	1:B:369:HIS:ND1	2.39	0.42
1:A:115:TYR:OH	1:A:220:GLU:HG2	2.20	0.42
1:B:135:ARG:NH1	1:B:146:SER:HB3	2.34	0.41
1:B:66[A]:ARG:HD2	1:B:244:TYR:OH	2.20	0.41
1:A:7:ILE:C	1:B:2:ASN:HB2	2.45	0.41
1:B:193:SER:O	1:B:197:VAL:HG23	2.20	0.41
1:B:263:ASP:O	1:B:266:ASP:OD1	2.38	0.41
1:A:26[A]:LEU:CD2	1:A:298:VAL:HG21	2.51	0.41
1:B:348:PRO:O	1:B:352:GLU:HG3	2.20	0.41
1:A:193:SER:O	1:A:197:VAL:HG23	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	458/462 (99%)	448 (98%)	8 (2%)	2 (0%)	30	27
1	B	458/462 (99%)	449 (98%)	9 (2%)	0	100	100
All	All	916/924 (99%)	897 (98%)	17 (2%)	2 (0%)	43	42

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	440	CYS
1	A	267	GLY

5.3.2 Protein sidechains ⓘ

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	374/376 (100%)	366 (98%)	8 (2%)	47	52
1	B	375/376 (100%)	364 (97%)	11 (3%)	37	40
All	All	749/752 (100%)	730 (98%)	19 (2%)	40	45

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	THR
1	A	202	ARG
1	A	237	ASN
1	A	238	LEU
1	A	284	ARG
1	A	337	ARG
1	A	368	MET
1	A	446	THR
1	B	1	MET
1	B	18	LEU
1	B	40	TRP
1	B	177	SER
1	B	237	ASN
1	B	239	GLN
1	B	240	THR
1	B	337	ARG
1	B	368	MET
1	B	415	GLU
1	B	446	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	GLN
1	A	239	GLN
1	A	297	GLN
1	A	373	HIS
1	B	239	GLN
1	B	326	HIS
1	B	360	GLN
1	B	373	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 2 are monoatomic - leaving 13 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	$\# Z > 2$
4	GOL	B	804	-	5,5,5	0.25	0	5,5,5	0.31	0
3	SO4	B	806	-	4,4,4	0.57	0	6,6,6	1.07	1 (16%)
3	SO4	A	800	-	4,4,4	0.39	0	6,6,6	0.47	0
4	GOL	A	802	-	5,5,5	0.84	0	5,5,5	0.67	0
3	SO4	B	802	-	4,4,4	0.49	0	6,6,6	0.41	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GOL	B	805	-	5,5,5	0.99	0	5,5,5	1.78	1 (20%)
5	PHE	A	900	-	11,12,12	0.61	0	11,15,15	0.98	1 (9%)
3	SO4	A	803	-	4,4,4	0.97	0	6,6,6	0.76	0
3	SO4	B	801	-	4,4,4	1.14	0	6,6,6	0.82	0
4	GOL	A	804	-	5,5,5	0.23	0	5,5,5	0.37	0
4	GOL	A	801	-	5,5,5	1.12	1 (20%)	5,5,5	1.28	0
4	GOL	B	803	-	5,5,5	0.77	0	5,5,5	1.20	1 (20%)
5	PHE	B	900	-	11,12,12	1.47	2 (18%)	11,15,15	0.86	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	B	804	-	-	0/4/4/4	-
4	GOL	A	802	-	-	2/4/4/4	-
4	GOL	B	805	-	-	2/4/4/4	-
5	PHE	A	900	-	-	1/8/8/8	0/1/1/1
4	GOL	A	804	-	-	1/4/4/4	-
4	GOL	A	801	-	-	2/4/4/4	-
4	GOL	B	803	-	-	1/4/4/4	-
5	PHE	B	900	-	-	0/8/8/8	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	900	PHE	O-C	2.88	1.30	1.22
5	B	900	PHE	CB-CG	2.60	1.57	1.51
4	A	801	GOL	O3-C3	-2.35	1.32	1.42

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	805	GOL	C3-C2-C1	-2.95	100.99	111.80
5	A	900	PHE	OXT-C-O	-2.84	117.65	124.08
4	B	803	GOL	O2-C2-C1	2.26	118.53	109.18
3	B	806	SO4	O3-S-O2	2.26	121.36	109.56

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	801	GOL	C1-C2-C3-O3
4	A	802	GOL	C1-C2-C3-O3
4	B	805	GOL	O1-C1-C2-C3
4	A	802	GOL	O2-C2-C3-O3
4	B	805	GOL	O1-C1-C2-O2
4	A	801	GOL	O2-C2-C3-O3
4	B	803	GOL	O1-C1-C2-O2
4	A	804	GOL	O1-C1-C2-O2
5	A	900	PHE	OXT-C-CA-N

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	804	GOL	2	0
3	B	806	SO4	1	0
4	A	801	GOL	2	0
4	B	803	GOL	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, 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	458/462 (99%)	0.31	47 (10%)	12 11	12, 26, 57, 104	6 (1%)
1	B	457/462 (98%)	0.10	40 (8%)	15 14	12, 21, 52, 103	5 (1%)
All	All	915/924 (99%)	0.20	87 (9%)	14 12	12, 24, 57, 104	11 (1%)

All (87) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	238	LEU	9.2
1	B	9	ILE	8.9
1	B	15	LEU	7.3
1	A	240	THR	6.9
1	B	238	LEU	6.8
1	A	267	GLY	6.5
1	A	1	MET	6.3
1	A	3	TRP	6.3
1	A	136	SER	5.9
1	A	15	LEU	5.7
1	B	1	MET	5.6
1	B	264	GLY	5.3
1	B	197	VAL	5.3
1	A	239	GLN	4.8
1	B	266	ASP	4.8
1	B	267	GLY	4.7
1	B	263	ASP	4.7
1	A	399	ARG	4.5
1	A	16	PRO	4.5
1	B	237	ASN	4.4
1	A	268	GLU	4.4
1	A	241	ALA	4.4
1	A	10	ASP	4.3
1	B	16	PRO	4.2

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Mol	Chain	Res	Type	RSRZ
1	B	462	ASP	4.2
1	B	195	HIS	4.2
1	A	339	GLY	4.2
1	B	236	ARG	4.2
1	A	235	ASP	4.1
1	A	135	ARG	4.1
1	B	239	GLN	4.1
1	A	234	ALA	4.0
1	A	236	ARG	4.0
1	A	137	ALA	3.9
1	A	462	ASP	3.7
1	B	200	TRP	3.6
1	A	195	HIS	3.6
1	B	202	ARG	3.6
1	A	189	SER	3.6
1	A	266	ASP	3.5
1	B	265	ASP	3.5
1	B	401	LEU	3.4
1	A	264	GLY	3.4
1	B	198	HIS	3.3
1	A	378	GLY	3.2
1	B	400	ALA	3.2
1	A	188	SER	3.2
1	A	198	HIS	3.1
1	B	336	SER	3.1
1	A	17	PRO	3.0
1	B	367	PRO	2.9
1	B	240	THR	2.9
1	B	235	ASP	2.9
1	A	237	ASN	2.9
1	A	367	PRO	2.8
1	A	22	LEU	2.7
1	A	200	TRP	2.7
1	B	378	GLY	2.7
1	B	206	ARG	2.6
1	B	2	ASN	2.6
1	B	377	THR	2.6
1	B	207	THR	2.6
1	A	138	ASP	2.6
1	A	197	VAL	2.5
1	A	2	ASN	2.5
1	A	369	HIS	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	265	ASP	2.4
1	B	199	ASP	2.4
1	B	369	HIS	2.4
1	A	263	ASP	2.3
1	A	206	ARG	2.3
1	B	374	GLU	2.3
1	B	234	ALA	2.3
1	A	233	VAL	2.3
1	A	376	SER	2.2
1	A	430	THR	2.2
1	A	6	ASP	2.2
1	A	7	ILE	2.2
1	A	377	THR	2.2
1	B	416	ASN	2.2
1	B	268	GLU	2.1
1	B	201	ASN	2.1
1	B	4	THR	2.1
1	B	399	ARG	2.1
1	A	190	GLY	2.1
1	B	189	SER	2.1
1	A	196	LEU	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 oligosaccharides 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($\text{\AA}^2$)	Q<0.9
3	SO4	A	803	5/5	0.79	0.40	34,39,46,49	5
4	GOL	B	804	6/6	0.81	0.22	39,42,45,50	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	GOL	A	804	6/6	0.83	0.22	20,20,20,20	0
3	SO4	B	806	5/5	0.84	0.11	57,59,62,64	0
4	GOL	A	802	6/6	0.85	0.18	41,49,52,54	0
4	GOL	A	801	6/6	0.90	0.16	31,39,41,46	0
4	GOL	B	805	6/6	0.93	0.11	22,28,30,31	0
5	PHE	A	900	12/12	0.94	0.08	18,21,24,25	0
5	PHE	B	900	12/12	0.94	0.08	19,24,26,29	0
4	GOL	B	803	6/6	0.95	0.09	22,27,35,36	0
3	SO4	B	801	5/5	0.98	0.09	24,25,30,34	0
2	MN	B	700	1/1	0.98	0.05	25,25,25,25	0
2	MN	A	700	1/1	0.99	0.03	25,25,25,25	0
3	SO4	A	800	5/5	0.99	0.07	24,25,27,30	0
3	SO4	B	802	5/5	0.99	0.04	20,22,23,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.